

# An Efficient Algorithm for the Detection of Eden

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In this paper, a polynomial time algorithm is presented for solving the Eden problem for graph cellular automata. The algorithm is based on our neighborhood elimination operation, which removes local neighborhood configurations that cannot be used in a preimage of the given configuration. This paper presents a detailed derivation of our algorithm from first principles, and a detailed complexity and accuracy analysis is also given. In the case of time complexity, it is shown that the average-case time complexity of the algorithm is  $\Theta(n^2)$ , and the best and worst cases are  $\Omega(n)$  and  $O(n^3)$ , respectively. This represents a vast improvement in the upper bound over current methods, without compromising average-case performance.

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## 1. Introduction

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Cellular automata and, more generally, discrete dynamical systems are powerful tools for modeling of complex phenomena [1]. This includes applications from physics, biology, and computer science [2]. Some have even speculated that the study of cellular automata may lead to a Grand Unified Theory of everything [3].

The study of the global dynamics of cellular automata (i.e., the study of automata configuration transition graphs) can provide unique insight into complex systems [4]. Efficient construction of a

configuration transition graph typically requires a method to determine if the given configuration is located on a leaf node of this graph [5].

This problem is known as the Eden problem, and has been shown to be computationally intractable for  $d$ -dimensional systems when  $d > 1$ . This is reflected in the worst-case computational complexity of algorithms that solve the Eden problem for higher dimensions (e.g., Wuensche's general reverse algorithm [6]).

We present a new algorithm for approximately solving the Eden problem for graph cellular automata (i.e., cellular automata on graphs [7, 8]), the most general form of deterministic cellular automata. Although there exist rare instances in which the algorithm will fail to identify the nonexistence of a preimage, this is made up for by its asymptotic complexity class, which is  $O(n^3)$  for the worst case and  $\Theta(n^2)$  for the average case. This provides a method that is more computationally feasible in the worst case than approaches based on Wuensche and Lesser's reverse algorithm [4] and Wuensche's general reverse algorithm [6] for the study of the global dynamics of higher-dimensional discrete dynamical systems with potentially a large number of cells.

## 2. Background

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### 2.1 Discrete Dynamical Systems

A regular cellular automaton can be defined as a lattice of finite state automata, typically referred to as *cells* or *sites*. A state transition function defines how a cell updates its state based on its current state and the state of its neighbors. Cells update synchronously in discrete time intervals. The sequence of all cell states at a given time is referred to as the automaton's *configuration*.

Random Boolean networks are binary cellular automata with one critical difference: there is no requirement that cells be located on a regular lattice [6]. Instead, neighborhoods are constructed via a random wiring. This random wiring makes random Boolean networks useful for theoretical biological models of genetic regulatory networks [9, 10].

Graph cellular automata (also referred to as generalized automata networks [11]) are a generalization of both cellular automata and random Boolean networks. For a graph cellular automaton, cell connectivity is defined by a connected graph. The class of graph cellular automata contains regular cellular automata and random Boolean networks as subclasses. Cellular automata and random Boolean networks can be considered as discrete dynamical systems. Despite their

simple construction, discrete dynamical systems have been shown to be capable of very complex behavior [12–14]. Furthermore, computationally intractable and formally undecidable problems relating to discrete dynamical systems have been shown to exist [15, 16].

## ■ 2.2 The Eden Problem

A particular problem of interest in the study of the global dynamics of discrete dynamical systems is the so-called Eden problem (also called the predecessor existence problem [17, 18]). The Eden problem attempts to determine, for a given automaton, if there exists a configuration (i.e., preimage) that will evolve to the given configuration in the next time step. If the Eden problem is resolved to be false, then the configuration is called a Garden of Eden configuration (i.e., it has no preimage). Wuensche and Lesser studied the Eden problem in depth and developed a reverse algorithm for one-dimensional regular cellular automata [4]. Wuensche further generalized this approach to the case of random Boolean networks, which may also be applied to graph cellular automata [6, 5]. While Wuensche and Lesser's method performs very well for small cellular automata, this method's upper bound is  $O(2^n)$  (as we will show in Section 5), which prevents exploration of large discrete dynamical systems.

For one-dimensional finite cellular automata, the Eden problem is in  $P$ ; however, for multi-dimensional finite cellular automata, the Eden problem has been shown to be  $NP$ -Complete [17]. Even certain variants of the Eden problem in one dimension (such as the constrained Eden problem [16]) have been shown to be  $NP$ -Complete. Assuming that  $P \neq NP$ , then there does not exist a polynomial time algorithm to solve the Eden problem for graph cellular automata.

If we assume  $P \neq NP$ , then a complete solution to the Eden problem for graph cellular automata is computationally intractable. However, this does not exclude the possibility of a good solution (i.e., one that can identify most Garden of Eden configurations) being achievable in polynomial time. In this paper, we present an algorithm that provides a good solution to the Eden problem for graph cellular automata in cubic time. By solving the problem for graph cellular automata we, by extension, solve the problem for regular cellular automata and random Boolean networks. Furthermore, we can show that our algorithm solves the Eden problem exactly when the topology of the graph cellular automaton is equivalent to a one-dimensional finite cellular automaton with periodic boundary conditions.

## ■ 2.3 Formal Definition of Graph Cellular Automata

In this section, we provide a formal definition of graph cellular automata. Our formalism is based heavily on the work of Fates [19], Marr et al. [7, 8], and Tomassini [11].

We consider a graph cellular automaton to be defined as a 4-tuple consisting of a connected graph, a set of states, a set of neighborhood mappings, and a set of state transition functions. This is given formally in Definition 1.

**Definition 1.** Let  $\mathbf{A} = (G, \Sigma, U, \Gamma)$  define a graph cellular automaton, where  $G = (V, E)$  is a graph with vertices  $V \subset \mathbb{Z}$  and edges  $E \subseteq V \times V$ ,  $\Sigma$  is a finite set of symbols referred to as the alphabet,  $U = \{b_i : i \in V\}$  is the set of neighborhoods  $b_i = \{i\} \cup \{j : (i, j) \in E \vee (j, i) \in E\}$ , and  $\Gamma = \{g_i : i \in V\}$  is the set of all state transition functions  $g_i : \Sigma^{|b_i|} \rightarrow \Sigma$ .

In Definition 1, the vertices of the graph  $G$  represent the cells of the automaton  $\mathbf{A}$ . Note that the neighborhood  $b_i$  of each cell  $i$  is effectively the set of cells that are connected to cell  $i$  via the set of edges  $E$ , including  $i$  itself. Note that the construction of  $b_i$  in Definition 1 assumes an undirected graph; the definition for a directed graph would be  $b_i = \{i\} \cup \{j : (j, i) \in E\}$ .

At any time  $t$ , each cell is associated with a state  $\sigma$ . For this we define the mapping in Definition 2. From this we can construct the global configuration of the automaton in Definition 3.

**Definition 2.** Let  $C : V \rightarrow \Sigma$  be a mapping from a cell  $i \in V$  to a state  $\sigma \in \Sigma$  such that  $C^t(i)$  represents the state of cell  $i$  at time  $t$ . Let  $C^t(b_i) \in \Sigma^{|b_i|}$  be the neighborhood configuration of  $i$ .

**Definition 3.** Let  $\phi^t = \{C^t(i) : i \in V\}$  be the configuration of the automaton  $\mathbf{A}$  at time  $t$ .  $\phi^t \in \Phi$ , where  $\Phi$  is the set of all possible configurations of  $\mathbf{A}$ .

Finally, we define the evolution of a graph cellular automaton as the sequence of configurations generated by repeated synchronous application of the local state transition functions. This is given as a recurrence relation expressed in terms of the global configuration transition function. This is given in Definition 4.

**Definition 4.** Let the recurrence relation  $\phi^{t+1} = f(\phi^t)$ ,  $t \geq 0$  be the evolution of  $\mathbf{A}$ , where  $f : \Phi \rightarrow \Phi$  is the global configuration transition function

$$f(\phi^t) = \{(\phi^t, \phi^{t+1}) : \phi^t = \{C^t(i) : i \in V\} \wedge \phi^{t+1} = \{g_i(C^t(b_i)) : i \in V\}\}.$$

We can now define formally an instance of the Eden problem.

**Definition 5.** Let EDEN denote the Eden problem, with instances consisting of a graph cellular automaton  $\mathbf{A}$  and configuration  $\phi \in \Sigma^{|V|}$ , and

with the question, does there exist an initial condition  $\phi^0$  such that  $\phi = f(\phi^0)$  under the evolution of  $\mathbf{A}$ ?

In Section 3, we will rely on the formalism given in this section to derive a polynomial time algorithm that provides the solution to  $\text{EDEN}(\mathbf{A}, \phi)$  in all but rare circumstances.

### 3. The Algorithm

In this section, we present a detailed derivation of our Eden detection algorithm, denoted by  $\text{EDEN-DET}(\mathbf{A}, \phi)$ . There are a number of steps involved in this derivation. First, some new mathematical constructions are defined. Then the fundamental operation of  $\text{EDEN-DET}(\mathbf{A}, \phi)$ , the neighborhood elimination operation, denoted by  $\text{NH-ELIM}(\mathbf{A}, H)$ , is derived. After presenting  $\text{NH-ELIM}(\mathbf{A}, H)$ , a simple Eden detection algorithm is provided, denoted by  $\text{S-EDEN-DET}(\mathbf{A}, \phi)$ . Using  $\text{S-EDEN-DET}(\mathbf{A}, \phi)$  as a starting point, we then derive a two-phase construction of  $\text{EDEN-DET}(\mathbf{A}, \phi)$ .

#### 3.1 Preliminaries

The graph cellular automata formalism given in Section 2.3 is not quite sufficient for us to express our Eden detection algorithm clearly. In this section, we present the definitions and notations that form the mathematical foundations of the algorithm. All definitions, notations, and theorems in this section assume the formalism in Section 2.3 to be given, and hence symbols used from Section 2.3 will not be redefined.

We will assume, without loss of generality, that  $\forall i \in V, |h_i| = k$ . This is done purely for notational convenience. All of the concepts applied in the construction of our algorithm can be extended trivially to nonuniform  $|h_i|$ . Note that we do not assume a uniform update rule across all cells  $\forall i, j \in V, g_i = g_j$ .

To describe our algorithm, we need a method of consistently referring to a specific neighborhood configuration (see Section 3.2). The notation for this reference is given in the following definition.

**Definition 6.** Assume that some ordering scheme has been applied to the set of all neighborhood configurations  $\Sigma^k$ . Subject to this ordering, the  $n^{\text{th}}$  neighborhood configuration is denoted by  $\psi_n \in \Sigma^k$ .

Note that the actual ordering scheme is arbitrary; all we require is an index into the possible neighborhood configuration space. For our

implementation, we simply map each configuration to its raw binary representation.

It is necessary for us to specify a set that contains all the cells that join adjacent neighborhoods. We refer to this set using the notation  ${}^i\Xi^j$ , and it is defined in Definition 7.

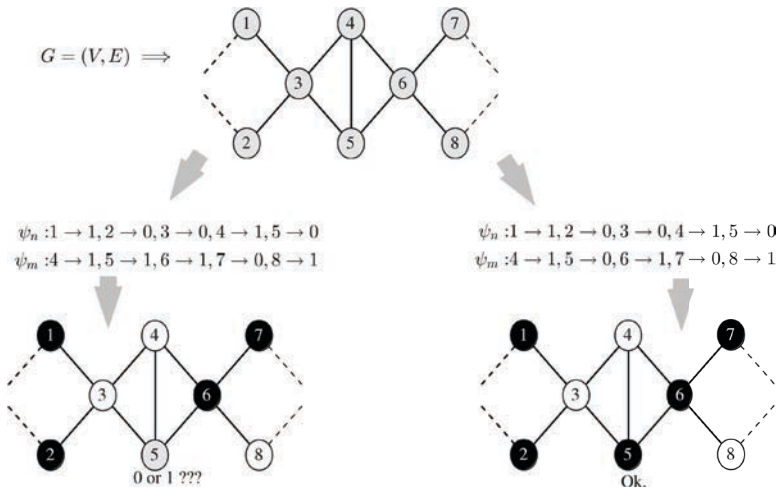
**Definition 7.** Let

${}^i\Xi^j = \{i\} \cup \{j\} \cup \{x : ((x, i) \in E \vee (i, x) \in E) \wedge ((x, j) \in E \vee (j, x) \in E)\}$  denote the boundary set of  $h_i$  and  $h_j$ . The  $n^{\text{th}}$  boundary cell,  $x \in V$ , is denoted by  $x = {}^i\Xi_n^j$ .

The basis of our algorithm is the detection and removal of neighborhood configurations that cannot exist in any preimage of  $\phi^t$  due to an inconsistency across boundary sets.

**Definition 8.** If there exists an initial configuration  $\phi^0$  such that  $C^0(h_i) = \psi_n$  and  $C^0(h_j) = \psi_m$ , then  $\psi_n$  is said to be  $i, j$ -consistent with respect to  $\psi_m$ .

The concept of  $i, j$ -consistency is readily visualized as shown in Figure 1. However, it would be preferable if a direct method of evaluating the  $i, j$ -consistency of two neighborhood configurations could be found. The function we require is given in Definition 9.



**Figure 1.** Example of  $i, j$ -consistency where  $i = 3$ ,  $j = 6$ , and  ${}^i\Xi^j = \{3, 4, 5, 6\}$ . Left:  $\psi_n$  is not  $i, j$ -consistent with respect to  $\psi_m$ , since they cause an inconsistent state in the boundary set (i.e., cell 5). Right: A modification to  $\psi_m$  allows consistency across the boundary set; hence,  $\psi_n$  is now  $i, j$ -consistent to  $\psi_m$ .

**Definition 9.** Let  $\theta_i^{i\Xi^j} : \Sigma^k \rightarrow \Sigma^{|i\Xi^j|}$  be a function that maps neighborhood configurations of  $h_i$  to the configuration of the boundary set  $i\Xi^j$ . The function is defined as

$$\theta_i^{i\Xi^j} = \left\{ (\psi_n, \psi'_n) : \psi_{n,s} = \psi_{n,q} \wedge y = h_{i,q} \wedge y = i\Xi^j_s \wedge s \in [0, i\Xi^j] \right\}.$$

The definition of  $\theta$  given in Definition 9 may seem strange, but it leads us into Theorem 1, which is a vital component of our algorithm.

**Theorem 1.** If  $\theta_i^{i\Xi^j}(\psi_n) = \theta_j^{i\Xi^j}(\psi_m)$ , then  $\psi_n$  is  $i,j$ -consistent with respect to  $\psi_m$ .

A formal proof is given in Appendix A.

### 3.2 Neighborhood Elimination

We can now formulate the core operation of our Eden detection algorithm. This core operation we call neighborhood elimination and denote it as  $\text{NH-ELIM}(\mathbf{A}, H)$ . As the name may suggest, its function is to eliminate neighborhood configurations that cannot be a component of any preimage of the automaton configuration in question.

To explain how we perform this operation, we first consider the matrix  $H \in \{0, 1\}^{|\Sigma|^k \times |V|}$ , where

$$H_{i,j} = \begin{cases} 0, & \text{impossible for } \psi_i = C^0(h_j) \\ 1, & \text{otherwise.} \end{cases} \tag{1}$$

It is important to note in equation (1) that  $H_{i,j} = 1$  should not be interpreted as  $\psi_i = C^0(h_j)$  in at least one preimage. Instead,  $H_{i,j} = 1$  means we cannot yet determine if  $\psi_i = C^0(h_j)$  or not. This is not the case for  $H_{i,j} = 0$ , which indicates that we have proven that there is no preimage such that  $\psi_i = C^0(h_j)$ .

The algorithm can be described as follows: Consider the case in which we have already determined that  $H_{i,j} = 0$  for specific  $i,j$  by the techniques described in Section 3.3. If we start with an arbitrary cell neighborhood  $h_i$ , then the column vector  $H_{*,i}$  provides us with the neighborhood configurations still under consideration. If  $H_{n,i} = 1$ , but the neighborhood configuration  $\psi_n$  is not  $i,j$ -consistent with respect to *any* candidate configurations in one or more connected neighborhoods  $h_j$ , then  $\psi_n$  can be excluded from the realm of possible configurations for  $h_i$ , as at least one boundary cell state cannot be satisfied consistently. By updating  $H_{*,i}$ , this will affect the validity of

other configurations, so we repeat the process for every neighborhood.

Theorem 1 provides us with a comparison operation for testing the  $i, j$ -consistency of two neighborhood configurations. With the function  $\theta_i^{j \exists!}$  as defined in Section 3.1, we arrive at  $\text{NH-ELIM}(\mathbf{A}, H)$  (i.e., Algorithm 1).

```

for all  $i \in V$  do
  for all  $j \in h_i - \{i\}$  do
     $\Theta_i \leftarrow \{x : x \in \theta_i^{j \exists!}(\psi_p) \wedge H_{p,i} = 1\}$ 
     $\Theta_j \leftarrow \{x : x \in \theta_i^{j \exists!}(\psi_q) \wedge H_{q,j} = 1\}$ 
     $\zeta_i \leftarrow \{x : x \in \Theta_i \wedge x \notin \Theta_j\}$ 
     $\forall p, H_{p,i} \leftarrow 0$  if  $\theta_i^{j \exists!}(\psi_p) \in \zeta_i$ 
  end for
end for

```

**Algorithm 1.**  $\text{NH-ELIM}(\mathbf{A}, H)$ : neighborhood elimination.

One step of  $\text{NH-ELIM}(\mathbf{A}, H)$  is shown in Figure 2, which displays contents of the data structures  $\Theta_i$ ,  $\Theta_j$ , and  $\zeta_i$ , along with the effect on the state of  $H$ . It should be noted that although the example in Figure 2 is for a small, one-dimensional cellular automaton with  $k = 3$ ,  $\text{NH-ELIM}(\mathbf{A}, H)$  is general enough to operate on graph cellular automata.

One particularly useful property of  $\text{NH-ELIM}(\mathbf{A}, H)$  is that the number of zero elements in  $H$  can never decrease. Therefore, repeating  $\text{NH-ELIM}(\mathbf{A}, H)$  on  $H$  in an iterative fashion will eventually result in an array  $H$  in which only configurations  $i, j$ -consistent with respect to all neighbors are candidates for preimage construction. This property also enables us to put an upper bound on the number of iterations required, which aids us in our complexity analysis (see Section 4).



$$\Sigma^3 = \{000, 001, 010, 011, 100, 101, 110, 111\}$$

$$H = \begin{matrix} & k & k+1 \\ \begin{bmatrix} \dots & 0 & 0 & \dots \\ \dots & 1 & 0 & \dots \\ \dots & 0 & 1 & \dots \\ \dots & 1 & 1 & \dots \\ \dots & 0 & 1 & \dots \\ \dots & 0 & 0 & \dots \\ \dots & 0 & 0 & \dots \\ \dots & 0 & 0 & \dots \end{bmatrix} & \begin{matrix} \Theta_i \leftarrow \{x : x \in \theta_i^{\exists'}(\psi_p) \wedge H_{p,i} = 1\} \\ \Theta_i = \{01, 11\} \\ \Theta_j \leftarrow \{x : x \in \theta_i^{\exists'}(\psi_q) \wedge H_{q,j} = 1\} \\ \Theta_j = \{01, 10\} \\ \zeta_i \leftarrow \{x : x \in \Theta_i \wedge x \notin \Theta_j\} \\ \zeta_i = \{11\} \end{matrix} \end{matrix}$$

$$H' = \begin{matrix} \begin{bmatrix} \dots & 0 & 0 & \dots \\ \dots & 1 & 0 & \dots \\ \dots & 0 & 1 & \dots \\ \dots & 0 & 1 & \dots \\ \dots & 0 & 1 & \dots \\ \dots & 0 & 0 & \dots \\ \dots & 0 & 0 & \dots \\ \dots & 0 & 0 & \dots \end{bmatrix} & \begin{matrix} H_{p,i} \leftarrow 0 \forall p, (\theta_i^{\exists'} \in \zeta_i) \\ H_{4,k} = 0 \end{matrix} \end{matrix}$$

**Figure 2.** One step of NH-ELIM(A,H). The upper-left matrix is H at the start of the iteration. After the iteration is completed, H<sub>4,k</sub> is set to 0, resulting in the lower-left matrix H'.

**3.3 Garden of Eden Detection**

In this section, we will describe our Eden detection algorithm (EDEN-DET(A,ϕ)) in full. Throughout this description we rely heavily on the formalism in Sections 2.3 and 3.1.

So far we have assumed that H is not all ones or all zeros, but we have not mentioned how H is initialized. If we are given an instance of EDEN(A,ϕ), we can prove the impossibility of some neighborhood configurations explicitly by using ϕ and the state transition functions g<sub>i</sub> ∈ Γ; that is,

$$H_{i,j} = \begin{cases} 0, & g_j(\psi_i) \neq \phi_j \\ 1, & g_j(\psi_i) = \phi_j. \end{cases} \tag{2}$$

In Section 3.2, it was stated for NH-ELIM(A,H) (Algorithm 1) that the number of zeros in H can never decrease. Therefore, repeated invoking of NH-ELIM(A,H) is guaranteed to converge to a steady state.

Once H is initialized, we can repeatedly operate the neighborhood elimination algorithm on H. Clearly, if for any column ∀ i, H<sub>i,j</sub> = 0 during an iteration, then there is no possible ψ<sub>i</sub> that can be selected for h<sub>j</sub> in any preimage. Furthermore, the steady state that H will con-

verge to in this case will be  $\forall i, j, H_{i,j} = 0$ . Therefore, we can conclude that  $\phi$  is a Garden of Eden configuration.

We might also assume that all Garden of Eden configurations will cause the condition  $\forall i, H_{i,j} = 0$ . Therefore, we could simply iterate until a steady state is reached and then look at the elements in  $H$  for any nonzero elements. This leads us to derive our initial algorithm for Eden detection, which we call simple Eden detection and denote as S-EDEN-DET( $\mathbf{A}, \phi$ ) (Algorithm 2).

```

 $\forall i, j, H_{i,j} \leftarrow 0$  if  $i, j, (g_j(\psi_i) \neq \phi_j)$ 
 $\forall i, j, H_{i,j} \leftarrow 1$  if  $i, j, (g_j(\psi_i) = \phi_j)$ 
while  $H \neq H'$  do
   $H' \leftarrow H$ 
   $H \leftarrow \text{NH-ELIM}(\mathbf{A}, H')$ 
end while
if  $\forall i, j, H_{i,j} = 0$  then
  GoE  $\leftarrow$  true
else
  GoE  $\leftarrow$  false
end if
return GoE

```

**Algorithm 2.** S-EDEN-DET( $\mathbf{A}, \phi$ ): simple Eden detection.

Unfortunately, S-EDEN-DET( $\mathbf{A}, \phi$ ) is not quite complete (hence the name “simple Eden detection”). It can be shown that  $\forall i, H_{i,j} = 0$  is a sufficient but not necessary condition of Eden. It is possible for cells within a cycle of  $G$  to have  $i, j$ -consistent neighbors, but there does not exist a combination of possible neighborhood configurations that can form a consistent chain. Figure 3 gives an example of such a case; note that for a one-dimensional cellular automaton the topology graph  $G$  contains one cycle that includes all cells. Clearly, more processing is required once S-EDEN-DET( $\mathbf{A}, \phi$ ) has converged, and there does not exist a  $j \in V$  such that  $\forall i, H_{i,j} = 0$ .

If for any possible neighborhood configuration (i.e.,  $H_{i,j} = 1$ ) we can construct at least one preimage, then we can conclude that  $\phi$  is not a Garden of Eden configuration. However, if it is found that a valid preimage cannot be constructed with  $C^0(b_j) = \psi_i$ , then we can set  $H_{i,j} = 0$  and repeat S-EDEN-DET( $\mathbf{A}, \phi$ ) until a new steady state is reached. This leads us to a second and more complete approach: EDEN-DET( $\mathbf{A}, H$ ).



```

while  $H \neq H'$  do
   $H' \leftarrow H$ 
   $H \leftarrow \text{NH-ELIM}(\mathbf{A}, H')$ 
end while

```

**Algorithm 3.** PH1( $\mathbf{A}, H$ ): Eden detection phase 1.

```

for all  $i \in V$  do
  for all  $j \in \Sigma^{|k|}$  do
    if  $H_{i,j} = 1$  then
       $H^{\text{tmp}} \leftarrow H$ 
       $\forall s, (s \neq j), H_{i,s}^{\text{tmp}} \leftarrow 0$ 
       $H^{\text{tmp}} \leftarrow \text{NH-ELIM}(\mathbf{A}, H^{\text{tmp}})$ 
      if  $H_{i,j}^{\text{tmp}} = 0$  then
         $H_{i,j} \leftarrow 0$ 
        return
      end if
    end if
  end for
end for

```

**Algorithm 4.** PH2( $\mathbf{A}, H$ ): Eden detection phase 2.

```

 $\forall i, j, H_{i,j} \leftarrow 0$  if  $i, j, (g_j(\psi_i) \neq \phi_j)$ 
 $\forall i, j, H_{i,j} \leftarrow 1$  if  $i, j, (g_j(\psi_i) = \phi_j)$ 
repeat
  CALL PH1( $\mathbf{A}, H$ )
  if  $\forall i, j, H_{i,j} = 0$  then
    GoE  $\leftarrow$  true
    return GoE
  else
    CALL PH2( $\mathbf{A}, H$ )
    GoE  $\leftarrow$  false
  end if
until  $H' = H$ 
return GoE

```

**Algorithm 5.** EDEN-DET( $\mathbf{A}, \phi$ ): Eden detection.

Leaving the details to Section 4, we simply claim that EDEN-DET( $\mathbf{A}, \phi$ ) is guaranteed to complete in polynomial time. More specifically, it can be shown to have a cubic worst-case time efficiency. Furthermore, when EDEN-DET( $\mathbf{A}, \phi$ ) returns GoE = false, then  $H$  encodes the complete set of preimages to  $\phi^t$  (except for rare cases when GoE = false is a false negative, as shown in Section 6).

#### 4. Time Complexity Analysis

In this section, we present the time complexity analysis for EDEN-DET( $\mathbf{A}, \phi$ ) (Algorithm 5). We show that the number of operations for the best case is a linear function of the number of cells; the worst case is shown to be cubic, and the average case is shown to be quadratic. Experimental results are also presented to reinforce theory with practice.

##### 4.1 Time Complexity of NH-ELIM( $\mathbf{A}, H$ )

The fundamental operation of EDEN-DET( $\mathbf{A}, \phi$ ) is clearly NH-ELIM( $\mathbf{A}, H$ ). From the pseudocode for NH-ELIM( $\mathbf{A}, H$ ) (Algorithm 1), it is also clear that the number of operations executed by NH-ELIM( $\mathbf{A}, H$ ) is a function of the number of cells  $n = |V|$ . We will show that this operation is in  $\Theta(n)$ .

The four lines within the innermost loop of NH-ELIM( $\mathbf{A}, H$ ) are only dependent on the number of neighborhood configurations. Without loss of generality, we assume  $\forall i, |b_i| = k$ ; thus the construction of  $\Theta_i$  and  $\Theta_j$  requires searching only a single column of  $H$ . That is,  $C_\Theta = c_0 |\Sigma^k|$ , where  $c_0 \approx k$  is the number of operations to evaluate  $\theta_i^{|\exists|}$ . The construction of  $\zeta_i$  is dependent only on the size of the  $\Theta$ ; hence,  $C_\zeta \leq C_\Theta$ . Furthermore, the number of elements in  $H$  is equal to the number of elements in  $\zeta_i \leq |\Sigma^k|$ . Thus the total operation count within the inner loop is given by

$$C_{\text{inner}} = 2 C_\Theta + C_\zeta + |\zeta| \approx 3 |\Sigma^k|. \tag{3}$$

Given equation (3), we can derive the total operation count for NH-ELIM( $\mathbf{A}, H$ ):

$$C_{NH}(n) = \sum_{i=1}^n \sum_{j=1}^k C_{\text{inner}} = k C_{\text{inner}} n \approx 3k |\Sigma^k| n. \tag{4}$$

Therefore  $C_{NH}(n) \in \Theta(n)$ .

**4.2 Best Case**

We now consider the best-case time complexity of EDEN-DET(A,ϕ). The best case occurs when there exist few possible *i,j*-consistent pairs for some subsequence in ϕ. This is very common in cellular automata in which Langton’s λ [14] is small. An example of this is when A is the elementary cellular automaton rule 2, and ϕ has a contiguous sequence of ones.

In this special case, only PH1(A,H) (Algorithm 3) will be required. Furthermore, a column of zeros will develop very quickly, as each iteration will eliminate at least one possible configuration from the unnatural area (due to few or no *i,j*-consistent neighborhood pairs); that is,  $I < |\Sigma^k|$ , where *I* is the number of iterations of the while loop. Using the results from equation (4), we have

$$C_{\text{best}}(n) = \sum_{i=1}^{|\Sigma^k|} C_{NH}(n) = |\Sigma^k| C_{NH}(n) \approx 3k |\Sigma^k|^2 n. \tag{5}$$

Therefore  $C_{\text{best}} \in \Omega(n)$ .

**4.3 Worst Case**

For the worst case, we must consider the full expression for the number of operations executed by EDEN-DET(A,ϕ). This is given by

$$C_{\text{Ops}}(n) = \sum_{t=1}^J \left( \underbrace{\sum_{i=1}^I C_{NH}(n)}_{\text{PH1(A,H)}} + \underbrace{\sum_{j=1}^{|V|} \sum_{i=1}^{|\Sigma^k|} C_{NH}(n)}_{\text{PH2(A,H)}} \right), \tag{6}$$

where *J* and *I* simply denote the number of iterations taken by the conditional loops. We require an upper bound on these loops.

In Section 3.3 we noted that the number of zeros in *H* can never decrease. Now we also note that if the number of zeros in *H* does not increase after an execution of PH2(A,H) (Algorithm 4), then EDEN-DET(A,ϕ) terminates with GoE = false. Hence for the worst case we must assume that the number of zeros decreases by exactly one. Furthermore, every iteration of PH1(A,H) will increase the number of zeros, terminate EDEN-DET(A,ϕ) with GoE = true, or continue to an iteration of PH2(A,H). Since  $H \in \{0, 1\}^{|\Sigma^k| \times n}$ , it must hold that

$$J(I + 1) \leq |\Sigma^k| n. \tag{7}$$

We want to maximize the value of *J*, as it has the greater effect on the total number of calls to NH-ELIM(A,H). If we assume the upper

bound is reachable, then as  $I \rightarrow 1$  we have  $J \rightarrow |\Sigma^k|/2n$ . We can apply this result to equation (6) to obtain an upper bound on  $C_{\text{ops}}(n)$ ,

$$C_{\text{ops}}(n) \leq \sum_{t=1}^{|\Sigma^k|/2n} \left( C_{NH}(n) + \sum_{j=1}^{|V|} \sum_{i=1}^{|\Sigma^k|} C_{NH}(n) \right) =$$

$$\frac{|\Sigma^k|}{2} n(C_{NH}(n) + |V| |\Sigma^k| C_{NH}(n)) = \frac{|\Sigma^k|}{2} (|\Sigma^k| n^2 + n) C_{NH}(n).$$

Furthermore, we have already shown that  $C_{NH} \in \Theta(n)$ . Therefore  $C_{\text{worst}} \in O(n^3)$ .

**4.4 Average Case**

Best- and worst-case bounds are important, but of limited practical use without an indication of the likelihood of EDEN(A,ϕ) instances that cause these bounds to occur. In this section we will show, using empirical data, that the average case is quadratic in time.

Consider equation (6): the values affecting the computational complexity are the number of iterations taken by the while loop of PH1(A,H), the number of iterations taken by the repeat-until loop, and the number of times PH2(A,H) needs to be executed. As in Section 4.3, we will denote the number of outer loops as  $J$  and the number of PH1(A,H) loops as  $I$ . Furthermore, we denote the number of iterations in which PH2(A,H) is executed as  $K$ .

We took random EDEN(A,ϕ) instances for  $|V| = n = 2^i, 2 \leq i \leq 13$ , where  $G$  is a single circuit. For each value of  $n$ , over 1000 samples were taken. The values of  $I, J$ , and  $K$  were counted for each sample. The expected values computed from these samples are shown in Figure 4.

From Figure 4 we can derive the overall expected values  $E(I) = 2.88, E(J) = 1.06$ , and  $E(K) = 0.25$ . So it is reasonable to approximate the average case as follows,

$$C_{\text{average}}(n) \approx \left( \sum_{i=1}^3 C_{NH}(n) \right) \times \Pr(\neg K) +$$

$$\left( \sum_{i=1}^3 C_{NH}(n) + \sum_{j=1}^{|V|} \sum_{i=1}^{|\Sigma^k|} C_{NH}(n) \right) \times \Pr(K) =$$

$$(3 C_{NH}(n)) \times \frac{3}{4} + (3 C_{NH}(n) + |\Sigma^k| n C_{NH}(n)) \times \frac{1}{4} = \frac{|\Sigma^k|}{4} n C_{NH}(n) + 3 C_{NH}(n).$$

Since  $C_{NH}(n) \in \Theta(n)$ , the approximate overall expected time complexity is  $C_{average}(n) \in \Theta(n^2)$ . Section 4.5 provides further experimentation to validate this approximation.

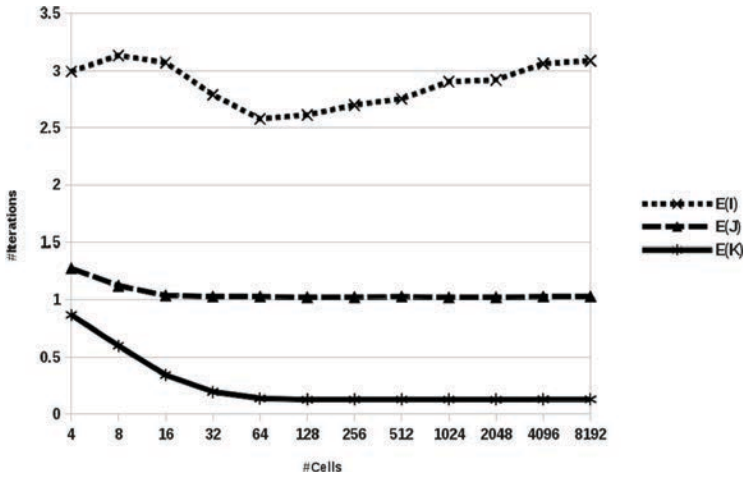


Figure 4. Expected iteration values. Based on 1000 random samples for each number of cells.

### 4.5 Experimental Results

For validation of the average case, we took a new random sample of 1000 instances of EDEN(A,ϕ) for  $|V| = n = 2^i, 2 \leq i \leq 13$ . For each sample, the average runtime of five separate runs was taken. Results were separated into two groups, based on whether EDEN-DET(A,ϕ) returned with GoE = true or GoE = false. The resulting average runtimes are shown in Figure 5.

Note that on average the runtime when GoE = false is approximately 16 times the runtime when GoE = true. This is because only a Garden of Eden configuration  $\phi_e$  can cause EDEN-DET(A,ϕ<sub>e</sub>) to return before Phase 2 is executed, which will complete in  $O(n)$  operations.

To confirm that the curves in Figure 5 are in fact quadratic, we can take the ratio  $R = C(2^{i+1}) / C(2^i)$ , where  $C(n)$  is the average runtime, as a function of the number of cells  $n$ . We would expect  $R \approx 4$  for a



quadratic (i.e., doubling the input takes four times longer). Figure 6 shows the  $R$  for the samples taken for Figure 5.

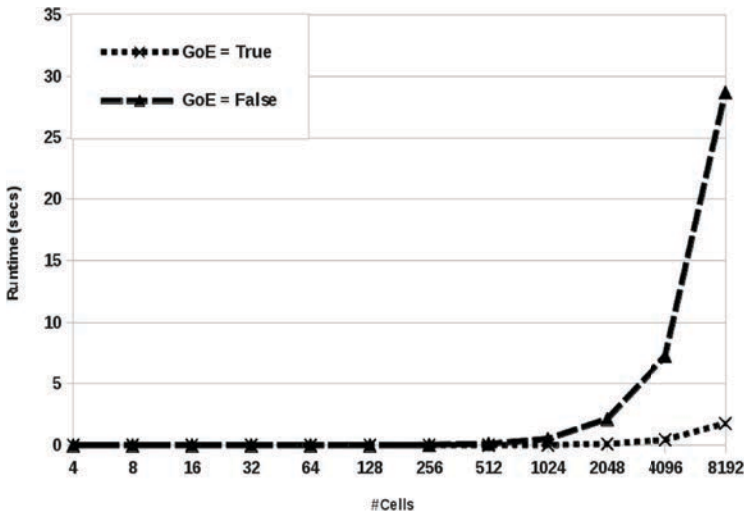


Figure 5. Runtimes for EDEN-DET (Algorithm 5).

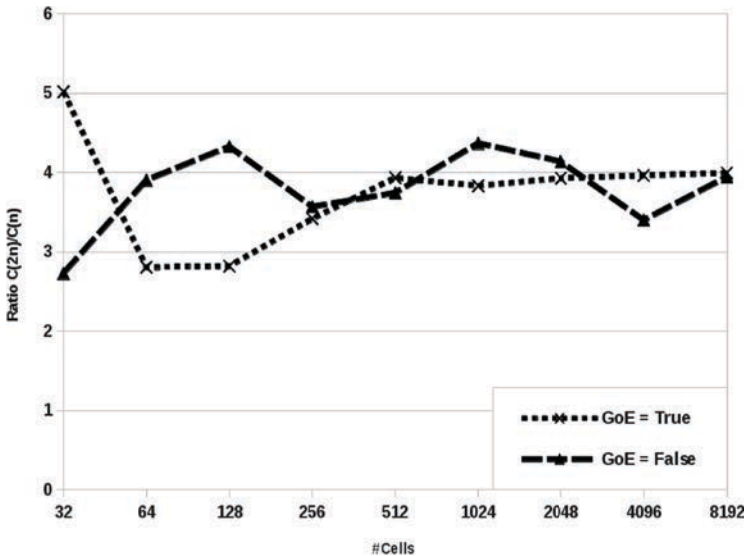


Figure 6. Ratio of runtimes  $C(2n) / C(n)$ .

From Figure 6 it is clear that  $R \approx 4$  (considering that  $R \approx 2$  for linear and  $R \approx 8$  for cubic). This provides support for our approximate average time complexity for EDEN-DET( $\mathbf{A}, \phi$ ) that we provided in Section 4.4.

## 5. Comparison with Wuensche and Lesser's Reverse Algorithm

In this section, we will compare the performance of EDEN-DET against the reverse algorithm developed by Wuensche and Lesser [4]. For the sake of simplicity, we will restrict the comparison to the simplest form of cellular automata: that of finite elementary cellular automata with periodic boundary conditions. As a result, we must emphasize that the following discussion and analysis relate specifically to Wuensche and Lesser's one-dimensional reverse algorithm [4] and not Wuensche's more general reverse algorithm, which applies to random Boolean networks and graph cellular automata [6, 5]. The results of this analysis, however, can certainly be generalized to the graph cellular automata case.

For a finite elementary cellular automaton with periodic boundary conditions  $\mathbf{A}$ , a configuration  $\phi^t$ , and a partial preimage  $\phi^{t-1}$  in which the first  $i$  cell states are known, Wuensche and Lesser's method is described as follows [4]:

1. If  $g(\phi_{i-1}^{t-1}, \phi_i^{t-1}, 0) = g(\phi_{i-1}^t - 1, \phi_i^t, 1) \neq \phi_i^t$ , then abandon the partial preimage. Resume derivation of next partial preimage (go to step 5).
2. If  $g(\phi_{i-1}^{t-1}, \phi_i^{t-1}, 0) \neq g(\phi_{i-1}^t - 1, \phi_i^t, 1)$ , then  $\phi_{i+1}^{t-1}$  can be uniquely determined. Proceed with next cell (go to step 1).
3. If  $g(\phi_{i-1}^{t-1}, \phi_i^{t-1}, 0) = g(\phi_{i-1}^t - 1, \phi_i^t, 1) = \phi_i^t$ , then  $\phi_{i+1}^{t-1}$  could be 0 or 1. Push the partial preimage  $(\phi_0^{t-1}, \phi_1^{t-1}, \dots, \phi_i^{t-1}, 1)$  onto the preimage queue to be processed later and continue with  $\phi_{i+1}^{t-1} = 0$ .
4. When  $i = n - 1$ , check that  $g(\phi_{n-2}^{t-1}, \phi_{n-1}^{t-1}, \phi_0^{t-1}) \neq g(\phi_{n-1}^{t-1}, \phi_0^{t-1}, \phi_1^{t-1})$ , then abandon this preimage; otherwise, add to the valid preimage list.
5. Take a new partial preimage from the queue and continue processing (step 1).
6. When the partial preimage queue is empty, all possible preimages starting with the start values of  $\phi_0^{t-1}, \phi_1^{t-1}$  are derived. Repeat for all possible  $\phi_0^{t-1}, \phi_1^{t-1}$ .

Note that the primary purpose of Wuensche and Lesser's method is the construction of all valid preimages, but it can be utilized directly to compute the solution to the Eden problem. Clearly, we can assert

GoE = false as soon as a valid preimage is found. We need not compute all of them. GoE = true will be asserted when no preimages are found.

### ■ 5.1 Worst-Case Complexity Analysis

With a brief description of Wuensche and Lesser's one-dimensional reverse algorithm, we can now show that the worst-case computation time is not bounded by a polynomial in the number of cells  $n$ . Consider Algorithm 6, which depicts Wuensche and Lesser's method modified for solving the Eden problem without computing all preimages.

```

GoE ← true
for all  $(p_1, p_2) \in \{(0, 0), (0, 1), (1, 0), (1, 1)\}$  do
   $\phi^{t-1} \leftarrow (p_1, p_2)$ 
   $Q \leftarrow \{\phi^{t-1}\}$ 
  while  $Q \neq \{\}$  do
     $\phi^{t-1} \leftarrow \text{pop}(Q)$ 
     $x = |\phi^{t-1}| - 1$ 
    for all  $i \in [x, n]$  do
       $T_0 \leftarrow g(\phi_{i-1}^{t-1}, \phi_i^{t-1}, 0)$ 
       $T_1 \leftarrow g(\phi_{i-1}^{t-1}, \phi_i^{t-1}, 1)$ 
      if  $T_0 = T_1 \neq \phi_i^t$  then
        break for loop
      else
        if  $T_0 \neq T_1$  then
          if  $T_0 = \phi_i^{t-1}$  then
             $\phi^{t-1} \leftarrow \phi^{t-1} \cup \{0\}$ 
          else
             $\phi^{t-1} \leftarrow \phi^{t-1} \cup \{1\}$ 
          end if
        else
          push( $Q, \phi^{t-1} \cup \{1\}$ )
           $\phi^{t-1} \leftarrow \phi^{t-1} \cup \{0\}$ 
        end if
      end if
    end for
  end while
   $T_n \leftarrow g(\phi_{n-1}^{t-1}, \phi_n^{t-1}, \phi_1^{t-1})$ 
   $T_1 \leftarrow g(\phi_n^{t-1}, \phi_1^{t-1}, \phi_2^{t-1})$ 
  if  $T_n = T_1$  then
    GoE ← false
    return GoE
  end if

```

```

end while
end for
return GoE

```

**Algorithm 6.** REVERSE( $\mathbf{A}, \phi$ ): Wuensche and Lesser's reverse algorithm.

Let  $C_{\text{inner}}$  denote the number of operations performed on a single iteration of the innermost loop in REVERSE( $\mathbf{A}, \phi$ ). Without loss of generality, we will assume  $C_{\text{inner}}$  is a constant. Clearly this is not true in reality, but instead  $3 \leq C_{\text{inner}} \leq 6$ .

Now, let  $C_x(n)$  denote the number of operations required to complete ignoring partial preimages already in  $\mathcal{Q}$  before reaching the  $x^{\text{th}}$  cell in the current preimage. We can express  $C_x(n)$  as the following recurrence relation,

$$C_x(n) = \sum_{i=x}^n C_{\text{inner}} + e(x) \sum_{i=x+1}^n a(i) C_i(n),$$

where  $e(x) = 0$  if  $T_0 = T_1 \neq \phi_x^{t-i}$ ; otherwise  $e(x) = 1$ , and  $a(x) = 1$  if  $T_0 = T_1 = \phi_x^{t-1}$ ; otherwise  $a(x) = 0$ .

The worst case for  $C_x(n)$  occurs when the number of partial preimages being pushed onto the queue is every iteration. In this case, we have  $\forall i > x, (a(i) = 1 \wedge e(i) = 1)$ , and the recurrence relation becomes

$$C_x(n) = \sum_{i=x}^n C_{\text{inner}} + \sum_{i=x+1}^n C_i(n).$$

We can now solve this recurrence relation. First consider expanding the  $C_{x+1}(n)$  term in the summation,

$$\begin{aligned} C_x(n) &= \sum_{i=x}^n C_{\text{inner}} + \sum_{i=x+1}^n C_i(n) = \\ &\sum_{i=x}^n C_{\text{inner}} + C_{x+1}(n) + \sum_{i=x+2}^n C_i(n) = C_{\text{inner}} + \\ &\sum_{i=x+1}^n C_{\text{inner}} + \left( \sum_{i=x+1}^n C_{\text{inner}} + \sum_{i=x+2}^n C_i(n) \right) + \sum_{i=x+2}^n C_i(n) = \\ &\sum_{i=x}^n C_{\text{inner}} + \left( \sum_{i=x+1}^n C_{\text{inner}} + \sum_{i=x+2}^n C_i(n) \right) + \sum_{i=x+2}^n C_i(n) = \end{aligned}$$

$$C_{\text{inner}} + 2 \sum_{i=x+1}^n C_{\text{inner}} + 2 \sum_{i=x+2}^n C_i(n).$$

Now, expanding the  $C_{x+2}(n)$  term,

$$\begin{aligned} C_x(n) &= C_{\text{inner}} + 2 \sum_{i=x+1}^n C_{\text{inner}} + 2 \sum_{i=x+2}^n C_i(n) = \\ &C_{\text{inner}} + 2 \sum_{i=x+1}^n C_{\text{inner}} + 2 C_{x+2}(n) + 2 \sum_{i=x+3}^n C_i(n) = \\ &C_{\text{inner}} + 2 C_{\text{inner}} + 2 \sum_{i=x+2}^n C_{\text{inner}} + \\ &2 \left( \sum_{i=x+2}^n C_{\text{inner}} + \sum_{x+3}^m C_i(n) \right) + 2 \sum_{i=x+3}^n C_i(n) = \\ &C_{\text{inner}} + 2 C_{\text{inner}} + 4 \sum_{i=x+2}^n C_{\text{inner}} + 4 \sum_{i=x+3}^n C_i(n). \end{aligned}$$

Repeating this process yields

$$\begin{aligned} C_x(n) &= C_{\text{inner}} + 2 C_{\text{inner}} + 4 C_{\text{inner}} + \dots + 2^{n-x} C_{\text{inner}} = \\ &C_{\text{inner}} \sum_{i=x}^n 2^{i-x}. \end{aligned}$$

The worst case for  $\text{REVERSE}(\mathbf{A}, \phi)$  requires that  $C_2(n)$  operations be executed four times,

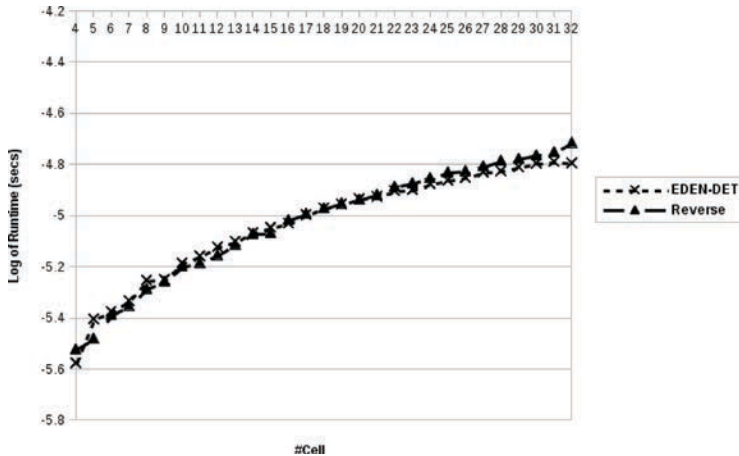
$$C_{\text{worst}} = 4 C_2(n) = 4 C_{\text{inner}} \sum_{i=2}^n 2^{i-2} = C_{\text{inner}} \sum_{i=2}^n 2^i,$$

hence  $\text{REVERSE}(\mathbf{A}, \phi)$  is in  $O(2^n)$ . It is worth noting that this worst case can only be achieved if the  $\phi$  is a Garden of Eden configuration, and the cell that determines this is the  $n^{\text{th}}$  cell. For example,  $\phi = (0, 0, \dots, 0, 1, 1)$  for the elementary cellular automaton rule 2. However, according to Wuensche and Lesser [4] the average case is orders of magnitude better. We confirm this experimentally in Section 5.2.

### 5.2 Experimental Comparison

We benchmarked  $\text{EDEN-DET}(\mathbf{A}, \phi)$  against  $\text{REVERSE}(\mathbf{A}, \phi)$ . Each experiment consisted of solving the Eden problem for 1000 random configu-

rations. Experiments were performed for both EDEN-DET( $\mathbf{A},\phi$ ) and REVERSE( $\mathbf{A},\phi$ ), using all the elementary cellular automata with cell counts ranging from 4 to 32. As shown in Figure 7, the benchmark average case is effectively the same order of magnitude for both methods.



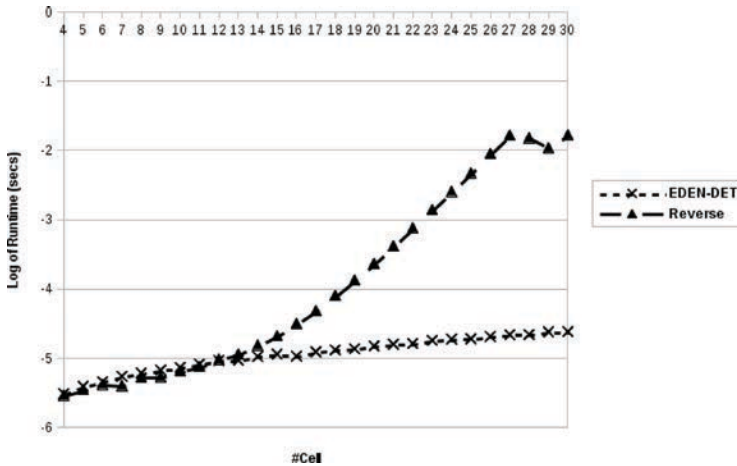
**Figure 7.** Comparison of EDEN-DET( $\mathbf{A},\phi$ ) against REVERSE( $\mathbf{A},\phi$ ), using a random sampling of configurations.

The worst case for REVERSE( $\mathbf{A},\phi$ ) is only approached for Garden of Eden configurations that are nearly identical to a non-Garden of Eden configuration, only differing in the last few cells. This is more likely to be possible with sparse configurations (i.e., very few 1 states compared with 0 states). If we restrict the random sample of test configurations to that of sparse configuration, then the probability of selecting a configuration that degrades the performance of REVERSE( $\mathbf{A},\phi$ ) increases.

Figure 8 indicates that the benchmark results are very different when we restrict the configuration sample this way. Such cases place a limitation on the usability of REVERSE( $\mathbf{A},\phi$ ) for large cell counts (as the cell count increases, any configuration with a relatively small sparse subsequence could render the Eden problem computationally intractable for REVERSE( $\mathbf{A},\phi$ )). The performance of EDEN-DET( $\mathbf{A},\phi$ ), however, is hardly affected by such sparse configurations.

The main difference in our approach, which provides such a large improvement in the worst-case performance, is the neighborhood elimination step. This operation performance is not affected by shifts (or rotations in a higher dimension) in the same configuration, because it treats each cell neighborhood independently of each other. As

a result,  $\text{EDEN-DET}(\mathbf{A}, \phi)$  provides a solution to the Eden problem that is scalable to very large cellular automata.  $\text{EDEN-DET}(\mathbf{A}, \phi)$  could be considered as a more stable alternative to Wuesnche and Lesser's  $\text{REVERSE}(\mathbf{A}, \phi)$ , as the worst case is vastly improved without degrading the average case.



**Figure 8.** Comparison of  $\text{EDEN-DET}(\mathbf{A}, \phi)$  against  $\text{REVERSE}(\mathbf{A}, \phi)$ , using a random sampling of sparse configurations.

## 6. Algorithm Correctness

In this section, we discuss the correctness of  $\text{EDEN-DET}(\mathbf{A}, \phi)$  in solving the Eden problem for graph cellular automata. We are able to show that  $\text{EDEN-DET}(\mathbf{A}, \phi)$  is completely correct for graphs with a single cycle. For graphs with more than one cycle, it is possible for incorrect results to be returned (i.e., false negatives); however, we show that these cases are rare.

It is first worth discussing the correctness of the solution when  $\text{EDEN-DET}(\mathbf{A}, \phi)$  returns with  $\text{GoE} = \text{true}$ . This result will never occur if  $\phi$  has a preimage (i.e., false positives cannot occur). This is because elements in  $H$  are only ever set to 0 when there is no  $i, j$ -consistent pair in a neighbor cell. If  $\text{GoE} = \text{true}$  is returned, then at some point there must have existed an  $i$  such that  $\forall j, H_{j,i} = 0$  (i.e., a cell has no possible  $i, j$ -consistent neighborhood configurations). For  $\phi$  to have a preimage, each cell must have at least one  $i, j$ -consistent neighborhood configuration. Therefore, only a true Garden of Eden configuration can cause  $\text{GoE} = \text{true}$  to be returned.

When  $\text{EDEN-DET}(\mathbf{A},\phi)$  returns with  $\text{GoE} = \text{false}$ , there are possible false identifications. That is, it is possible for a Garden of Eden configuration to cause  $\text{GoE} = \text{false}$  to be returned. However, this rare case is only a possibility when the graph  $G$  has more than one cycle.

We will now show that for  $\text{EDEN-DET}(\mathbf{A},\phi)$  to return  $\text{GoE} = \text{false}$  is always correct if  $G$  contains only one cycle. Consider  $H$ , which has reached a nonzero steady state after executing  $\text{PH1}(\mathbf{A},H)$ . If we assume a neighborhood configuration  $H_{j,i} = 1$  (see Section 3.3) and carry out an iteration of  $\text{NH-ELIM}(\mathbf{A},H)$ , we are essentially propagating the assumption around the cycle of  $G$ . When this propagation returns to  $i$ , then there are only two possibilities: (1) the assumed  $H_{j,i}$  is not eliminated, meaning a chain of  $i,j$ -consistent pairs can be constructed (i.e., a preimage can exist under this assumption); and (2) the assumed  $H_{j,i}$  is eliminated; hence,  $\psi_j$  cannot contribute to any preimage. Since  $\text{EDEN-DET}(\mathbf{A},\phi)$  only returns  $\text{GoE} = \text{false}$  when every element in  $H$  has passed assumption testing, we can conclude that this can only occur if  $\phi$  does in fact have a possible preimage. Therefore  $\text{EDEN-DET}(\mathbf{A},\phi)$  is completely correct for  $G$  with a single cycle.

These correctness results for the single-cycle (i.e., one-dimensional) case have also been supported by experimental results. We executed  $\text{EDEN-DET}(\mathbf{A},\phi)$  on the entire configuration space for all elementary cellular automata where  $n = \{4, 8, 16\}$ . Each return value was validated via a brute-force search for a preimage. This resulted in a 100% success rate.

Unfortunately, things are not so easy for  $G$  with multiple cycles. The assumption-testing method we apply in  $\text{PH2}(\mathbf{A},H)$  is really only powerful enough to test consistency within a single cycle. It may be possible for every  $H_{j,i}$  to pass the assumption test, but any choice made from one cycle breaks consistency in another. Hence a complete solution would require looking at pairs of cycles, triples of cycles, and so forth (we do not have a rigorous proof of this). This is likely the result of the  $NP$ -complete nature of the Eden problem in more than one dimension.

Again, we look to empirical data to show that in the majority of cases the single cycle accuracy is all we need. This time, over 170 000 random instances of  $\text{EDEN}(\mathbf{A},\phi)$  (for a fixed choice of rules representing Wolfram classes 1, 2, and 3 [13]) were taken as inputs. The topology of the graph  $G$  was equivalent to a dodecahedron. Since  $|V| = 20$ , the complete configuration space is  $2^{20}$ . Every result was compared to a brute-force approach.

We found that for class 1 cellular automata (i.e., point attractors), no false negatives ever seem to occur. Rules that fall under class 2 (i.e., simple structures, maybe periodic) had a low number of false neg-



atives, around 0.01%. Class 3 cellular automata (i.e., chaotic) are a different story: around 16% of cases in which  $\text{EDEN-DET}(\mathbf{A},\phi)$  returned  $\text{GoE} = \text{false}$  were incorrect. Over all samples, the false negative rate was around 10%.

It is interesting to note that it is only class 3 cellular automata that seem to cause  $\text{PH2}(\mathbf{A},H)$  to be executed in the one-dimensional case.

False negatives can be detected without resorting to a brute-force sweep. As previously stated in Section 3.3, the final state of  $H$  completely encodes all possible preimages. Neighborhoods in  $H$  can be stitched together using a method similar to Wuensche’s general reverse algorithm [6]; if no preimage can be constructed, then we have detected a false negative. In light of this, our algorithm could also be considered as a search-reduction step to be used prior to invoking Wuensche’s general method. Combined, these would provide a completely correct and more efficient method for constructing configuration transition graphs.

## 7. Conclusion

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In this paper we have presented an efficient algorithm (i.e., average case in  $\Theta(n^2)$ ),  $\text{EDEN-DET}(\mathbf{A},H)$ , for solving the Eden problem for graph cellular automata. By changing the topology of the graph  $G$ , the Eden problem can be solved for all classes of deterministic discrete dynamical systems (e.g., regular cellular automata and random Boolean networks). This analysis provides a firm foundation for further study of the global dynamics of discrete dynamical systems.

## Appendix

### A. Proof of Theorem 1

---

First consider the equality

$$\theta_i^{\exists^j}(\psi_n) = \theta_j^{\exists^i}(\psi_m).$$

Given Definition 9, we can expand the above expression. This yields

$$\begin{aligned} \equiv \forall s, \left( \psi'_{n,s} = \psi'_{m,s} \wedge \exists p, \left( \psi_{n,p} = \psi'_{n,s} \wedge y = h_{i,p} \wedge y = {}^i \Xi_s^j \right) \wedge \right. \\ \left. \exists q, \left( \psi_{m,q} = \psi'_{m,s} \wedge z = h_{j,q} \wedge z = {}^i \Xi_s^j \right) \right). \end{aligned}$$

This can be reduced using predicate calculus,

$$\begin{aligned} &\equiv \forall s, \left( \psi'_{n,s} = \psi'_{m,s} \wedge \exists p, \left( \psi_{n,p} = \psi'_{n,s} \wedge h_{i,p} = {}^i \Xi_s^j \right) \wedge \right. \\ &\quad \left. \exists q, \left( \psi_{m,q} = \psi'_{m,s} \wedge h_{j,q} = {}^i \Xi_s^j \right) \right) \equiv \\ &\quad \forall s, \left( \psi'_{n,s} = \psi'_{m,s} \wedge \exists p, q \left( \psi_{n,p} = \psi'_{n,s} \wedge h_{i,p} = \right. \right. \\ &\quad \quad \left. \left. {}^i \Xi_s^j \wedge \psi_{m,q} = \psi'_{m,s} \wedge h_{j,q} = {}^i \Xi_s^j \right) \right) \equiv \\ &\quad \forall s, \left( \exists p, q \left( \psi'_{n,s} = \psi'_{m,s} \wedge \psi_{n,p} = \psi'_{n,s} \wedge h_{i,p} = \right. \right. \\ &\quad \quad \left. \left. {}^i \Xi_s^j \wedge \psi_{m,q} = \psi'_{m,s} \wedge h_{j,q} = {}^i \Xi_s^j \right) \right) \equiv \forall s, \\ &\quad \left( \exists p, q \left( \psi_{n,p} = \psi_{m,q} \wedge h_{i,p} = {}^i \Xi_s^j \wedge h_{j,q} = {}^i \Xi_s^j \right) \right). \end{aligned}$$

Now let  $C^0(h_i) = \psi_n$  and  $C^0(h_j) = \psi_m$ ; hence,  $C^0(h_{i,p}) = \psi_{n,p}$  and  $C^0(h_{j,q}) = \psi_{m,q}$ :

$$\begin{aligned} &\equiv \forall s, \left( \exists p, q \left( \psi_{n,p} = \psi_{m,q} \wedge h_{i,p} = {}^i \Xi_s^j \wedge h_{j,q} = {}^i \Xi_s^j \right) \wedge C^0(h_{i,p}) = \right. \\ &\quad \left. \psi_{n,p} \wedge C^0(h_{j,q}) = \psi_{m,q} \right) \equiv \\ &\quad \forall s, \left( \exists p, q \left( h_{i,p} = {}^i \Xi_s^j \wedge h_{j,q} = {}^i \Xi_s^j \right) \wedge C^0(h_{i,p}) = C^0(h_{j,q}) \right). \end{aligned}$$

This satisfies our definition of  $i,j$ -consistency (i.e., Definition 8). Therefore  $\psi_n$  and  $\psi_m$  are  $i,j$ -consistent.  $\square$

## References

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