Part IV

Games and design of randomized algorithms

In this chapter we present a new way how to see randomized algorithms and several basic techniques how to design and analyse randomized algorithms:

- Especially we deal with:
 - Application of linearity of expectations
 - Game theory based lower bounds methods for randomized algorithms.

A way to see basics of deterministic, randomized and quantum computations and their differences.

Let us consider an *n* bits strings set $S \subset \{0, 1\}^n$.

To describe a **deterministic computation** on S we need to specify

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an initial state - by an n-bit string - say s_0
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and an evolution (computation mapping), $E: S \to S$ which can be described by a vector of the length 2^n , the elements and indices of which are *n*-bit strings.

A computation step is then an application of the evolution mapping E to the current state represented by an *n*-bit string *s*.

However, for any a bit significant task, the number of bits needed to describe such an evolution mapping, $n2^n$, is much too big. The task of programming is then/therefore to replace an application of such an enormously huge mapping by an application of a much shorter circuit/program.

To describe a randomized computation we need;

1:) to specify an **initial probability distribution** on all *n*-bit strings, what can be done by a vector of length 2^n , indexed by *n*-bit strings, the elements of which are non-negative numbers that sum up to 1

2:) to specify a randomized evolution, which has to be done, in case of a homogeneous evolution, by a $2^n \times 2^n$ matrix A of conditional probabilities for obtaining a new state/string from an old state/string.

The matrix A has to be **stochastic** - all columns have to sum up to one and A[i, j] is a probability of going from a string representing j to a string representing i.

To perform a computation step, one then needs to multiply by A the 2^n -elements vector specifying the current probability distribution on 2^n states.

However, for any nontrivial problem the number 2^n is larger than the number of particles in the universe. Therefore, the task of programming is to design a small circuit/program that can implement such a multiplication by a matrix of an enormous size.

In case of **quantum computation** on *n* quantum bits:

1:) Initial state has to be given by an 2^n vector of complex numbers (probability amplitudes) the sum of the squares of which is one.

2:) Homogeneous quantum evolution has to be described by an $2^n \times 2^n$ unitary matrix of complex numbers - at which inner products of any two different columns and any two different rows are $0.^1$

Concerning a **computation step**, this has to be again a multiplication of a vector of the probability amplitudes, representing the current state, by a very huge $2^n \times 2^n$ unitary matrix which has to be realized by a "small" quantum circuit (program).

 $^{^{1}}$ A matrix A is usually called unitary if its inverse matrix can be obtained from A by transposition around the main diagonal and replacement of each element by its complex conjugate.

LINEARITY OF EXPECTATIONS

A very simple, but very often very useful, fact is that for any random variables X_1, X_2, \ldots it holds

$$\mathbf{E}[\sum_{i} X_{i}] = \sum_{i} \mathbf{E}[X_{i}].$$

even if X_i are dependent and dependencies among X_i 's are very complex.

Example: A ship arrives at a port, and all 40 sailors on board go ashore to have fun. At night, all sailors return to the ship and, being drunk, each chooses randomly a cabin to sleep in. Now comes the **question**: *What is the expected number of sailors sleeping in their own cabins?*

Solution Let X_i be a random variable, so called *(indicator variable)*, which has value 1 if the *i*-th sailor chooses his own cabin, and 0 otherwise.

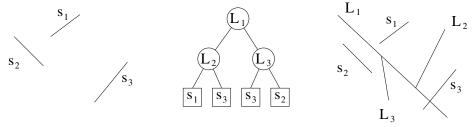
Expected number of sailors who get to their own cabin is

$$\mathsf{E}[\sum_{i=1}^{40} X_i] = \sum_{i=1}^{40} \mathsf{E}[X_i]$$

Since cabins are chosen randomly $\mathbf{E}[X_i] = \frac{1}{40}$ and therefore, $\mathbf{E}[\sum_{i=1}^{40} X_i] = 40.\frac{1}{40} = 1.$

EXAMPLE - **BINARY PARTITION** of a **SET** of **LINE SEGMENTS** 1/3

Problem Given a set $S = \{s_1, \ldots, s_n\}$ of non-intersecting line segments, find a partition of the plane such that every region will contain at most one line segment (or at most a part of a line segment).



A (binary) partition will be described by a binary tree + additional information (about nodes). With each node v a region r_v of the plane will be associated (the whole plane will be represented by the root) and also a line L_v intersecting r_v .

Each line L_v will partition the region r_v into two regions $r_{l,v}$ and $r_{r,v}$ which correspond to two children of v - to the left and right one.

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Notation: $l(s_i)$ will denote a **line-extension of the segment** s_i .

autopartitions will use only line-extensions of given segments. Algorithm RandAuto:

Input: A set $S = \{s_1, \ldots, s_n\}$ of non-intersecting line segments.

Output: A binary autopartition P_{Π} of S.

1: Pick a permutation Π of $\{1, \ldots, n\}$ uniformly and randomly.

2: While there is a region R that contains more than one segment, choose one of them randomly and cut it with $l(s_i)$ where i is the first element in the ordering induced by Π such that $l(s_i)$ cuts the region R. Theorem: The expected size of the autopartition P_{Π} of S, produced by the above RandAuto algorithm is $\theta(n \ln n)$. **Proof:** Notation (for line segments u, v).

 $index(u, v) = {i \ if \ l(u) intersects i - 1 segments before hitting v;} \\ \infty \ if l(u) does not hit v.$

 $u \dashv v$ will be an event that l(u) cuts v in the constructed (autopartition) tree.

EXAMPLE - BINARY PARTITION of a SET of LINE SEGMENTS 3/3

Probability: Let u and v be segments, index(u, v) = i and let u_1, \ldots, u_{i-1} be segments the line I(u) intersects before hitting v.

The event $u \dashv v$ happens, during an execution of RandPart, only if u occurs before any of $\{u_1, \ldots, u_{i-1}, v\}$ in the permutation Π . Therefore the probability that event $u \dashv v$ happens is $\frac{1}{i+1} = \frac{1}{index(u,v)+1}$.

Notation: Let $C_{u,v}$ be the indicator variable that has value 1 if $u \dashv v$ and 0 otherwise.

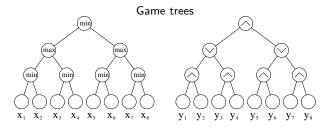
$$\mathbf{E}[C_{u,v}] = \Pr[u \dashv v] = \frac{1}{index(u,v)+1}.$$

Clearly, the size of the created partition P_{Π} equals n plus the number of intersections due to cuts. Its expectation value is therefore

$$n+E[\sum_{u}\sum_{v\neq u}C_{u,v}]=n+\sum_{u}\sum_{v\neq u}Pr[u\dashv v]=n+\sum_{u}\sum_{v\neq u}\frac{1}{index(u,v)+1}.$$

For any line segment u and integer i there are at most two v, w such that index(u, v) = index(u, w) = i. Hence $\sum_{v \neq u} \frac{1}{index(u, v)+1} \leq \sum_{i=1}^{n-1} \frac{2}{i+1}$ and therefore $n + \mathbf{E}[\sum_{u} \sum_{v \neq u} C_{u,v}] \le n + \sum_{u} \sum_{i=1}^{n-1} \frac{2}{i+1} \le n + 2nH_n.$

GAME TREE EVALUATION - I.

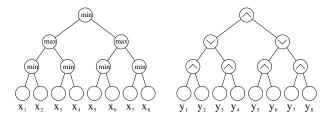


Game trees are trees with operations max and min alternating in internal nodes and values assigned to their leaves. In case all such values are Boolean - 0 or 1 Boolean operation OR and AND are considered instead of max and min.



 T_k – binary game tree of depth 2k. Goal is to evaluate the tree - the root.

GAME TREE EVALUATION - II.

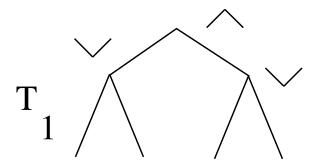


Evaluation of game trees plays a crucial role in AI, in various game playing programs.

Assumption: An evaluation algorithm chooses at each step (somehow) a leaf, reads its value and performs all evaluations of internal nodes it can perform. Cost of an evaluation algorithm is the number of leaves inspected. Determine the total number of such steps needed.

WORST CASE COMPLEXITY

 T_k – will denote the binary game tree of depth 2k.



Every deterministic algorithm can be forced to inspect all leaves. The worst-case complexity of a deterministic algorithm is therefore:

$$n=4^k=2^{2k}.$$

To evaluate an AND-node v, the algorithm chooses randomly one of its children and evaluates it.

If 1 is returned, algorithm proceeds to evaluate other children subtree and returns as the value of v the value of that subtree. If 0 is returned, algorithm returns immediately 0 for v (without evaluating other subtree).

To evaluate an OR-node v, algorithm chooses randomly one of its children and evaluates it.

If 0 is returned, algorithm proceeds to evaluate other subtree and returns as the value of v the value of the subtree. If 1 is returned, the algorithm returns 1 for v.

- Start at the root and in order to evaluate a node evaluate (recursively) a random child of the current node.
- If this does not determine the value of the current node, evaluate the node of other child.

Theorem: Given any instance of T_k , the expected number of steps for the above randomized algorithm is at most 3^k .

Proof by induction: **Base step:** Case k = 1 easy - verify by computations for all cases. **Inductive step:** Assume that the expected cost of the evaluation of any instance of T_{k-1} is at most 3^{k-1} .

Consider an OR-node tree T with both children being T_{k-1} -trees. If the root of T were to return 1, at least one of its T_{k-1} -subtrees has to return 1. With probability $\frac{1}{2}$ this child is chosen first, given in average at most 3^{k-1} leaf-evaluations. With probability $\frac{1}{2}$ both subtrees are to be evaluated. The expected cost of determining the value of T is therefore:

$$\frac{1}{2} \times 3^{k-1} + \frac{1}{2} \times 2 \times 3^{k-1} = \frac{1}{2} \times 3^k = \frac{3}{2} \times 3^{k-1}.$$

If the root of T were to return 0 both subtrees have to be evaluated, giving the cost $2 \times 3^{k-1}$.

Consider now the root of T_k .

If the root evaluates to 1, both of its OR-subtrees have to evaluate to 1. The expected cost is therefore

$$2\times\frac{3}{2}\times3^{k-1}=3^k.$$

If the root evaluates to 0, at least one of the subtrees evaluates to 0. The expected cost is therefore

$$\frac{1}{2} \times 2 \times 2 \times 3^{k-1} + \frac{1}{2} \times \frac{3}{2} \times 3^{k-1} \le 3^k = n^{\lg_4 3} = n^{0.793}$$

Our algorithm is therefore a *Las Vegas algorithm*. Its running time (number of leaves evaluations) is: $n^{0.793}$.

CLASSICAL GAMES THEORY BRIEFLY

We will consider **games with two players**, Alice and Bob.*X* and *Y* will be nonempty sets of their game (pure) strategies -*X* of Alice, *Y* of Bob. Mappings $p_X : X \times Y \to \mathbf{R}$ and $p_Y : X \times Y \to \mathbf{R}$ will be called **payoff functions** of Alice and Bob. The quadruple (X, Y, p_X, p_Y) will be called a (mathematical) game.

A mixed strategy will be a probability distribution on pure strategies.

An element $(x, y) \in X \times Y$ is said to be a **Nash equilibrium** of the game (X, Y, p_X, p_Y) iff $p_X(x', y) \le p_X(x, y)$ for any $x' \in X$, and $p_Y(x, y') \le p_Y(x, y)$ for all $y' \in Y$.

Informally, Nash equilibrium is such a pair of strategies that none of the players gains by changing his/her strategy.

A game is called **zero-sum game** if $p_X(x, y) + p_Y(x, y) = 0$ for all $x \in X$ and $y \in Y$.

One of the basic result of the classical game theory is that not every two-players zero-sum game has a Nash equilibrium in the set of pure strategies, but there is always a Nash equilibrium if players follow mixed strategies. It has been shown, for several zero-sum games, that if one of the players can use quantum tools and thereby **quantum strategies**, then he/she can increase his/her chance to win the game.

This way, from a fair game, in which both players have the same chance to win if only classical computation and communication tools are used, an unfair game can arise, or from an unfair game a fair one. Alice and Bob play with a box and a penny as follows:

- Alice places a penny head up in a box.
- Bob flips or does not flip the coin
- Alice flips or does not flip the coin
- Bob flips or does not flip the coin

After the "game" is over, they open the box and Bob wins if the penny is head up.

It is easy to check that using pure strategies chances to win are $\frac{1}{2}$ for each player and there is no (Nash) equilibrium in the case of pure classical strategies.

However, there is equilibrium if Alice chooses its strategy with probability $\frac{1}{2}$ and Bob chooses each of the four possible strategies with probability $\frac{1}{4}$.

Two members of a gang are imprisoned, each in a separate cell, without possibility to communicate. However, police has not enough evidence to convict them on the principal charge and therefore police intends to put both of them for one year to jail on a lesser charge.

Simultaneously police offer both of them so called Faustian bargain. Each prisoner gets a chance either to betray the other one by testifying that he committed the crime, or to cooperate with the other one by remaining silent. pause Here are payoffs they are offered:.

- If both betray, they will get into jail for 2 years.
- If one betrays and second decides to cooperate, then first will get free and second will go to jail for 3 years.
- If both cooperate they will go to jail for 1 year.

What is the best way for them to behave? This game is a model for a variety of real-life situations involving cooperative behaviour. Game was originally framed in 1950 by M. Flood and M. Dresher

Two prisoners, Alice and Bob, can use, independently, any of the following two strategies: **to cooperate** or **to defect** (not to cooperate).

The problem is that the payoff function (p_A, p_B) , in millions, is a very special one (first (second) value is payoff of Alice (of Bob):

What is the best way for Alice and Bob to proceed in order to maximize their payoffs?

A strategy s_A is called **dominant** for Alice if for any other strategy s'_A of Alice and s_B of Bob, it holds

$$P_A(s_A, s_B) \geq P_A(s'_A, s_B).$$

Clearly, defection is the dominant strategy of Alice (and also of Bob) in the case of Prisoners Dilemma game.

Prisoners Dilemma game has therefore dominant-strategy equilibrium

$$\begin{array}{cccc} {\it Alice} & C_A & D_A \ {\it Bob} & (3,3) & (5,0) \ D_B & (0,5) & (1,1) \end{array}$$

BATTLE of SEX GAME

Alice and Bob have to decide, independently of each other, where to spent the evening.

Alice prefers to go to opera (O), Bob wants to watch TV (T) - tennis.

However, at the same time both of them prefer to be together than to be apart.

Pay-off function is given by the matrix (columns are for Alice) (columns are for Bob)

$$egin{array}{ccc} O & T \ O & (lpha,eta) & (\gamma,\gamma) \ T & (\gamma,\gamma) & (eta,lpha) \end{array}$$

where $\alpha > \beta > \gamma$.

What kind of strategy they should choose?

The two Nash equilibria are (O, O) and (T, T), but players are faced with tactics dilemma, because these equilibria bring them different payoffs.

There are three coins: one fair, with both sides different, and two unfair, one with two heads and one with two tails.

The game proceeds as follows.

- Alice puts coins into a black box and shakes the box.
- Bob picks up one coin.
- Alice wins if coin is unfair, otherwise Bob wins

Clearly, in the classical case, the probability that Alice wins is $\frac{2}{3}$.

- Next goal is to present, using zero-sum games theory, a method how to prove lower bounds for the average running time of randomized algorithms.
- This techniques can be applied to algorithms that terminate for all inputs and all random choices.

A two players zero-sum game is represented by an $n \times m$ payoff-matrix M with all rows and columns summing up to 0.

Payoffs for n possible strategies of Alice are given in rows of M.

Payoffs for m possible strategies of Bob are given in columns of M.

 $M_{i,j}$ is the amount paid by Bob to Alice if Alice chooses strategy *i* and Bob's choice is strategy *j*.

The goal of Alice (Bob) is to maximize (minimize) her payoff.

D 1

Example - stone-scissors-paper game

	Bob				
		Scissors	Paper	Stone	
Alice	Scissors	0	1	-1	
	Paper	-1	0	1	
	Stone	1	-1	0	

PAYOFF-MATRIX

 $\begin{array}{ccc} _ & \to \mathsf{Table} & \mathsf{shows} & \mathsf{how} \\ & & \mathsf{much} & \mathsf{Bob} & \mathsf{has} & \mathsf{to} & \mathsf{pay} \\ & & & \mathsf{to} & \mathsf{Alice} \end{array}$

(Games with players having no information about their opponents' strategies.)

Observe that if Alice chooses a strategy *i*, then she is guaranteed a payoff of $\min_j M_{ij}$ regardless of Bob's strategy.

An optimal strategy O_A for Alice is such an *i* that maximises min_j M_{ij} .

 $O_A = \max_i \min_j M_{ij}$

denotes therefore the lower bound on the value of the payoff Alice gains (from Bob) when she uses an optimal strategy.

An optimal strategy O_B for Bob is such a *j* that minimizes max_i M_{ij} . Bob's optimal strategy ensures therefore that his payoff is at least

$$O_B = \min_i \max_i M_{ij}$$

Theorem

$$O_A = \max_i \min_j M_{ij} \leq \min_i \max_i M_{ij} = O_B$$

Often $O_A < O_B$. In our last (scissors-...) example, $-1 = O_A < O_B = +1$.

If $O_B = O_A$ we say that the game has a solution – a specific choice of strategies that leads to this solution.

 ϱ and γ are so called optional strategies for Alice and Bob if

 $O_A = O_B = M_{\varrho\gamma}$

Example of the game which has a solution $(O_A = O_B = 0)$

0 1 2 -1 0 1 -2 -1 0

What happens if a game has no solution ? There is no clear-cut strategy for any player. Way out: to use randomized strategies.

Alice chooses strategies according to a probability vector $p = (p_1, \ldots, p_n)$; p_i is probability that Alice chooses strategy $s_{A,i}$

Bob chooses strategies according to a probability vector $q = (q_1, \ldots, q_n)$; q_j is a probability that Bob chooses strategy $s_{B,j}$.

Payoff is now a random variable – if p, q are taken as column vectors then

$$E[\mathsf{payoff}] = p^{ op} M q = \sum_{i=1}^n \sum_{j=1}^m p_i M_{ij} q_j$$

Let $O_A(O_B)$ denote the best possible (optimal) lower (upper) bound on the expected payoff of Alice (Bob). Then it holds:

$$O_A = \max_p \min_q p^T M q$$
 $O_B = \min_q \max_p p^T M q$

Theorem (von Neumann Minimax theorem) For any two-person zero-sum game specified by a payoff matrix *M* it holds

$$\max_{p} \min_{q} p^{T} M q = \min_{q} \max_{p} p^{T} M q$$

Observe that once p is fixed, $\max_p \min_q p^T Mq = \min_q \max_p p^T Mq$ is a linear function and is minimized by setting to 1 the q_j with the smallest coefficient in this linear function.

This has interesting/important implications:

If Bob knows the distribution p used by Alice, then his optimal strategy is a pure strategy.

A similar comment applies in the opposite direction. This leads to a simplified version of the minimax theorem, where e_k denotes a unit vector with 1 at the *k*-th position and 0 elsewhere.

Theorem (Loomis' Theorem) For any two-persons zero-sum game

$$\max_{p} \min_{j} p^{T} M e_{j} = \min_{q} \max_{i} e_{i}^{T} M q$$

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Yao's technique provides an application of the game-theoretic results to the establishment of lower bounds for randomized algorithms. For a given algorithmic problem $\mathcal P$ let us consider the following payoff matrix.

		deterministic algorithms $\mathcal{A}_1 \mathcal{A}_2 \mathcal{A}_3$		
1	C 1		Bop –	a designer
NI	-			choosing good algorithms
Ν	<i>c</i> ₂	entries		
Р	C 3	=	Alico	an adversary
U	C 4	recourses	Alice -	an adversary
т	••	resources		choosing bad inputs
1		(i.e. used computation time)		
S				

Pure strategy for Bob corresponds to the choice of a deterministic algorithm. **Optimal pure strategy** for Bob corresponds to a choice of an optimal deterministic algorithm.

YAO'S TECHNIQUE 2/3

 V_B is the worst-case running time of any deterministic algorithm

Problem: How to interpret mixed strategies ?

A *mixed strategy* for Bob is a probability distribution over (always correct) deterministic algorithms—so it is a Las Vegas randomized algorithm.

An optimal mixed strategy for Bob is an optimal Las Vegas algorithm. **Distributional complexity** of a problem is an expected running time of the best deterministic algorithm for the worst distribution on the inputs.

Loomis theorem implies that distributional complexity equals to the least possible time achievable by any randomized algorithm

Reformulation of von Neumann+Loomis' theorem in the language of algorithms

Corollary Let Π be a problem with a finite set I of input instances and \mathcal{A} be a finite set od deterministic algorithms for Π . For any input $i \in I$ and any algorithm $A \in \mathcal{A}$, let T(i, A) denote computation time of A on input i. For probability distributions p over I and q over \mathcal{A} , let i_p denote random input chosen according to p and A_q a random algorithm chosen according to q. Then

$$\max_{p} \min_{q} E\left[T(i_{p}, A_{q})\right] = \min_{q} \max_{p} E\left[T(i_{p}, A_{q})\right]$$
$$\max_{p} \min_{A \in \mathcal{A}} E\left[T(i_{p}, A)\right] = \min_{q} \max_{i \in I} E\left[T(i, A_{q})\right]$$

Consequence:

Theorem(Yao's Minimax Principle) For all distributions p over l and q over A.

 $\min_{A \in \mathcal{A}} \mathbf{E}[T(i_{p}, A)] \leq \max_{i \in I} \mathbf{E}[T(i, A_{q})]$

Interpretation: Expected running time of the optimal deterministic algorithm for an arbitrarily chosen input distribution p for a problem Π is a lower bound on the expected running time of the optimal (Las Vegas) randomized algorithm for Π .

In other words, to determine a lower bound on the performance of all randomized algorithms for a problem P, derive instead a lower bound for any deterministic algorithm for P when its inputs are drawn from a specific probability distribution (of your choice).

Interpretation again Expected running time of the optimal deterministic algorithm for an arbitrarily chosen input distribution p for a problem Π is a lower bound on the expected running time of the optimal (Las Vegas) randomized algorithm for Π .

Consequence:

In order to prove a lower bound on the randomized complexity of an algorithmic problem, it suffices to choose *any* probability distribution p on the input and prove a lower bound on the expected running time of deterministic algorithms for that distribution.

The power of this technique lies in

- \blacksquare the flexibility at the choice of p
- the reduction of the task to determine lower bounds for randomized algorithms to the task to determine lower bounds for deterministic algorithms.

(It is important to remember that we can expect that the deterministic algorithm "knows" the chosen distribution p.)

The above discussion holds for Las Vegas algorithms only!

THE CASE OF MONTE CARLO ALGORITHMS

Let us consider Monte Carlo algorithms with error probability $0 < \varepsilon < \frac{1}{2}$.

Let us define the **distributional complexity with error** ε , notation

$$\min_{A\in\mathcal{A}}E\left[T_{\varepsilon}(I_{p},A)\right],$$

to be the minimum expected time of any deterministic algorithm that errs with probability at most ε under the input I_p with distribution p.

Let us denote by

$$\max_{i\in\mathcal{I}}E\left[\left(T_{\varepsilon}(i,A_{q})\right)\right]$$

the expected time (under the worst input) of any randomized algorithm A_q that errs with probability at most ε .

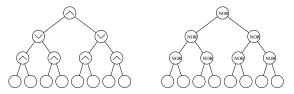
Theorem For all distributions p over inputs and q over Algorithms, and any $\varepsilon \in [0, 1/2]$, it holds

$$\frac{1}{2}(\min_{A\in\mathcal{A}}\mathsf{E}[T_{2\varepsilon}(i_{p},A)]) \leq \max_{i\in\mathcal{I}}\mathsf{E}[T_{\varepsilon}(i,A_{q})]$$

A randomized algorithm for a game-tree T evaluations can be viewed as a probability distribution over deterministic algorithms for T, because the length of computation and the number of choices at each step are finite.

Instead of AND–OR trees of depth 2k we can consider NOR–trees of depth 2k. Indeed, it holds:

 $(a \lor b) \land (c \lor d) \equiv (a \text{ NOR } b) \text{NOR}(c \text{ NOR } d)$



Note: It's important to distinguish between:

- the expected running time of the randomized algorithm with a fixed input (where probability is considered over all random choices made by the algorithm)
- and
- the expected running time of the deterministic algorithm when proving the lower bound (the average time is taken over all random input instances).

LOWER BOUND FOR GAME TREE EVALUATION - I

Assume now that each leaf of a NOR-tree is set up to have value 1 with probability $p = \frac{3-\sqrt{5}}{2}$ (observe that $(1-p)^2 = p$ for such a p).

Observe that if inputs of a NOR-gate have value 1 with probability p then its output value is also 1 with probability (1-p)(1-p) = p.

Consider now only *depth-first pruning algorithms* for tree evaluation. (They are such depth-first algorithms that make use of the knowledge that subtrees that provide no additional useful information can be "pruned away".)

Of importance for the overall analysis is the following technical lemma:

Lemma Let T be a NOR-tree each leaf of which is set to 1 with a fixed probability. Let W(T) denote the minimum, over all deterministic algorithms, of the expected number of steps to evaluate T. Then there is a depth-first pruning algorithm whose expected number of steps to evaluate T is W(T).

The last lemma tells us that for the purposes of our lower bound, we may restrict our attention to the depth-first pruning algorithms.

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For a depth-first pruning algorithm evaluating a NOR-tree, let W(h) be the expected number of leaves the algorithm inspects in determining the value of a node at distance h from the leaves.

It holds

$$W(h) = pW(h-1) + (1-p)2W(h-1) = (2-p)W(h-1)$$

because with the probability 1 - p the first subtree produces 0 and therefore also the second tree has to be evaluated. If $h = \lg_2 n$, then the above recursion has a solution

$$W(h) \geq n^{0.694}$$

This implies:

Theorem The expected running time of any randomized algorithm that always evaluates an instance of T_k correctly is at least $n^{0.694}$, where $n = 2^{2k}$ is the number of leaves.

The upper bound for randomized game tree evaluation algorithms already shown, at the beginning of this chapter was $n^{0.79}$, what is more than the lower bound n^{694} just shown.

It was therefore natural to ask what does the previous theorem really says?

For example, is our lower bound technique weak? ?

No, the above result just says that in order to get a better lower bound another probability distribution on inputs may be needed.

Two recent results put more light on the Game tree evaluation problem.

- It has been shown that for our game tree evaluation problem the upper bound presented at the beginning is the best possible and therefore that $\theta(n^{0.79})$ is indeed the classical (query) complexity of the problem.
- It has also been shown, by Farhi et al. (2009), that the upper bound for the case quantum computation tools can be used is O(n^{0.5}).

The concept of the number of wisdom introduced in the following and related results helped to show that randomness is deeply rooted even in arithmetic.

In order to define numbers of wisdom the concept of self-delimiting programs is needed.

A program represented by a binary word p, is self-delimiting for a computer C, if for any input pw the computer C can recognize where p ends after reading p only.

Another way to see self-delimiting programs is to consider only such programming languages L that no program in L is a prefix of another program in L.

For a universal computer C with only self-delimiting programs, the number of wisdom Ω_C is the probability that randomly constructed program for C halts. More formally

$$\Omega_C = \sum_{p \text{ halts}} 2^{-|p|}$$

where p are (self-delimiting) halting programs for C.

 Ω_C is therefore the probability that a self-delimiting computer program for C generated at random, by choosing each of its bits using an independent toss of a fair coin, will eventually halt.

- \bullet 0 \leq $\Omega_C \leq$ 1
- \blacksquare Ω_C is an uncomputable and random real number.
- At least *n*-bits long theory is needed to determine *n* bits of Ω_C .
- At least *n* bits long program is needed to determine *n* bits of Ω_C
- Bits of Ω can be seen as mathematical facts that are true for no reason.

- Greg Chaitin, who introduced numbers of wisdom, designed a specific universal computer C and a two hundred pages long Diophantine equation E, with 17,000 variables and with one parameter k, such that for a given k the equation E has a finite (infinite) number of solutions if and only if the k-th bit of Ω_C is 0 (is 1). { As a consequence, we have that randomness, unpredictability and uncertainty occur even in the theory of Diophantine equations of elementary arithmetic.}
- Knowing the value of Ω_C with *n* bits of precision allows to decide which programs for *C* with at most *n* bits halt.