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### An Improved Activation Function for Energy Minimization

Gadi Pinkas and Rina Dechter

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**An Improved Activation Function for Energy  
Minimization**

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**WUCS-92-02**

**August, 1992**

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# An Improved Activation Function for Energy Minimization

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WUCS-92-2

## Abstract

Symmetric networks<sup>a</sup> that are based on energy minimization, such as Boltzmann machines or Hopfield nets, are used extensively for optimization, constraint satisfaction, and approximation of NP-hard problems. Nevertheless, finding a global minimum for the energy function is not guaranteed, and even a local minimum may take an exponential number of steps. We propose an improvement to the standard activation function used for such networks. The improved algorithm guarantees that a global minimum is found in linear time for tree-like subnetworks. The algorithm is uniform and *does not* assume that the network is a tree. It performs no worse than the standard algorithms for any network topology. In the case where there are trees growing from a cyclic subnetwork, the new algorithm performs better than the standard algorithms by avoiding local minima along the trees and by optimizing the energy of these trees in linear time. The algorithm is self-stabilizing for trees (cycle-free undirected graphs) and remains correct under various scheduling demons. However, no uniform protocol exists to optimize trees under a pure distributed demon and no such protocol exists for cyclic networks under central demon.

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<sup>a</sup>A shorter version of the paper appears in the Proceedings of The American Association for Artificial Intelligence (AAAI), 1992.

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# An Improved Activation Function for Energy Minimization

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## 1. Introduction

Symmetric networks, such as Hopfield nets, and Boltzmann machines, mean-field and harmony networks are widely used for optimization, constraint satisfaction, and approximation of NP-hard problems [Hopfield 82], [Hopfield 84], [Hinton, Sejnowski 86], [Peterson, Hartman 89], [Smolensky 86, page 259].

These models are characterized by a symmetric matrix of weights and a quadratic energy function that should be minimized. Usually, each unit computes the gradient of the energy function and updates its own activation value so that the energy decreases gradually. Convergence to a *local* minimum is guaranteed, although in the worst case it is exponential in the number of units [Kasif et al. 89], [Papadimitriou et al. 90].

In many cases, the problem at hand is formulated as an energy minimization problem and the best solutions (sometimes the *only* solutions) are the global minima [Hopfield, Tank 85], [Ballard et al. 86], [Pinkas 91], [Pinkas 92]. The desired connectionist algorithm is, therefore, one that reduces the impact of shallow local minima and improves the chances of finding a global minimum. Models such as Boltzmann machines and Harmony nets use simulated annealing to escape from local minima. These models asymptotically converge to a global minimum, meaning that if the annealing schedule is slow enough, a global minimum is found. Nevertheless, such a schedule is hard to find and therefore, practically, finding a global minimum for such networks is not guaranteed even in exponential time (note that the problem is NP-hard).

In this paper, we look at the *topology* of symmetric neural networks. We present an algorithm that optimizes tree-like subnetworks<sup>1</sup> in linear time. It is based on a dynamic programming algorithm presented in [Dechter et al. 90]. Our adaptation is connectionist in style; that is, the algorithm can be stated as a simple, uniform activation function [Rumelhart et al. 86], [Feldman, Ballard 82]. It does not assume the desired topology (tree) and performs no worse than the standard algorithms for all topologies. In fact, it may be integrated with many of the standard algorithms in such a way that if the network happens to have tree-like subnetworks, the new algorithm out-performs the standard algorithms. From the negative side, we show that once cycles are introduced it is impossible to have a uniform algorithm that guarantees optimal solution.

The paper is organized as follows: Section 2 defines the problem of energy minimization and gives examples of standard techniques. Section 3 presents the new algorithm and gives examples where it out-performs the standard algorithms. Section 4 discusses negative results concerning the feasibility of uniform optimization activation functions. Section 5 summarizes.

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<sup>1</sup>The network is characterized by an undirected graph without cycles; i.e., only one path exists between any two nodes. The terms cycle-free or unrooted tree are synonymous in this context.

## 2. Connectionist energy minimization

### 2.1. The problem

Suppose a quadratic energy function of the form

$$E(X_1, \dots, X_n) = - \sum_{i < j}^n w_{i,j} X_i X_j + \sum_i^n -\theta_i X_i.$$

Each of the variables  $X_i$  may have a value of zero or one (called the activation value), and the task is to find a zero/one assignment to the variables  $X_1, \dots, X_n$  that minimizes the energy function. To avoid confusion with signs, we will consider the equivalent problem of maximizing the goodness function:

$$G(X_1, \dots, X_n) = -E(X_1, \dots, X_n) = \sum_{i < j} w_{i,j} X_i X_j + \sum_i \theta_i X_i \quad (1)$$

In connectionist approaches, we look at the network that is generated by assigning a node  $i$  for every variable  $X_i$  in the function and by creating a symmetric weighted arc (with weight  $w_{i,j} = w_{j,i}$ ) between node  $i$  and node  $j$  for every term  $w_{i,j} X_i X_j$ . Similarly, a bias  $\theta_i$  is given to unit  $i$  if the term  $\theta_i X_i$  is in the function. For example, Figure 2-b shows the network that corresponds to the goodness function  $E(X_1, \dots, X_7) = 3X_2X_3 - X_1X_3 + 2X_3X_4 - 2X_4X_5 - 3X_3 - X_2 + 2X_1$ . Each of the nodes is assigned a processing unit and the network collectively searches for an assignment that maximizes the goodness. The algorithm that is repeatedly executed in each unit/node is called the *protocol* or the *activation function*. A protocol is *uniform* if all the units execute it.

### 2.2. Examples of standard connectionist activation functions

We give examples for the discrete Hopfield network [Hopfield 82] and the Boltzmann machine [Hinton, Sejnowski 86], which are two of the most popular models for connectionist energy minimization.

In the discrete Hopfield model, each unit computes its activation value using the formula

$$X_i = \begin{cases} 1 & \text{iff } \sum_j w_{i,j} X_j \geq -\theta_i, \\ 0 & \text{otherwise.} \end{cases}$$

For Boltzmann machines the determination of the activation value is stochastic and the probability of setting the activation value of a unit to one is

$$P(X_i = 1) = \frac{1}{1 + e^{-(\sum_j w_{i,j} X_j + \theta_i)/T}}$$

where  $T$  is the annealing temperature.

Both approaches may be integrated with our topology-based algorithm; i.e., nodes that cannot be identified as parts of a tree-like topology, use one of the standard algorithms.

## 3. The algorithm

### 3.1. Assumptions

We assume that the model of communication is shared memory, multi-reader/single-writer, that scheduling is under a central demon, and that execution is fair. In a *shared memory, multi-reader/single-writer*,



each unit has a shared register called the activation register. A unit may read the content of the registers of all its neighbors, but write only its own. *Central demon* means that the units are activated one at a time in an arbitrary order.<sup>2</sup> An execution is said to be *fair* if every unit is activated infinitely often.

### 3.2. Key idea

The algorithm identifies parts of the network that have no cycles (tree-like subnetworks) and optimizes the energy on these subnetworks. Once a tree is identified, it is optimized using an adaptation of a constraint optimization algorithm for trees presented in [Dechter et al. 90]. The algorithm belongs to the family of nonserial dynamic programming methods [Bertelé, Brioschi 72].

Let us assume first that the network is an unrooted tree (cycle-free). Any such network may be directed into a rooted tree. The algorithm is based on the observation that given an activation value (0/1) for a node in a tree, the optimal assignments for all its adjacent nodes are independent of each other. In particular, the optimal assignment to the node's descendants are independent of the assignments for its ancestors. Therefore, each node  $i$  in the tree may compute two values:  $G_i^0$  and  $G_i^1$ .  $G_i^1$  is the maximal goodness contribution of the subtree rooted at  $i$ , including the connection to  $i$ 's parent whose activation is *one*. Similarly,  $G_i^0$  is the maximal goodness of the subtree, including the connection to  $i$ 's parent whose activation value is *zero*. The acyclicity property will allow us to compute each node's  $G_i^1$  and  $G_i^0$  as a simple function of its children's values, implemented as a propagation algorithm initiated by the leaves.

Knowing the activation value of its parent and the values  $G_j^0, G_j^1$  of all its children, a node can compute the maximal goodness of its subtree. When the information reaches the root, it can assign a value (0/1) that maximizes the goodness of the whole network. The assignment information propagates now towards the leaves: knowing the activation value of its parent, a node can compute the preferred activation value for itself. At termination (at stable state), the tree is optimized.

The algorithm has three basic steps:

- 1) **Directing a tree.** Knowledge is propagated from leaves towards the center of the network, so that after a linear number of steps, every unit in the tree knows its parent and children.
- 2) **Propagation of goodness values.** The values  $G_i^1, G_i^0$  are propagated from leaves to the root. At termination, every node knows the maximal goodness of its subtree, and the appropriate activation value it should assign, given that of its parent. In particular, the root can now decide its own activation value so as to maximize the whole tree.
- 3) **Propagation of activation values.** Starting with the root, each node in turn determines its activation value. After  $O(\text{depth of tree})$  steps, the units are in a stable state, which globally maximizes the goodness.

Each unit contains an *activation register* that consists of the following fields:

- $X_i$ : The activation value, which contains zero or one (or in between for some connectionist models like [Peterson, Hartman 89] and [Hopfield 84]).
- $G_i^0$ : the maximal goodness of the subtree rooted by  $i$  providing its parent's activation value is zero.
- $G_i^1$ : the maximal goodness of the subtree rooted by  $i$  providing its parent's activation value is one.
- $(P_i^1, \dots, P_i^j)$ : a bit for each of the  $j$  neighbors of  $i$  that indicates which neighbor  $k$  of  $i$  is the parent; i.e.,  $P_i^k = 1$  iff node  $k$  is the parent of  $i$  (the rest are zeros).

<sup>2</sup>Standard algorithms need to assume the same condition in order to guarantee convergence to a *local* minimum (see [Hopfield 82]). This condition can be relaxed by restricting only that adjacent nodes may not be activated at the same time.

### 3.3. Directing a tree

The goal of this algorithm is to inform every node of its role in the network and of its child-parent relationships. Nodes with a single neighbor identify themselves as leaves first and then identify their neighbor as a parent (point to it). A node whose neighbors all point towards it identifies itself as a root. A node whose neighbors all but one point towards it selects the one as a parent. Finally, a node that has at least two neighbors *not* pointing towards it identifies itself as being outside the tree.

Each unit uses one bit per neighbor to keep the pointing information:  $P_i^j = 1$  indicates that node  $i$  sees its  $j$ th neighbor as its parent. By looking at  $P_i^j$ , node  $i$  knows whether  $j$  is pointing to it.

Identifying tree-like subnetworks in a general network may be done by the following algorithm:

#### Tree Directing (for unit $i$ ):

1. Initialization: If first time, then for all neighbors  $j$ ,  $P_i^j = 0$ . /\* Start with clear pointers (this step is not needed for trees) \*/
2. If there is only a single neighbor ( $j$ ) and  $P_i^j = 0$ , then  $P_i^j = 1$ . /\* A leaf selects its neighbor as parent if that neighbor doesn't point to it \*/
3. Else, if one and only one neighbor ( $k$ ) does not point to  $i$  ( $P_i^k = 0$ ), then  $P_i^k = 1$ , and, for the rest of the neighbors,  $P_i^j = 0$ . /\*  $k$  is the parent \*/
4. Else, for all neighbors  $j$ ,  $P_i^j = 0$ . /\* Node is either a root or outside the tree \*/

In Figure 1-(a), we see a cycle-free network after the tree-directing phase. The numbers on the edges represent the values of the  $P_i^j$  bits. In Figure 1-(b), a tree-like subnetwork is identified inside a cyclic network; note that node 5 is not a root, since not all its neighbors are pointing towards it.

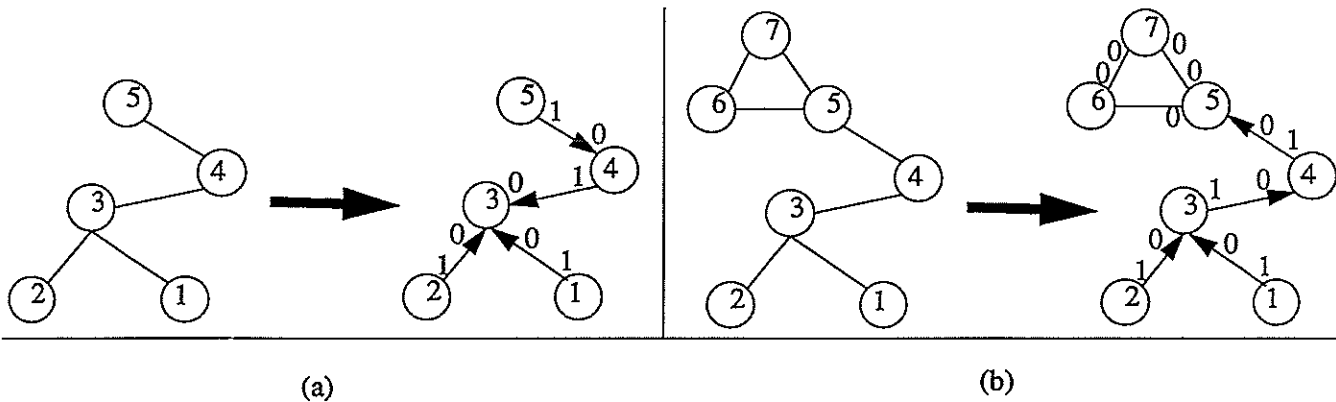


Figure 1: Directing a tree: (a) A tree, (b) A cyclic graph

### 3.4. Propagation of goodness values

In this phase, every node  $i$  computes its goodness values  $G_i^1$ ,  $G_i^0$  by propagating these two values from the leaves to the root (see figure 2).

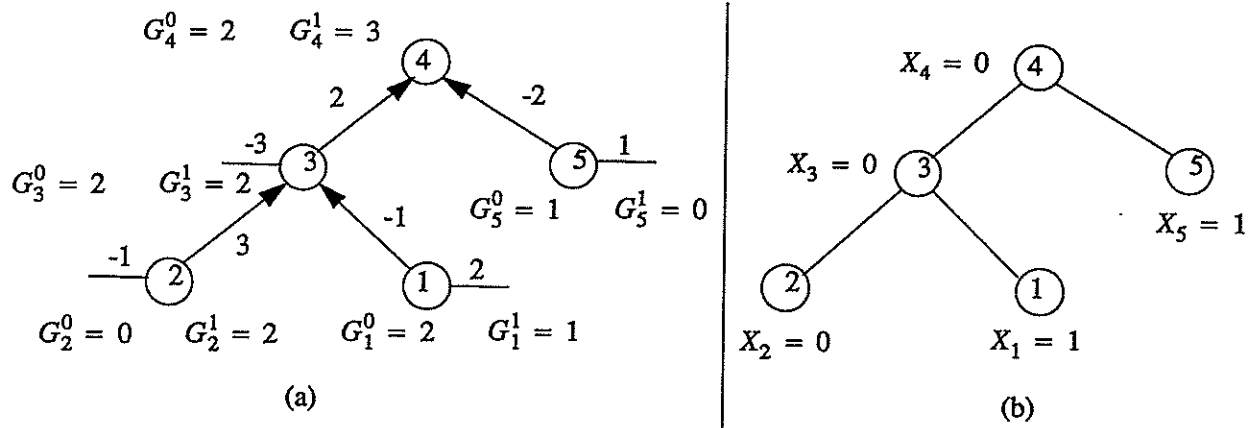


Figure 2: (a) Propagating goodness values, (b) Propagating activation values

Given a node  $i$ , its parent  $k$ , and its children  $child(i)$  in the tree, it can be shown, based on the energy function (1), that the goodness values obey the following recurrence:

$$G_i^{X_k} = \max\left\{ \sum_{j \in child(i)} G_j^{X_i} + w_{i,k} X_i X_k + \theta_i X_i \right\}.$$

Consequently, a nonleaf node  $i$  computes its goodness values using the goodness values of its children as follows: If  $X_k = 0$ , then  $i$  must decide between setting  $X_i = 0$ , obtaining a goodness of  $\sum_j G_j^0$ , or setting  $X_i = 1$ , obtaining a goodness of  $\sum_j G_j^1 + \theta_i$ . This yields

$$G_i^0 = \max\left\{ \sum_{j \in child(i)} G_j^0, \sum_{j \in child(i)} G_j^1 + \theta_i \right\}.$$

Similarly, when  $X_k = 1$ , the choice between  $X_i = 0$  and  $X_i = 1$  yields

$$G_i^1 = \max\left\{ \sum_{j \in child(i)} G_j^0, \sum_{j \in child(i)} G_j^1 + w_{i,k} + \theta_i \right\}.$$

The initial goodness values for leaf nodes can be obtained from the above (no children). Thus,  $G_i^0 = \max\{0, \theta_i\}$ ,  $G_i^1 = \{0, w_{i,k} + \theta_i\}$ .

For example: If unit 3 in Figure 2 is zero, then the maximal goodness contributed by node 1 is  $G_1^0 = \max_{X_1 \in \{0,1\}} \{2X_1\} = 2$  and it is obtained at  $X_1 = 1$ . Unit 2 (when  $X_3 = 0$ ) contributes  $G_2^0 = \max_{X_2 \in \{0,1\}} \{-X_2\} = 0$ , obtained at  $X_2 = 0$ , while  $G_2^1 = \max_{X_2 \in \{0,1\}} \{3X_2 - X_2\} = 2$  is obtained at  $X_2 = 1$ . As for nonleaf nodes, if  $X_4 = 0$ , then when  $X_3 = 0$ , the goodness contribution is  $\sum_k G_k^0 = 2 + 0 = 2$ , while if  $X_3 = 1$ , the contribution is  $-3 + \sum_k G_k^1 = -3 + 1 + 2 = 0$ . The maximal contribution  $G_3^0 = 2$  is achieved at  $X_3 = 0$ .

#### Propagating Energy Alternatives (unit $i$ ):

1. If node  $i$  is a leaf and  $k$  is its parent:  $G_i^0 = \max\{\theta_i, 0\}$ ;  $G_i^1 = \max\{w_{i,k} + \theta_i, 0\}$
2. If node  $i$  is not a leaf and has a parent  $k$  and children  $j$ :  
 $G_i^0 = \max\{\sum_{j \in children(i)} G_j^0, \sum_{j \in children(i)} G_j^1 + \theta_i\}$ ;  
 $G_i^1 = \max\{\sum_{j \in children(i)} G_j^0, \sum_{j \in children(i)} G_j^1 + w_{i,k} + \theta_i\}$ ;

### 3.5. Propagation of activation values

Once a node is assigned an activation value, all its children can activate themselves so as to maximize the goodness of the subtrees they control. When such a value is chosen for a node, its children can evaluate their own activation values, and the process continues until the whole tree is assigned.

There are two kinds of nodes that may start the process: a root which will choose an activation value to optimize the entire tree, and a non-tree node which uses a standard activation function.

When a root  $X_i$  is identified, it chooses the value 0 if the maximal goodness is  $\sum_j G_j^0$ , while it chooses 1 if the maximal goodness is  $\sum_j G_j^1 + \theta_i$ . In summary, the root chooses its value according to

$$X_i = \begin{cases} 1 & \text{iff } \sum_j G_j^1 + \theta_i \geq \sum_j G_j^0, \\ 0 & \text{otherwise.} \end{cases}$$

In Figure 2, for example,  $G_5^1 + G_3^1 + 0 = 2 < G_5^0 + G_3^0 = 3$  and therefore  $X_4 = 0$ .

An internal node whose parent is  $k$  chooses an activation value that maximizes

$$\sum_j G_j^{x_i} + w_{i,k} X_i X_k + \theta_i X_i.$$

The choice therefore, is between  $\sum_j G_j^0$  (when  $X_i = 0$ ) and  $\sum_j G_j^1 + w_{i,k} X_k + \theta_i$  (when  $X_i = 1$ ), yielding:

$$X_i = \begin{cases} 1 & \text{if } \sum_j G_j^1 + w_{i,k} X_k + \theta_i \geq \sum_j G_j^0 \\ 0 & \text{otherwise.} \end{cases}$$

As a special case, a leaf  $i$ , chooses  $X_i = 1$  iff  $w_{i,k} X_k \geq -\theta_i$ , which is exactly the discrete Hopfield activation function for a node with a single neighbor. For example, in Figure 2,  $X_5 = 1$  since  $w_{4,5} X_4 = 0 > -\theta_5 = -1$ , and  $X_3 = 0$  since  $G_1^1 + G_2^1 + 2X_4 + \theta_3 = 1 + 2 + 0 - 3 = 0 < G_2^0 + G_1^0 = 2$ . Figure 2-(b) shows the activation values obtained by propagating them from the root to the leaves.

#### Propagation of Activation Values (unit $i$ ):

1. If for all neighbors ( $j$ ),  $P_j^i = 1$ , then /\* a root \*/

$$X_i = \begin{cases} 1 & \text{if } \sum_j (G_j^1 - G_j^0) \geq -\theta_i \\ 0 & \text{otherwise} \end{cases}$$

2. else, if there exist one and only one neighbor ( $k$ ) such that  $P_k^i = 0$ , and the rest of the neighbors (if exist) have  $P_j^i = 1$ , then /\* A non-root node in the tree \*/

$$X_i = \begin{cases} 1 & \text{if } \sum_j (G_j^1 - G_j^0) + w_{i,k} X_k \geq -\theta_i \\ 0 & \text{otherwise} \end{cases}$$

3. else,  $X_i = \langle \text{standard-activation-function} \rangle$ . /\* Not part of a tree \*/

### 3.6. A complete activation function

Interleaving the three algorithms described earlier achieves the goal of identifying tree-like subnetworks and maximizes their goodness. In this subsection, we present the complete algorithm, combining the three phases while simplifying the computation. The algorithm is integrated with the discrete Hopfield

activation function; however it can be integrated also with other activation function (eg. Boltzmann machine).<sup>3</sup>

Let  $i$  be the executing unit,  $j$  a non-parent neighbor of  $i$ , and  $k$  the parent of  $i$ :

**Optimizing on Tree-like Subnetworks (unit  $i$ ):**

1. Initialization: If first time, then  $(\forall j) P_j^i = 0$ . /\*Clear pointers (needed only for cyclic nets)\*/
2. Tree directing: If there exists a single neighbor  $k$ , such that  $P_k^i = 0$ , then  $P_k^i = 1$ , and for all other neighbors  $j$ ,  $P_j^i = 0$ ; else, for all neighbors,  $P_j^i = 0$ .
3. Computing goodness values:  
 $G_i^0 = \max\{\sum_{j \in \text{child}(i)} G_j^0 P_j^i, \sum_{j \in \text{child}(i)} G_j^1 P_j^i + \theta_i\}$ .  
 $G_i^1 = \max\{\sum_{j \in \text{child}(i)} G_j^0 P_j^i, \sum_{j \in \text{child}(i)} (G_j^1 P_j^i + w_{i,j} P_j^j) + \theta_i\}$ .
4. Assigning activation values:  
 If at least two neighbors are not pointing to  $i$ , then  
 /\*use standard activation function (Hopfield) \*/  

$$X_i = \begin{cases} 1 & \text{if } \sum_j w_{i,j} X_j \geq -\theta_i, \\ 0 & \text{otherwise;} \end{cases}$$
 else, /\* Node in a tree (including root and leaves) \*/  

$$X_i = \begin{cases} 1 & \text{if } \sum_j ((G_j^1 - G_j^0) P_j^i + w_{i,j} X_j P_j^j) \geq -\theta_i, \\ 0 & \text{otherwise.} \end{cases}$$

For proof of convergence see the appendix.

### 3.7. An example

The example illustrated in Figure 3 demonstrates a case where a local minimum of the standard algorithms is avoided. Standard algorithms may enter such local minimum and stay in a stable state that is clearly wrong.

The example is a variation on a harmony network [Smolensky 86, page 259] and an example from [McClelland et al. 86, page 22]. The task of the network is to identify words from low-level line segments. Certain patterns of line segments excite units that represent characters, and certain patterns of characters excite units that represent words. The line strokes used to draw the characters are the input units: L1,..., L5. The units "N," "S," "A," and "T" represent characters. The units "able," "nose," "time," and "cart" represent words, and Hn, Hs, Ha, Ht, H1,...,H4 are hidden units required by the Harmony model. For example, given the line segments of the character S, unit L4 is activated (input), and this causes units Hs and "S" to be activated. Since "NOSE" is the only word that contains the character "S," both H2 and the unit "nose" are also activated and the word "NOSE" is identified.

The network has feedback cycles (symmetric weights) so that ambiguity among characters or line-segments may be resolved as a result of identifying a word. For example, assume that the line segments required to recognize the word "NOSE" appear, but the character "N" in the input is blurred and therefore the setting of unit L2 is ambiguous. Given the rest of the line segments (e.g., those of the character "S"), the network identifies the word "NOSE" and activates units "nose" and H2. This causes unit "N" to be activated and so are all of its line segments. Thus the ambiguity of L2 is resolved.

The network is indeed designed to have a global minimum when L2, Hn, "N," H2, and "nose" are all activated; however, standard connectionist algorithms may fall into a local minimum when all these

<sup>3</sup>Note how similar the new activation function is to the original Hopfield function.

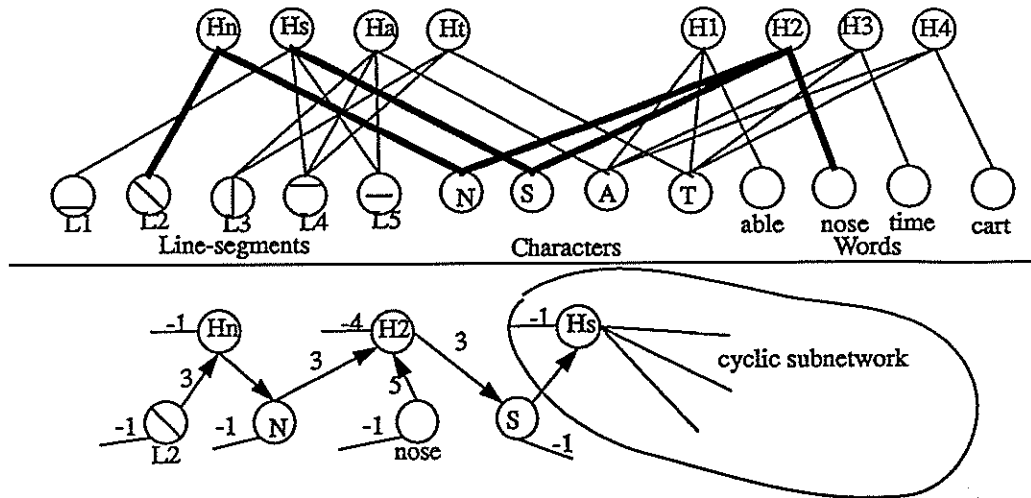


Figure 3: A Harmony network for recognizing words: Local minima along the subtrees are avoided

units are zero, generating goodness of  $5 - 4 = 1$ . The correct setting (global minimum) is found by our tree-optimization protocol (with goodness:  $3-1+3-1+3-1+5-1-4+3-1+5=13$ ). The thick arcs in the upper network of Figure 3 mark the arcs of a tree-like subnetwork. This tree-like subnetwork is drawn with pointers and weights in the lower part of the figure. Node "S" is not part of the tree and its activation value is set to one because the line-segments of "S" are activated. Once "S" is set, the units along the tree are optimized (by setting them all to one) and the local minimum is avoided.

#### 4. Limitations and extensions

We have shown a way to enhance the performance of connectionist energy minimization networks without losing much of the simplicity of the standard approaches. Our simple algorithm is limited in two ways, however. First, the central demon (or atomicity of the protocol) is not a realistic restriction. We would like the network to work correctly also under a *distributed demon*, where any subset of units may be scheduled for execution at the same time. Second, we would like the algorithm to be *self-stabilizing*. It should converge to a legal, stable state given enough time, even after noisy fluctuations that cause the units to execute an arbitrary program state and the registers to have arbitrary content.

##### 4.1. Distributed demon

Following [Dijkstra 74] and [Collin et al. 91], we show that no deterministic uniform algorithm exists that guarantees a stable global minimum under a distributed demon, even for simple chain-like trees.

Proof: Consider the network of figure 4. There are two global minima possible :  $(11\dots1101\dots11)$  and  $(11\dots1011\dots11)$ . If the network is initialized such that all units have the same register values, and all units start with the same program state, than there exists a fair execution under a distributed demon, such that in every step all units are activated. The units left of the center  $(1, 2, 3, \dots, i)$  "see" the same input as those units right of the center  $(2i, 2i-1, 2i-2, \dots, i+1)$  respectively. Because of the uniformity and the determinism, the units in each pair  $(i, i+1), (i-1, i+2), \dots, (1, 2i)$  must transfer to the same program state and produce the same output on the activation register.<sup>4</sup> Thus, after every step of that

<sup>4</sup>We assume that when execution begins, all units are initialized with same values and that all of them starts in the same program state.

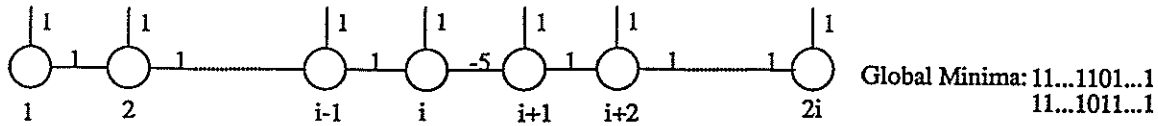


Figure 4: No uniform algorithm exists to optimize chains under distributed demon.

execution, units  $i$  and  $i+1$  will always have the same activation value and a global minimum (where the two units have different values) will never be obtained.

This negative result should not discourage us, since it relies on obscure infinite sequence of executions which is unlikely to occur under a truly random demon. Our algorithm will converge to a global minimum under a distributed demon in each of the following cases: (1) If step 2 of the protocol in section 3.6 is atomic; (2) if for every node  $i$  and every neighbor  $j$ , node  $i$  is executed without  $j$  infinitely often; (3) if one node is unique and acts as a root, that is, does not execute step 2 (an almost uniform protocol); and (4) if the network is cyclic (one node will be acting as a root).

Another negative results similar to [Collin et al. 91], is that if the network is cyclic, no uniform algorithm exists for its optimization, even under central demon. This may be proved even for cyclic networks simple as rings: In figure 5, we see a ring-like network whose global minima is (010101) and (101010). Consider a fair execution under central demon that activates the units 1,4,2,5,3,6 in order, and repeats this order indefinitely. Starting with the same program state and same inputs, the two units in every pair of (1,4), (2,5), (3,6) "see" the same input, and therefore have the same output and transfer to the same program state. As a result, these units never output different values and a global minimum is not obtained.

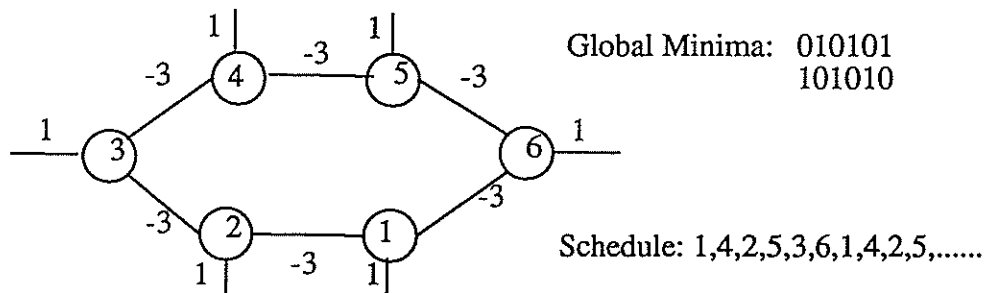


Figure 5: No uniform algorithm exists that guarantees to optimize rings even under central demon.

## 4.2. Self-stabilization

A protocol is self-stabilizing if in any fair execution, starting from any input configuration and any program state (of the units), the system reaches a valid stable configuration. The motivation here is that even if hardware fluctuations occur and change the content of the registers (as well as the state of the program), the algorithm is guaranteed to recover and to find a correct stable state.

The algorithm in section 3.6 is self-stabilizing for cycle free networks (trees), and it remains self-stabilizing under distributed demon with the same weak restrictions as in the previous section; i.e., executes without a neighbor infinitely often or is almost uniform (has a root). For proofs see the

appendix. However, the algorithm is not self-stabilizing for general cyclic topologies. For example, consider the configuration of the pointers in the ring of figure 6. It is in stable state, although clearly not a valid tree.<sup>5</sup> To solve the problem of self-stabilization in cyclic networks, we need to make our

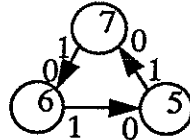


Figure 6: The simple protocol is not self-stabilizing in cyclic networks.

algorithm more complex. For example, we may use a variation of the self-stabilizing tree-directing protocol of [Collin et al. 91]. This algorithm remains self-stabilizing even in cyclic networks. Thus, self-stabilization may be obtained in the general case, but at the expense of more complexity and more space requirements.

## 5. Summary

We have shown a uniform self-stabilizing connectionist activation function that guarantees to find a global minimum of acyclic symmetric networks in linear time. The algorithm optimizes tree-like subnetworks in general (cyclic) networks. The simple version of the algorithm loses its self-stabilization property in cyclic networks; however, we can extend it to be self-stabilizing for all network topologies at the expense of more space requirements.

We proved two negative results: (1) Under a distributed demon, no *uniform* algorithm exists to optimize even simple chains, and (2) no uniform algorithm exists to optimize simple cyclic networks (rings) even under a central demon.

We conjecture that these negative results are not of significant practical importance, since in truly random scheduling demons the probability of having such pathological executions approaches zero. Our algorithm remains correct under a distributed demon if some weak assumptions are made on the demon or if we assume almost uniformity or atomicity.

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<sup>5</sup>Such configuration will never occur, if all units start at the *starting point*; i.e., clearing the bits of  $P_i$ . It may only happen due to some noise or hardware fluctuation.



## A. Convergence and Self-Stabilization of the Activation Function for Trees

To show: The algorithm converges and is self-stabilizing in networks with tree topology under a distributed demon with fair exclusion<sup>6</sup>. The algorithm also converges (but is not self-stabilizing) when the network is an arbitrary graph (e.g., has cycles) and is self-stabilizing for any topology if an almost uniform<sup>7</sup> version of the algorithm is applied. The almost uniform algorithm remains self-stabilized for unrestricted networks even under pure distributed demon<sup>8</sup>

*Proof: (sketch)*

The second and third phases of the algorithm are adaptations of an existing dynamic programming algorithm [Bertelé, Brioschi 72] and their correctness therefore are not proved here. The self-stabilization of these steps is obvious because no variables are initialized. The proof therefore concentrates on the tree directing phase.

Let us first assume that the scheduler is a distributed demon with fair exclusion and that the network is cycle-free (a tree). We want to show that the algorithm converges, it is self-stabilized and that the final stable result is that the pointers  $P_i^j$  represent a tree.

Define:

A node is *legal* if it either a root (i.e., all its neighbors are legal, point to it and it doesn't point to any of them), or an intermediate node (i.e., it points to one of the neighbors and the rest of its neighbors are all legal and point back).

A node is a *candidate* if it is illegal and have all its neighbors but one pointing to it and legal.

To show:

- 1) The property of being legal is stable; i.e., once a node becomes legal it will stay legal.
- 2) A state where the number of illegal nodes is  $k > 0$ , leads to a state where the number of illegal nodes is less than  $k$ ; i.e. the number of illegals decreases and eventually all nodes turn legal.
- 3) If all the nodes are legal then the graph is marked as a tree.
- 4) The algorithm is self-stabilized for trees.
- 5) The algorithm converge even if the graph has cycles.
- 6) The algorithm is self-stabilized in arbitrary networks if an almost uniform version is used (even under a distributed demon).

Proof:

1) To show that a legal state is stable. Assume a legal node  $i$  becomes illegal. It is either a root node which one of its children became illegal, or an intermediate node whose one of its children became illegal (it cannot be that its parent suddenly points to  $i$  or that one of the children stopped pointing and still is legal). Therefore, there must be a chain of  $i_1, i_2, \dots, i_k$  of nodes that became illegal. Since there are no cycles, there must be a leaf that was legal and turned illegal. This cannot be true since a leaf does not have children. Contradiction.

2) To show that if there are illegal nodes, their number is reduced. To prove this claim we need three steps: 2.1) Eventually a state is reached where if there is at least one illegal node then there is also a candidate node among the illegals. 2.2) This candidacy is stable. 2.3) eventually the candidate will become legal (therefore the number of illegals is reduced).

2.1) Because of the fair execution, eventually a state is reached where each node was executed at least once. Assume that at least one node is illegal and all the illegal nodes are not candidates. If a node is illegal and not candidate, then either it is a root-type (all point to it) but at least one of its children is illegal or there are at least two of its neighbors that are illegal. Suppose there are no root-type

<sup>6</sup>Fair exclusion is when for every node  $i$  and its neighbor  $j$ ,  $i$  is executed without  $j$  infinitely often.

<sup>7</sup>The roots are marked and the algorithm specifically handle them by constantly setting the pointers to zero.

<sup>8</sup>Any random subset of nodes may be chosen at one time for synchronous execution.

illegal nodes, then all illegal nodes have at least two illegal neighbors. Therefore there must be a cycle that connects illegal nodes (contradiction). Therefore, one of the illegals must be root-type. Suppose  $i$  is a root-type illegal node. It must have a neighbor  $j$  which is illegal. Consider the subtree of  $j$  that does not include  $i$ : it must contain illegal nodes, If there are no root-type illegals we get contradiction again; however, if there is a root-type node, we eliminate it and look at the subtree of some illegal  $j'$  that does not include  $j$ . Eventually, since the network is finite we obtain a subtree with no root-like illegals but which includes other illegal nodes. This leads to a contradiction. The conclusion is that there must be candidates if there are illegals.

2.2) To show that a candidate is stable unless it becomes legal:

If a node  $i$  is a candidate, all its legal children remain legal. There are three types of candidate nodes (node  $j$  is an illegal neighbor of  $i$ ): 1) Node  $j$  points to  $i$ ; 2) the pointer goes in both directions; 3) no pointer from  $i$  to  $j$  or wise-versa. All possible changes in the pointers  $P_j^i$  or  $P_i^j$  will cause  $i$  to remain a candidate or to turn legal (the rest of the pointers will not be changed).

2.3) To show that every candidate node will eventually turn legal: Assume  $j$  is the illegal neighbor of the candidate  $i$ . In the next execution of  $i$  without  $j$ , if  $P_j^i = 0$  then  $i$  becomes legal by pointing to  $j$ ; otherwise,  $i$  becomes a root-type candidate (all its neighbors point to it) but  $j$  is illegal. We'll prove now that if an illegal node  $j$  points to  $i$  then eventually a state is reached where either  $j$  is legal or  $P_j^i = 0$  and this proposition is stable once it holds. If this statement is true, then  $i$  is executed eventually: if  $j$  is legal then all of  $i$ 's neighbors are legal and therefor  $i$  turns legal. If  $j$  is illegal then  $P_j^i = 0$  and  $i$  will point to it ( $P_i^j = 1$ ) making itself legal.

To prove: if  $j$  is an illegal node pointing to  $i$  then there will be a state where either  $j$  is legal or  $P_j^i = 0$  and this state is stable.

By induction on the size of the subtree of  $j$  that does not include  $i$ .

Base: If  $j$  is a leaf and  $j$  points to  $i$  then if at the time  $j$  is executed (without  $i$ )  $P_j^i = 0$ , then node  $j$  points to  $i$  and turns legal; otherwise,  $j$  updates  $P_j^i = 0$ . This status is stable because the legal state is stable and since a leaf will point to a node only if it turns legal.

Induction step: Assume hypothesis is true for trees of size less than  $n$ . Suppose  $j$  the illegal neighbor of  $i$ . Node  $j$  points to  $i$  and it has  $j_1, \dots, j_k$  other neighbors. Because we assume that all nodes were executed at least one time, since  $j$  points to  $i$  we assume that at the last execution of  $j$  all the other neighbors  $j_1, \dots, j_k$  pointed to  $j$ . The subtrees rooted by  $j_l$  (not including  $j$ ) are of size  $n$  and therefore by hypothesis, there will be a state where all the nodes  $j_1, \dots, j_k$  are either legal or  $P_{j_l}^j = 0$ . This state is stable so when eventually  $j$  is executed it will either point to  $i$  turning legal (if all  $j_1, \dots, j_k$  are pointing to it), or it will make  $P_j^i = 0$  (if some of its neighbors do not point to it). Since the status of  $j_1, \dots, j_k$  is stable at that point, whenever  $j$  is executed it will either become legal or its pointers become zero.

3) To show that if all the nodes are legal then the graph is marked as a tree: If a node is legal, then all its children are legal and point to it. Therefore each node represents a subtree (if not a leaf) and have one parent at the most. To show that there is only one root: If several roots exist, then because of connectivity, there is one node that is shared between at least two subtrees and has therefore two parents (contradiction).

4) The algorithm is self-stabilizing for cycle-free networks since no initialization is needed (in the proof we haven't use the first initialization step; i.e.,  $P_i^i = 0$ ). In the case where no cycles exist, we do not need this step. The pointers can get any initial values and the algorithm still converges.

5) The algorithm (with  $P_i^i = 0$  initialization) converges even if the graph has cycles: Since all the nodes start with zero pointers, a (pseudo) root of a tree-like subnetwork will never point towards any of its neighbors (since it is part of a cycle and all of its neighbors but one must be legal).

6) To show that the algorithm is self-stabilizing in arbitrary networks if an almost uniform version is used, even under pure distributed demon: We need to show that a candidate will eventually turn legal even if its neighbors are executed in the same time.

Suppose node  $i$  is a candidate and node  $j$  is its illegal neighbor:

1) If  $j$  is a root, then it will never point to  $i$  and therefore  $i$  will eventually turn legal by pointing to  $j$ .

2) If  $i$  is the root, then  $P_i^j = 0$ , and if  $j$  becomes legal it will point to  $i$  making  $i$  legal. Node  $j$  will turn eventually legal using the following induction (on the size of the subtree of  $j$ ).

Hypothesis: In a subtree without a node that acts as a root, all illegal nodes will eventually turn legal.

Base: If  $j$  is a leaf, it will point eventually to its neighbor  $i$  which in its turn will make  $j$  legal by  $P_i^j = 0$ .

Induction: If  $j_1, \dots, j_k$  are other neighbors of  $j$ , then they will eventually turn legal (induction hypothesis) while pointing to  $j$ . Eventually  $j$  is executed and also turns legal.

3) Suppose, neither  $i$  nor  $j$  are roots, but suppose one of them is not part of a cycle (and therefore is part of a subtree that does not include a node marked as a root. Using the above induction, all the nodes in the subtree will eventually turn legal. As a result either  $i$  or  $j$  eventually turns legal and therefore  $i$  will eventually turn legal as well.

□

## References

- [Ballard et al. 86] D. H. Ballard, P. C. Gardner, M. A. Srinivas, "Graph problems and connectionist architectures," Department of Computer Science, University of Rochester, Technical Report 167, 1986.
- [Bertelé, Brioschi 72] U. Bertelé, F. Brioschi, *"Nonserial dynamic programming,"* Academic Press, 1972.
- [Collin et al. 91] Z. Collin, R. Dechter, S. Katz, "On the feasibility of distributed constraint satisfaction," In IJCAI-91: *Proceedings of 12th International Conference on Artificial Intelligence*, Sydney, Australia, 1991.
- [Dechter, Pearl 88] R. Dechter, J. Pearl, "Network-based heuristics for constraint-satisfaction problems," *Artificial Intelligence* 34, pp. 1-38, 1988.
- [Dechter et al. 90] R. Dechter, A. Dechter, J. Pearl, "Optimization in constraint networks," in R.M. Oliver and J.Q. Smith (editors), *Influence diagrams, belief nets and decision analysis*, John Wiley, 1990.
- [Dijkstra 74] E.W. Dijkstra, "Self-stabilizing systems in spite of distributed control," *Communications of the ACM* 17, pp. 643-644, 1974.
- [Feldman, Ballard 82] J.A. Feldman, D.H. Ballard, "Connectionist models and their properties," *Cognitive Science* 6, 1982.
- [Hinton, Sejnowski 86] G.E. Hinton, T.J. Sejnowski, "Learning and re-learning in Boltzman machines," in J. L. McClelland, D. E. Rumelhart, *Parallel Distributed Processing: Explorations in the Microstructure of Cognition I*, MIT Press, 1986.
- [Hopfield 82] J. J. Hopfield, "Neural networks and physical systems with emergent collective computational abilities," *Proceedings of the National Academy of Sciences* 79, pp. 2554-2558, 1982.
- [Hopfield 84] J. J. Hopfield, "Neurons with graded response have collective computational properties like those of two-state neurons," *Proceedings of the National Academy of Sciences*, pp. 3088-3092, 1984.

- [Hopfield, Tank 85] J.J. Hopfield, D.W. Tank, "Neural computation of decisions in optimization problems," *Biological Cybernetics* 52, pp. 144-152, 1985.
- [Kasif et al. 89] S. Kasif, S. Banerjee, A. Delcher, G. Sullivan, "Some results on the computational complexity of symmetric connectionist networks," Department of Computer Science, The John Hopkins University, Technical Report JHU/CS-89/10, 1989.
- [McClelland et al. 86, page 22] J. L. McClelland, D. E. Rumelhart, G.E Hinton, J.L. McClelland, "The appeal of PDP," in J. L. McClelland, D. E. Rumelhart, *Parallel Distributed Processing: Explorations in the Microstructure of Cognition I*, MIT Press, 1986.
- [Papadimitriou et al. 90] C. Papadimitriou, A. Shaffer, M. Yannakakis, "On the complexity of local search," in *ACM Symposium on the Theory of Computation*, pp. 438-445, 1990.
- [Peterson, Hartman 89] C. Peterson, E. Hartman, "Explorations of mean field theory learning algorithm," *Neural Networks* 2, no. 6, 1989.
- [Pinkas 91] G. Pinkas, "Energy minimization and the satisfiability of propositional calculus," *Neural Computation* 3, no. 2, 1991.
- [Pinkas 92] G. Pinkas, "Constructing proofs in symmetric networks," to appear in J. E Moody, S. j. Hanson, R. P. Lipmann (eds.), *Advances in Neural Information Processing Systems 4*, (NIPS-91), 1992.
- [Rumelhart et al. 86] D.E. Rumelhart, G.E Hinton, J.L. McClelland, "A general framework for parallel distributed processing," in J. L. McClelland, D. E. Rumelhart, *Parallel Distributed Processing: Explorations in the Microstructure of Cognition I*, MIT Press, 1986.
- [Smolensky 86, page 259] P. Smolensky, "Information processing in dynamic systems: Foundations of harmony theory," in J. L. McClelland, D. E. Rumelhart, *Parallel Distributed Processing: Explorations in the Microstructure of Cognition I*, MIT Press, 1986.