Putting Humpty-Dumpty together again: Reconstructing functions from their projections.

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ABSTRACT:
We present a problem decomposition approach to reduce neural net training times. The basic idea is to train neural nets in parallel on marginal distributions obtained from the original distribution (via projection), and then reconstruct the original table from the marginals (via a procedure similar to the join operator in database theory). A function is said to be reconstructible, if it may be recovered without error from its projections. Most distributions are non-reconstructible. The main result of this paper is the Reconstruction theorem, which enables non-reconstructible functions to be expressed in terms of reconstructible ones, and thus facilitates the application of decomposition methods.

Main Category: Algorithms & Architectures,
Sub-Category: Constructive & Pruning Algorithms.
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Reconstructing functions from their projections.

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1 Introduction

Suppose a neural network is to be taught a complicated, real valued multivariable function \( f : \mathbb{R}^n \to \mathbb{R} \), represented as a table. Real-time calculation of \( f \) is assumed to be infeasible or undesirable. The decomposition-reconstruction problem is the problem of reconstructing a function from its projections. Metaphorically, it is the problem of putting Humpty-Dumpty together again, without having to call in all the King's horses or men. We present a solution to the decomposition-reconstruction problem, using tools and techniques from database theory.

The main idea underlying our solution is quite simple. We start with a multivariable distribution and obtain a set of marginals. Obviously, it is not always possible to reconstruct the original distribution from the marginals. One may construct a "best" reconstruction in many different ways. A particularly appealing criterion is to obtain a reconstruction that has the maximum Shannon entropy. In probabilistic database theory such a maximum entropy reconstruction is referred to as the join distribution. It is possible then, to compute an error distribution; basically nothing more than the difference between the original and the join distributions. It may be shown
(the Reconstruction theorem) that after some minor processing, the error distribution is reconstructible. In other words, the error distribution after some modification, is also a maximum entropy reconstruction of a set of marginals. The original distribution may then be expressed as a linear combination of the join distribution and the error distribution. The marginals of each of these two distributions are taught to a disjoint set of neural nets in parallel. For other approaches to the problem of reducing the training time, of neural networks, (for example, by reducing the size of the network), see [5, 15, 17].

The decomposition-reconstruction problem has of course, already been studied in great depth from various points of view, including Reconstructability theory, Database theory, Statistics and Information theory. Many of the fundamental issues involved were first worked out by Ross Ashby in the early 60's, in a Systems theoretic context [1]. The joint work of George Klir and Roger Cavallo laid the foundations for the notion of system reconstructability [6, 7], much of it destined to be rediscovered in the then emerging field of relational database theory. In recent years there has been an increasing interest in a generalization of relational databases, the theory of probabilistic databases [2, 8].

We believe probabilistic database theory offers a fresh, simple, powerful and elegant approach to the problem of training neural networks on tables of data. It is possible to develop a much more systematic and logical approach to this problem, free of the rampant ad hocery present in current approaches.

Section 2 defines the basic tools underlying this approach, drawn from the theory of probabilistic databases [8]. In Section 3 we present the Reconstruction theorem, the central theorem of this paper. Section 4 is concerned with how ideas from probabilistic database theory may be used to advantage in neural network theory. Finally, in Section 5, we illustrate these ideas by applying them to a control problem. There are many kinds of neural networks. For concreteness, the semilinear feedforward network may be assumed to be the basic underlying neural architecture for this paper (though the results are architecture independent).

2 Probabilistic Systems

Definition 2.1 A probabilistic system (PS) is a 2-tuple \( D = (V, p) \) where, \( V = \{v_1, \ldots, v_n\} \), referred to as the scheme of the PS, is a non-empty set
of variables, each \( v_i \) taking values from a finite set \( S_i \), and,

\[
p : T(V) \rightarrow [0,1] \\
\sum_{t \in T(V)} p(t) = 1;
\]

where, the finite product space \( T(V) = \prod_i S_i \), is the set of tuples of the system \( D \). A model of a scheme \( V \) is a set \( X = \{V_1, \ldots, V_n\} \) such that \( \cup_i V_i \subseteq V \) and \( V_i \not\subseteq V_j, \forall i, j \in \{1, \ldots, n\} \). A model \( X \), of a scheme \( V \), is non-trivial iff \( X \neq \emptyset \) and \( X \neq \{V\} \). A collection of probabilistic systems \( D_1, \ldots, D_k \) is a probabilistic database (PD). A collection of probabilistic systems \( D_i = (V_i, p_i) \) together define a probabilistic database (PD).

Notation:
The tuples of a PS will be referred to by small greek letters \( \alpha, \beta, \) etc. The \( k \)th component of a tuple \( \alpha \), will be denoted by \( \alpha_k \). Given two schemes \( V \) and \( V' \) with \( V' \subseteq V \), and tuple \( \alpha \in T(V) \) then \( \alpha[V'] \) is the restriction of \( \alpha \) to variables in \( V' \).

Probabilistic systems closely resemble contingency tables. However, viewing them as generalizations of relational databases leads to important conceptual and computational advantages. It is fortunate that the development of probabilistic database theory was largely motivated by problems peculiar to database theory. A wealth of new ideas (at least, most of them are only a decade old) lie waiting to be interpreted in contexts and applications, other than database theory. We will consider this issue in greater detail in Section 4. The next two definitions formalize the notions of decomposition and reconstruction.

**Definition 2.2 (Projections):** Let \( D = (V, p) \) be a probabilistic system. Let \( V = \{v_1, \ldots, v_n\} \), where variables \( v_i \) take values from finite sets \( S_i \). Let \( V' = \{v_1, \ldots, v_k\} \subseteq V, k \leq n \) and \( T(V') = \prod_{i=1}^k S_i \). The projection of \( p \) onto \( V' \) yields a new distribution, \( p' = \pi_{V'}(p) \) such that,

\[
p' : T(V') \rightarrow [0,1] \\
p'(\beta) = \sum_{\alpha \in T(V), \alpha[V]=\beta} p(\alpha)
\]

The projection of a distribution \( p \) onto a model, \( X = \{V_1, \ldots, V_k\} \) is,

\[
\pi_X(p) = \{\pi_{V_1}(p), \ldots, \pi_{V_k}(p)\}.
\]
Definition 2.3 (Extensions): Let $P_V$ denote the set of all probability distributions over the finite product space $T(V)$. Then if $V'$ is a scheme with distribution $p$, $V' \subseteq V$, the extension of $p'$ to the scheme $V$, is the set $E^V(p')$ of all preimages of $p'$ under the mapping $\pi_{V'}$. Formally, $E^V(p') = \{ p \in P_V | \pi_{V'}(p) = p' \}$. Similarly, the extension of a set of distributions $P = \{ p_1, \ldots, p_r \}$, is defined to be the intersection of the extensions of its constituent distributions. Formally, $E^V(P) = \cap_{p \in P} E^V(p)$. In particular, if $X$ is a model for a scheme $V$, the extension of $V$ relative to $X$, $E^V(\pi_X(p))$, is seen to be the set $E^V(\pi_X(p)) = \{ p' \in P_V | \pi_X(p') = \pi_X(p) \}$.

Example 2.1 Consider the PS $D = (V = \{v_1, v_2, v_3\}, p)$:

<table>
<thead>
<tr>
<th>$v_1$</th>
<th>$v_2$</th>
<th>$v_3$</th>
<th>$p(\cdot)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>0</td>
<td>0</td>
<td>0.12 = $p_0$</td>
</tr>
<tr>
<td>0</td>
<td>0</td>
<td>1</td>
<td>0.04 = $p_1$</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>0</td>
<td>0.32 = $p_2$</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>1</td>
<td>0.11 = $p_3$</td>
</tr>
<tr>
<td>1</td>
<td>0</td>
<td>0</td>
<td>0.01 = $p_4$</td>
</tr>
<tr>
<td>1</td>
<td>0</td>
<td>1</td>
<td>0.00 = $p_5$</td>
</tr>
<tr>
<td>1</td>
<td>1</td>
<td>0</td>
<td>0.29 = $p_6$</td>
</tr>
<tr>
<td>1</td>
<td>1</td>
<td>1</td>
<td>0.11 = $p_7$</td>
</tr>
</tbody>
</table>

Choose a model $X$ of $V$ to be $\{ \{v_1, v_2\}, \{v_2, v_3\}, \{v_1, v_3\} \}$. Then $P = \pi_X(p) = \{ \pi_{\{v_1, v_2\}}, \pi_{\{v_2, v_3\}}, \pi_{\{v_1, v_3\}} \}$ is:

<table>
<thead>
<tr>
<th>$v_1$</th>
<th>$v_2$</th>
<th>$\pi_{{v_1, v_2}}$</th>
<th>$v_2$</th>
<th>$v_3$</th>
<th>$\pi_{{v_2, v_3}}$</th>
<th>$v_1$</th>
<th>$v_3$</th>
<th>$\pi_{{v_1, v_3}}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>0</td>
<td>0.16</td>
<td>0</td>
<td>0</td>
<td>0.13</td>
<td>0</td>
<td>0</td>
<td>0.44</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>0.43</td>
<td>0</td>
<td>1</td>
<td>0.04</td>
<td>0</td>
<td>1</td>
<td>0.15</td>
</tr>
<tr>
<td>1</td>
<td>0</td>
<td>0.01</td>
<td>1</td>
<td>0</td>
<td>0.61</td>
<td>1</td>
<td>0</td>
<td>0.30</td>
</tr>
<tr>
<td>1</td>
<td>1</td>
<td>0.40</td>
<td>1</td>
<td>1</td>
<td>1.22</td>
<td>1</td>
<td>1</td>
<td>0.11</td>
</tr>
</tbody>
</table>

It is clear that there are many such $p$'s that would have produced precisely the same sub-tables. In fact any solution to the set of equations below:

- $p_0 + p_1 = 0.16$
- $p_0 + p_4 = 0.13$
- $p_0 + p_2 = 0.44$
- $p_2 + p_3 = 0.43$
- $p_1 + p_5 = 0.04$
- $p_1 + p_3 = 0.15$
- $p_4 + p_5 = 0.01$
- $p_2 + p_6 = 0.61$
- $p_4 + p_6 = 0.30$
- $p_6 + p_7 = 0.40$
- $p_3 + p_7 = 0.22$
- $p_5 + p_7 = 0.11$
will be a member of $E^V(P) = E^V(\pi_X(p))$. The above twelve equations in eight unknowns characterize $E^V$ completely. Note that the system of equations is overdetermined; it is by no means guaranteed that solutions will always exist. When solutions do exist, the set of projections are said to be \textit{consistent}. Equivalently, $E^V(P) = E^V(\pi_X(p)) \neq \emptyset$. 

A model $X$ of $V$, partitions $P_v$ into classes $E^V(\pi_X(p))$ equivalent with respect to projections onto $X$. The polyhedral structure of $E^V$, can be used to show that there exists a unique representative for each class, the maximum entropy distribution, or the join distribution [13]. Formally:

**Definition 2.4 (Join):** Let $\{D_i = (V_i, p_i)\}$ be a set of probabilistic systems. Let $V = \cup_i V_i$ and $X = \{V_1, \ldots, V_n\}$ be a model for $V$. Let $P = \{p_1, \ldots, p_n\}$. If $E^V(P) \neq \emptyset$ (the distributions are consistent), then define the \textit{join} of the systems $\{D_i\}$, $D_w = (V, p_w)$, such that,

$$H(p_w) = \max \{H(p') \mid p' \in E(\pi_X(P))\}$$

where $H()$ is the Shannon entropy function. In particular, $p_w$ will be referred to as the \textit{probabilistic join} of the distributions $\{p_1, \ldots, p_n\}$ and is denoted by $p_w = \oplus (\{p_1, \ldots, p_n\})$. If $p = \oplus (\pi_X(p))$ then $D = (V, p)$ is said to be \textit{reconstructible} relative to $X$.

**Remark 2.1** To reconstruct via the join operator, is to reconstruct adopting a maximum entropy philosophy. The pros and cons of this approach have been argued extensively in the literature [13, 14, 18, 19]. The $\oplus$ operator may be thought of as a binary operator acting on a pair of probabilistic systems, and defined in a manner analogous to the classical relational database join operator. It is a remarkable fact that the generalization of the database join operator to probabilistic systems is equivalent to maximum entropy reconstruction.

### 3 Function Reconstruction

A relational database is a set $RD = \{D_1, D_2, \ldots, D_n\}$ of \textit{relational systems} $D_i = (V_i, \tau_i)$, where the range of each $\tau_i$ is the set $\{0, 1\}$. These functions are relatively trivial to reconstruct owing to the fact that, at the cost of introducing one additional variable, for any relational system, a non-trivial model exists such that, the system is reconstructible relative to that model, with respect to relational join. Formally,
**Theorem 3.1** Let $D = (V, r)$ be a relational system and $X = \{V_1, V_2\}$ be a model for the scheme $V$. Let $v_g$ be a variable such that $v_g \notin V$, and let $v_g$ take values from a finite set $S_g$. For the scheme $V' = V \cup \{v_h\}$, let $X' = \{V_1 \cup \{v_h\}, V_2 \cup \{v_h\}\}$ be a model. Then, there exists a relational system $D' = (V', r')$ such that,

$$
egin{align*}
   r_1' &= \pi_{\{V_1\}}(r') = \pi_{\{V_1\}}(r) = r_1 \\
   r_2' &= \pi_{\{V_2\}}(r') = \pi_{\{V_2\}}(r) = r_2 \\
   r' &= \otimes (r_1', r_2')
\end{align*}
$$

A proof for the above theorem is given in Appendix A. We adopt the position that $0-1$ functions are intrinsically easy to learn and are interested primarily in the general case of real valued functions. It is assumed that efficient methods exist for the computation of a given $0-1$ function. It may even be the case that a methodology other than that of neural networks be more appropriate for computing functions that take only either of two values.

The situation for probabilistic systems is embodied in the Reconstruction theorem. This theorem asserts that the error distribution, obtained as a difference between the original distribution and the maximum entropy reconstruction of a set of marginals, is always reconstructible. This is done by separating the magnitude of the error distribution ($p_1$) from its sign ($p_2$). The former is reconstructible with respect to a non-trivial model, and the latter is converted to a simple $0-1$ function and is trivially learnable. The practical implications are that we may always parallelize the training process by projecting the original function onto some appropriate model. The sum of the pre-calculated error and the join enables us to retrieve the original distribution. The following definition aids the statement of the theorem. A proof of Theorem 3.2 is given in Appendix B.

**Definition 3.1** Let $D = (V, p)$ be a PS, $|V| \geq 2$, $X$ be a model for $V$, $T$ be the set of tuples of $D$, and let $p_\omega = \min (\pi_X(p))$. $err : T \rightarrow \mathbb{R}$ is the error function of $D$, where $err(\alpha) = p(\alpha) - p_\omega(\alpha)$. Further, if $K = \sum_{\gamma \in T} |p(\gamma) - p_\omega(\gamma)| = \sum_{\gamma \in T} |err(\gamma)|$, define $p_\varepsilon : T \rightarrow [0, 1]$, and $p_\sigma : T \rightarrow \{0, 1\}$ such that, $p_\varepsilon(\alpha) = |err(\alpha)|/K$, and $p_\sigma(\alpha) = 0$ if $err(\alpha) \leq 0$, otherwise $p_\sigma(\alpha) = 1$. 

Theorem 3.2 (Reconstruction theorem) Let $D = (V, p)$ be a PS, $|V| \geq 2$, $X$ be a model for $V$, $p_X = \pi_X(p)$, and let $p_c$, $p_s$, and $K$ be as defined above. Then, $p_c = \pi_X(p)$ (i.e. $p_c$ is reconstructible). Hence $p_c$ is reconstructible and $p(\alpha) = p_X(\alpha) + K(2p_s(\alpha) - 1) \times p_c(\alpha) \forall \alpha \in T$.

When is a function reconstructible? An answer to this query is provided by the concept of functional and multivalued dependency. If one variable $y$, functionally dependent on another $x$, then given $x$, there is no uncertainty regarding the $y$. On the other hand if $x$ multidetermines $y$, knowledge of $x$ may not remove all our uncertainty regarding $y$, but it is definitely the case that knowledge of any other variable, say $z$, will not remove any further uncertainty regarding $y$. Formally,

Definition 3.2 For a probabilistic system $D = (V, p)$, with $X, Y \subseteq V$, and with the distributions over $X$ and $Y$ obtained by projection of $p$, we say that $X$ is functionally dependent on $Y$, denoted $X \rightarrow Y$ iff. $H(Y \mid X) = 0$. Additionally, we say that $X$ multidetermines $Y$, denoted $X \rightarrow\rightarrow Y$, iff. $H(Y \mid X) = H(Y \mid V - (Y \cup X))$. Let $Z = V - (X \cup Y)$. The degrees of functional and multivalued dependencies of $Y$ on $X$, $FD(D; X, Y)$, $MVD(D; X, Y)$ respectively, are given by:

$$FD(D; X, Y) = \begin{cases} 1 & \text{if } H(Y) = 0 \\ \frac{H(Y) - H(Y \mid X)}{H(Y)} & \text{otherwise} \end{cases}$$

$$MVD(D; X, Y) = \begin{cases} 1 & \text{if } H(Y \mid X) = 0 \\ \frac{H(Y \mid \{X \cup Z\})}{H(Y \mid X)} & \text{otherwise} \end{cases}$$

Clearly, if $X \rightarrow Y$ then, $X \rightarrow\rightarrow Y$. In general however, the converse is not true. The answer to our query about the reconstructability of a function with respect to a model, lies in Theorem 3.3, first proved for relational databases by Fagin [9], for probabilistic databases by Cavallo and Pittarelli [8], and may be generalized for models other than binary ones.

Theorem 3.3 Let $D = (V, p)$ be a PS, $X, Y \subseteq V$ and $X \rightarrow\rightarrow Y$. Then $p$ is reconstructible with respect to the binary model $\{X, Y\}$. Formally,

$$p = \pi_X(p), \pi_Y(p)$$
4 Choosing a Model

It is clear that given a $PS \ D = (V, p)$, a model $X$ for $V$ has to be chosen judiciously, so that effective use can be made of parallelization. We illustrate how basic concepts from database theory suggest guidelines for such a choice.

A model $X$ may be viewed as a reduced hypergraph, basically a collection of some of the incomparable subsets of a set [4]. An important class of reduced hypergraphs is that of the acyclic hypergraphs. Hypergraph acyclicity is a non-trivial generalization of the familiar concept of a tree (in a graph), and has deep connections with many other fields [16]. In hypergraphs, there are three degrees of acyclicity, $\alpha$, $\beta$ and $\gamma$ [11]. The main guideline for choosing a model, is to select one that is also acyclic ($\alpha$, $\beta$ or $\gamma$) [3]. The main reason for choosing an acyclic model is that its presence guarantees many nice properties for the join operator such as monotonicity and consistency, [10, 11] and convergence [7, 18].

Faced with two or more acyclic models, a choice may be made on the basis of their entropies relative to a standard distribution, such as the uniform distribution. Given two distributions $p_1$ and $p_2$, define $H(p_1 \parallel p_2)$ to be the directed divergence (cross-entropy, relative entropy) of $p_1$ with respect to $p_2$ [12]. Then, if we have two acyclic models, $X_i$ and $X_j$, and $p_i = \pi (\pi_{X_i}(p))$, $p_j = \pi (\pi_{X_j}(p))$, select $X_i$ over $X_j$ iff, $H(p_i \parallel u) \geq H(p_j \parallel u)$, where $u$ is the uniform distribution. Choosing the smaller value ensures that information loss is minimized in the decomposition process.

For a multivariable function, it may be the case that some variables group together naturally into interdependent blocks. Models that give decompositions along these "natural fissures" are better than those that ignore such interdependencies. These qualitative statements may be made precise by using the dependency measures defined in the last section.

1. If $X, Y \subseteq V$, and if $X \rightarrow Y$, then decompose the system $D = (V, p)$ onto the model $\{X, Y\}$.

2. If such a multivalued dependency does not exist for a scheme $V$, then choose a decomposition of $V$ into subsets $X$ and $Y$, such that $MVD(D; X, Y)$ is maximum.
5 Example

The methods outlined earlier are particularly suited for control problems. For example, consider the well known Gantry Crane system control problem. A gantry crane is used to move large parts and assemblies from one location to another on a factory floor; the control system is responsible for the horizontal motion of crane and load. Four key variables may be identified, (see Figure 5), $x_1$, $x_2$, the positions of the load and crane, with respect to the origin, and $v_1$ and $v_2$, the velocities of the load and crane respectively. The problem is to compute the value of the force, for a given set of values of $x_1$, $x_2$, $v_1$ and $v_2$. It is assumed that it is infeasible to compute $F$ in real-time using the above two equations. Essentially two data sets (one for testing and the other for training, were generated, using the system control equations. All values were normalized to lie in closed-open interval $[0, 1)$. The domains of the input variables were quantized into ten states, by using the following rule for classification:

$$x/10 \leq y < (x + 1)/10, \quad x \in \{0, \ldots, 9\}$$
i.e. \( y \) was said to belong to "class \( x \)" if it satisfied the above equation for the given \( x \). Thus if \( x_1 = 0.56 \) then \( x_1 \) would belong to class \( 0 \) etc. Note that quantization was done for the input variables alone and not for \( F \). In what follows, the training table will be referred to as the master table or Table 0. A typical set of entries in the master table is shown below.

<table>
<thead>
<tr>
<th>( x_1 )</th>
<th>( v_1 )</th>
<th>( x_2 )</th>
<th>( v_2 )</th>
<th>( F )</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>3</td>
<td>1</td>
<td>1</td>
<td>0.0068925912</td>
</tr>
<tr>
<td>1</td>
<td>3</td>
<td>1</td>
<td>1</td>
<td>0.0073849196</td>
</tr>
<tr>
<td>\ldots</td>
<td>\ldots</td>
<td>\ldots</td>
<td>\ldots</td>
<td>\ldots</td>
</tr>
<tr>
<td>6</td>
<td>3</td>
<td>6</td>
<td>5</td>
<td>0.0046771155</td>
</tr>
<tr>
<td>6</td>
<td>4</td>
<td>6</td>
<td>5</td>
<td>0.0085336845</td>
</tr>
<tr>
<td>\ldots</td>
<td>\ldots</td>
<td>\ldots</td>
<td>\ldots</td>
<td>\ldots</td>
</tr>
</tbody>
</table>

The training procedure is schematically outlined in Figure 5. Table 0 was decomposed into Tables 1, 2, 5, 7 and 8. A separate 1-2-1 layer backpropagation network was trained to learn the relation in each of these tables. The testing procedure involved the following,

- For each input \((x_1, v_1, x_2, v_2)\), in testing sample do:
  1. Present \((x_1, v_1, x_2)\) to the Table 1 and Table 7 networks\(^1\); present \((x_2, v_2, v_1)\) to the Table 2 and Table 8 networks; and similarly present \((x_1, v_1, x_2, v_2)\) to the Table 0 and Table 5 networks.
  2. Compute \( p_r = p_o + K (2p_o - 1) \times \sigma (p'_o, p''_o) \), the reconstructed output, where \( K \) is a normalization factor obtained while training. This gives the output obtained after decomposition.
  3. For each network, the deviation with the actual output value is noted.

The results are shown below. The last row lists the net mean square error of the reconstructed output compares very favorably with that for the Table 0 network.

\(^1\) By "Table 1" network, we mean the "the network originally trained on Table 1" etc.
After Quantization

Master Table
Table 0

Projecting onto $V_1$
Table 1

Projecting onto $V_2$
Table 2

Join Table
Table 3

Table 4

Error table = Table 0 - Table 3

0 - 1 table, Derived error table
Table 5

Table 6

Table 7

Table 8

Figure 2: Tables Generated during Training Phase
### Tables

<table>
<thead>
<tr>
<th>Trained/Tested</th>
<th>Training</th>
<th>Testing</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>Samples</td>
<td>M.S.E</td>
</tr>
<tr>
<td>Table 0 = (V, p)</td>
<td>120</td>
<td>0.000005</td>
</tr>
<tr>
<td>Table 1 = (V', p')</td>
<td>120</td>
<td>0.000043</td>
</tr>
<tr>
<td>Table 2 = (V'', p'')</td>
<td>322</td>
<td>0.000003</td>
</tr>
<tr>
<td>Table 7 = (V', p'_i)</td>
<td>120</td>
<td>0.000013</td>
</tr>
<tr>
<td>Table 8 = (V'', p''_i)</td>
<td>222</td>
<td>0.000008</td>
</tr>
<tr>
<td>Reconstructed table = (V, p_r)</td>
<td>846</td>
<td>0.0000101</td>
</tr>
</tbody>
</table>

### 6 Conclusion

The Divide-and-Conquer paradigm has long been recognized as a useful tool for conquering complexity. Our work address the question of when problem decomposition is possible. The Reconstruction theorem (Section 3), asserts that any non reconstructible function may be expressed as a linear combination of reconstructible functions. This suggests a natural input space decomposition method for training neural networks on functions expressed as large tables, drawing on ideas from probabilistic database theory. This approach will be most useful in solving problems characterized by a large number of inputs.

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Appendix A

Theorem 4.1 Let $D = (V, r)$ be a relational system and $X = \{V_1, V_2\}$ be a model for the scheme $V$. Let $v_g$ be a variable such that $v_g \not\in V$, and let $v_g$ take values from a finite set $S_g$. For the scheme $V' = V \cup \{v_g\}$, let $X' = \{V_1 \cup \{v_g\}, V_2 \cup \{v_g\}\}$ be a model. Then, there exists a relational system $D' = (V', r')$ such that,

\[
\begin{align*}
    r_1' &= \pi_{\{V_1\}}(r') = \pi_{\{V_1\}}(r) = r_1 \\
    r_2' &= \pi_{\{V_2\}}(r') = \pi_{\{V_2\}}(r) = r_2 \\
    r' &= \Join (r_1', r_2')
\end{align*}
\]

Proof:
The basic idea is to extend the set of variables $V$, by one, and construct a reconstructible relational system $D' = (V', r')$, which yields the same projections as $D$, including $D$ itself (i.e. $\pi_{\{V_1\}}(r') = r$).

Let $D_1 = (V_1, \pi_{V_1}(r))$ and $D_2 = (V_2, \pi_{V_2}(r))$ be the projections of $D$ onto the schemes $V_1$ and $V_2$ respectively, and let $D_\wedge$ be their relational join. Also, let $T_1, T_2, T_3$, and $T_\wedge$ be the set of tuples in $D, D_1, D_2$ and $D_\wedge$ respectively. If $D = (V, r)$ is reconstructible with respect to $X = \{V_1, V_2\}$, then there is nothing to prove and we are done. So we may assume that $D$ is not reconstructible with respect to $X$, relative to relational join. From the definition of relational join, it then follows that $T_1 \subseteq T_2$. Let $\Delta T$ denote the set of such tuples found in $T_\wedge$, but not in $T$.

Every tuple $t$ in $\Delta T$ is formed by the join of a pair, consisting of one tuple, say $t_1 \in T_1$, from $D_1$, and one from $D_2$, say $t_2 \in T_2$. Tuples $t_1$ and $t_2$ agree on the set of attributes belonging to the set $V_1 \cap V_2$. If they didn't so agree, then the tuple $t$ would not have been created while taking the join of $D_1$ and $D_2$. This suggests a simple method to prevent the creation of unwanted tuples in $D_\wedge$.

Construct the systems $D'_1 = (V_1 \cup \{v_g\}, r'_1)$ and $D'_2 = (V_2 \cup \{v_g\}, r'_2)$ and let $D' = (V \cup \{v_g\}, r')$ be the join of the systems $D'_1$ and $D'_2$. Let $t \not\in \Delta T$, i.e. $t \in T_\wedge \cap T$. Now $t = t_1 \Join t_2$ for some $t_1 \in T_1$, and $t_2 \in T_2$. Now, for every such $t_1$ and $t_2$, there exists $t'_1 \in T'_1, t'_2 \in T'_2$ with $t'_1[V_1] = t_1, t'_2[V_2] = t_2$, and $t'_1[\{v_g\}] = t'_2[\{v_g\}]$ (see Section 2 for notation). In other words, the tuples $t'_1$ and $t'_2$ agree on the value of $v_g$, and $t'_1$ agrees with $t_1$ on all other variables. Similarly, $t'_2$ agrees with $t_2$ on all variables.
other than \( v_2 \). By choosing a different value for \( v_2 \) for each other pair \( t_1 \) and \( t_2 \), we can ensure that, \( \pi_{(V_1)}(r_1') = r_1, \pi_{(V_2)}(r_2') = r_2 \) and \( \pi_{(V)}(r') = r \). 

**Example A.1** This example illustrates the procedure described in the proof. Consider the relational database \( D = (V, r) \) given by the table below.

<table>
<thead>
<tr>
<th>( x )</th>
<th>( y )</th>
<th>( z )</th>
<th>( r )</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>0</td>
<td>0</td>
<td>1</td>
</tr>
<tr>
<td>0</td>
<td>0</td>
<td>1</td>
<td>0</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>0</td>
<td>1</td>
</tr>
<tr>
<td>0</td>
<td>1</td>
<td>1</td>
<td>1</td>
</tr>
<tr>
<td>1</td>
<td>0</td>
<td>0</td>
<td>1</td>
</tr>
<tr>
<td>1</td>
<td>0</td>
<td>1</td>
<td>0</td>
</tr>
<tr>
<td>1</td>
<td>1</td>
<td>0</td>
<td>0</td>
</tr>
<tr>
<td>1</td>
<td>1</td>
<td>1</td>
<td>1</td>
</tr>
</tbody>
</table>

Upon projecting the above system onto the model \( X = \{ V_1, V_2 \} \), where \( V_1 = \{ x, y \} \) and \( V_2 = \{ y, z \} \) we get,

\[
\begin{array}{c|c}
  x & y \\
  \hline
  0 & 0 \\
  0 & 1 \\
  1 & 0 \\
  1 & 1 \\
\end{array}
\quad
\begin{array}{c|c}
  y & z \\
  \hline
  0 & 0 \\
  0 & 1 \\
  1 & 0 \\
  1 & 1 \\
\end{array}
\]

The relational join of \( D_1 = (\{ x, y \}, r_1) \) and \( D_2 = (\{ y, z \}, r_2) \) yields,

\[
\begin{array}{c|c|c}
  x & y & z \\
  \hline
  0 & 0 & 0 \\
  0 & 0 & 1 \\
  0 & 1 & 0 \\
  0 & 1 & 1 \\
  1 & 0 & 0 \\
  1 & 0 & 1 \\
  1 & 1 & 0 \\
  1 & 1 & 1 \\
\end{array}
\]

The tuple marked with an arrow is not present in the original system \( D \). Hence \( D \) is not reconstructible. We note that the problem stems from the fact that the tuple \( t_1 = 11 \) in \( D_1 \) joins with tuple \( t_2 = 10 \) in \( D_2 \). All the following combinations however are required viz.

![Image](image-url)
Introduce a new variable \( w \), such that,

\[
\begin{array}{cccc}
  x & y & w & w \\
  0 & 0 & 0 & 0 \\
  0 & 1 & 1 & 1 \\
  0 & 1 & 2 & 2 \\
  1 & 1 & 3 & 3 \\
\end{array}
\]

Or, defining two new systems \( D_1' = (\{x, y, w\}, r_1') \) and \( D_2' = (\{y, z, w\}, r_2') \),

\[
\begin{array}{cccc|c}
  x & y & w & r_1' & w & y & z & r_2' \\
  0 & 0 & 0 & 1 & 0 & 0 & 0 & 1 \\
  0 & 1 & 1 & 1 & 1 & 1 & 0 & 1 \\
  0 & 1 & 2 & 1 & 2 & 1 & 1 & 1 \\
  1 & 1 & 3 & 1 & 3 & 1 & 1 & 1 \\
\end{array}
\]

We have shown only the nonzero tuples in both \( D_1' \) and \( D_2' \). The join of \( D_1' \) and \( D_2' \), \( D'_w = (\{x, y, w, z\}, r'_w) \), yields,

\[
\begin{array}{cccc|c}
  x & y & w & z & r'_w \\
  0 & 0 & 0 & 0 & 1 \\
  0 & 1 & 1 & 0 & 1 \\
  0 & 1 & 2 & 1 & 1 \\
  1 & 1 & 3 & 1 & 1 \\
\end{array}
\]

Again all other tuples \( \alpha \), have \( r'_w(\alpha) = 0 \). It is easy to see that \( D' = D_1' \uplus D_2' \) yields \( D \) when projected onto \( \{V\} \) and also yields \( D_1 \) and \( D_2 \) when projected onto \( V_1 \) and \( V_2 \). By construction \( D' \) is reconstructible and may be used in place of \( D \). Note that \( w \) here takes values from a set \( \{0, 1, 2, 3\} \). It is not too difficult to see that in this case, we could have done with \( w \) taking values from the smaller set \( \{0, 1, 2\} \). It is fairly straightforward to extend the procedure. \( \blacksquare \).
Appendix B

Theorem 5.1 (Reconstruction theorem) Let $D = (V, p)$ be a PS, $|V| \geq 2$, $X$ be a model for $V$, $p^* = \pi_X(p)$, and let $p_\varepsilon$, $p_\sigma$, and $K$ be the derived error functions, and the normalization factor of $D$ respectively (see Definition 3.1). Then,

$$p_\varepsilon = \pi_X(p_\varepsilon)$$

i.e. $p_\varepsilon$ is reconstructible with relative to $X$. In particular,

$$p(\alpha) = p^*(\alpha) + K(2p_\sigma(\alpha) - 1) \times p_\varepsilon(\alpha) \quad \forall \alpha \in T$$

Proof:
From Theorem 3.1, we know that $p_\sigma$, being a simple $0-1$ relational function is reconstructible, although an additional variable may have to be introduced to effect this. The issue at hand is whether $p_\varepsilon$ is also reconstructible. This is quite easy to demonstrate. Suppose it is the case that

$$(6.1) \quad \pi_X(p_\varepsilon) = p'_\varepsilon \neq p_\varepsilon$$

Since $p'_\varepsilon$ is the join of the probabilistic database $\pi_X(p_\varepsilon)$, it must satisfy the consistency requirement that,

$$(6.2) \quad \pi_X(p'_\varepsilon) = \pi_X(p_\varepsilon)$$

This in turn implies that there exists at least one distribution $p' \neq p^*$, such that,

$$(6.3) \quad p'_\varepsilon = k_1 |p' - p|$$

where $k_1 = \sum_{\alpha \in T} |p'(\alpha) - p(\alpha)|$. Since $\pi_X(p'_\varepsilon) = p'_\varepsilon$, we have (from Eqns. 6.1, 6.3),

$$(6.4) \quad \pi_X(p_\varepsilon) = \pi_X(p'_\varepsilon) = \pi_X(k_1 |p' - p|)$$

Noting that, $(1) \ p_\varepsilon = k_2 |p^* - p|$, $(2) \ p'_\varepsilon$ is a maximum entropy distribution (from Eqns. 6.1, 6.2), we get,

$$(6.5) \quad H(k_1 |p' - p|) > H(k_2 |p^* - p|)$$

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It is well known that [12], if \( q \) is an arbitrary finite probability distribution of say \( n \) variables, and \( f() \) a function, with domain and range the set of all finite probability distributions on \( n \) variables (so that \( f(q) \) is also a distribution on \( n \) variables), then,

\[
H(q) = H(f(q)) + H(q|f(q))
\]

(Roughly, this translates to the statement that in the absence of no new observation, uncertainty cannot be removed by mere transformation of the distribution alone).

Define \( f_i(q) = |k_i(q - p)| \) where \( q \) is some arbitrary distribution on \( n \) variables, \( p \) is the original non reconstructible function, and \( k_i \) is a normalizing factor and depends on both \( q \) as well as \( p \). Then,

\[
H(p') = H(f_1(p')) + H(p'|f_1(p'))
= H(|k_1(p' - p)|) + H(p'|(|k_1(p' - p)|))
\]

\[
H(p_w) = H(f_2(p_w)) + H(p_w|f_2(p_w))
= H(|k_2(p_w - p)|) + H(p_w|(|k_2(p_w - p)|))
\]

But knowledge of \( k_1 |p' - p| \) determines \( p' \) (\( p \) is fixed and \( p' \) has to be non negative),

\[
H(p'|f_1(p')) = 0
\]

and similarly,

\[
H(p_w|f_2(p_w)) = 0
\]

Then Eqns. 6.6,6.7 reduce to,

\[
H(p') = H(f_1(p')) = H(|k_1(p' - p)|)
H(p_w) = H(f_2(p_w)) = H(|k_2(p_w - p)|)
\]

Comparing Eqn 6.10 with Eqn. 6.5, it is seen that,

\[
H(p_w) < H(p')
\]

This immediately implies a contradiction, since we must have:

\[
H(p_w) > H(p')
\]

(because of the definitions of \( p_w \)), Hence such a distribution \( p' \) cannot exist and the theorem follows.  

\[17\]
References


