Coherent Functions and Program Checkers *

(extended abstract)

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ABSTRACT

Functions whose value at one point can be interpolated from values at other points have occupied a special place in cryptography. In this paper we will develop a general concept of "coherence" to capture this property and show its applications to several models of program checking. The concept and techniques also have applications in Hellman's timememory tradeoff for inverting random functions.

1 Introduction

Functions with the property that the value at one point can be interpolated from values at other points have occupied a special place in cryptography. Concepts like "self-reducibility" (Angluin and Lichtenstein [AL], Abadi, Feigenbaum, and Kilian [AFK]), "randomly testable" (Lipton [L]), and other related notions (Blum, Luby, and Rubinfeld [BLR]) which were raised for various purposes have captured some aspects of this quality. In this paper we will develop a general concept of "coherence", which is in some sense the minimum requirement for a function to have the property mentioned above, and show its applications

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to program checking and other subjects.

In Section 2, we illustrate the concept of coherence by considering a simple puzzle. In Section 3 a direct application of this concept is given for a time-memory tradeoff problem of Hellman [H] for inverting one-way functions. In Sections 4 and 5, we extend and develop the coherence concept to Turing machine models and comparison tree models, and show the connection with program checking (Blum and Kannan [BK]); the connection gives an approach to prove that certain computations do not have efficient program checkers. Section 6 concludes with some open problems.

2 A Puzzle

Let N, m be positive integers. Consider the following game to be played by A and B. There are N boxes with lids $BOX_1, BOX_2, \dots, BOX_N$ each containing a boolean bit. In the preprocessing stage, Player A will inspect the bits and take notes using an m-bit pad. Afterwards, Player B will ask Player A a question of the form "What is in BOX_i ?". Before answering the question, Player A is allowed to consult the m-bit pad and take off the lids of an adaptively chosen sequence of boxes not including BOX_i . The puzzle is, what is the minimum number of boxes A needs to examine in order to find the answer?

One strategy that A may adopt is the following. Divide the boxes into m consecutive groups each containing no more than $\lceil N/m \rceil$ members. In the preprocessing stage, A records for each $1 \le j \le m$ a bit

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 a_j which is the parity of all the bits in the j-th group. To answer a query "what is in BOX_i ", A can lift the lids of all the boxes (except BOX_i) in the group containing BOX_i , say group j, and compute the parity b of these bits; clearly, the bit in BOX_i must be equal to $a_j \oplus b$. With this strategy, A never needs to lift more than $\lceil N/m \rceil - 1$ lids. Is this the best possible strategy?

Let us define the model precisely. Let F_N be the set of all functions $f: \{1, 2, \dots, N\} \rightarrow \{0, 1\}$. For convenience we represent each of the 2^N possible bit patterns of the boxes as a function $f \in F_N$; f(i) = 1 if and only if BOX_i contains 1. An algorithm G consists of a partition $\{D_{\alpha} \mid \alpha \in \{0,1\}^m\}$ of F_N and a collection of $N2^m$ decision trees $T_{\alpha,i}$, $\alpha \in \{0,1\}^m, 1 \leq i \leq N$. Each $T_{\alpha,i}$ is a binary tree, in which every internal node is associated with a query $BOX_k = ?$ $(k \neq i)$ and has its two outgoing edges labeled by 0 and 1; each leaf contains a boolean output bit. We further require $T_{\alpha,i}$ to yield the correct output when it is used in the standard way to answer Player B's question "what is in BOX_i ?" when α is the m-bit string recorded on the pad. Define the cost of the algorithm, cost(G), to be the maximum height of all the trees $T_{\alpha,i}$. Let $s(m,N) = \min_G \operatorname{cost}(G)$. The following theorem shows that the strategy described earlier is an optimal one.

Theorem 1 Any algorithm must have cost at least $\lceil N/m \rceil - 1$.

Corollary $s(m, N) = \lceil N/m \rceil - 1$.

Proof Without loss of generality we can assume that $1 \leq m < N$. Consider any algorithm $\{D_{\alpha}, T_{\alpha,i}\}$. We will prove that for each $\alpha \in \{0, 1\}^m$

$$|D_{\alpha}| \le 2^{Nt/(t+1)},\tag{1}$$

where t is the cost of the algorithm.

Let $\ell = \lfloor N/(t+1) \rfloor t + \max(0, N - \lfloor N/(t+1) \rfloor (t+1) - 1)$. It is easy to verify that $\ell \leq Nt/(t+1)$. To prove (1), we only need to demonstrate that every $f \in D_{\alpha}$ can be uniquely specified by ℓ bits. Set $J \leftarrow \emptyset$. Repeat the following step $\lfloor N/(t+1) \rfloor$ times: take the smallest $i \notin J$, use the tree $T_{\alpha,i}$ to ask exactly t

queries f(j) = ? where $j \notin J \cup \{i\}$ and find the value of f(i), then enlarge J to include i and all these j's. At this point $t \cdot \lfloor N/(t+1) \rfloor$ queries have been answered and at least $(t+1) \cdot \lfloor N/(t+1) \rfloor$ values of f(j) have been determined; it takes at most $\max(0, N - (t+1) \lfloor N/(t+1) \rfloor - 1)$ more queries to find out the rest of the values of f(j). Since the answer to the queries determines f completely, ℓ bits suffice to specify f. This proves (1).

It follows from (1) that

$$2^N \le \sum_{\alpha} |D_{\alpha}| \le 2^m \cdot 2^{Nt/(t+1)},$$

and hence $N/m \le t+1$. This proves Theorem 1. The corollary follows immediately. \square

Let us view the above discussion more abstractly. Call a function f coherent if, when given some hints (m bits) about f, one can interpolate its value at any w from the value of f at other points efficiently (t points). There is an implicit assumption that one can discuss how to interpret hints before f is chosen. (See Beaver and Feigenbaum [BeaF] for a discussion of this assumption in a different context.) Which functions are coherent depend not only on m, t, but also the interpretation agreed upon.

We will extend this concept to the Turing machine model in Section 4. A function f will be called coherent if, when given some hints (a finite number of bits to specify a machine S), one can interpolate any f(w) from the value of f at other points efficiently (using S). In this case, the different possible interpretations of hints correspond to the different ways of specifying a universal Turing machine. The new ingredients to be added are the Turing-uniformity condition imposed on algorithms and the probabilistic setting needed for application purposes. Before doing that, we first discuss in the next section a question for which the simple model defined in the present section can be directly applied.

3 Inverting Random Functions

In the context of cryptanalysis, Hellman [H] developed a method for finding the inverse of any func-

tion. We will be interested in the question whether his method is the best possible among procedures of similar nature.

To explain Hellman's method, suppose we are given a one-cycle permutation $f: \{1, 2, \dots, N\} \rightarrow$ $\{1, 2, \dots, N\}$ as a "blackbox"; that is, the only way to get information about f is to feed the blackbox inputs x and get the value f(x) as outputs. The goal is to preprocess the information about f so as to be able to find the inverse $f^{-1}(y)$ fast for inputs y. His strategy is as follows. Write $\{1, 2, \dots, N\}$ as $\{x_0, x_1, \dots, x_{N-1}\}$ such that $f(x_i) = x_{i+1 \mod N}$. In the preprocessing stage, we experiment with the blackbox and store the sequence $x_{jN/m}$, $0 \le j \le m-1$ using $m \log_2 N$ bits. Afterwards, if we are given y, we can find $f^{-1}(y)$ by first computing $y_1 = f(y), y_2 =$ $f(y_1), \cdots$ until we recognize some y_r to be x_{i+1} , and then compute $z_1 = f(x_i), z_2 = f(z_1), z_3 = f(z_2), \cdots$ until we notice some $z_{r'+1}$ to be the same as y; clearly, $z_{r'}$ is $f^{-1}(y)$. Note that r + r' = N/m. Thus, using $m \log_2 N$ bits of memory, we can invert f in time N/m.

If we allow f to be arbitrary functions from $\{1, 2, \dots, N\}$ into $\{1, 2, \dots, N\}$, then it is not always possible to partition the points into disjoint chains of uniform length. Hellman [H] considered a certain strategy of storing 2m numbers $x_i, x_i', 1 \le i \le m$ such that there is a chain $z_{j+1} = f(z_j), j = 1, 2, \cdots$ leading from x_i to x_i' for each i, and a way of trying to find $f^{-1}(y)$ with this information. He showed that, if f is a random function uniformly chosen from all functions and if x is uniformly chosen from $\{1, 2, \dots, N\}$, then the probability of this algorithm successfully finding x from the value y = f(x) in time $O((N/m)^2)$ is at least 0.63.

Are these memory-time tradeoffs the best possible? If we restrict algorithms to those based on storing endpoints of chains, Shamir and Spencer [SS] showed that t = O(N/m) and $t = O((N/m)^2)$ are the best achievable for random 1-cycle permutations and random functions. What if we remove the restriction of basing algorithms on the chain-chasing principle?

Let us consider a general model for answering this question. The problem is similar to that considered in Section 2 except for the probabilistic setting. Let H_N denote the set of all functions $f:\{1,2,\cdots,N\}\to$ $\{1, 2, \dots, N\}$, and q_N be a probability distribution over H_N . An algorithm G consists of a partition of $\{D_{\alpha} \mid \alpha \in \{0,1\}^m\}$ of H_N and a collection of N-ary decision trees $T_{\alpha,i}$, $\alpha \in \{0,1\}^m, 1 \leq i \leq N$; the internal nodes of $T_{\alpha,i}$ ask queries of the form f(j) = ? $(j \neq i)$, and the leaves contain answers $1 \leq k \leq N$. Let cost(G) be the maximum height of the trees. Denote by $\eta(G,q_N)$ the probability that, for a random $x \in \{1, 2, \dots, N\}$ and a random f distributed according to q_N , the algorithm G with y = f(x) as input will return z satisfying f(z) = y. The following theorem can be obtained by an extension of the proof of Theorem 1. Let q_N be either the uniform distribution over all 1-cycle permutations or the uniform distribution over H_N . Let 0 < c < 1 be any fixed constant.

Theorem 2 Any algorithm G with $\eta(G, q_N) \geq c$ must satisfy $\cos(G) = \Omega((N \log N)/m)$.

Proof Omitted from this extended abstract.

Thus, Hellman's scheme is nearly optimal for random 1-cycle permutations, even allowing probabilistic algorithms. It remains an interesting open question to bridge the gap between the upper and lower bounds for the random function case.

4 Coherent Functions

Let f be a function, and \mathcal{M} be a class of algorithms, called examiners, which for every input w try to guess the value of f(w) by asking a series of questions $f(x_1) =?$, $f(x_2) =?$, \cdots , where the x_i 's are not allowed to include w. Informally, f is coherent with respect to \mathcal{M} if there exists an examiner in \mathcal{M} which can correctly predict f(w) for all w. The precise nature of the functions f and the model of algorithms will depend on the specific applications. Note that whether a function is coherent, i.e. intuitively easy to interpolate, is different from the difficulty of computing the function.

We will define below in detail the concept of coherent functions in the Turing machine model. Extensions to other models will be discussed in Section 5.2.

An examiner S is a multitape probabilistic polynomial time Turing machine with a special query tape. Given an input string $w \in \{0,1\}^*$, the oracle machine S^A with any oracle set A operates in the standard way, except that oracle A will never be used to query whether $w \in A$.

Let $f: \{0,1\}^* \to \{0,1\}$ be any boolean function. Define A_f to be the set of w such that f(w) = 1. Write S(f, w) as the random variable whose value is 1 if S^{A_f} accepts w and 0 otherwise. Let $\alpha(S, f, w) = \Pr\{S(f, w) \neq f(w)\}$, and $\alpha(S, f) = \max_w \alpha(S, f, w)$.

Definition Let $\epsilon \geq 0$. We say that S is $(1 - \epsilon)$ -successful on f if $\alpha(S, f) \leq \epsilon$.

Definition A boolean function f is said to be coherent (with respect to probabilistic polynomial time Turing machines) if for every $\epsilon > 0$ there exists an examiner S that is $(1 - \epsilon)$ -successful on f.

Remark A main difference between between the definitions for coherent functions and self-reducible functions is that in the former case we allow an adaptive choice of queries in the interpolation process, while for self-reducible functions the choice of points for interpolation purpose is made in one pass.

Theorem 3 There exists a language in $DSPACE(2^{n^{\log \log n}})$ that is not coherent.

Sketch of Proof To make the main idea clear, we will assume that checkers only ask adversary programs P to return P(x) for strings x having the same length as the input string w. To simplify the matter further, we will only prove the existence of f which have no efficient checkers. We will indicate at the end how to modify the arguments to remove these simplifying assumptions.

The general approach is as follows. For any examiner S, integer n, let $h_S(n)$ be the number of n-variable boolean functions for which S interpolates successfully (for all 2^n possible inputs). If we can

prove that $h_S(n)$ is $o(2^n)$, then we can use the diagonalization technique to construct an f which will avoid being interpolated by any M. Now the arguments used in proving Theorem 1 almost tell us that $h_S(n)$ is $o(2^n)$, except that we now allow probabilistic algorithms with errors, which entails a more involved argument.

To study the behavior of S on input strings of length n, we introduce some notations. An n-boolean function is a mapping $b: \{0,1\}^n \to \{0,1\}$. For any boolean function $f: \{0,1\}^* \to \{0,1\}$, let $f^{(n)}$ denote the induced n-boolean function when f is restricted to n-bit inputs. An (n,m)-tree is a rooted decision tree K such that the following conditions are met:

- (a) The root has 2^n edges labeled by distinct *n*-bit strings,
- (b) If v is an internal node in the w-th subtree from the root (where $w \in \{0,1\}^n$), then v contains a query of the form b(x) = ? where $x \neq w$,
- (c) Each leaf ℓ contains an output $\mu(\ell)$, and
- (d) The height of the tree is m+1.

For any n-boolean function b, we can regard K as a deterministic algorithm for guessing the value b(w) for any input w. Given w, we first take the edge labeled by w from the root, and then trace a path by asking queries and branching according to the answers until a leaf ℓ is reached; we then make the guess $b(w) = \mu(\ell)$. Clearly, for any w, the number of queries asked does not exceed m. Let $\mathcal{K}_{n,m}$ denote the set of all (n,m)-trees. We now define randomized algorithms.

An (n,m)-tree examiner T is specified by a probability distribution τ over $\mathcal{K}_{n,m}$. As an algorithm for guessing values b(w), we first pick a random K according to τ and execute K. For any w,b, let $\beta(T,b,w)$ denote the probability that T will guess b(w) incorrectly; let $\beta(T,b) = \max_{w} \beta(T,b,w)$. For any ϵ , let $\mathcal{B}_{T,\epsilon}$ denote the set of all n-boolean functions b such that $\beta(T,b) \leq \epsilon$.

CLAIM 1 Let $0 < \epsilon < 1/2$. If there exists an examiner S which is $(1-\epsilon)$ -successful on f, then there exists a k > 0 and a sequence of (n, n^k) -tree examiners T_n

such that $\beta(T_n, f^{(n)}) < 4^{-n}$ for all n.

To prove the claim, let S' be the examiner which, for any input (w, P), employs $S \ 10 \lceil n/(1/2 - \epsilon)^2 \rceil + 1$ times and takes the majority vote as the verdict. It is easy to verify that the probability that S' has made a mistake is at most 4^{-n} for any input (w, P) with |w| = n. Choose integer k large enough so that, for all n, the running time of S' is at most n^k for input (w, P) with |w| = n. The behavior of S' when |w| = n naturally gives an n-tree examiner T_n satisfying the claim. We have thus proved CLAIM 1.

The next claim extends the argument used in Theorem 1 to the probabilistic case, and is central to the proof of Theorem 3. Let k > 0 be fixed. We will prove the following claim for all large enough n.

CLAIM 2 Let $\epsilon = 4^{-n}$. If T is an (n, n^k) -tree examiner, then $|\mathcal{B}_{T,\epsilon}| \leq 2^{2^n(1-\frac{1}{2n^k})}$.

To prove CLAIM 2, let us label the elements of $\mathcal{B}_{T,\epsilon}$ as K_1, K_2, K_3, \cdots . Let $M = \lceil 2^n/(n^k+1) \rceil$, and I_n be the set of all M-tuples of positive integers. For each $\tilde{a} = (a_1, a_2, \cdots, a_M) \in I_n$, consider the following deterministic procedure which, for any n-boolean function b, generates a binary string $\sigma(\tilde{a}, b)$.

Procedure encode

begin Set $J \leftarrow \emptyset$;

for j = 1 to M do

begin take smallest $i \notin J$;

use $K_{a,j}$ to ask exactly n^k queries b(t)=? $(t \neq i)$ and then make a guess of b(i); Set $\gamma_j \in \{0,1\}^{n^k}$ to be the sequence of answers to the n^k queries asked;

enlarge J to include i and all the t's

end

if all the M guesses of b(i) are correct then $\sigma(\tilde{a}, b) \leftarrow 0\gamma_1\gamma_2\cdots\gamma_M$; if any of the M guesses of b(i) is incorrect then $\sigma(\tilde{a}, b) \leftarrow 1$ followed by the 2^n values of b(w)

end encode

In the above procedure it is understood that i, t are n-bit strings, and the phrase "smallest i" means "lexicographically smallest".

Let τ be the probability distribution describing T such that $\tau(i)$ is the probability that T takes on the value K_i . We extend the notation τ by defining $\tau(\tilde{a})$ to be $\prod_{1 \le j \le M} \tau(a_j)$, when $\tilde{a} = (a_1, a_2, \dots, a_M) \in I_n$.

For any \tilde{a} , $\sigma(\tilde{a}, b)$ gives an encoding of all the *n*-boolean functions b. Let $\mu(\tilde{a})$ denote the average length of the encoding for a random $b \in \mathcal{B}_{T,\epsilon}$. Then

$$\sum_{\tilde{a}} \tau(\tilde{a}) \mu(\tilde{a}) = \sum_{\tilde{a}} \tau(\tilde{a}) \frac{1}{|\mathcal{B}_{T,\epsilon}|} \sum_{b \in \mathcal{B}_{T,\epsilon}} |\sigma(\tilde{a}, b)|$$
$$= \frac{1}{|\mathcal{B}_{T,\epsilon}|} \sum_{b \in \mathcal{B}_{T,\epsilon}} \sum_{\tilde{a}} \tau(\tilde{a}) |\sigma(\tilde{a}, b)|(2)$$

Now, for any fixed $b \in \mathcal{B}_{T,\epsilon}$, the probability for any of the M guesses to be incorrect is bounded by $M \cdot \epsilon \leq 2^{-n}$, and hence for large n

$$\sum_{\tilde{a}} \tau(\tilde{a}) |\sigma(\tilde{a}, b)| \leq (1 + Mn^{k}) + 2^{-n} (1 + 2^{n})$$

$$\leq 2^{n} (1 - \frac{1}{2n^{k}}). \tag{3}$$

It follows from (2) and (3) that there exists \tilde{a} such that

$$\mu(\tilde{a}) \leq 2^n \left(1 - \frac{1}{2n^k}\right).$$

Since any encoding in binary of the element in the set $\mathcal{B}_{T,\epsilon}$ must have an average length of $\log_2 |\mathcal{B}_{T,\epsilon}|$ or more, it follows that

$$|\mathcal{B}_{T,\epsilon}| \leq 2^{2^n(1-\frac{1}{2n^k})}.$$

This completes the proof of CLAIM 2.

CLAIMs 1 and 2 allow us to construct by diagonalization a function f which is not coherent, as we will show below. Let S_1, S_2, \cdots be the sequence of all examiners. Let $\epsilon = 1/4$, and let $T_1^{[i]}, T_2^{[i]}, \cdots, T_n^{[i]}, \cdots$ be the sequence of tree examiners mentioned in CLAIM 1 for S_i . Let $\epsilon_n = 4^{-n}$. We can define a sequence of integers Λ_i such that $\Lambda_1 < \Lambda_2 < \cdots < \Lambda_i < \cdots$ and, for each j,

$$\left| \bigcup_{1 \leq i \leq j} |\mathcal{B}_{T_n^{[i]}, \epsilon_n} \right| \leq j \cdot 2^{2^n (1 - 1/n^{\log \log n})} < 2^{2^n}$$

for all $n \geq \Lambda_j$. As there are 2^{2^n} n-boolean functions, there exists a function f such that $f^{(n)} \notin \bigcup_{1 \leq i \leq j} \mathcal{B}_{T_n^{[i]}, \epsilon_n}$ for all $n \geq \Lambda_j$. This means, for any i, $T_n^{[i]}$ will fail to interpolate $f^{(n)}$ with error ϵ_n for all sufficiently large n. By CLAIM 2 no examiner S_i can be $(1 - \epsilon)$ -successful on f.

This completes the proof of Theorem 3, with the assumptions stated at the beginning of the proof. To show that the function f can be constructed in space $2^{n^{\log \log n}}$, one only needs to study the resources needed in carrying out the diagonalization argument. To remove the restriction that checkers can ask queries only of the same length as the input, we observe that the proof presented remains valid even if n does not take on all possible integer values. Let $\mathcal L$ denote the family of languages that contain no strings of length not of the form $2^{2^{2^i}}$. Then the checkers have no need to ask queries longer than the input. The present proof then shows the existence of a language in $\mathcal L$ that is not coherent. The details will be described in the complete paper. \square

Remark Theorem 3 remains true if we replace $\log \log n$ by any exponential-space-constructible monotonic function $k(n) \to \infty$.

Remark A result related to Theorem 3 was proved by Feigenbaum, Kannan, and Nisan [FKN] for selfreducible functions.

5 Program Checkers

5.1 Turing Machine Model

Program checking, which is a novel approach for checking the correctness of programs, was recently raised by Blum and Kannan [BK] (see also Blum, Luby, and Rubinfeld [BLR], Lipton [L]). Seemingly intractable computations such as graph isomorphism have simple *checkers* which run in polynomial time. The limitation of this approach is not well understood; in fact no computations have been shown to be hard to check.

Let \mathcal{M} be the class of probabilistic polynomial time Turing machines. We will show that any booleanvalued function with efficient checkers must be coherent with respect to \mathcal{M} . As a corollary, this shows that there exist functions in $DSPACE(2^{n^{\log \log n}})$ for which no efficient checker exists.

Let f be a computational problem. Let P be a deterministic Turing machine that halts for all input strings w; call the output P(w). Following the terminology in [BK], we say that P has a bug (with respect to f) if $P(w) \neq f(w)$ for some w. An efficent checker for f is a probabilistic polynomial time oracle Turing machine which, for any P as the oracle and any w, will output either CORRECT or BUGGY such that the following conditions are satisfied:

- (a) If P has no bugs, then with probability at least2/3, M will output CORRECT;
- (b) If $P(w) \neq f(w)$, then with probability at least 2/3, M will output BUGGY.

Clearly, the quantity 2/3 in the above definition can be substituted by any fixed constant 1/2 < c < 1.

Theorem 4 If f has an efficient checker, then f is coherent.

Proof Let $\epsilon > 0$, and V be an efficient checker for f with error probability at most ϵ . For any $w \in \{0,1\}^*$, let P_w denote the program which returns $P_w(w) = 0$ and $P_w(x) = f(x)$ for all $x \neq w$. Suppose (w, P) is input to V. From the definition of program checker, if we make the guess f(w) = 0 when V gives the CORRECT verdict, and 1 when V outputs BUGGY, then the probability $r_V(w)$ of the guess being correct is at least $1 - \epsilon$.

Now turn V into an examiner S_V by simulating V, except omitting queries f(w) = ? where w is the input and behaving as if an adversary returns the value f(w) = 0; at the end if V gives a CORRECT verdict, then S outputs the value 0, and if V gives a BUGGY verdict, then S outputs the value 1. Clearly, the probability for S_V to output the correct value of f(w) is exactly $r_V(w)$, which is at least $1 - \epsilon$. This proves the theorem. \square

Theorem 5 There exists $L \in DSPACE(2^{n^{\log \log n}})$ for which there is no efficient checker.

Proof Follows immediately from Theorems 3 and 4. \square

Remark Theorem 5 can be strengthened to $DSPACE(2^{cn})$ for any c > 1 by a direct proof from the definition of program checkers.

5.2 Comparison Tree Models

Although program checking was defined originally for the Turing machine model in [BK], it is of interest to consider program checking in other models such as the comparison-based model, the bounded-degree algebraic tree model, or the arithmetic computation tree model. The merits are (a) these models have more realistic cost measure than the Turing machine model for certain types of problems such as geometric optimization, and (b) nonlinear lower bounds are known in these models, which might make it possible to compare the cost of optimal checkers versus optimal algorithms for a problem.

We take a step in this direction by considering two classes of ordering problems in a comparison-base model, namely, sorting-like problems as defined by Fredman [F] and partial order production problems as given by Schönhage [Sch] (also see [Y2]). We will show that these two classes of problems behave quite differently with respect to program checking. In fact, for almost all sorting-like problems, checking is as expensive as direct computing; while every partial order production problem has a linear time checker and hence checking is always cheap.

Before proceeding, we give an example of a well-known problem that has an efficient checker in the comparison tree model.

ELEMENT DISINCTNESS: Given real numbers x_1, x_2, \dots, x_n , determine whether they are all distinct.

In the comparison tree model, a program can perform a sequence of tests of the form $x_i : x_j$ with $\langle , =, \text{ or } \rangle$ as possible outcomes, and the cost of

the program is the number of comparisons needed. It is well known that the cost of any program is at least $\Omega(n \log n)$, even for probabilistic algorithms allowing error (Manber and Tompa [MT]). Let us consider probabilistic procedures V which, when given a set W of numbers and a program P, want to decide if the answer of P to W is reliable. Aside from performing comparisons, V can call on P to determine whether a subset (allowing repeated entries) of the input numbers has the all-distinctness property. We will say that V is a program checker if for any W and P, the final judgment made has probability at least 2/3 to be right. The cost of the program checker is the maximum expected number of comparisons plus the number of calls to P. Since we will describe an algorithm and not prove lower bounds for this problem, we will not describe the complexity model in detail. The model will become clear from the later discussion of sorting-like problems.

Theorem 6 ELEMENT DISTINCTNESS has a program checker with O(n) cost in the comparison tree model.

Proof We describe below a program checker V. Given an input (W, P), where $W = \{x_1, x_2, \dots, x_n\}$ is a set of real numbers and P is a program, possibly faulty, for the element distinctness problem. Depending on the answer of P to W, we distinguish two cases.

- (A) If P returns "no", then for $j=n,n-1,\cdots$, ask P whether $\{x_1,x_2,\cdots,x_{j-1}\}$ contains all distinct elements, until either (a) in the round $j=k\geq 2$, P returns a "yes" answer, or (b) up to and including round j=2, P has not returned a "yes" answer. In case (b) V says BUGGY. In case (a) V compares x_k against every x_i , $1\leq i< k$; if $x_k=x_i$ for some i then says CORRECT, and otherwise gives the BUGGY verdict.
- (B) If P returns "yes", then V randomly permutes x_1, x_2, \dots, x_n and partitions it into two sets $\{y_1, y_2, \dots, y_{n/2}\}$ and $\{y_{n/2+1}, \dots, y_n\}$. Then it constructs n sets $W_j = \{y_1, y_2, \dots, y_{n/2}\} \cup \{j\}, 1 \leq j \leq n$, and presents to P in random order these sets. If P

returns "yes" for any W_j with $1 \le j \le n/2$ or if P returns "no" for any W_j with $1 \le j \le n/2$, then V says BUGGY; otherwise V says CORRECT.

The above V clearly runs in time O(n). It is easy to see that V always gives the right verdict in situation (A). In situation (B) if P did not lie, M clearly will give the right verdict. It remains to show that, when x_1, x_2, \dots, x_n have some $x_i = x_j$ but P returns "yes", M will catch P with a BUGGY verdict with probability at least 1/4. However, this is true since with probability 1/2, x_i and x_j will be in different halves of the y's list, and P cannot distinguish between S_i and S_j . This proves Theorem 6. \square

Remark In Blum and Kannan [BK] and Lipton [L], O(n)-time program checkers were given for the element distinctness problem, but not in the comparison model.

We now turn to the discussion of sorting-like problems and partial order production problems. A sorting-like problem g is a function from the set of all permutations on n or fewer elements into $\{1, 2, \dots, r\}$ where $n, r \geq 2$ are integers. Let J_m denote the set of all m-tuples of distinct real numbers. For any $\tilde{x} = (x_1, x_2, \dots, x_m) \in J_m$, let $\sigma_{\tilde{x}}$ be the permutation (i_1, i_2, \dots, i_m) such that $x_{i_1} < x_{i_2} < \dots < x_{i_m}$. Given an input $\tilde{x} \in J_m$ where $m \leq n$, we want to determine the value of $g(\sigma_{\tilde{x}})$. For example, the standard sorting problem corresponds to the case $r = 1! + 2! + \dots + n!$ with each integer $1 \leq j \leq r$ representing a linear ordering of n or fewer elements. Let $G_{n,r}$ be the set of all sorting-like problems.

Below we give a rough description of the model. A detailed specification is given in the Appendix.

Let $\mathcal{F}_{n,r}$ be the set of all decision trees using binary comparisons $x_i : x_j$ as primitives. A decision tree $F \in \mathcal{F}_{n,r}$ is an algorithm for g if it outputs the value of $g(\sigma_{\tilde{x}})$ for all inputs \tilde{x} .

A deterministic checker H is a decision tree for which the inputs are triplets (\tilde{x}, F, j) ; $F \in \mathcal{F}_{n,r}$ and j is the output of F when \tilde{x} is given to F. The deterministic checker H needs to either endorse the value

j as $g(\sigma_{\tilde{x}})$ or returns a BUGGY verdict. In addition to performing comparisons $x_i:x_j$, H can make subroutine calls of F, each time supplying F with a permuted tuple $(x_{i_1}, x_{i_2}, \dots, x_{i_k})$ and then running F. In the spirit of [BK], each call is counted as one step. A checker V is a randomized decision tree defined in the standard way [Y1]. We say that V is a checker with error ϵ if, for every \tilde{x} and $F \in \mathcal{F}_{n,r}$, the probability of V returning the wrong verdict is at most ϵ .

Clearly, for the standard sorting problem, there is a checker V with error 0 and running time n-1. In fact, if the input is (\tilde{x}, F, j) where j is the encoding of (i_1, i_2, \dots, i_m) , then V can simply check the m-1 inequalities $x_{i_1} < x_{i_2} < \dots < x_{i_m}$. (Our model is different from the one used in [BK] for sorting and has no need to check set equalities.) Since it is clear that any algorithm (even randomized and with error, say 1/3) for sorting must use $\Omega(n \log n)$ comparisons, it is easier to check than to compute for the sorting problem. We will show that this is an anomaly for sorting-like problems, and in fact almost all sorting-like problems are as hard to check as to compute. Before doing that, we define the concept of coherence in this model.

An examiner S is a randomized decision tree for which the inputs are pairs (\tilde{x},g) , where $\tilde{x} \in J_m, m \leq n$ and $g \in G_{n,r}$; the outputs are integers $1 \leq j \leq r$. It can make comparisons $x_i : x_j$ and ask queries $g(\sigma_{\tilde{y}}) = r$? where $\tilde{y} = (x_{i_1}, x_{i_2}, \cdots, x_{i_k}) \neq \tilde{x}$; each comparison or query is counted as one step for the cost. Let $\alpha(S,g,\tilde{x})$ denote the probability that the output of S is different from $g(\sigma_{\tilde{x}})$; let $\alpha(S,g) = \max_{\tilde{x}} \alpha(S,g,\tilde{x})$. For any S and g, we will say that S is $(1-\epsilon)$ -successful on g if $\alpha(S,g) \leq \epsilon$. Let $S_{n,r}$ denote the set of all examiners.

Definition Let r > 1 be fixed. A sequence of sorting-like problems $\hat{g} = g_1, g_2, \cdots$ where $g_n \in G_{n,r}$ is said to be *coherent* (with respect to O(n)-cost decision trees) if, for every $\epsilon > 0$, there exists a sequence of examiners S_1, S_2, \cdots , where $S_n \in \mathcal{S}_{n,r}$, such that S_n has O(n) cost and is $(1 - \epsilon)$ -successful on g_n for

every n.

Theorem 7 Let $\hat{g} = g_1, g_2, \cdots$ where $g_n \in G_{n,r}$. If there exist a sequence of checkers V_n for g_n with cost O(n) and error 1/3, then \hat{g} must be coherent.

Proof Similar to the proof of Theorem 4. □

Theorem 8 Take a random sequence of sorting-like problems g_1, g_2, \cdots where g_n is uniformly chosen from $G_{n,r}$. Then with probability $1 \hat{g}$ is not coherent.

Proof Omitted. □

It follows from Theorems 7 and 8 that, if we choose a random g_n uniformly from $G_{n,r}$ for each n, then with probability 1, there cannot be any sequence of O(n)-cost checkers V_n for g_n with error 1/3. A stronger version of the result is given below.

Theorem 9 Let r > 1 be fixed. Take a random g uniformly chosen from $G_{n,r}$. Then with probability 1 - o(1) for large n, the following is true: any checker for g with error probability less than 1/3 must have $\Omega(n \log n)$ running time.

Proof Omitted, □

The other class of decision tree problems we consider is the partial order production problem ([Sch][Y2]). Let Q be a partial order on n elements z_1, z_2, \dots, z_n . Given n distinct numbers x_1, x_2, \dots, x_n , we want to find a permutation i_1, i_2, \dots, i_n such that $z_j <_Q z_k$ implies $x_{i_j} < x_{i_k}$. Define checkers as in the sorting-like problems case. The following result shows that, for any Q with only a moderate number #Q of linear extensions, checking is easier than computing. The proof is an easy consequence of results in the literature originally considered in other contexts.

Theorem 10 Any randomized algorithm for Q-production with error 1/3 must have running time $\Omega(\log(n!/\#Q))$. There is a checker for Q with error 1/3 and running time O(n).

The first statement is a straightforward extension of Schönhage's result [Sch] that any deterministic algorithm for Q-production must have running time $\Omega(\log(n!/\#Q))$. The second statement follows imme-

diately from a recent result of Valerie King [K] that the partial order verification problem on n elements can be done in O(n) time probabilistically (improving on a previous result by Kenyon-Mathieu and King [KK]).

6 Concluding Remarks

There are many open problems in this subject. The broad direction is to seek other basic properties for classes of functions, study their relationships, and use them to resolve cryptographic questions. It seems that the randomness concept for functions is now well understood (see Goldreich, Goldwasser, and Micali [GGM]), but concepts concerning various nonrandomness properties are not yet fully explored. Our development of the "coherence" notion is one step in this direction.

We list below a few concrete open questions which are immediately suggested by this work.

- (a) Is there any language in the polynomial time hierarchy that is not coherent, under reasonable assumptions such as the noncollapsing of certain complexity hierarchies? A result of this nature for the encrypted data computation was given by coherence, Feigenbaum, and Kilian [AFK]. How about assuming the existence of one-way functions? Some interesting results along this line have recently been obtained by Beigel and Feigenbaum [BeiF].
- (b) Investigate program checking in the algebraic computation tree model. For example, can one prove that any checker for the problem of finding the closest pair of points in the plane must take $\Omega(n \log n)$ time?
- (c) Is the time-memory tradeoff for inverting random functions achieved by Hellman's algorithm optimal?

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Appendix: Model for Sorting-Like Problems

Let n, r be positive integers. Let $\Delta_n = \bigcup_{1 \leq m \leq n} \Gamma_m$, where Γ_m is the set of all permutations of $\{1, 2, \dots, m\}$. Following Fredman [F] with minor modifications, a sorting-like problem Q is a family of disjoint sets (Y_1, Y_2, \dots, Y_r) that partition Δ_n . Given an m-tuple of distinct real numbers $\tilde{x} = (x_1, x_2, \dots, x_m)$, we want to determine $f_Q(\tilde{x})$, the integer j such that Y_j contains the linear ordering satisfied by the $x_i's$.

A comparison forest F is a collection of n comparison-based decision trees A_1, A_2, \dots, A_n , where A_m uses only comparisons of the form $x_i : x_j$ with $1 \le i, j \le m$; each leaf of A_m contains an output $s_\ell \in \{1, 2, \dots, r\}$. Let \mathcal{F} denote the set of all comparison forests.

For $\tilde{x} = (x_1, x_2, \dots, x_m)$ and $F \in \mathcal{F}$, let $cost(F, \tilde{x})$ denote the number of comparisons used when A_m is used for the input \tilde{x} . We will say that F correctly computes Q for input $\tilde{x} = (x_1, x_2, \dots, x_m)$, if for A_m the leaf ℓ reached by \tilde{x} satisfies $s_{\ell} = f_Q(\tilde{x})$; otherwise, F incorrectly computes Q for \tilde{x} . Write $\nu(F, \tilde{x}) = 0$ in the former case and 1 in the latter case. Let \mathcal{F}_Q denote the set of F that correctly compute Q for all inputs \tilde{x} .

A randomized algorithm B for Q is a probability distribution p on \mathcal{F} . For input \tilde{x} , the cost of B is defined to be $c(B,\tilde{x}) = \sum_{F \in \mathcal{F}} p(F) \cos(F,\tilde{x})$; the error probability $e_Q(B,\tilde{x})$ is, for a random F distributed according to p, the probability that F incorrectly computes \tilde{x} . Let $c(B) = \max_{\tilde{x}} c(B,\tilde{x})$. For any $0 \le \epsilon \le 1$, let $\mathcal{B}_{Q,\epsilon}$ denote the set of all randomized algorithms B satisfying $e_Q(B,\tilde{x}) \le \epsilon$. The randomized complexity with error ϵ is defined as $C_{Q,\epsilon} = \min \{c(B) \mid B \in \mathcal{B}_{Q,\epsilon}\}$.

Clearly, every sorting-like problem Q on n elements has a comparison forest F such that F computes Q correctly for every \tilde{x} and uses $O(n\log n)$ comparisons. As a trivial corollary, $C_{Q,\epsilon} = O(n\log n)$ for all $0 \le \epsilon < 1$.

A deterministic checker H is a collection of rn rooted trees $D_{k,1}, D_{k,2}, \cdots, D_{k,n}, 1 \leq k \leq r$, where $D_{k,m}$ has two kinds of internal nodes: comparison nodes with tests $x_i: x_j$, and subroutine nodes specified with a j-tuple $\tilde{x}' = (x_{i_1}, x_{i_2}, \cdots, x_{i_j})$ where $1 \leq j \leq n$. Out of every comparison node there are two edges labeled with < and >; out of each subroutine node, there are up to r edges labeled with distinct integers between 1 and r. At each leaf ℓ , a value $\mu(\ell) \in \{\text{CORRECT}, \text{BUGGY}\}$ is stored. Let \mathcal{H} denote the set of all comparison forest checkers.

As a deterministic checker, $H \in \mathcal{H}$ works as follows. Given any adversary program, i.e. a comparison forest $F \in \mathcal{F}_n$ and $\tilde{x} = (x_1, x_2, \dots, x_m)$, suppose for input \tilde{x} the output of F is k. Then H will try to check the result by tracing a path in $D_{k,m}$ until a leaf ℓ is reached; the rules for tracing the path are (a) when a comparsion node with test $x_i : x_j$ is encoun-

tered, perform the test and take the branch according to the result, and (b) when a subroutine node with $\tilde{x}'=(x_{i_1},x_{i_2},\cdots,x_{i_j})$ is encountered, ask F to perform a computation with input \tilde{x}' , and take the edge labeled by t if t is the output from F. The value $\mu(\ell)$ is the result to be returned by the checker H. Define $cost(H,F,\tilde{x})$ as the path length from the root to ℓ . Let $\delta(H,F,\tilde{x})$ be the boolean function whose value is 1 if and only if H fails in checking the validity of F's computing Q for input \tilde{x} . Precisely, let $\delta(H,F,\tilde{x})=1$ if and only if either of the following conditions is true: (a) $(F\in\mathcal{F}_Q) \wedge (\mu(\ell)=\mathrm{BUGGY})$, (b) $(\nu(F,\tilde{x})=1) \wedge (\mu(\ell)=\mathrm{CORRECT})$.

A checker V for Q is a probability distribution p on \mathcal{H} . For input (\tilde{x},F) , the expected number of operations used by V is clearly $c(V,F,\tilde{x})=\sum_{H\in\mathcal{H}}p(H)\cos(H,F,\tilde{x});$ the error probability $\delta_Q(V,F,\tilde{x})$ is, for a random H distributed according to p, the probability that H fails as a checker, i.e. $\sum_{H\in\mathcal{H}}p(H)\delta(H,F,\tilde{x}).$ The cost of V is defined to be $\max_{F,\tilde{x}}c(V,F,\tilde{x}).$ We say that V is a checker for Q with error ϵ , if $\delta_Q(V,F,\tilde{x})\leq \epsilon$ for all $F,\tilde{x}.$