Probabilistic Analysis for Randomized Game Tree Evaluation

Tämur Ali Khan and Ralph Neininger

ABSTRACT: We give a probabilistic analysis for the randomized game tree evaluation algorithm of Snir. We first show that there exists an input such that the running time, measured as the number of external nodes read by the algorithm, on that input is maximal in stochastic order among all possible inputs. For this worst case input we identify the exact expectation of the number of external nodes read by the algorithm, give the asymptotic order of the variance including the leading constant, provide a limit law for an appropriate normalization as well as a tail bound estimating large deviations. Our tail bound improves upon the exponent of an earlier bound due to Karp and Zhang, where subgaussian tails were shown based on an approach using multitype branching processes and Azuma's inequality. Our approach rests on a direct, inductive estimate of the moment generating function.

1 Introduction

In this note we analyze the performance of the randomized algorithm to evaluate Boolean decision trees proposed by Snir (1985). Given is a complete binary tree of height 2k, $k \ge 1$, where the root (at depth 0) is labeled \land as are all internal nodes with even depth, all internal nodes with odd depth are labeled \lor . The $n=2^{2k}$ external nodes are labeled either 0 or 1 and the objective is to calculate the value of the root. For each node its value is given as the value of the operation labeled at that node applied to the values of its children. The cost for evaluating the Boolean decision tree is measured as the number of external nodes read by the algorithm.

Snir proposed and analyzed the following randomized algorithm to evaluate a Boolean decision tree: At each node one chooses randomly (with probability 1/2) one of its children and calculates its value recursively. If the result allows to identify the value of the node (that is a 0 for a \land -labeled node and a 1 for a \lor -labeled node, respectively) one is done, otherwise also the other child's value has to be calculated recursively in order to obtain the value of the node. Applying this to the root of the tree yields the value of the Boolean decision tree.

The advantage of this algorithm over any deterministic algorithm is that for any input at the external nodes its expected cost is sublinear in n, whereas any deterministic algorithm has linear worst case cost. More precisely, Saks and Wigderson (1986) obtained that the maximum expected cost is of the order $\Theta(n^{\alpha})$ with $\alpha = \log_2((1+\sqrt{33})/4) \doteq 0.753$ and showed that this is also a lower bound on the maximum expected cost for any other randomized algorithm to evaluate a Boolean decision tree; see also Motwani and Raghavan (1995, Chapter 2) for an account on this subject. Further analysis was given by Karp and Zhang (1995). For certain regular inputs at the external nodes the cost of the algorithm can be represented via 2-type Galton-Watson processes. Karp and Zhang showed that the normalized cost has subgaussian tails. That argument was based on Azuma's inequality.

We denote the input of 0's and 1's at the external nodes as a vector $v \in \{0, 1\}^n$ and the number of external nodes read by the algorithm on input v by C(v). We

will see subsequently that for particular $v^* \in \{0,1\}^n$ not only the expectation of the cost of the algorithm is maximized, i.e., $\mathbb{E} C(v^*) = \max_{v \in \{0,1\}^n} \mathbb{E} C(v)$, but also that $C(v^*)$ is maximal in stochastic order, $C(v) \leq C(v^*)$ for all $v \in \{0,1\}^n$. Here, $X \leq Y$ for random variables X, Y denotes that the corresponding distribution functions F_X, F_Y satisfy $F_X(x) \geq F_Y(x)$ for all $x \in \mathbb{R}$, or, equivalently, that there are realizations X', Y' of the distributions $\mathcal{L}(X), \mathcal{L}(Y)$ of X, Y on a joint probability space such that we pointwise have $X' \leq Y'$.

From this perspective it is reasonable to consider $C(v^*)$ as the worst case complexity of the randomized algorithm and to analyze its asymptotic probabilistic behavior. Our results for the exact mean of $C(v^*)$, the asymptotic growth of its variance including the evaluation of the leading constant, a limit law for $C(v^*)$ after normalization as $k \to \infty$ together with an explicit tail estimate are based on a recursive description of the problem. Since v^* is a regular input in the sense of Karp and Zhang, also their 2-type Galton-Watson approach applies.

Our main finding is an improvement of the tail bound $\exp(-\text{const }t^2)$ for t > 0, to $\exp(-\text{const }t^{\kappa})$, with $1 < \kappa < 1/(1-\alpha) \doteq 4.06$, see Theorem 3.6. This is based on a direct, inductive estimate of the moment generating function. Our approach is also applicable to any regular input as well as to other related problems.

The paper is organized as follows: In section 2 we explain, how a worst case input v^* is obtained. Section 3 contains the statements of the results. In sections 4 and 5 the 2-type branching process of Karp and Zhang (1995) is recalled and the recursive description of the quantities, that our analysis is based on, is introduced. Section 6 contains the proofs of our results and section 7 has extensions to m-ary Boolean decision trees.

2 Worst case input

In this section we explain how a worst case input v^* is constructed. We first have a look at the case k=1 and $v\in\{0,1\}^4$ such that the decision tree is evaluated to 1 at the root. Clearly both children of the root have to lead to an evaluation of 1. Now each pair of external nodes attached to the children needs to have at least one value 1. Note that the algorithm reads in both pairs of external nodes until it finds the first one. Hence there will in total be read two 1's no matter how $v\in\{0,1\}^4$ is drawn among the choices that lead to an evaluation of 1 for the decision tree. Clearly, to maximize the number of 0's being read we choose in each pair of external nodes one 0 and one 1. Then both 0's are being read independently with probability 1/2. Hence, $v_1=(0,1,0,1)$ stochastically maximizes C(v) for all $v\in\{0,1\}^4$ such that the decision tree evaluates 1, see Figure 1.

Analogously look at the case k=1 and $v \in \{0,1\}^4$ such that the decision tree is evaluated to 0. Clearly, one child of the root has to have the value 0, whose external nodes attached need to have both values 0. If we choose also value 0 for the other child of the root, we are lead to v=(0,0,0,0), and the algorithm reads exactly 2 external nodes with values both 0. Therefore, to stochastically maximize C(v) we choose the second child of the root with value 1 and again its external nodes attached with values 0 and 1. Then, $v_0=(0,0,0,1)$ stochastically maximizes C(v) for all $v \in \{0,1\}^4$ for which the decision tree evaluates to 0, see Figure 1.

Since we have $C(v_0) \leq C(v_1)$, it follows that $v^* = (0, 1, 0, 1)$ is a choice with $C(v) \leq C(v^*)$ for all $v \in \{0, 1\}^4$. For general $k \geq 2$ a corresponding $v^* = v^*(k)$ can recursively be constructed from $v^*(k-1)$ as follows: Each component 0 in

 $v^*(k-1)$ is replaced by the block 0, 0, 0, 1, whereas each 1 is replaced by the block 0, 1, 0, 1. For example, for k = 3, this yields

$$\begin{array}{lll} v^{\star} & = & (0,0,0,1,0,0,0,1,0,0,0,1,0,1,0,1,\\ & & 0,0,0,1,0,1,0,1,0,0,0,1,0,1,0,1,\\ & & 0,0,0,1,0,0,0,1,0,0,0,1,0,1,0,1,\\ & & 0,0,0,1,0,1,0,1,0,0,0,1,0,1,0,1,0,1. \end{array}$$

In Proposition 3.1 we show that this construction yields a v^* with $C(v) \leq C(v^*)$ for all $v \in \{0,1\}^n$ and $k \geq 1$.

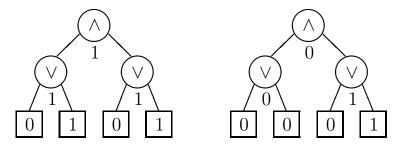


Figure 1: Shown are decision trees for k = 1 evaluating at the root to 1 and 0, respectively, together with a choice for the external nodes that stochastically maximizes the number of external nodes read by the algorithm.

If we would only want to stochastically maximize the cost over all $v \in R_0(n) \subset \{0,1\}^n$ that evaluate to a 0 at the root, the same recursive construction of replacing digits by corresponding blocks, starting with $v_0 = (0,0,0,1)$, yields a $v_* \in R_0(n)$ such that $C(v) \preceq C(v_*)$ for all $v \in R_0(n)$.

3 Results

We assume that we have $n=2^{2k}$ with $k \ge 1$ and denote by $v^* \in \{0,1\}^n$ an input as constructed in section 2.

Proposition 3.1 For $v^* \in \{0,1\}^n$ as defined in section 2 we have $C(v) \leq C(v^*)$ for all $v \in \{0,1\}^n$.

The stochastic worst case behavior $C(v^*)$ of the randomized game tree evaluation algorithm has the following asymptotic properties: The subsequent theorems describe the behavior of mean, variance, limit distribution, and large deviations of $C(v^*)$. For the mean we have:

Theorem 3.2 The expectation of $C(v^*)$ is given by $\mathbb{E} C(v^*) = c_1 n^{\alpha} - c_2 n^{\beta}$, with

$$\alpha = \log_2 \frac{1 + \sqrt{33}}{4}, \quad \beta = \log_2 \frac{1 - \sqrt{33}}{4}, \quad c_1 = \frac{1}{2} + \frac{7}{2\sqrt{33}}, \quad c_2 = c_1 - 1.$$

We denote for sequences (a_k) , (b_k) by $a_k \sim b_k$ asymptotic equivalence, i.e., $a_k/b_k \to 1$ as $k \to \infty$. Then we have for the variance of $C(v^*)$:

Theorem 3.3 The variance of $C(v^*)$ satisfies asymptotically $\operatorname{Var} C(v^*) \sim d n^{2\alpha}$ as $k \to \infty$, where $d \doteq 0.0938$. The constant d can also be given in closed form.

For random variables X, Y we denote by $X \stackrel{d}{=} Y$ equality in distribution, i.e., $\mathcal{L}(X) = \mathcal{L}(Y)$. Then we have the following limit law for $C(v^*)$:

Theorem 3.4 For $C(v^*)$ we have after normalization convergence in distribution,

$$\frac{C(v^{\star})}{n^{\alpha}} \longrightarrow C, \quad k \to \infty,$$

where the distribution of C is given as $\mathcal{L}(C) = \mathcal{L}(G_1)$ and $\mathcal{L}(G) = \mathcal{L}(G_0, G_1)$ is characterized by $\mathbb{E} ||G||^2 < \infty$, $\mathbb{E} G = (c_0, c_1)$, with $c_0 = 1/2 + 5/(2\sqrt{33})$, and

$$G \stackrel{d}{=} \frac{1}{4^{\alpha}} \left\{ G^{(1)} + G^{(2)} + \begin{bmatrix} B_1 B_2 & 0 \\ 1 - B_2 & 0 \end{bmatrix} G^{(3)} + \begin{bmatrix} 0 & B_1 \\ B_1 & 0 \end{bmatrix} G^{(4)} \right\},\,$$

with $G^{(1)}, \ldots, G^{(4)}$, B_1, B_2 independent with $\mathcal{L}(G^{(r)}) = \mathcal{L}(G)$, for $r = 1, \ldots, 4$, and $\mathcal{L}(B_1) = \mathcal{L}(B_2) = B(1/2)$. Here, B(1/2) denotes the Bernoulli(1/2) distribution.

For the estimate of large deviations we rely on Chernoff's bounding technique. We need to follow a bivariate setting for the vector $(C(v^*), C(v_*))$ as introduced in section 5. The following bound on the moment generating function is obtained:

Proposition 3.5 It exists a sequence $(Y_k)_{k\geq 0} = (Y_{k,0}, Y_{k,1})_{k\geq 0}$ of bivariate random variables with marginal distributions $\mathcal{L}((C(v^*) - \mathbb{E}C(v^*))/n^{\alpha})$, $\mathcal{L}((C(v_*) - \mathbb{E}C(v_*))/n^{\alpha})$ such that for all $q > 1/\alpha \doteq 1.33$ there is a K > 0 with

$$\mathbb{E} \exp\langle s, Y_k \rangle \le \exp(K \|s\|^q) \tag{1}$$

for all $s \in \mathbb{R}^2$ and $k \geq 0$. An explicit value for $K = K_q$ is given in (4).

The bound on the moment generating function in the previous proposition implies a large deviation estimate via Chernoff bounds:

Theorem 3.6 For all $1 < \kappa < 1/(1 - \alpha) \doteq 4.06$ there exists an L > 0 such that for any t > 0 and $n = 2^{2k}$

$$\mathbb{P}\left(\frac{C(v^*) - \mathbb{E}C(v^*)}{n^{\alpha}} > t\right) \le \exp(-Lt^{\kappa}). \tag{2}$$

An explicit value for L is given in (5). The same bound applies to the left tail.

The approach of Karp and Zhang (1995) based on Azuma's inequality gives the tail bound $\exp(-L't^2)$ for an explicitly known L'. For $\kappa=2$ the prefactor $L=L_2$ in Theorem 3.6 can also be evaluated and satisfies $L_2>11L'$.

4 Karp and Zhang's 2-type branching process

For the analysis of $C(v^*)$ note that whenever the algorithm has to evaluate the value of a node at a certain depth that yields a 1, according to the discussion of section 2, the algorithm has to evaluate the values of two nodes of depths two levels

below that each yield a 1, and $B_3 + B_4$ nodes of depths two levels below that each yield a 0, cf. Figure 1. Here, B_3 , B_4 are independent Bernoulli B(1/2) distributed random variables. Analogously, when the algorithm has to evaluate the value of a node at a certain depth that yields 0, two levels below it has to evaluate B_1 nodes yielding a 1 and $2 + B_1B_2$ nodes yielding a 0, where B_1 , B_2 are independent B(1/2) distributed random variables. Here, the event $\{B_1 = 1\}$ corresponds to the algorithm first checking the right child of the node to be evaluated and $\{B_2 = 1\}$ to first checking the left child of that child, cf. Figure 1. Since at each node the child being evaluated first is independently drawn from all other choices, this gives rise to the following 2-type Galton-Watson branching process.

We have individuals of type 0 and 1 where the population of the k-th generation corresponds to the number of nodes at depth 2k that are read by the algorithm. We consider processes starting either with an individual of type 1 or type 0 and assume that the algorithm is applied to the worst case inputs v^* and v_* , respectively. Then we have the following offspring distributions: An individual of type 1 has an offspring of 2 individuals of type 1 and $B_3 + B_4$ individuals of type 0. An individual of type 0 has an offspring of B_1 individuals of type 1 and $2 + B_1B_2$ individuals of type 0. We denote the number of individuals of type 0 and 1 in generation k by $(V_n^{(i)}, W_n^{(i)})$, when starting with an individual of type i = 0, 1, where $n = 2^{2k}$. Note that for $v^*, v_* \in \{0, 1\}^n$ we have the representations

$$C(v^{\star}) \stackrel{d}{=} V_n^{(1)} + W_n^{(1)}, \quad C(v_{\star}) \stackrel{d}{=} V_n^{(0)} + W_n^{(0)}.$$

This is the approach of Karp and Zhang (1995) for regular inputs like v^* , v_* . Hence, part of the analysis of $C(v^*)$ can be reduced to the application of the theory of multitype branching processes; see for general reference Harris (1963) and Athreya and Ney (1972), and for a survey on the application of branching processes to tree structures and tree algorithms see Devroye (1998).

However, we will also use a recursive description of the problem. This will be given in the next section and enables to use as well results from the probabilistic analysis of recursive algorithms by the contraction method.

5 The recursive point of view

It is convenient to work as well with a recursive description of the distributions $\mathcal{L}(C(v_{\star}))$ and $\mathcal{L}(C(v^{\star}))$. For this, we define the distributions of a bivariate random sequence $(Z_n) = (Z_{n,0}, Z_{n,1})$ for all $n = 2^{2k}$, $k \ge 0$ by $Z_1 = (1,1)$ and, for $k \ge 1$,

$$Z_n \stackrel{d}{=} Z_{n/4}^{(1)} + Z_{n/4}^{(2)} + \left[\begin{array}{cc} B_1 B_2 & 0 \\ 1 - B_2 & 0 \end{array} \right] Z_{n/4}^{(3)} + \left[\begin{array}{cc} 0 & B_1 \\ B_1 & 0 \end{array} \right] Z_{n/4}^{(4)},$$

where $Z_{n/4}^{(1)}, \ldots, Z_{n/4}^{(4)}$, B_1, B_2 are independent, B_1, B_2 are Bernoulli B(1/2) distributed and $\mathcal{L}(Z_{n/4}^{(1)}) = \cdots = \mathcal{L}(Z_{n/4}^{(4)}) = \mathcal{L}(Z_{n/4})$. It can directly be checked by induction that the marginals of Z_n satisfy $\mathcal{L}(Z_{n,0}) = \mathcal{L}(C(v_*))$ and $\mathcal{L}(Z_{n,1}) = \mathcal{L}(C(v^*))$. Note that $Z_{n,0}$ and $Z_{n,1}$ become dependent, firstly, since we have coupled the offspring distributions using for the second component of Z_n again B_1 and $1 - B_2$ instead of B_3 and B_4 , cf. section 4, and, secondly, since the first component of $Z_{n/4}^{(3)}$ contributes to both components of Z_n . Sequences satisfying

recursive equations as (Z_n) are being dealt with in a probabilistic framework, the so called contraction method; see Rösler (1991, 1992), Rachev and Rüschendorf (1995), Rösler and Rüschendorf (2001), and Neininger and Rüschendorf (2004).

6 Proofs

In this section we sketch the proofs of the results stated in section 3.

Proof of Proposition 3.1: (Sketch) We denote by $R_0(n)$, $R_1(n) \subset \{0,1\}^n$ the sets of vectors at the external nodes at depth 2k that yield an evaluation at the root of the decision tree of value 0 and 1, respectively. From the discussion in section 2 we have

$$C(v) \prec C(v_{\star}), \quad v \in R_0(n), \quad \text{and} \quad C(v) \prec C(v^{\star}), \quad v \in R_1(n).$$

Hence, it remains to show that $C(v_\star) \leq C(v^\star)$. This is shown by induction on $k \geq 1$. For k=1 this can directly be checked. For the step $k-1 \to k$ assume that we have $C(v_\star(k-1)) \leq C(v^\star(k-1))$. It suffices to find realizations of the quantities $(V_n^{(1)}, W_n^{(1)})$ and $(V_n^{(0)}, W_n^{(0)})$ on a joint probability space with $V_n^{(0)} + W_n^{(0)} \leq V_n^{(1)} + W_n^{(1)}$ almost surely, $n=2^{2k}$. For this we use B, B', $(V_{n/4}^{(i),j}, W_{n/4}^{(i),j})$ for $i=1,2,j=1,\ldots,4$ being independictly in the sum of the sum

For this we use $B, B', (V_{n/4}^{(i),j}, W_{n/4}^{(i),j})$ for $i=1,2, j=1,\ldots,4$ being independent for each i=0,1 and with B, B' Bernoulli B(1/2) distributed, $\mathcal{L}(V_{n/4}^{(i),j}) = \mathcal{L}(V_{n/4}^{(i)})$, $\mathcal{L}(W_{n/4}^{(i),j}) = \mathcal{L}(W_{n/4}^{(i)})$ for i=1,2 and $j=1,\ldots,4$. By the induction hypothesis we may assume that we have versions of these random variates with $V_{n/4}^{(0),j} + W_{n/4}^{(0),j} \leq V_{n/4}^{(1),j} + W_{n/4}^{(1),j}$ for $j=1,\ldots,4$. With this coupling we define $(V_n^{(1)}, W_n^{(1)})$ and $(V_n^{(0)}, W_n^{(0)})$ according to the values of B, B': On $\{B=1, B'=0\}$ we set

$$\begin{pmatrix} V_{n}^{(0)} \\ W_{n}^{(0)} \end{pmatrix} := \begin{pmatrix} V_{n/4}^{(0),2} \\ W_{n/4}^{(0),2} \end{pmatrix} + \begin{pmatrix} V_{n/4}^{(0),3} \\ W_{n/4}^{(0),3} \end{pmatrix} + BB' \begin{pmatrix} V_{n/4}^{(0),4} \\ W_{n/4}^{(0),4} \end{pmatrix} + B \begin{pmatrix} V_{n/4}^{(1),1} \\ W_{n/4}^{(1),1} \end{pmatrix},$$

$$\begin{pmatrix} V_{n}^{(1)} \\ W_{n}^{(1)} \end{pmatrix} := B \begin{pmatrix} V_{n/4}^{(0),3} \\ W_{n/4}^{(0),3} \end{pmatrix} + \begin{pmatrix} V_{n/4}^{(1),1} \\ W_{n/4}^{(1),1} \end{pmatrix} + B' \begin{pmatrix} V_{n/4}^{(0),4} \\ W_{n/4}^{(0),4} \end{pmatrix} + \begin{pmatrix} V_{n/4}^{(1),2} \\ W_{n/4}^{(1),2} \end{pmatrix},$$

and obtain $V_n^{(0)} + W_n^{(0)} \leq V_n^{(1)} + W_n^{(1)}$. On the remaining sets $\{B=0, B'=0\}$, $\{B=0, B'=1\}$, and $\{B=1, B'=1\}$ similar couplings of $(V_n^{(0)}, W_n^{(0)})$, $(V_n^{(1)}, W_n^{(1)})$ can be defined with $V_n^{(0)} + W_n^{(0)} \leq V_n^{(1)} + W_n^{(1)}$.

Proof of Theorem 3.2: (Sketch) Assume that a generation has (w_0, w_1) individuals of type 0 and 1. Then, by the definition on the offspring distribution in section 4, the expected number of individuals in the subsequent generation is given by

$$M \begin{pmatrix} w_0 \\ w_1 \end{pmatrix}, \quad M := \left[egin{array}{cc} 9/4 & 1 \\ 1/2 & 2 \end{array} \right].$$

Since $C(v^*) = C(v^*(k))$ is the sum of the individuals at generation k for the process started with an individual of type 1 we obtain

$$\mathbb{E} C(v^*) = (1,1)M^k \begin{pmatrix} 0 \\ 1 \end{pmatrix}.$$

The matrix M has the eigenvalues $\lambda_1 = (17 + \sqrt{33})/8$ and $\lambda_2 = (17 - \sqrt{33})/8$ and its k-th power can be evaluated to

$$M^k = \frac{1}{2\sqrt{33}} \left[\begin{array}{cc} (\sqrt{33} + 1)\lambda_1^k + (\sqrt{33} - 1)\lambda_2^k & 8(\lambda_1^k - \lambda_2^k) \\ 4(\lambda_1^k - \lambda_2^k) & (\sqrt{33} - 1)\lambda_1^k + (\sqrt{33} + 1)\lambda_2^k \end{array} \right].$$

From this, $\mathbb{E} C(v^*)$ and various constants needed subsequently can be read off. Note, that $\lambda_1^k = n^{\alpha}$ with α given in Theorem 3.2 and $n = 2^{2k}$.

Before proving Theorem 3.3 it is convenient to first prove Theorem 3.4.

Proof of Theorem 3.4: (Sketch) The 2-type branching process defined in section 4 is supercritical, nonsingular, and positive regular. Hence, a theorem of Harris (1963) implies that

$$\frac{1}{n^{\alpha}} \begin{pmatrix} V_n^{(1)} \\ W_n^{(1)} \end{pmatrix} \longrightarrow Y \begin{pmatrix} \nu_1 \\ \nu_2 \end{pmatrix}$$

almost surely, as $k \to \infty$, where Y is a nonnegative random variable and (ν_1, ν_2) a deterministic vector that could also be further specified. Thus we obtain

$$\frac{C(v^{\star})}{n^{\alpha}} \longrightarrow C$$

in distribution, as $k \to \infty$, with $\mathcal{L}(C) = \mathcal{L}((\nu_1 + \nu_2)Y)$.

On the other hand the recursive formulation of section 5 leads after the normalization $X_n := Z_n/n^{\alpha}$ to

$$X_n \stackrel{d}{=} \sum_{r=1}^4 A_r X_{n/4}^{(r)},$$

for $k \geq 1$, where $A_1 = A_2 = (1/4^{\alpha})I_2$, with the 2×2 identity matrix I_2 , and

$$A_3 = \frac{1}{4^{\alpha}} \begin{bmatrix} B_1 B_2 & 0\\ 1 - B_2 & 0 \end{bmatrix}, \quad A_4 = \frac{1}{4^{\alpha}} \begin{bmatrix} 0 & B_1\\ B_1 & 0 \end{bmatrix}, \tag{3}$$

where $X_{n/4}^{(1)}, \ldots, X_{n/4}^{(4)}$, B_1, B_2 are independent with $\mathcal{L}(X_{n/4}^{(r)}) = \mathcal{L}(X_{n/4})$ for $r = 1, \ldots, 4$ and $\mathcal{L}(B_1) = \mathcal{L}(B_2) = B(1/2)$. It follows from the contraction method that X_n converges weakly and with all mixed second moments to some G, that can be characterized as in Theorem 3.4. For details, how to apply the contraction method, see Theorem 4.1 in Neininger (2001). Thus, we have $C(v^*)/n^{\alpha} \to G_1$ in distribution. \blacksquare

Proof of Theorem 3.3: (Sketch) As shown in the proof of Theorem 3.4 we have the convergence $X_n = Z_n/n^\alpha \to G$ for all mixed second moments. This, in particular, implies $\operatorname{Var} X_{n,1} \to \operatorname{Var} G_1$. The variances of G_1 can be obtained from the distributional identity for G stated in Theorem 3.4. Then we obtain $\operatorname{Var} C(v^\star) = \operatorname{Var}(n^\alpha X_{n,1}) \sim dn^{2\alpha}$ with $d = \operatorname{Var} G_1$.

Proof of Proposition 3.5: For $Y_n = (1/n^{\alpha})(Z_n - \mathbb{E} Z_n)$ we have marginals $\mathcal{L}(Y_{n,1}) = \mathcal{L}((C(v^{\star}) - \mathbb{E} C(v^{\star}))/n^{\alpha})$ and $\mathcal{L}(Y_{n,0}) = \mathcal{L}((C(v_{\star}) - \mathbb{E} C(v_{\star}))/n^{\alpha})$. The distributional recurrence for Z_n from section 5 implies the relation

$$Y_n \stackrel{d}{=} \sum_{r=1}^4 A_r Y_{n/4}^{(r)} + b_n, \quad k \ge 1,$$

with $Y_{n/4}^{(1)}, \ldots, Y_{n/4}^{(4)}, B_1, B_2$ independent, $\mathcal{L}(Y_{n/4}^{(r)}) = \mathcal{L}(Y_{n/4})$, for $r = 1, \ldots, 4, \mathcal{L}(B_1) = \mathcal{L}(B_2) = B(1/2)$ and $b_n = (1/n^{\alpha})(4^{\alpha} \sum_{r=1}^{4} (A_r \mathbb{E} Z_{n/4}) - \mathbb{E} Z_n)$. The matrices A_r are given in (3).

We prove the assertion by induction on k. For k=0 we have $Y_0=0$, thus the assertion is true. Assume the assertion is true for some $n/4=2^{2(k-1)}$. Then, conditioning on (A_1,\ldots,A_4,b_n) , denoting the distribution of this vector by σ_n , and using the induction hypothesis, we obtain

$$\mathbb{E} \exp\langle s, Y_n \rangle = \int \exp\langle s, \beta_n \rangle \prod_{r=1}^4 \mathbb{E} \exp\langle s, a_r Y_{n/4} \rangle d\sigma_n(a_1, \dots, a_4, \beta_n)$$

$$\leq \int \exp\langle s, \beta_n \rangle \prod_{r=1}^4 \exp(K \| a_r^T s \|^q) d\sigma_n(a_1, \dots, a_4, \beta_n)$$

$$\leq \int \exp\left(\langle s, \beta_n \rangle + K \| s \|^q \sum_{r=1}^4 \| a_r \|_{\text{op}}^q \right) d\sigma_n(a_1, \dots, a_4, \beta_n)$$

$$= \mathbb{E} \exp(\langle s, b_n \rangle + K \| s \|^q U) \exp(K \| s \|^q),$$

with $U := \sum_{r=1}^{4} (\|A_r\|_{\text{op}}^q) - 1 = 4^{-\alpha q} (2 + B_1 B_2 + (1 - B_2) + B_1) - 1$ and $\|A\|_{\text{op}} = \sup_{\|x\|=1} \|Ax\|$ for matrices A. Hence, the proof is completed by showing

$$\sup_{k>1} \mathbb{E} \exp(\langle s, b_n \rangle + K ||s||^q U) \le 1,$$

for some appropriate K > 0. We denote $\xi := -\operatorname{ess\,sup} U = 1 - 4^{1-\alpha q}$, thus $q > 1/\alpha$ implies $\xi > 0$.

Small ||s||: First we consider small ||s|| with $||s|| \le c/\sup_{k\ge 1} ||b_n||_{2,\infty}$ for some c>0, where $||b_n||_{2,\infty}:=|||b_n|||_{\infty}$, the inner norm being the Euclidean norm. Note that throughout we have $n=n(k)=2^{2k}$. For these small ||s|| we have

$$\mathbb{E} \exp((\langle s, b_n \rangle + K \| s \|^q U) \le \exp(-K \| s \|^q \xi) \mathbb{E} \exp(s, b_n)$$

and, with $\mathbb{E}\langle s, b_n \rangle = 0$,

$$\mathbb{E} \exp\langle s, b_n \rangle = \mathbb{E} \left[1 + \langle s, b_n \rangle + \sum_{k=2}^{\infty} \frac{\langle s, b_n \rangle^k}{k!} \right]$$

$$= 1 + \mathbb{E} \langle s, b_n \rangle^2 \sum_{k=2}^{\infty} \frac{\langle s, b_n \rangle^{k-2}}{k!}$$

$$\leq 1 + \|s\|^2 \mathbb{E} \|b_n\|^2 \sum_{k=2}^{\infty} \frac{c^{k-2}}{k!}$$

$$= 1 + \|s\|^2 \mathbb{E} \|b_n\|^2 \frac{e^c - 1 - c}{c^2}.$$

Using $\exp(-K\|s\|^q\xi) \le 1/(1+K\|s\|^q\xi)$ and with $\Psi(c) = (e^c-1-c)/c^2$ we obtain

$$\mathbb{E} \exp(\langle s, b_n \rangle + K \|s\|^q U) \le \frac{1 + \|s\|^2 \mathbb{E} \|b_n\|^2 \Psi(c)}{1 + K \|s\|^q \xi}.$$

Hence, we have to choose K with

$$K \ge \frac{\|s\|^{2-q}\Psi(c)}{\xi} \sup_{k\ge 1} \mathbb{E} \|b_n\|^2.$$

With $\|s\| \le c/\sup_{k \ge 1} \|b_n\|_{2,\infty}$ a possible choice is

$$K = \frac{\sup_{k \ge 1} \mathbb{E} \|b_n\|^2}{\sup_{k > 1} \|b_n\|_{2,\infty}^{2-q}} \frac{\Psi_q(c)}{\xi},$$

with $\Psi_q(c) = (e^c - 1 - c)/c^q$.

Large ||s||: For general $s \in \mathbb{R}^2$ we have

$$\langle s, b_n \rangle + K \|s\|^q U \le \|s\| \|b_n\| - \|s\|^q K \xi \le \|s\| \|b_n\|_{2,\infty} - \|s\|^q K \xi,$$

and this is less than zero if

$$||s||^{q-1} \ge \frac{\sup_{k \ge 1} ||b_n||_{2,\infty}}{K\xi} = \frac{\sup_{k \ge 1} ||b_n||_{2,\infty}^{3-q}}{\sup_{k \ge 1} \mathbb{E} ||b_n||^2 \Psi_q(c)}.$$

If ||s|| satisfies the latter inequality we call it large. Thus, for large ||s|| we have $\sup_{k\geq 1}\mathbb{E}\exp(\langle s,b_n\rangle+K||s||^qU)\leq 1.$

In order to overlap the regions for small and large ||s|| we need

$$\Psi_1(c) \ge \frac{\sup_{k \ge 1} \|b_n\|_{2,\infty}^2}{\sup_{k > 1} \mathbb{E} \|b_n\|^2}.$$

The right hand side of the latter display can be evaluated explicitly for our problem and equals 104/77. Thus, this inequality is true for, e.g., c=1.53. Hence, with the explicit value

$$K := K_q = \frac{\sup_{k \ge 1} \mathbb{E} \|b_n\|^2}{\sup_{k \ge 1} \|b_n\|_{2,\infty}^{2-q}} \frac{e^{1.53} - 2.53}{1.53^q (1 - 4^{1-q\alpha})}$$
(4)

the proof is completed. ■

Proof of Theorem 3.6: By Chernoff's bounding technique we have, for u > 0 and with Proposition 3.5,

$$\mathbb{P}\left(\frac{C(v^{\star}) - \mathbb{E}C(v^{\star})}{n^{\alpha}} > t\right) = \mathbb{P}(\exp(uY_{n,1}) > \exp(ut)))$$

$$\leq \mathbb{E}\exp(uY_{n,1} - ut)$$

$$= \mathbb{E}\exp(\langle (0, u), Y_n \rangle - ut)$$

$$\leq \exp(K_q u^q - ut),$$

for all q, K_q as in Proposition 3.5 and (4). Minimizing over u>0 we obtain the bound

$$\mathbb{P}\left(\frac{C(v^*) - \mathbb{E}C(v^*)}{n^{\alpha}} > t\right) \le \exp(-Lt^{\kappa}),$$

for $1 < \kappa < 1/(1 - \alpha)$, with

$$L = L_{\kappa} = K_{\kappa/(\kappa - 1)}^{1 - \kappa} \frac{(\kappa - 1)^{\kappa - 1}}{\kappa^{\kappa}}$$
 (5)

and $K_{\kappa/(\kappa-1)}$ given in (4). This completes the tail bound.

7 m-ary Boolean decision trees

The analysis can be carried over to the case of m-ary Boolean decision trees. The algorithm visits randomly chosen children and evaluates recursively their values until the value of the root can be identified, the remaining children are discarded afterwards. A worst case input $v^* \in \{0,1\}^n$ with $n=m^{2k}$ can be constructed similarly. Then we have similar results for $C(v^*)$:

Theorem 7.1 For the worst case complexity $C(v^*)$ of evaluating an m-ary Boolean decision tree we have the following asymptotics:

$$\mathbb{E} C(v^{\star}) = c_1^{(m)} n^{\alpha_m} + c_2^{(m)} n^{\beta_m},$$

$$\operatorname{Var} C(v^{\star}) \sim d_m n^{2\alpha_m},$$

$$\frac{C(v^{\star})}{n^{\alpha_m}} \rightarrow C_m,$$

$$\mathbb{P} \left(\frac{C(v^{\star}) - \mathbb{E} C(v^{\star})}{n^{\alpha_m}} > t \right) \leq \exp(-L^{(m)} t^{\kappa}), \quad t > 0,$$

with constants $c_1^{(m)}$, α_m , β_m , d_m , $L^{(m)} > 0$, $c_2^{(m)} \in \mathbb{R}$, and $1 < \kappa < \kappa_m = 1/(1-\alpha_m)$.

Numerical values for α_m , d_m and κ_m are listed in Table 1. The distribution of C_m is given as $\mathcal{L}(C_m) = \mathcal{L}(G_1)$ and $\mathcal{L}(G) = \mathcal{L}(G_0, G_1)$ is characterized by $\mathbb{E} \|G\|^2 < \infty$,

$$\mathbb{E} G = (c_0^{(m)}, c_1^{(m)})$$
 and

$$G \stackrel{d}{=} \frac{1}{m^{2\alpha_m}} \left\{ \sum_{r=1}^m G^{(r)} + \sum_{r=1}^{m-1} \begin{bmatrix} 0 & \mathbf{1}_r(U_0) \\ \mathbf{1}_r(U_0) & 0 \end{bmatrix} \bar{G}^{(r)} + \sum_{r,\ell=1}^{m-1} \begin{bmatrix} \mathbf{1}_r(U_0)\mathbf{1}_\ell(U_r) & 0 \\ 1 - \mathbf{1}_\ell(U_r) & 0 \end{bmatrix} G^{(r,\ell)} \right\},$$

with $\mathcal{L}(G^{(r)}) = \mathcal{L}(\bar{G}^{(r)}) = \mathcal{L}(G^{(r,\ell)}) = \mathcal{L}(G)$ and $G^{(r)}$, $\bar{G}^{(r)}$, $G^{(r,\ell)}$, U_r independent with $\mathcal{L}(U_r) = \text{unif}\{0,\ldots,m-1\}$ for all r,ℓ . Here, we denote $\mathbf{1}_i(Y) := \mathbf{1}_{\{i \leq Y\}}$ for integer i and a random variable Y, and we have

$$c_0^{(m)} = \frac{1}{2} + \frac{m+3}{2\sqrt{16m+(m-1)^2}}, \quad c_1^{(m)} = \frac{1}{2} + \frac{3m+1}{2\sqrt{16m+(m-1)^2}}.$$

m	2	3	4	5	6	7	8
α_m	0.754	0.759	0.765	0.769	0.774	0.778	0.781
d_m	0.0938	0.0847	0.0782	0.0731	0.0689	0.0652	0.0619
κ_m	4.060	4.154	4.247	4.336	4.419	4.497	4.571
m	9	10	11	12	13	14	15
α_m	0.785	0.788	0.790	0.793	0.795	0.798	0.800
d_m	0.0590	0.0564	0.0541	0.0519	0.0499	0.0481	0.0464
-	0.0590 4.641	0.0564 4.707		0.0519 4.829	0.0499 4.886	0.0481 4.940	0.0464 4.993
d_m			0.0541				
d_m			0.0541				
$\frac{d_m}{\kappa_m}$	4.641	4.707	0.0541 4.769	4.829	4.886	4.940	4.993
d_m κ_m	4.641	4.707	0.0541 4.769 20	4.829	4.886	4.940	4.993

Table 1: Numerical values of the quantities α_m , d_m and κ_m appearing in Theorem 7.1 for various values of m.

References

- [1] Athreya, K. B. and Ney, P. (1972) Branching processes. Die Grundlehren der mathematischen Wissenschaften, Bd. 196, Springer-Verlag, New York-Heidelberg.
- [2] Devroye, L. (1998) Branching processes and their applications in the analysis of tree structures and tree algorithms. *Probabilistic methods for algorithmic discrete mathematics*, 249–314, Algorithms Combin., 16, Springer, Berlin.
- [3] Harris, T. E. (1963) The theory of branching processes. Die Grundlehren der Mathematischen Wissenschaften, Bd. 119, Springer-Verlag, Berlin; Prentice-Hall, Inc., Englewood Cliffs, N.J.

- [4] Karp, R. and Zhang, Y. (1995) Bounded branching process and AND/OR tree evaluation. *Random Structures Algorithms* 7, 97–116.
- [5] Motwani, R. and Raghavan, P. (1995) Randomized algorithms. Cambridge University Press, Cambridge.
- [6] Neininger, R. (2001) On a multivariate contraction method for random recursive structures with applications to Quicksort. Random Structures Algorithms 19, 498–524.
- [7] Neininger, R. and Rüschendorf, L. (2004) A general limit theorem for recursive algorithms and combinatorial structures. *Ann. Appl. Probab.* 14, 378–418.
- [8] Rachev, S. T. and Rüschendorf, L. (1995). Probability metrics and recursive algorithms. Adv. in Appl. Probab. 27, 770–799.
- [9] Rösler, U. (1991). A limit theorem for "Quicksort". RAIRO Inform. Théor. Appl. 25, 85–100.
- [10] Rösler, U. (1992). A fixed point theorem for distributions. *Stochastic Process.* Appl. 42, 195–214.
- [11] Rösler, U. and Rüschendorf, L. (2001). The contraction method for recursive algorithms. Algorithmica 29, 3–33.
- [12] Saks, M. and Wigderson, A. (1986) Probabilistic boolean decision trees and the complexity of evaluating game trees. *Proceedings of the 27th Annual IEEE Symposium on Foundations of Computer Science*, 29–38, Toronto, Ontario.
- [13] Snir, M. (1985) Lower bounds on probabilistic linear decision trees. Theoret. Comput. Sci. 38, 69–82.

Tämur Ali Khan and Ralph Neininger

Department of Mathematics
J.W. Goethe University
Robert-Mayer-Str. 10
60325 Frankfurt a.M.
Germany
{alikhan, neiningr}@ismi.math.uni-frankfurt.de