Robust Algorithms for the Secretary Problem

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— Abstract

In classical secretary problems, a sequence of n elements arrive in a uniformly random order, and we want to choose a single item, or a set of size K. The random order model allows us to escape from the strong lower bounds for the adversarial order setting, and excellent algorithms are known in this setting. However, one worrying aspect of these results is that the algorithms overfit to the model: they are not very robust. Indeed, if a few "outlier" arrivals are adversarially placed in the arrival sequence, the algorithms perform poorly. E.g., Dynkin's popular 1/e-secretary algorithm is sensitive to even a single adversarial arrival: if the adversary gives one large bid at the beginning of the stream, the algorithm does not select any element at all.

We investigate a robust version of the secretary problem. In the *Byzantine Secretary* model, we have two kinds of elements: green (good) and red (rogue). The values of all elements are chosen by the adversary. The green elements arrive at times uniformly randomly drawn from [0, 1]. The red elements, however, arrive at adversarially chosen times. Naturally, the algorithm does not see these colors: how well can it solve secretary problems?

We show that selecting the highest value red set, or the single largest green element is not possible with even a small fraction of red items. However, on the positive side, we show that these are the only bad cases, by giving algorithms which get value comparable to the value of the optimal green set *minus the largest green item*. (This benchmark reminds us of regret minimization and digital auctions, where we subtract an additive term depending on the "scale" of the problem.) Specifically, we give an algorithm to pick K elements, which gets within $(1 - \varepsilon)$ factor of the above benchmark, as long as $K \ge \text{poly}(\varepsilon^{-1} \log n)$. We extend this to the knapsack secretary problem, for large knapsack size K.

For the single-item case, an analogous benchmark is the value of the second-largest green item. For value-maximization, we give a poly $\log^* n$ -competitive algorithm, using a multi-layered bucketing scheme that adaptively refines our estimates of second-max over time. For probability-maximization, we show the *existence* of a good randomized algorithm, using the minimax principle.

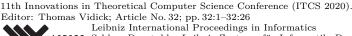
We hope that this work will spur further research on robust algorithms for the secretary problem, and for other problems in sequential decision-making, where the existing algorithms are not robust and often tend to overfit to the model.

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1 Introduction

In sequential decision-making, we have to serve a sequence of requests *online*, i.e., we must serve each request before seeing the next one. E.g., in online auctions and advertising, given a sequence of arriving buyers, we want to choose a high bidder. Equivalently, given a sequence of n numbers, we want to choose the highest of these. The worst-case bounds for this problem are bleak: choosing a random buyer is the best we can do. So we make (hopefully reasonable) stochastic assumptions about the input stream, and give algorithms that work well under those assumptions.

A popular assumption is that the values/bids are chosen by an adversary, but presented to the algorithm in a uniformly random order. This gives the *secretary* or the *random-order model*, under which we can get much better results. E.g., Dynkin's secretary algorithm that selects the first prefix-maximum bidder after discarding the first 1/e-fraction of the bids, selects the highest bid with probability 1/e [13]. The underlying idea – of fixing one or more thresholds after seeing some prefix of the elements – can be generalized to solve classes of packing linear programs near-optimally [8, 9, 25, 21], and to get $O(\log \log n)$ -competitive algorithms for matroids [28, 16] in the random-order model.

However, the assumption that we see the elements in a uniformly random order is quite strong, and most current algorithms are not robust to small perturbations to the model. E.g., Dynkin's algorithm is sensitive to even a single adversarial corruption: if the adversary gives one large bid at the beginning of the stream, the algorithm does not select any buyer at all, even if the rest of the stream is perfectly random! Many other algorithms in the secretary model suffer from similar deficiencies, which suggests that we may be over-fitting to the assumptions of the model.

We propose the *Byzantine secretary model*, where the goal is to design algorithms robust to outliers and adversarial changes. The use of the term "Byzantine" parallels its use in distributed systems, where some of the input is well-behaved while the rest is arbitrarily corrupted by an adversary. Alternatively, our model can be called *semi-random* or *robust*: these other terms are used in the literature which inspires our work. Indeed, there is much interest currently in designing stochastic algorithms that are robust to adversarial noise (see [10, 34, 29, 3, 34, 11, 12, 14, 30] and references therein). Our work seeks to extend robustness to online problems. Our work is also related in spirit to investigations into how much randomness in the stream is necessary and sufficient to get competitive algorithms [5, 23].

1.1 Our Model

In the secretary problem, n elements arrive one-by-one. Each item has a value that is revealed upon its arrival, which happens at a time chosen independently and *uniformly at random* in [0, 1]. (We choose the continuous time model, instead of the *uniformly random* arrival order model, since the independence allows us to get clean proofs.) When we see an item, we must either *select* it or discard it before we see the next item. Our decisions are *irrevocable*. We can select at most K elements, where K = 1 for the classical version of the problem. We typically want to maximize the expected total value of the selected elements where the value of a set is simply the sum of values of individual elements. (For the single-item case we may also want to maximize the probability of selecting the highest-value item, which is called the *ordinal case*.) Given its generality and wide applicability, this model and its extensions are widely studied; see §1.3.

The difference between the classical and Byzantine secretary models is in how the sequence is generated. In both models, the adversary chooses the values of all n elements. In the classical model, these are then permuted in a random order (say by choosing the arrival times independently and uniformly at random (u.a.r.) from [0,1]). In the Byzantine model, the elements are divided into two groups: the green (or good) elements/items G, and the red (or rogue/bad) elements/items R. This partition and the colors are not visible to the algorithm. Now elements in G arrive at independently chosen u.a.r. times between [0, 1], but those in R arrive at times chosen by the adversary. Faced with this sequence, the algorithm must select some subset of elements (say, having size at most K, or more generally belonging to some down-closed family). The precise order of actions is important:

- First, the adversary chooses values of elements in $R \cup G$, and the arrival times of elements in R.
- Then each element $e \in G$ is independently assigned a uniformly random arrival time $t_e \sim U[0, 1]$.

Hence the adversary is powerful and strategic, and can "stuff" the sequence with values in an order that fools our algorithms the most. The green elements are non-strategic (hence are in random order) and beyond the adversary's control. When an element is presented, the algorithm does not see the color (green vs. red): it just sees the value and the time of arrival. We assume that the algorithm knows n := |R| + |G|, but not |R| or |G|; see Appendix B on how to relax this assumption. The green elements are denoted $G = \{g_{\max} = g_1, g_2, \ldots, g_{|G|}\}$ in non-increasing order of values.

What results can we hope to get in this model? Here are two cautionary examples:

- Since the red elements behave completely arbitrarily, the adversary can give non-zero values to only the reds, and plant a bad example for the adversarial order using them. Hence, we cannot hope to get the value of the optimal red set in general, and should aim to get value only from the greens.
- Moreover, suppose essentially all the value among the greens is concentrated in a single item g_{max} . Here's a bad example: the adversary gives a sequence of increasing reds, all having value much smaller than g_{max} , but values which are very far from each other. When the algorithm does see the green item, it will not be able to distinguish it from the next red, and hence will fail. This is formalized in Observation 19. Hence, to succeed, the green value must be spread among more than one item.

Given these examples, here is the "leave-one-out" benchmark we propose:

 $V^* :=$ value of the best feasible green set from $G \setminus g_{\text{max}}$.

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This benchmark is at least as strong as the following guarantee:

(value of best feasible green set from G) $-v(g_{\text{max}})$. (2)

The advantage of (1) over (2) is that V^* is interesting even when we want to select a single item, since it asks for value v_{g_2} or higher.

We draw two parallels to other commonly used benchmarks. Firstly, the perspective (2) suggests the regret-type guarantees, where we seek the best solution in hindsight, minus the "scale of the problem instance". The value of g_{max} is the scale of the instance here. Secondly, think of the benchmark (1) as assuming the existence of at least two high bids, then the second-largest element is almost as good a benchmark as the top element. This is a commonly used assumption, e.g., in digital goods auctions [4].

Finally, if we really care about a benchmark that includes g_{max} , our main results for selecting multiple items (Theorem 1 and Theorem 2) continue to hold, under the (mild?) assumption that the algorithm starts with a polynomial approximation to $v(g_{\text{max}})$.

1.2 Our Results

We first consider the setting where we want to select at most K elements to maximize the expected total value. In order to get within $(1 + \varepsilon)$ factor of the benchmark V^* defined in (1), we need to assume that we have a "large budget", i.e., we are selecting a sufficiently large number of elements. Indeed, having a larger budget K allows us to make some mistakes and yet get a good expected value.

▶ **Theorem 1** (Uniform Matroids). There is an algorithm for Byzantine secretary on uniform matroids of rank $K \ge poly(\varepsilon^{-1} \log n)$ that is $(1 + \varepsilon)$ -competitive with the benchmark V^* .

For the standard version of the problem, i.e. without any red elements, [26] gave an algorithm that achieves the same competitiveness when $K \ge \Omega(1/\varepsilon^2)$. The algorithm from [26] uses a single threshold, that it updates dynamically; we extend this idea to having several thresholds/budgets that "cascade down" over time; we sketch the main ideas in §2.2. In fact, we give a more general result – an algorithm for the *knapsack* setting where each element has a size in [0, 1], and the total size of elements we can select is at most K. (The uniform-matroids case corresponds to all sizes being one.) Whereas the main intuition remain unchanged, the presence of non-uniform sizes requires a little more care.

▶ **Theorem 2** (Knapsack). There is an algorithm for Byzantine secretary on knapsacks with size at least $K \ge \text{poly}(\varepsilon^{-1} \log n)$ (and elements of at most unit size) that is $(1+\varepsilon)$ -competitive with the benchmark V^* .

As mentioned earlier, under mild assumptions the guarantee in Theorem 2 can be extended against the stronger benchmark that includes g_{max} . Formally, assuming the algorithm starts with a poly(m)-approximation to the value of g_{max} , we get a $(1 + \varepsilon)$ -competitive algorithm for $K \geq \text{poly}(\varepsilon^{-1} \log(mn))$ against the stronger benchmark.

Selecting a Single Item. What if we want to select a single item, to maximize its expected value? Note that the benchmark V^* is now the value of g_2 , the second-largest green item. Our main result for this setting is the following, where $\log^* n$ denotes the iterated logarithm:

▶ **Theorem 3** (Value Maximization Single-Item). There is a randomized algorithm for the value-maximization (single-item) Byzantine secretary problem which gets an expected value at least $(\log^* n)^{-2} \cdot V^*$.

Interestingly, our result is unaffected by the corruption level, and works even if just two elements g_{max}, g_2 are green, and every other item is red. This is in contrast to many other robustness models where the algorithm's performance decays with the fraction of bad elements [14, 3, 12, 30]. Moreover, our algorithms do not depend on the fraction of bad items. Intuitively, we obtain such strong guarantees because the adversary has no incentive to present too many bad elements with large values, as otherwise an algorithm that selects a random element would have a good performance.

In the classical setting, the proofs for the value-maximization proceed via showing that the best item itself is chosen with constant probability. Indeed, in that setting, the competitiveness of value-maximization and probability-maximization versions is the same. We do not know of such a result in the Byzantine model. However, we can show a non-trivial performance for the probability-maximization (ordinal) problem:

► Theorem 4 (Ordinal Single-item Algorithm). There is a randomized algorithm for the ordinal Byzantine secretary which selects an element of value at least the second-largest green item with probability $\Omega(1/\log^2 n)$.

Other Settings. Finally, we consider some other constraint sets given by matroids. In (simple) partition matroids, the universe U is partitioned into r groups, and the goal is to select one item from each group to maximize the total value. If we were to set the benchmark to be the sum of second-largest green items from each group, we can just run the single-item algorithm from Theorem 1 on each group independently. But our benchmark V^* is much higher: among the items $g_2, \ldots, g_{|G|}$, the set V^* selects the largest one from each group. Hence, we need to get the largest green item from r - 1 groups! Still, we do much better than random guessing.

▶ **Theorem 5** (Partition Matroids). There is an algorithm for Byzantine secretary on partition matroids that is $O(\log \log n)^2$ -competitive with the benchmark V^* .

Finally, we record a simple but useful logarithmic competitive ratio for arbitrary matroids (proof in §6.2), showing how to extend the corresponding result from [1] for the non-robust case.

▶ **Observation 6** (General Matroids). There is an algorithm for Byzantine secretary on general matroids that is $O(\log n)$ -competitive with the benchmark V^* .

Our results show how to get robust algorithms for the widely-studied secretary problems, and we hope it will generate further interest in robust algorithm design. Interesting next directions include improving the quantitative bounds in our results (which are almost certainly not optimal), and understanding tradeoffs between competitiveness and robustness.

1.3 Related Work

The secretary problem has a long history, see [17] for a discussion. The papers [1, 28, 16] studied generalizations of the secretary problem to matroids, [20, 27, 24, 22] studied extensions to matchings, and [38, 39] studied extensions to arbitrary packing constraints. More generally, the random-order model has been considered, both as a tool to speed up algorithms (see [6, 41]), and to go beyond the worst-case in online algorithms (see [32, 18, 19]). E.g., we can solve *linear programs* online if the columns of the constraint matrix arrive in a random order [8, 9, 25, 21], and its entries are small compared to the bounds. In online algorithms, the random-order model provides one way of modeling benign nature, as opposed to an adversary hand-crafting a worst-case input sequence; this model is at least as general as i.i.d. draws from an unknown distribution.

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Both the random-order model and the Byzantine model are *semi-random* models [2, 15], with different levels of power to the adversary. Other restrictions of the random-order model have been studied, via lower-bounding the entropy of the input stream, to ensure the permutations are "random enough" [5, 23]. These papers give sufficient conditions for the classical algorithms to perform well, whereas we take the dual approach of permitting outliers and then asking for new robust algorithms. There are works (e.g., [31, 33, 35]) that give algorithms which have a worst-case adversarial bound, and which work better when the input is purely stochastic; most of these do not study the performance on mixed arrival sequences. One exception is the work [14] who study online matching for mixed arrivals, under the assumption that the "magnitude of noise" is bounded. Another exception is a recent (unpublished) work of Kesselheim and Molinaro, who define a robust K-secretary problem similar to ours. They assume the corruptions have a bursty pattern, and get 1 - f(K)-competitive algorithms. Our model is directly inspired by theirs.

2 Preliminaries and Techniques

By $[a \dots b]$ we denote the set of integers $\{a, a + 1, \dots, b - 1, b\}$. The universe $U := R \cup G$ consists of red/corrupted elements R and greed/good elements G. Let v(e) denote the value of element e: in the ordinal case v(e) merely defines a total ordering on the elements, whereas in the value-maximization case $v(e) \in \mathbb{R}_{\geq 0}$. Similarly, let $v(\mathcal{A})$ be the random variable denoting the value of the elements selected by algorithm \mathcal{A} . Let $t_e \in [0, 1]$ be the arrival time of element e. Let $R = \{r_1, r_2, \dots, r_{|R|}\}$ and $G = \{g_{\max}, g_2, g_3, \dots, g_{|G|}\}$; the elements in each set are ordered in non-increasing values. Let V^* be the benchmark to which we compare our algorithm. Note that V^* is some function of $G \setminus \{g_{\max}\}$, depending on the setting. We sometimes refer to elements $\{e \in U \mid v(e) \geq v(g_2)\}$ as big.

2.1 Two Useful Subroutines

Here are two useful subroutines.

Select a Random Element. The subroutine is simple: select an element uniformly at random. The algorithm can implement this in an online fashion since it knows the total number of elements n. An important property of this subroutine is that, in the value case, if any element in U has value at least nV^* , this subroutine gets at least V^* in expectation since this highest value element is selected with probability 1/n.

Two-Checkpoints Secretary. The subroutine is defined on two checkpoints $T_1, T_2 \in [0, 1]$, and let $\mathbb{I} := [T_1, T_2]$ be the interval between them. The subroutine ignores the input up to time T_1 , observes it during \mathbb{I} by setting threshold τ to be the highest value seen in the interval \mathbb{I} , i.e., $\tau \leftarrow \max\{v(e) \mid t_e \in \mathbb{I}\}$. Finally, during $\langle T_2, 1 \rangle$ the subroutine selects the first element e with value $v(e) \geq \tau$.

We use the subroutine in the single-item setting where the goal is to find a "big" element, i.e., an element with value at least $v(g_2)$. Suppose that there are no big red elements in I. Now, if g_2 lands in I, and also g_{\max} lands in $\langle T_2, 1 \rangle$, we surely select some element with value at least g_2 . Indeed, if there are no big items, threshold $\tau \leftarrow g_2$, and because g_{\max} lands after I, it or some other element will be selected. Hence, with probability $\Pr[t_{g_2} \in I] \Pr[t_{g_{\max}} \in \langle T_2, 1 \rangle] = (T_2 - T_1) \cdot (1 - T_2)$, we select an element of value at least $v(g_2)$.

2.2 Our Techniques

A common theme of our approaches is to prove a "good-or-learnable" lemma for each problem. Our algorithms begin by putting down a small number of *checkpoints* $\{T_i\}_i$ to partition the time horizon [0, 1] – and the arriving items – into disjoint *intervals* $\{\mathbb{I}_i\}_i$. We maintain *thresholds* in each interval to decide whether to select the next element. Now a "good-orlearnable" lemma says that either the setting of the thresholds in the current interval \mathbb{I}_i will give us a "good" performance, or we can "learn" that this is not the case and update the thresholds for the next interval \mathbb{I}_{i+1} . Next we give details for each of our problems.

Uniform Matroid Value Maximization (§3). Recall that here we want to pick K elements (in particular, all elements have size 1, unlike the knapsack case where sizes are in the range [0,1]). For simplicity, suppose the algorithm knows that the benchmark V^* lies in [1,n]; we remove this assumption later. We define $O(\varepsilon^{-1} \log n)$ levels, where level $\ell \geq 0$ corresponds to values in the range $[n/(1+\varepsilon)^{\ell+1}, n/(1+\varepsilon)^{\ell})$. For each interval \mathbb{I}_i and level ℓ , we maintain a budget $B_{\ell,i}$. Within this interval \mathbb{I}_i , we select the next arriving element having a value in some level ℓ only if the budget $B_{\ell,i}$ has not been used up. How should we set these budgets? If there are $1/\delta$ intervals of equal size, we expect to select δK elements in this interval. So we have a total of δK budget to distribute among the various levels. We start off optimistically, giving all the budget to the highest-value level. Now this budget gradually cascades from a level ℓ to the next (lower-value) level $\ell + 1$, if level ℓ is not selecting elements at a "good" enough" rate. The intuition is that for the "heavy" levels (i.e., those that contain many elements from the benchmark-achieving set S^*), we will roughly see the right number of them arriving in each interval. This allows us to prove a good-or-learnable lemma, that either we select elements at a "good enough" rate in the current interval, or this is not the case and we "learn" that the budgets should cascade to lower value levels. There are many details to be handled: e.g., this reasoning is only possible for levels with many benchmark elements, and so we need to define a dedicated budget to handle the "light" levels.

Single-Item Value-Maximization (§5). We want to maximize the expected value of the selected element, compared to $V^* := v(g_2)$, the value of the second-max green. With some small constant probability our algorithm selects a uniformly random element. This allows us to assume that every element has value less than nV^* , as otherwise the expected value of a random guess is $\Omega(V^*)$. We now describe how applying the above "good-or-learnable" paradigm in a natural way guarantees an expected value of $\Omega(V^*/\log n)$. Running the two-checkpoint secretary (with constant probability) during $T_1 = 0, T_2 = 1/2$ we know that it gets value $\Omega(V^*)$ and we are done, or failing that, there exist a red element of value at least V^* in [0, 1/2]. But then we can use this red element (highest value in the first half) to get a factor n estimate on the value of V^* . So by grouping elements into buckets if their values are within a factor 2, and randomly guessing the bucket that contains $v(g_2)$, gives us an expected value of $\Omega(V^*/\log n)$. To obtain the stronger factor of poly log* n in Theorem 3, we now define log* n checkpoints. We prove a "good-or-learnable" lemma that either selecting a random element from one of the current buckets has a good value, or we can learn a tighter estimate on V^* and reduce the number of buckets.

Ordinal Single-Item Secretary (§4). We now want to maximize the probability of selecting an element whose value is as large as the green second-max; this is more challenging than value-maximization since there is no notion of values for bucketing. Our approach is crucially different. Indeed, we use the *minimax principle* in the "hard" direction: we give an algorithm

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that does well when the input distribution is *known to the algorithm* (i.e., where the algorithm can adapt to the distribution), and hence infer the existence of a (randomized) algorithm that does well on worst-case inputs.

The known-distribution algorithm uses $O(\log n)$ intervals. Again, we can guarantee there is a "big" (larger than g_2) red element within each interval, as otherwise running Dynkin's algorithm on a random interval with a small probability already gives a "good" approximation. This implies that even if the algorithm "learns" a good estimate of the second-max just before the last interval, it will win. This is because the algorithm can set this estimate of second-max as a threshold, and it wins by selecting the big red element of the last interval. Finally, to learn a good estimate on the second-max, we again prove a "good-or-learnable" lemma. Its proof crucially relies on the algorithm knowing the arrival distribution, since that allows us to set "median" of the conditional distribution as a threshold.

Other Results (§6). We also give $O(\log \log n)^2$ -competitive algorithms for Partition matroids, where the difficulty is that we cannot afford to lose the max-green element in every part. Our idea is to only lose one element globally to get a very rough scale of the problem, and then exploit this scale in every part. We also show why other potential benchmarks are too optimistic in §A, and how to relax the assumption that n is known in §B. See those sections for details.

3 Knapsack Byzantine Secretary

Consider a knapsack of size K; for all the results in this section we assume that $K \ge poly(\varepsilon^{-1}\log n)$. Each arriving element e has a size $s(e) \in [0, 1]$ and a value $v(e) \ge 0$. Let $g_{\max}, g_2, g_3, \ldots$, denote the green elements G with decreasing values and let

$$V^* := \max\left\{\sum_{e \in S} v(e) \mid S \subseteq G \setminus g_{\max} \text{ and } \sum_{e \in S} s(e) \le K\right\}$$
(3)

be the value of the benchmark solution, i.e., the optimal solution obtained after discarding the top green element g_{max} . Let S^* be the set of green elements corresponding to this benchmark.

In §3.1 we give a $(1 + \varepsilon)$ -competitive algorithm assuming we have a factor poly(n)-approximation to the benchmark value V^* . (In fact, given this poly(n)-approximation, we can even get within a $(1 + \varepsilon)$ -factor of the optimal set *including* g_{max} .) Then in §3.2 we remove the assumption, but now our value is only comparable to V^* (i.e., excluding g_{max}).

Intuition. The main idea of the regular (non-robust) multiple-secretary problem (where we pick at most K items) is to observe a small ε fraction of the input, estimate the value of the K^{th} largest element, and then select elements with value exceeding this estimate. (A better algorithm revises these estimates over time, but let us ignore this optimization for now.) In the Byzantine case, there may be an arbitrary number of red items, so strategies that try to estimate some statistics (like the K^{th} largest) to use for the rest of the algorithm are susceptible to adversarial attacks.

For now, suppose we know that all items of S^* have values in $[1, n^c]$ for some constant c. The density of an item to be its value divided by its size. We define $O(\log n)$ density levels, where elements in the same level have roughly the same density, so our algorithm does not distinguish between them. The main idea of our algorithm is to use cascading budgets. At the beginning we allocate all our budget to picking only the highest-density level items. If we find that we are not picking items at a rate that is "good enough", we re-allocate parts of

our budget to lower-density levels. The hope is that if the benchmark solution S^* selects many elements from a certain density level, we can get a good estimate of the "right" rate at which to pick up items from this level. Moreover, since our budgets trickle from higher to lower densities, the only way the adversary can confuse us is by giving dense red elements, in which case we will select them.

Such an idea works only for the value levels that contain many elements of S^* . For the remaining value levels, we allocate *dedicated budgets* whose sole purpose is to pick a "few" elements from that level, irrespective of whether they are from S^* . By making the total number of levels logarithmic, we argue that the total amount of dedicated budgets is only o(K), so it does not affect the analysis for the cascading budget.

3.1 An Algorithm Assuming a Polynomial Approximation

Suppose we know the benchmark V^* to within a polynomial factor: by rescaling, assume that V^* lies in the range $[1 \dots n^c]$ for some constant c. This allows us to simplify the instance structure as follows: Firstly, we can pick all elements of size at most 1/n, since the total space usage is at most $n \cdot 1/n = 1$ (recall, $K \ge \text{poly}(\varepsilon^{-1} \log n)$). Next, we can ignore all elements with value less than $1/n^2$ because their total value is at most $1/n \ll 1 \le V^*$. If the *density* of an element is defined to be the ratio v(e)/s(e), then all remaining elements have density between n^{c+1} and n^{-2} . The main result of this section is the following:

▶ Lemma 7. If V^* lies between 1 and n^c for some constant c, each element has size at least 1/n and value at least $1/n^2$, and $K \ge \text{poly}(\varepsilon^{-1} \log n)$, then there exists a $(1 + O(\varepsilon))$ -competitive algorithm.

The idea of our algorithm is to partition the input into $1/\delta$ disjoint pieces (δ is a small parameter that will be chosen later) and try to solve $1/\delta$ "similar-looking" instances of the knapsack problem, each with a knapsack of size δK .

The Algorithm. Define checkpoints $T_i := \delta i$ and corresponding intervals $\mathbb{I}_i := \langle T_{i-1}, T_i]$ for all $i \in [1 \dots 1/\delta]$. Define $L := (1 + \frac{c+3}{\varepsilon} \log n)$ density levels as follows: for each integer $\ell \in [0 \dots L)$, density value $\rho_{\ell} := n^{c+1}/(1+\varepsilon)^{\ell}$. Now density level ℓ corresponds to all densities lying in the range $(\rho_{\ell+1}, \rho_{\ell}]$. Note that densities decrease as ℓ increases. We later show that the setting of parameters $K \ge \Omega(\frac{L^2 \log L/\varepsilon}{\varepsilon^4})$ and $1/\delta = \Omega(L/\varepsilon)$ suffices.

We maintain two kinds of *budgets*:

- Cascading budgets: We maintain a budget $B_{\ell,i}$ for each density level ℓ and each interval \mathbb{I}_i . For the first interval \mathbb{I}_1 , define $B_{0,1} := \delta K$, and $B_{\ell,1} := 0$ for $\ell > 0$. For the subsequent intervals, we will set $B_{\ell,i}$ in an online way as described later.
- Dedicated budgets: We maintain a dedicated budget $\hat{B}_{\ell} := H$ for each density level ℓ ; we will later show that setting $H := \Omega(\frac{L \log L/\varepsilon}{\varepsilon^3})$ suffices.

Suppose we are in the interval \mathbb{I}_i , and the arriving element e has density v(e)/s(e) in level ℓ . 1. If the remaining cascading budget $B_{\ell',i}$ of one of the density levels $\ell' \geq \ell$ is positive then

- select e. For the smallest $\ell' \geq \ell$ satisfying this condition, update $B_{\ell',i} \leftarrow B_{\ell',i} s(e)$.
- 2. Else, if the remaining dedicated budget \tilde{B}_{ℓ} for level ℓ is positive, select e and update $\tilde{B}_{\ell} \leftarrow \tilde{B}_{\ell} s(e)$.

Finally, for i > 1, we define the cascading budgets $B_{\ell,i}$ for this interval \mathbb{I}_i based on how much of the budgets at levels ℓ and $\ell - 1$ are consumed in the previous interval \mathbb{I}_{i-1} as follows. The amount of budget $B_{\ell-1,i-1}$ at level $\ell - 1$ that is not consumed in interval \mathbb{I}_{i-1} is moved to level ℓ (which has lower density), and the budget that gets consumed in \mathbb{I}_{i-1} is restored

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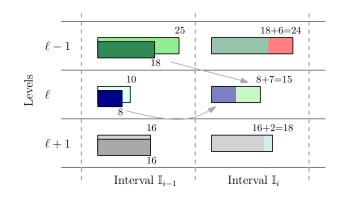


Figure 1 The bars for \mathbb{I}_{i-1} show the budget, and (in darker colors) the amount consumed. The consumed budget (in dark blue) at level ℓ in interval \mathbb{I}_{i-1} is restored at level ℓ in \mathbb{I}_i ; the unconsumed budget at level $\ell - 1$ in interval \mathbb{I}_{i-1} is then added to it.

at level ℓ (see Figure 1). Formally, if $C_{\ell,i-1}$ is the amount of *consumed* cascading budget for level ℓ in interval \mathbb{I}_{i-1} and $R_{\ell,i-1}$ is the amount of *remaining* budget at level ℓ at the end of interval i-1 (i.e., the value of $B_{\ell,i-1}$ at the time corresponding to the end of \mathbb{I}_{i-1}), then we define the initial budget for level ℓ at the start of interval i to be

 $B_{\ell,i} := C_{\ell,i-1} + R_{\ell-1,i-1}.$

It is easy to see that we can compute these cascading budgets online.

A Note about Budget Violations. The total budget, summed over both categories and over all the intervals for the cascading budgets, is $K' := ((1/\delta) \cdot \delta K) + L H > K$. If we use up all this budget, we would violate the knapsack capacity. Moreover, we select an element as long as the corresponding budget is positive, and hence may exceed each budget by the size of the last element. However, since $K' \leq (1 + \varepsilon)K$ and K is much larger than individual element sizes, the violations is a small fraction of K, so we can reject each element originally selected by the algorithm with some small probability (e.g., $p = 2\varepsilon$) to guarantee that the non-rejected selected elements have size at most K with high probability (i.e., at least $1 - 1/n^c$, for an arbitrary constant c > 0). Henceforth, we will not worry about violating the budget.

The Analysis. Recall the benchmark V^* from (3), and let S^* be a set that achieves this value. All the elements have value in $[1/n^2, n^c]$ and size at least [1/n, 1], so each element $e \in S^*$ has a corresponding density level $\ell(e) \in [0 \dots L)$ based on its density v(e)/s(e). We need the notion of "heavy" and "light" levels. For any level $\ell \in [0 \dots L)$, define s^*_{ℓ} to be the total size of elements in S^* with density level ℓ :

$$s_{\ell}^* := \sum_{e \in S^* : \ell(e) = \ell} s(e).$$
 (4)

We say a level ℓ is heavy if $s_{\ell}^* \geq H$, else level ℓ is light. We refer to (green) elements of S^* at a heavy (resp., light) level as heavy-green (resp., light-green) elements. Note that elements not in S^* (some are red and others green) are left unclassified. If H is sufficiently large, a concentration-of-measure argument using the uniformly random arrival times for green items shows that each heavy level receives $(1 - \varepsilon)\delta H$ size during each interval with high probability. The idea of the proof is to argue that the cascading budget never "skips" a heavy level, and hence we get almost all the value of the heavy levels.

To avoid double-counting the values of the light levels, we separately account for the algorithm's value attained (a) on light levels using the dedicated budget or on light-green elements using the cascading budget, and (b) for elements that are not light-green (incl.

red elements) using the cascading budget. Note that (a) and (b) are disjoint, hence their contributions can be added up. We show that (a) exceeds the value of S^* restricted to the light levels, while (b) exceeds $(1 - \varepsilon)$ times the value of of S^* on the heavy levels. This is sufficient to prove our result. We start by arguing the former claim.

 \triangleright Claim 8 (Light-Green Elements). The sum of values of elements selected using the dedicated budget at light levels, and of light-green elements selected using the cascading budget, is at least $\sum_{\ell \text{ light }} s_{\ell}^* \cdot \rho_{\ell+1}$.

Proof. Our algorithm attempts to select each light-green element in S^* using the cascading budget, or failing that, by the dedicated budget at its density level. The only case in which a light-green element $e \in S^*$ is dropped is if all the dedicated budget at its level $\ell(e)$ has been exhausted. But this means the algorithm has already collected at least $s_{\ell}^* \cdot \rho_{\ell+1}$ from the dedicated budget at this light level ℓ .

Next, to prove that (b) exceeds the value on heavy levels (up to $1 - \varepsilon$), we need the following property of the cascading budget on the heavy levels.

 \triangleright Claim 9. For all intervals \mathbb{I}_i and levels ℓ , w.h.p. we have that if $B_{\ell,i} > 0$ then every heavy level $\ell' < \ell$ satisfies $B_{\ell',i} \ge \delta s_{\ell'}^* \cdot (1 - \varepsilon)$.

Proof. For a heavy level ℓ' , the expected size of heavy-green elements from S^* falling in any interval is $\delta s_{\ell'}^* \geq \delta H$. If $\delta H \geq \frac{\Omega(\log(L/(\delta \varepsilon)))}{\varepsilon^2}$ then with probability $1 - \varepsilon$ we get that for each interval *i* and each heavy level ℓ' , the total size of elements from S^* lying at level ℓ' and arriving in interval \mathbb{I}_i is at least $\delta s_{\ell'}^* \cdot (1 - \varepsilon)$, by a concentration bound. Henceforth, let us condition on this event happening for all heavy levels ℓ' .

Now if the cascading budget $B_{\ell,i} > 0$, this budget must have gradually come from levels $\ell' < \ell$ of higher densities. But this means $B_{\ell',i} \ge \delta s^*_{\ell'} \cdot (1-\varepsilon)$ because otherwise the cascading budget would never move to level $\ell' + 1$, since level ℓ' receives at least $\delta s^*_{\ell'} \cdot (1-\varepsilon)$ size of elements in every interval.

For a level τ let $h^*_{[0,\tau\rangle} := \sum_{\ell' \text{heavy}, \ \ell' < \tau} s^*_{\ell'}$ denote the total size of items in S^* restricted to heavy levels from $[0, \tau\rangle$. Similarly, let $h^{\mathcal{A}}_{[0,\tau\rangle}$ be the total size of non-light-green items collected by the algorithm in levels $[0, \tau\rangle$ and charged against the cascading budget.

 \triangleright Claim 10. For all levels τ we have that $h_{[0,\tau)}^{\mathcal{A}} \ge (1 - O(\varepsilon))h_{[0,\tau)}^{*}$.

Proof. Let t be the smallest index of an interval where $B_{\tau,t+1} > 0$. We partition the intervals into two groups: $\mathbb{I}_1, \mathbb{I}_2, \ldots, \mathbb{I}_t$ and $\mathbb{I}_{t+1}, \ldots, \mathbb{I}_{1/\delta}$. From Claim 9 we can conclude that for *each interval* in the latter group, the algorithm collects a total size of at least $\sum_{\ell \text{ heavy}, \ \ell < \tau} (1-\varepsilon) \, \delta \, s_{\ell}^* = (1-\varepsilon) \, \delta \, h_{[0,\tau)}^*$ from levels $[0, \tau)$. Hence the total contribution over all the intervals of the latter group is $(\frac{1}{\delta} - t)(1-\varepsilon) \, \delta \, h_{[0,\tau)}^* = (1-t\delta)(1-\varepsilon) \, h_{[0,\tau)}^*$.

We now consider the group $\mathbb{I}_1, \ldots, \mathbb{I}_t$. Let C_i, R_i and Q'_i be the total size of the consumed non-light-green, remaining budget, and consumed light-green elements charged to the cascading budget in interval i with levels $[0, \tau \rangle$. By definition, the total size of all light-green elements is at most LH, giving $\sum_{i=1}^t Q_i \leq LH$. Furthermore, since the full cascading budget is contained in $[0, \tau \rangle$, the algorithm construction guarantees $C_i + R_i + Q_i = \delta K$. Finally, we argue that $\sum_{i=1}^t R_i \leq \delta KL$: consider an infinitesimally small part dB of the budget. At the end of each interval, dB is either used to consume an element or it "moves" from level ℓ to $\ell + 1$, which can happen at most L times. Since the total amount of budget per interval is $\int dB = \delta K$, the total sum is at most δKL .

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This lower-bounds the total size contribution of the group $\mathbb{I}_{t+1}, \ldots, \mathbb{I}_{1/\delta}$.

$$\sum_{i=1}^{t} C_{i} = t\delta K - \sum_{i=1}^{t} R_{i} - \sum_{i=1}^{t} Q_{i} \ge t\delta K - \delta KL - LH \ge (t\delta - \varepsilon - \varepsilon)K$$
$$\ge (t\delta - O(\varepsilon))h_{[0,\tau)}^{*},$$

where we use $K \ge h^*_{[0,\tau\rangle}$ (since the total size of elements in S^* is at most K), $\delta L \le \varepsilon$, and $LH \le \varepsilon K$. Combining contributions from both groups we get:

$$(1-t\delta)(1-\varepsilon) h^*_{[0,\tau\rangle} + (t\delta - O(\varepsilon)) h^*_{[0,\tau\rangle} = [(1-\varepsilon) - t\delta(1-\varepsilon) + t\delta - O(\varepsilon)] h^*_{[0,\tau\rangle}$$
$$= (1-O(\varepsilon)) h^*_{[0,\tau\rangle}.$$

Hence, we conclude that $h_{[0,\tau\rangle}^{\mathcal{A}} \ge (1 - O(\varepsilon)) h_{[0,\tau\rangle}^{*}$.

Using the above claims we now prove Lemma 7.

Proof of Lemma 7. Our fine-grained discretization of densities gives us that

$$V^* \le (1+\varepsilon) \bigg(\sum_{\ell \text{ light}} s^*_{\ell} \rho_{\ell+1} + \sum_{\ell \text{ heavy}} s^*_{\ell} \rho_{\ell+1} \bigg).$$
(5)

From Claim 8, our algorithm accrues value at least $\sum_{\ell \text{ light}} s_{\ell}^* \cdot \rho_{\ell+1}$ due to the elements from light levels that were charged to the dedicated budget and light-green elements charged to the cascading budget. It is therefore sufficient to prove a similar bound on the value accrued on non-light-green elements charged to the cascading budget with respect to $\sum_{\ell \text{ heavy}} s_{\ell}^* \cdot \rho_{\ell+1}$, which we deduce from Claim 10.

Let $\ell'(x)$ be defined as the largest level ℓ' where $\rho_{\ell'} \ge x$, then

$$\sum_{\ell \text{ heavy}} s_{\ell}^* \rho_{\ell+1} = \int_0^\infty \sum_{\ell \text{ heavy, } \rho_{\ell+1} \ge x} s_{\ell}^* dx = \int_0^\infty h_{[0,\ell'(x))}^* dx$$
$$\leq (1+O(\varepsilon)) \int_0^\infty h_{[0,\ell'(x))}^{\mathcal{A}} dx$$

where the last inequality uses Claim 10. Notice the right-hand side is the value of non-lightgreen elements charged against the cascading budget. Thus, this part of the algorithm's value exceeds (up to $1 - O(\varepsilon)$) the value of heavy levels of S^* , finalizing our proof.

3.2 An Algorithm for the General Case

To remove the assumption that we know a polynomial approximation to V^* , the idea is to ignore the first ε fraction of the arrivals, and use the maximum value in this interval to get a poly(n) approximation to the benchmark. This strategy is easily seen to work if there are $\Omega(1/\varepsilon)$ elements with a non-negligible contribution to V^* . For the other case where most of the value in V^* comes from a small number of elements, we separately reserve some of the budget, and run a collection of algorithms to catch these elements when they arrive.

Formally, we define $(1/\varepsilon)$ checkpoints $T_i := i\varepsilon$ and corresponding intervals $\mathbb{I}_i := \langle T_{i-1}, T_i \rangle$ for all $i \in [1 \dots (1/\varepsilon)]$. We run the following three algorithms in parallel, and select the union of elements selected by them.

(i) Select one of the n elements uniformly at random; i.e., run Select-Random-Element from §2.1.

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- (ii) Ignore elements that arrive in times $[0, \varepsilon)$, and let \hat{v} denote the highest of their values. Run the algorithm from §3.1 during time $[1/\varepsilon, 1]$, assuming that $V^* \in [\hat{v}/n^2, \hat{v} n^2]$.
- (iii) At every checkpoint T_i , consider the largest value \hat{v}_i seen until then. Define $L := \frac{10}{\varepsilon} \log n$ value levels as follows: for $\ell \in (-L/2 \dots L/2)$ and $\tau_\ell := \hat{v}_i/(1+\varepsilon)^\ell$, define level $\ell(i)$ as corresponding to values in $(\tau_{\ell(i)+1}, \tau_{\ell(i)}]$. For each of these levels ℓ , keep $D := \frac{10}{\varepsilon} \log \frac{1}{\varepsilon}$ dedicated slots, and select any element having this value level and arriving after T_i , as long as there is an empty slot in its level.

The total space used by the there algorithms is at most

$$1 + K + (1/\varepsilon) \cdot L \cdot D = K + O\left(\frac{\log n \log 1/\varepsilon}{\varepsilon^3}\right) \leq (1+\varepsilon)K,$$

where the last inequality holds because $K \ge \Omega\left(\frac{L^2 \log(L/\varepsilon)}{\varepsilon^4}\right)$ from the size condition from §3.1. We can now fit this into our knapsack of size K w.h.p. by sub-sampling each selected element with probability $(1 - O(\varepsilon))$. To complete the proof of Theorem 2, we need to show that we get expected value $(1 - O(\varepsilon))V^*$.

Proof of Theorem 2. The proof considers several cases. Firstly, if there is any single element with value more than $n \cdot V^*$, then the algorithm in Step (i) will select it with probability 1/n, proving the claim. Hence, all elements have value at most nV^* .

Now suppose at least $D = \frac{10}{\varepsilon} \log \frac{1}{\varepsilon}$ elements in S^* (recall S^* has total value V^*) have individual values at least V^*/n^2 . In this case, at least one of these D elements arrives in the interval $[0, \varepsilon)$ with probability $1 - \varepsilon$, and that element gives us the desired n^2 -approximation to V^* . Moreover, the expected value of elements in S^* arriving in times $[\varepsilon, 1]$ is at least $(1 - O(\varepsilon)) V^*$, even conditioning on one of them arriving in $[0, \varepsilon)$.

Finally, consider the case where $D' \leq D$ elements of S^* have value more than V^*/n^2 . The idea of the algorithm in Step (iii) is to use the earliest arriving of these D' elements, or the element g_{\max} , to get a rough estimate of V^* , and from thereon use the dedicated slots to select the remaining elements. Indeed, if the first of these elements arrive in interval \mathbb{I}_i , the threshold \hat{v}_i lies in $[V^*/n^2, nV^*]$ (since we did not satisfy the first case above). Now the value levels and dedicated budgets set up at the end of this interval would pick the rest of these D' elements – except those that fall in this same interval \mathbb{I}_i . We argue that each of the remaining D' elements has at least $(1 - \varepsilon)$ probability of not being in \mathbb{I}_i , which gives us an expected value of $(1 - O(\varepsilon))V^*$ in this case as well. This is true because the expected number of these 1 + D' elements (including g_{\max}) that land in any interval that contains at least one of them is at most $1 + \varepsilon D'$ (even after we condition on the first arrival, each remaining element has ε chance of falling in this interval). Since any such interval has the same chance of being the first interval \mathbb{I}_i , and these 1 + D' elements have the same distribution, the expected number of additional elements in \mathbb{I}_i is $\varepsilon D'$.

This completes the proof of Theorem 2 for the knapsack case, where the size K of the knapsack is large enough compared to the largest size of any element. This generalizes the multiple-secretary problem, where all items have unit size. We have not optimized the value of K that suffices, opting for modularity and simplicity. It can certainly be improved further, though getting an algorithm that works under the assumption that $K \ge O(1/\varepsilon^2)$, like in the non-robust case, may require new ideas.

4 Single-Item Ordinal Case

In this section we give a proof of Theorem 4, showing that there exists an algorithm which selects an element with value no smaller than g_2 , with probability at least $\Omega(1/\log^2 n)$. Our proof for this theorem is non-constructive and uses (the hard direction of) the Minimax

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Theorem; hence we can currently only show the *existence* of this algorithm, and not give a compact description for it. Our main technical lemma furnishes an algorithm which, given a known (general) probability distribution \mathcal{B} over input instances, selects a big element with probability at least $\Omega(1/\log^2 n)$. Consequently, we use the Minimax Lemma to deduce that the known-distribution case is equivalent to the worst-case input setting and recover the analogous result.

Since our algorithms crucially argue about the input distribution \mathcal{B} and rely on the Minimax, we need to formally define these terms and establish notation connecting the Byzantine secretary problem with two-player zero-sum games. Suppose we want to maximize the probability of selecting a big element and to this end we choose an algorithm \mathcal{A} , while the adversary chooses a distribution \mathcal{B} over the input instances and there is an (infinite) payoff matrix K prescribing the outcomes. Its rows are indexed by different algorithms, and columns by input instances. Formally, a "pure" input instance is represented as an $|\mathcal{R}|$ -tuple of numbers in [0, 1], representing the arrival times t_e of the red elements; and a permutation $\pi \in S_n$ over U representing the total ordering of all values in $U = \mathcal{R} \cup G$. Recall that the green elements G choose their arrival times independently and uniformly at random in [0, 1], hence their t_e 's are not part of the input. A "mixed" input instance is a probability distribution \mathcal{B} over pure instances $[0, 1]^{|\mathcal{R}|} \times S_n$.

While we do not need the full formal specifications of algorithms, we will mention that a "mixed" algorithm \mathcal{A} is a distribution over deterministic algorithms. An algorithm \mathcal{A} on an input instance I gets a payoff of $K(\mathcal{A}, I) := \Pr[v(\mathcal{A}) \ge v(g_2) \mid I]$ where the probability is taken over the assignment of random arrival times to elements in G and the distribution of deterministic algorithms \mathcal{A} . The following Lemma states that for each \mathcal{B} there is an algorithm (that depends on \mathcal{B}) that selects a big elements with probability $\Omega(1/\log^2 n)$. We prove the result in §4.1 and §4.2.

▶ Lemma 11 (Known Distribution Ordinal Single-Item Algorithm). Given a distribution over input instances \mathcal{B} , there exists an algorithm \mathcal{A} that has an expected payoff of $\Omega(1/\log^2 n)$.

To deduce the general case from the known distribution setting, we use a minimax lemma for two-player games. We postpone the details to Appendix C and simply state the final result here.

► Theorem 4 (Ordinal Single-item Algorithm). There is a randomized algorithm for the ordinal Byzantine secretary which selects an element of value at least the second-largest green item with probability $\Omega(1/\log^2 n)$.

4.1 The Algorithm when \mathcal{B} is Known

In this section we give the algorithm for Lemma 11. We start with some preliminary notation. For each element e, let t_e denote the time at which it appears. Furthermore, for $t \in [0, 1]$, let $\mathcal{K}(t)$ denote the information seen by the algorithm up to and including time t, consisting of arrival times and relative values of elements appearing before t.

We define $\log n + 1$ time *checkpoints* as follows: set the initial checkpoint $T_0 := \frac{1}{4}$, and then subsequent checkpoints $T_i := \frac{1}{4} + \frac{i}{2 \cdot \log n}$ for all $i \in [1 \dots \log n]$. Note that the last checkpoint is $T_{\log n} = \frac{3}{4}$. Now the corresponding intervals are

$$\mathbb{I}_{0} := [0, 1/4] \quad , \quad \mathbb{I}_{i} := \langle T_{i-1}, T_{i}] \ \forall i \in [1 \dots \log n], \quad \text{and} \quad \mathbb{I}_{\log n+1} := \langle 3/4, 1].$$
(6)

Let $m_i := \max\{v(e) \mid e \in R \text{ and } t_e \in \mathbb{I}_i\}$ be the maximum value among the red elements that land in interval \mathbb{I}_i , and let $\mathcal{H} := \{m_i > v(g_2) \text{ for all } i \in [1 \dots \log n]\}$ be the event where the maximum value red item in all intervals is larger than the target g_2 , i.e., is "big". We call this event \mathcal{H} the *hard cases* and \mathcal{H}^c the *easy cases*; we will show the Two Checkpoints Secretary (from §2.1) achieves $\Omega(1/\log^2 n)$ winning probability for all input instances in \mathcal{H}^c . Finally, define

$$p_e^i := \Pr_{\mathcal{B}}[e = g_2 \mid \mathcal{H} \text{ and } \mathcal{K}(T_i)],$$

i.e., p_e^i is the probability that e is the second-highest green element conditioned on the information seen until checkpoint T_i and the current instance being hard. Importantly, the algorithm can compute p_e^i at T_i .

Now to solve the hard cases, at each checkpoint T_i the algorithm computes sets S_i satisfying $S_{i+1} \subseteq S_i$. These sets represent elements which are candidates for the second-max. In other words, at time T_i there is reasonable probability that second-max is in S_i . We start with defining $S_0 \leftarrow \{e \in U \mid t_e \in \mathbb{I}_0\}$, the elements the algorithm saw before T_0 . For $i \ge 0$, let c_i denote the *center* of S_i , i.e., the element of S_i such that there are exactly $\lfloor |S_i|/2 \rfloor$ elements smaller than it. Define $p^i(X) := \sum_{e \in X} p_e^i$ for a set X and index $i \ge 0$. Given S_{i-1} , we determine S_i as follows:

Define $\mathsf{bot}_{i-1} \leftarrow \{e \in S_{i-1} \mid v(e) \le v(c_{i-1})\}$, and note that $\mathsf{bot}_{i-1} \subseteq S_{i-1}$.

If $p^i(\mathsf{bot}_{i-1}) = p^{i-1}(S_{i-1})$ then $S_i \leftarrow \mathsf{bot}_{i-1}$, else $S_i \leftarrow S_{i-1} \setminus \mathsf{bot}_{i-1}$.

Our algorithm runs one of the following three algorithms uniformly at random:

- (i) Select a random $i \in [1 \dots \log n]$, define $\tau \leftarrow \max\{v(e) \mid t_e \in \mathbb{I}_i\}$ and select the first element larger than τ . I.e., run Two Checkpoints Secretary (from §2.1) with the checkpoints being the ends of interval \mathbb{I}_i .
- (ii) Select a random $i \in [0 \dots \log n]$, read input until checkpoint T_i , define $\tau \leftarrow v(c_i)$ and select the first element larger than it.
- (iii) Compute the sets S_i until $|S_k| \leq 10$ for some k: then define τ to be the value of a random element in S_k , and select the first element larger than it.

4.2 The Analysis

In this section we prove Lemma 11. Let us give some intuition. We can assume we have a hard case, else the first algorithm achieves $\Omega(1/\log^2 n)$ winning probability. For the other two algorithms, let us condition on g_2 falling in the first interval \mathbb{I}_0 , and then exploit the fact that there is a big red element in every interval \mathbb{I}_i . It may be useful to imagine that we are trying to guess, at each checkpoint, which of the elements in the past were actually g_2 . If we could do this, we would set a threshold at its value, and select the first subsequent element bigger than the threshold – and since there is a 1/4 chance that g_{\max} would fall in $\mathbb{I}_{\log n+1}$, we'd succeed! Of course, since there are red elements all around, guessing g_2 is not straightforward.

So suppose we are at checkpoint T_k , and suppose there is a reasonable probability that $v(g_2) \leq v(c_{k-1})$, but also still some nonzero probability that $v(g_2) > v(c_{k-1})$. In such a scenario, we claim that trying to choose an element in the interval \mathbb{I}_k larger than c_{k-1} will give us a reasonable probability of success. Indeed, we claim there would have been at least one red element in \mathbb{I}_k bigger than c_{k-1} (since there is still a non-zero probability that $v(g_2) > v(c_{k-1})$ even at the end of the interval \mathbb{I}_k , and since the case is hard), and $v(g_2) \leq v(c_{k-1})$ with reasonable probability. Of course, we only know this at the end of the interval, but the algorithm can randomly guess k with $\Omega(1/\log n)$ probability. Finally, if there is no such checkpoint, then in every interval we reduce the size of set $|S_i|$ by half while suffering a small loss in $p(S_i)$. In this case, both $|S_{\log n}| = O(1)$ and $p(S_{\log n}) = \Omega(1)$, so the third algorithm can guess g_2 with constant probability and select an element larger than it in the last interval.

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Formal Analysis. Let ALG be 1 if $v(\mathcal{A}) \geq v(g_2)$ and 0 otherwise, where \mathcal{A} is the algorithm from the last section. Suppose we're in an easy case, i.e., there is an interval \mathbb{I}_s such that all red elements in this interval are smaller than g_2 . Now if the first algorithm is chosen, suppose it selects the interval \mathbb{I}_s , suppose g_2 lands in \mathbb{I}_s , and g_{\max} lands in $\mathbb{I}_{\log n+1}$. Then the algorithm surely selects an element greater than g_2 , and it has expected value:

$$\mathbb{E}[ALG] \ge \frac{1}{3} \cdot \frac{1}{\log n} \cdot \frac{1}{2\log n} \cdot \frac{1}{4} = \Omega\left(\frac{1}{\log^2 n}\right).$$

Henceforth we can assume the case is hard, and hence each interval I_i contains a red element bigger than g_2 . We condition on the event that g_2 appears in \mathbb{I}_0 , which happens with constant probability. Define

$$k^* := \min \left\{ i \in [1 \dots \log n] \mid \frac{1}{\log n} \le \frac{p^i(\mathsf{bot}_{i-1})}{p^{i-1}(S_{i-1})} < 1 \right\},$$

and set $k^* = \log n + 1$ if the above set is empty.

 \triangleright Claim 12. For all $i < k^*$, the probability $p^i(S_i) = \Omega(1)$.

Proof. By definition, $p^0(S_0) = 1$. By our definition of the sets S_i , we know that if $p^i(\mathsf{bot}_{i-1}) = p^{i-1}(S_{i-1})$ then $p^i(S_i) = p^{i-1}(S_{i-1})$. Else since $i < k^*$, we have

$$p^{i}(S_{i}) = p^{i-1}(S_{i-1}) - p^{i}(\mathsf{bot}_{i-1}) \le p^{i-1}(S_{i-1})(1 - \frac{1}{\log n}).$$

Hence, $p^i(S_i) \ge (1 - \frac{1}{\log n})^{\log n} = \Omega(1)$, proving the claim.

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Now there are two cases, depending on the value of k^* . Suppose $k^* \leq \log n$. Condition on the event that the second algorithm is chosen, that it chooses the $i^{th} = (k^* - 1)^{th}$ checkpoint, and that $v(g_2) \leq v(c_i)$. By our choice of k^* , we get that $v(g_2) \leq \tau = v(c_i)$ with probability at least $p^i(S_i) \cdot \frac{1}{\log n}$, and by Claim 12 this is $\Omega(\frac{1}{\log n})$. Since the case we are considering is hard and $\Pr[v(g_2) > v(c_i) \mid \mathcal{H}$ and $\mathcal{K}(T_{i+1})] > 0$, there is a red element larger than $v(c_i) = \tau$ appearing in \mathbb{I}_i . Thus the algorithm will always select an element in this interval. The correct interval is chosen with probability $\frac{1}{\log n}$, so the algorithm's value is

$$\mathbb{E}[ALG] = \frac{1}{3} \cdot \frac{1}{\log n} \cdot \Omega\left(\frac{1}{\log n}\right) = \Omega\left(\frac{1}{\log^2 n}\right)$$

The other case is when $k^* = \log n + 1$. By definition $|S_0| \leq n$ and $|S_i| \leq \lceil |S_{i-1}|/2 \rceil$. Therefore $|S_{\log n}| \leq 10$. Let us condition on the event that the third algorithm is chosen, that g_{\max} appears in $\mathbb{I}_{\log n+1}$, and that the algorithm guesses g_2 correctly. The probability of this event is at least

$$\frac{1}{3} \cdot \frac{1}{4} \cdot p^{\log n}(S_{\log n}) \cdot \frac{1}{10} = \Omega(1).$$

where we use Claim 12 to bound the probability $p^{\log n}(S_{\log n})$. In this event, the algorithm selects an element larger than g_2 and has expected value $\mathbb{E}[ALG] = \Omega(1)$.

Putting all these cases together, we get that our algorithm selects an element with value at least $v(g_2)$ with probability at least $\Omega((\log n)^{-2})$. This finishes the proof of Lemma 11, and hence of Theorem 4. It remains an intriguing open question to get a direct algorithm that achieves similar guarantees.

5 Single-Item Value-Maximization

In this section, we give an algorithm for the problem of selecting an item to maximize the expected *value*, instead of maximizing the probability of selecting the second-largest green item (the ordinal problem considered in §4). In the classical secretary problem, both problems are well known to be equivalent, with Dynkin's algorithm giving a tight 1/e bound for both. But in the Byzantine case the problems thus far appear to have different levels of complexity: in §6.2 we present a simple $O(\log n)$ -competitive algorithm for the value-maximization byzantine secretary problem, which is already better than the poly log *n*-competitive of §4. We now substantially improve it to give a poly log* *n*-competitive ratio.

▶ **Theorem 3** (Value Maximization Single-Item). There is a randomized algorithm for the value-maximization (single-item) Byzantine secretary problem which gets an expected value at least $(\log^* n)^{-2} \cdot V^*$.

In the rest of this section, let $V^* := v(g_2)$ denote the benchmark, the value of the second-largest green element. The high level idea of our algorithm is to partition the input into $O(\log^* n)$ intervals and argue that every interval contains a red element of value $v_i > V^*$, as otherwise Dynkin's algorithm will be successful. Moreover, this v_i cannot be much larger than V^* , as otherwise we can just select a random element. This implies we can use the largest value in each interval to find a good estimate of V^* , and eventually set it as a threshold in the last interval to select a large value element.

5.1 The Algorithm

Define $\log^{(i)} n$ to be the *iterated logarithm* function: $\log^{(0)} n = n$ and $\log^{(i+1)} n = \log(\log^{(i)} n)$. We define $\log^* n + 1$ time *checkpoints* as follows: the initial checkpoint $T_0 = \frac{1}{2}$, and then subsequent checkpoints $T_i = \frac{1}{2} + \frac{i}{4 \cdot \log^* n}$ for all $i \in [1, \ldots, \log^* n]$. Note that the last checkpoint is $T_{\log^* n} = \frac{3}{4}$. Now the intervals are

$$\mathbb{I}_{0} := [0, T_{0}] \quad , \quad \mathbb{I}_{i} := \langle T_{i-1}, T_{i} \rangle \, \forall i \in [1 \dots \log^{*} n], \quad \text{and} \quad \mathbb{I}_{\log^{*} n+1} := \langle T_{\log^{*} n}, 1].$$
(7)

Our algorithm runs one of the following three algorithms chosen uniformly at random.

- (i) Select one of the *n* elements uniformly at random; i.e., run Select-Random-Element from §2.1.
- (ii) Select a random interval $i \in [1 \dots \log^* n]$ and run Dynkin's secretary algorithm on \mathbb{I}_i . Formally, run Two-Checkpoints-Secretary (from §2.1) with the interval being $[T_{i-1}, \frac{1}{2}(T_{i-1} + T_i)]$.
- (iii) Select a random index $i \in [0 \dots \log^* n]$ and observe the maximum value during the interval \mathbb{I}_i ; let this maximum value be v_i . Choose a uniformly random $s \in [0 \dots 2 \log^{(i)} n]$. Select the first element arriving after T_i that has value at least $\tau := (v_i \log^{(i)} n)/2^s$.

5.2 The Analysis

To prove Theorem 3, assume WLOG that there are only two green elements g_{max} and g_2 , and every other element is red (otherwise, we can condition on the arrival times of all other green elements). Let v_i be the value of the highest red element in \mathbb{I}_i , i.e., excluding g_{max} and g_2 .

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Proof of Theorem 3. We assume $\log^{(i)} n$ is an integer for all *i*; this is true with a constant factor loss. For sake of a contradiction, assume that the algorithm in §5.1 does not get expected value $\Omega((\log^* n)^{-2}V^*)$. Under this assumption, we first show that every interval contains a red element of value at least V^* .

 \triangleright Claim 13. For all $j \in [1 \dots \log^* n]$ we have $v_j \ge V^*$.

Proof. Suppose this is not the case. Let \mathcal{E}_1 be the event that the following three things happen simultaneously: that we select Algorithm ii in §5.1 with random variable i = j, that the second-highest green element g_2 falls in the interval $[T_{i-1}, \frac{1}{2}(T_{i-1} + T_i))$, and that the highest green element g_{\max} falls in $\mathbb{I}_{\log^* n+1}$. Note that $\Pr[\mathcal{E}_1] = \frac{1}{3\log^* n} \cdot \frac{1}{4\log^* n} \cdot \frac{1}{4} = \Omega((\log^* n)^{-2})$. Conditioned on this event \mathcal{E}_1 , our algorithm (or specifically, Algorithm ii on the interval \mathbb{I}_j) gets a value at least $v(g_2) = V^*$. Hence the algorithm has expected valuation $\Omega((\log^* n)^{-2}V^*)$, which is a contradiction to our assumption on its performance.

We now prove that these red elements with large values cannot be much larger than V^* . **Lemma 14.** For all $j \in [1 \dots \log^* n]$ we have $v_j \leq V^* \cdot \log^{(j-1)} n$.

Proof. We prove this lemma by induction. The base case j = 1 says $v_1 \leq nV^*$, i.e., the highest observed value in $\mathbb{I}_1 = [T_0, T_1\rangle$ is at most nV^* . Suppose this is not the case – there exists a red element e in \mathbb{I}_1 with value at least nV^* . Let \mathcal{E}_1 be the event that we select Algorithm i in §5.1 (i.e., Select-Random-Element) and that it selects e. Since $\Pr[\mathcal{E}_1] = \Omega(\frac{1}{n})$, we have a contradiction that the expected valuation is $\Omega(V^*)$.

Now suppose the statement is true until $j \ge 1$. We prove the inductive step j+1. Suppose not, i.e., $v_{j+1} > V^* \log^{(j)} n$. Let \mathcal{E}_2 be the event that we select Algorithm iii in §5.1 with parameter i = j and that the random $s \in [0 \dots 2 \log^{(j)} n]$ is such that $\frac{v_j}{2^{s+1}} \le V^* < \frac{v_j}{2^s}$ (it exists by induction hypothesis). This implies threshold $\tau := \frac{v_j \log^{(j)} n}{2^s}$ is between $\frac{1}{2}V^* \log^{(j)} n$ and $V^* \log^{(j)} n$. Note $\Pr[\mathcal{E}_2] \ge \frac{1}{3} \cdot \frac{1}{\log^* n} \cdot \frac{1}{2 \log^{(j)} n}$. Since event \mathcal{E}_2 implies the algorithm gets value at least $\tau \ge \frac{1}{2}V^* \log^{(j)} n$ (because $v_{j+1} > \tau$), its expected value is $\Omega((\log^* n)^{-1}V^*)$, a contradiction.

Now by Claim 13 and Lemma 14, we have $v_j \in [V^*, V^* \cdot \log^{(j)} n]$ for all $j \in [1 \dots \log^* n]$. We still get a contradiction. Let \mathcal{E}_3 be the event that the following three things happen simultaneously: that we select Algorithm iii in §5.1 with $i = \log^* n$, that the highest green element g_{\max} is in interval $\mathbb{I}_{\log^* n+1}$, and that we select s in Algorithm iii such that $\tau := (v_{\log^* n} \log^{(j)} n)/(2^s)$ is between $\frac{1}{2}V^*$ and V^* . Note $\Pr[\mathcal{E}_3] \ge \frac{1}{3\log^* n} \cdot \frac{1}{4} \cdot \frac{1}{2\log^* n}$. Since the event \mathcal{E}_2 implies the algorithm gets value at least $\tau \ge \frac{1}{2}V^*$ (because g_{\max} is in $\mathbb{I}_{\log^* n+1}$), its expected value is $\Omega((\log^* n)^{-2}V^*)$. Thus, we have a contradiction in every case, which means our assumption is incorrect and the algorithm has expected value $\Omega((\log^* n)^{-2}V^*)$.

6 Value Maximization for Matroids

In this section we discuss multiple-choice Byzantine secretary algorithms in the matroid setting.

▶ Definition 15 (Byzantine secretary problem on matroids). Let \mathcal{M} be a matroid over $U = R \cup G$, where elements in $G = \{g_{\max}, g_2, \ldots, g_{|G|}\}$ arrive uniformly at random in [0, 1]. When an element $e \in U$ arrives, the algorithm must irrevocably select or ignore e, while ensuring that the set of selected elements forms an independent set in \mathcal{M} . The leave-one-out benchmark V^* is the highest-value independent subset of $G \setminus \{g_{\max}\}$.

The knapsack results imply $(1 - \varepsilon)$ -competitiveness for uniform matroids as long as the rank r is large enough; we now consider other matroids.

6.1 $O(\log \log n)^2$ -competitiveness for Partition Matroids

A partition matroid is where the elements of the universe are partitioned into parts $\{P_1, P_2, \ldots\}$. Given some integers r_1, r_2, \ldots , a subset of elements is independent if for every *i* it contains at most r_i element from part P_i .

▶ **Theorem 5** (Partition Matroids). There is an algorithm for Byzantine secretary on partition matroids that is $O(\log \log n)^2$ -competitive with the benchmark V^* .

We prove Theorem 5 for simple partition matroids where all $r_i = 1$, i.e., we can select at most one element in each part. This is without loss of generality (up to O(1) approximation) because we can randomly partition each part P_i further into r_i parts and run the simple partition matroid algorithm.

Recall that our single item poly log^{*} n algorithm from §5 no longer works for partition matroids. This is because besides one part we want to get the highest green element in all the other parts. Formally, Claim 13 where we use Dynkin's secretary algorithm in the proof of Theorem 3 fails because it needs at least two green elements. So we need to overcome the lower bound to getting the highest-value green element $v(g_1)$ in Observation 19. We achieve this and design an $O(\log \log n)^2$ -approximation algorithm by making an assumption that the algorithm starts with a polynomial approximation to $v(g_1)$. Although in general this is a strong assumption, it turns out that for partition matroids this assumption is w.l.o.g. because the algorithm may lose the highest green element in one of the parts.

6.1.1 The Algorithm

We define $\log \log n + 1$ time *checkpoints* as follows: the initial checkpoint $T_0 = \frac{1}{2}$, and then subsequent checkpoints $T_i = \frac{1}{2} + \frac{i}{2 \cdot \log \log n}$ for all $i \in [1 \dots \log \log n]$. Now the corresponding intervals are

$$\mathbb{I}_0 := [0, T_0] \quad \text{and} \quad \mathbb{I}_i = \langle T_{i-1}, T_i \rangle \quad \forall i \in [1 \dots \log \log n]$$

$$\tag{8}$$

Let v_0 denote the value of the max element seen by the algorithm in \mathbb{I}_0 .

Now for every part P of the partition matroid, we execute the following algorithm separately. Let v_i for $i \in [1 \dots \log \log n]$ denote the value of the max element seen by the algorithm in part P during interval \mathbb{I}_i . Let V^* denote the element of our benchmark in P. Notice that $v_i \in P$ and V^* cannot be the overall highest green element as we exclude it. We define $4 \log^{1/i} n$ levels for \mathbb{I}_i where *level* j for $j \in [1 \dots 4 \log^{1/i} n]$ is given by elements with values in

$$\Big[\frac{v_{i-1} \cdot \log^{1/i} n}{2^j}, \frac{v_{i-1} \cdot \log^{1/i} n}{2^{j-1}}\Big].$$

We run one of the following algorithms uniformly at random.

- (a) Select an element uniformly at random as discussed in §2.1.
- (b) For every part P, select a random interval $i \in [1 \dots \log \log n]$ and select a random level $j \in [4 \log^{1/i} n]$. Select the first element above $\frac{v_{i-1} \cdot \log^{1/i} n}{2^i}$ in P.
- (c) For every part P, select a random interval $i \in [1 \dots \log \log n]$ and if there is an element with value more than $2^{\log^{1/i} n}$ times the max of all the already seen elements in \mathbb{I}_i , selects it with constant probability, say 1/100.

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6.1.2 The Analysis

Since with constant probability our algorithm selects one of the *n* elements uniformly at random (Algorithm a), we can assume that $v_0 \leq n^2 \cdot V^*$. We always condition on the event that g_{\max} arrives in the interval \mathbb{I}_0 , which happens with constant probability and implies $v_0 \geq v(g_{\max})$. Moreover, we ignore parts *P* where V^* is below $v(g_{\max})/n^2$ because they do not contribute significantly to the benchmark. So from now assume

 $V^*/n^2 \le v_0 \le n^2 \cdot V^*.$

We design an algorithm that gets value $\Omega(V^*/(\log \log n)^2)$ in each part P, which implies Theorem 5 by linearity of expectation over parts.

Let $v_i^{(\text{red})} \in P$ for $i \in [1 \dots \log \log n]$ denote the value of the max red element that the adversary presents in \mathbb{I}_i .

 \triangleright Claim 16. If there exists an $i \in [1 \dots \log \log n]$ with $v_i^{(\text{red})} > V^* \cdot \log^{1/i} n$ then the expected value of the algorithm is $\Omega(V^*/\log \log n)$.

Proof. With constant probability, our algorithm selects a random interval i and selects a random level element in it (Algorithm b). Since w.p. $1/\log \log n$ it selects this i, and w.p. $\frac{1}{4} \log^{1/i} n$ it selects the random level of $v_i^{(\text{red})}$ in \mathbb{I}_i , the algorithm has expected value at least

$$\frac{1}{\log\log n} \cdot \frac{1}{4\log^{1/i} n} \cdot v_i^{(\operatorname{red})} \geq \frac{1}{4\log\log n} \cdot V^*.$$

By the last claim we can assume for all $i \in [1 \dots \log \log n]$, we have $v_i^{(\text{red})} \leq V^* \cdot \log^{1/i} n$.

 \triangleright Claim 17. If there exists an $i \in [1 \dots \log \log n]$ with $v_i^{(\text{red})} < V^*/2^{\log^{1/i} n}$ then the expected value of the algorithm is $\Omega(V^*/(\log \log n)^2)$.

Proof. With constant probability the algorithm guesses one of the intervals i and if there is an element with value more than $2^{\log^{1/i} n}$ times the max of all the already seen elements in \mathbb{I}_i , selects it with constant probability (Algorithm c). With $1/\log\log n$ probability the algorithm selects this particular i and with $1/\log\log n$ probability V^* appears in this interval with value at least $2^{\log^{1/i} n}$ times the max seen element in this interval. Notice there can be at most $O\left(\frac{4\log^{1/i} n}{\log^{1/i} n}\right) = O(1)$ elements with such large jumps in value in this interval. In this case our algorithm selects V^* with constant probability.

Finally, we are only left with the case where for all $i \in [1 \dots \log \log n]$ value $\frac{V^*}{2^{\log^{1/i} n}} \leq v_i^{(\text{red})} \leq V^* \cdot \log^{1/i} n$, which we handle using Algorithm b.

 \triangleright Claim 18. If for all $i \in [1 \dots \log \log n]$ we have

$$\frac{V^*}{2^{\log^{1/i} n}} \le v_i^{(\text{red})} \le V^* \cdot \log^{1/i} n$$

then the expected value of the algorithm is $\Omega(V^*/(\log \log n)^2)$.

Proof. Consider Algorithm b. It selects $i = \log \log n - 1$ w.p. $1/\log \log n$. Moreover, suppose V^* appears in $\mathbb{I}_{\log \log n-1}$. Now since there are only a constant number of levels in this interval, our algorithm selects an element of value at least V^* with constant probability.

We have shown that in every case the algorithm has expected value $\Omega(V^*/(\log \log n)^2)$ for any fixed part *P*. This implies Theorem 5 by linearity of expectation over parts.

6.2 $O(\log n)$ -approx for General Matroids

▶ **Observation 6** (General Matroids). There is an algorithm for Byzantine secretary on general matroids that is $O(\log n)$ -competitive with the benchmark V^* .

Proof. Notice that no element can have weight more than nr times the second max-element because w.p. 1/n our algorithm selects one of the n elements uniformly at random. Given this, condition on the event that the max element with value v lands in the first half of the input. Define $2\log(nr)$ exponentially separated levels as follows:

 $\left[\frac{v}{2^{\log(nr)}}, \frac{v}{2^{\log(nr)-1}}\right\rangle, \left[\frac{v}{2^{\log(nr)-1}}, \frac{v}{2^{\log(nr)-2}}\right\rangle, \dots, \left[\frac{v}{2}, v\right\rangle, \dots, \left[v2^{\log(nr)-1}, v2^{\log(nr)}\right\rangle.$

Since at least one of these intervals contains at least $2\log(nr)$ fraction of OPT, we can guess that interval and run a greedy algorithm, i.e., accept any element with value in that interval or above if it is independent.

7 Conclusion

In this paper we defined a robust model for the secretary problem, one where some of the elements can arrive at adversarially chosen times, whereas the others arrive at random times. For this setting, we argue that a natural is the optimal solution on all but the highest-valued green item (or even simpler, the optimal solution on the green items, minus the single highest-value item). This benchmark reflects the fact that we cannot hope to compete with the red (adversarial) items, and also cannot do well if all the green value is concentrated in a single green item.

We show that for the case where we want to pick K items, or if we have a knapsack of size K, we can get within $(1 - \varepsilon)$ of this benchmark, assuming K is large enough. We can also get non-trivial results for the single-item case, where our benchmark is now the second-highest valued green item. In the ordinal setting where we only see the relative order of arriving elements and the goal is to maximize the probability of getting an element whose value is above the benchmark, we use the minimax principle to show existence of an $O(\log^2 n)$ -approximation algorithm in §4. In the value maximization setting, we give an $O(\log^* n)^2$ -approximation algorithm in §5. We also show $O(\log \log n)$ -competitiveness for partition matroids.

The results above suggest many question. Can we improve the lower bound on the size required for $(1 - \varepsilon)$ -competitiveness? Can we get a constant-competitive algorithm for the single-item case? For the probability-maximization problem, our proof only shows the existence of an algorithm; can we make this constructive? More generally, many of the algorithms for secretary problems seem to overfit to the model, at least in the presence of small adversarial changes: how can we make our algorithms robust?

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A Hard Benchmarks

We show that for the benchmark $V^* := v(g_{max})$, every algorithm has an approximation of at most O(1/n).

▶ **Observation 19** (Lower Bound for g_{max}). Any randomized algorithm for the single-item Byzantine secretary problem cannot select the highest-value good/green item with probability larger than 1/(|R|+1).

Proof. We use Yao's minimax lemma, so it is enough to construct an input distribution \mathcal{B} for which no deterministic algorithm can achieve an approximation better than $\frac{1}{|R|+1}$. The distribution is as follows. The red elements arrive at random times, that is $(t_{r_1}, t_{r_2}, ...t_{r_{|R|}}) \sim U[0, 1]^{|R|}$. The linear ordering among the elements is set such that the red elements are strictly increasing according to their arrival time, or formally: $t_{r_i} > t_{r_j} \implies v(r_i) > v(r_j)$. The maximum element is green and all the other green elements are smaller than all red elements. Formally: $v(g_{max}) > v(r_1) > v(r_2) ... > v(r_{|R|}) > v(g_2) > v(g_3) ... > v(g_{|G|})$. This fully defines the input distribution.

All the arrival times are distinct with probability 1. Let $\mathcal{K}(t)$ denote the information seen by the algorithm up to and including time t. Partition the probability space according to $S := \{t_{g_{max}}, t_{r_1}, t_{r_2}, ...t_{r_{|R|}}\}$ and $L := (t_{g_2}, t_{g_3}, ...t_{g_{|G|}})$. Let $s_1 < s_2 < ... < s_{|R|+1}$ be the elements of S. Let $M_i := \{t_{g_{max}} = s_i\}$. By definition, we have $Pr[M_i|S, L] = \frac{1}{|R|+1}$. Note that, since the red items arrive in increasing order of value and the green item has maximum value, we have $Pr[M_i|S, L, \mathcal{K}(t)] = Pr[M_j|S, L, \mathcal{K}(t)]$ for all $t \leq s_i, s_j$. Therefore, $Pr[M_i|S, L, \mathcal{K}(s_i)] \leq \frac{1}{|R|+2-i}$. In other words, there is no way to distinguish the maximum green element from the red elements before it is too late, that is at the time of the green element's arrival. Thus, by a simple inductive argument, the proof is finished.

Using techniques presented in [7], we can extend this result to the value case as well.

B Relaxing the Assumption that *n* is Known

In this section we extend our results to some settings where n is unknown. Most importantly, observe that all of the results in this paper hold even if n is known only up to a constant factor with at most a constant factor degradation in the quality of the result. As a simple example, note that picking a uniformly random element from an n-element sequence when the assumed number of elements is $\tilde{n} \in [n, 2n]$ will select an element $x \in U$ with probability $p_x \in [\frac{1}{2n}, \frac{1}{n}]$, leading to a degradation in the result by a factor of at most 2, which we typically ignore in this paper.

This still leaves us open to the possibility that we do not even know the scale of n. Surprisingly, it is still possible to "guess" \tilde{n} while only incurring a loss of $\tilde{O}(\log n)$ in the quality, even if there is no prior known upper limit on n.¹ The following claim formalizes this result.

 \triangleright Claim 20. There exists a distribution X over the integers such that for every $n \ge 1$ the probability that the sampled number $\tilde{n} \sim X$ is within a constant factor of n, is at least $1/\tilde{O}(\log n)$.

¹ By $\widetilde{O}(f(n))$ we mean $f(n) \cdot \operatorname{poly}(\log f(n))$.

Proof. Consider the sequence $a_k := \frac{1}{k(\log_2 k)(\log_2 \log_2 k)^2}$ defined for $k \ge 2$. It is well-known that this sequence converges, i.e., $\sum_{k=2}^{\infty} a_k = O(1)$. A simple way to see this is by noting that a non-negative decreasing sequence $(A_k)_k$ converges if and only if $(2^k A_{2^k})_k$ converges [40, Thm 3.27].

Let $\sum_{k}^{\infty} A_k \sim \sum_{k}^{\infty} B_k$ be the equivalence relation denoting that $(A_k)_k$ and $(B_k)_k$ either both converge or both diverge. Then the above fact implies that $\sum_{k}^{\infty} \frac{1}{k \log_2 k (\log_2 \log_2 k)^2} \sim \sum_{k}^{\infty} \frac{1}{k^2}$, where the last sequence clearly converges.

We can assume without loss of generality that $n \ge 100$ by handling those cases separately. The strategy for guessing the estimate \tilde{n} is now immediate: we sample \tilde{n} from $\mathbb{Z}_{\ge 2}$ according to the distribution $\Pr[\tilde{n} = k] = a_k/Z$ where $Z := \sum_{k=100}^{\infty} a_k = O(1)$. We observe that $\Pr[\tilde{n} \in \langle 2^{k-1} \dots 2^k]] \ge \frac{1}{2} 2^k a_{2^k}/Z = \tilde{\Omega}(1/k)$. Let k' be the unique index such that $n \in \langle 2^{k'-1} \dots 2^{k'}]$, hence $k' = \Theta(\log n)$. Then $\Pr[\tilde{n} \in \langle 2^{k'} \dots 2^{k'+1}]] = \tilde{\Omega}(1/k') = \tilde{\Omega}(1/\log n)$. But also in that case we have that $n \in [\tilde{n}/4, \tilde{n}]$ and we are done.

Finally, consider an important case where the fraction of red elements is bounded away from $1 - \Omega(1)$. This is a reasonable assumption for most applications, e.g., online auctions, where we do not expect that most of the arrivals will be chosen by an adversary. By simply observing the first half of the sequence, i.e., $[0, 1/2\rangle$, we can typically estimate n up to a constant while degrading the expected output of our algorithms by at most a constant factor.

 \triangleright Claim 21. If there is a constant $\varepsilon < 1$ such that the fraction of red elements $\frac{|R|}{|R|+|G|} \le \varepsilon$ then we can estimate n up to a constant factor by time t = 1/2.

Proof. We run a simple preprocessing step to estimate n up to a constant factor by t = 1/2. Notice that the expected number of green elements to arrive in the interval [0, 1/2) is $0.5 \cdot |G| = 0.5 \cdot n(1 - \varepsilon) = \Omega(n)$. Since by simple Chernoff bounds this means that w.h.p. we see $\Omega(n)$ elements in the first half, we run a simple algorithm that does not select any element till t = 1/2, and then use the number of elements that arrive in [0, 1/2) as an estimate of n.

C Minimax

In this section we argue that an α -payoff (i.e., the probability of selecting the second-max element or better is at least α) known distribution algorithm for the ordinal single-item Byzantine secretary implies an α -payoff algorithm for the general, worst-case input, setting. This can be directly modeled as a two-player game where player A chooses an algorithm \mathcal{A} and player B chooses a distribution over the input instances \mathcal{B} . Our coveted result would go along the lines of

$$\sup_{\mathcal{A}} \inf_{\mathcal{B}} K(a, b) = \inf_{\mathcal{B}} \sup_{\mathcal{A}} K(\mathcal{A}, \mathcal{B}),$$

where $K(\mathcal{A}, \mathcal{B})$ denotes the payoff when we run algorithm \mathcal{A} on the input distribution \mathcal{B} . The left-hand side denotes the worst-case input setting, while the right-hand side denotes the known distribution setting.

The main challenge in proving such a claim stems from the infiniteness of the set of algorithms and set of input distributions. Indeed, if one makes no finiteness assumption for either A or B, the Minimax property can fail even for relatively well-behaved two-player games [37]. On the other hand if both A and B would be finite, then the result would follow from the classic Von Neumann's Minimax [36].

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▶ Fact 22 (Von Neumann's Minimax). Let A and B be finite sets. Denote by $\mathcal{D}(A)$ and $\mathcal{D}(B)$ distributions over A and B, respectively. Then for any matrix of values $K : A \times B \to \mathbb{R}$ it holds that

$$\max_{a \in \mathcal{D}(A)} \min_{b \in \mathcal{D}(B)} K(a, b) = \min_{b \in \mathcal{D}(B)} \max_{a \in \mathcal{D}(A)} K(a, b).$$
(9)

The infiniteness of the sets stems from the arrival times being in the infinite set [0, 1]. To solve this issue, we slightly modify our algorithm by discretizing [0, 1]. Let $N = n^3, \mathcal{T} :=$ $\{\frac{0}{N}, \frac{1}{N}, \frac{2}{N}, \ldots, \frac{N}{N}\}$ and $\Phi : [0, 1] \to \mathcal{T}, \Phi(t) := \lfloor N \cdot t \rfloor / N$ be the discretizing function. We modify our algorithm in the following way: apply Φ to the input distribution \mathcal{B} , as well as to every arrival time. Note that the elements are presented to the algorithm exactly as before, it just *pretends* they arrive in discrete time steps. We can assume $\Phi(t_e) \neq \Phi(t_g)$ for every $e \in U, g \in G, e \neq g$ (otherwise, we say the algorithm loses), since this happens with at most $n^2/N = o(1)$ probability. Using completely analoguous techniques as in Section 4 we can show this algorithm is $\Omega(1/\log^2 n)$ -competitive.

We note that a randomized algorithm is simply a distribution over deterministic algorithms. Hence our second goal is to argue that the number of distinct deterministic algorithms is finite (i.e., bounded by a function of n). To this end we have to specify how we represent them with at least some formality. We identify a deterministic algorithm \mathcal{A} with a function that gets evaluated each time a new element arrives; its parameter is the information history $\mathcal{K}(t)$ (t being the current time) represented in any appropriate format; its output is $\{\perp, \top\}$ representing whether to select the current element. For concreteness, the information history consists of (t_e, π_e) for every element e that arrived before the function call, where $t_e \in \mathcal{T}$ is the discretized arrival time (after applying Φ) and $\pi_e \in [0 \dots n-1]$ is the relative value order of e with respect to prior arrived elements. The number of distinct histories is bounded by $((N+1)n)^n$, a function of n; therefore the set of deterministic algorithms, i.e., functions from the history to $\{\perp, \top\}$, is also bounded.

We remember that an input distribution is simply a distribution over "pure" inputs. Note that the payoff of a deterministic algorithm for a specific input depends only on the following: $\Phi(t_r)$ for every red element; $\pi \in S_n$, the permutation representing the total order among the elements; and $\pi_t \in S_{|R|}$, the permutation denoting the order in which the red elements arrive (since red elements can have the same discretized arrival time, but an arbitrary order in which they are presented to the algorithm). The above discretization makes the number of pure inputs at most $(n!)^2 \cdot (N+1)^n$, i.e., bounded by a function of n. The reader can refresh their memory about the representation of pure inputs by reviewing the introduction to Section 4.

Finally, for our discretized algorithm, we proved that the set of pure inputs with different payoffs, as well as the number of deterministic algorithms is bounded by a function of n. Therefore, for a fixed n, both numbers are finite. We invoke the Von Neumann's Minimax (Fact 22) to conclude that the best result in the known distribution setting and worst-case input setting are equivalent, recovering the following theorem.

► **Theorem 4** (Ordinal Single-item Algorithm). There is a randomized algorithm for the ordinal Byzantine secretary which selects an element of value at least the second-largest green item with probability $\Omega(1/\log^2 n)$.