
Adaptive Sampling for Clustered Ranking

Sumeet Katariya
katariya@wisc.edu

Lalit Jain
lalitj@umich.edu

Nandana Sengupta
nandana@uchicago.edu

James Evans
jevans@uchicago.edu

Robert Nowak
rdnowak@wisc.edu

Abstract

We consider the problem of sequential or active *clustered ranking*, where the goal is to sort items according to their means into clusters of pre-specified sizes, by adaptively sampling from their reward distributions. This setting is useful in many social science applications, where an *approximate* rank of *every* item is desired. In contrast, a complete ranking of the items requires a large number of samples if the means are close, while only finding the top few items as is done in recommender systems reveals no information about the ranks of other items. We propose a computationally efficient PAC algorithm **LUCBRank** to solve this problem, and derive an upper bound on its sample complexity. We also derive a nearly matching distribution-dependent lower bound. Experiments on synthetic as well as real-world data arising from assessing the safety scores of images show that **LUCBRank** performs better than state-of-the-art baseline methods, even when these methods have the advantage of knowing the underlying parametric model.

1 Introduction

We consider the problem of efficiently sorting items according to their means into clusters of pre-specified sizes, which we refer to as *clustered ranking*. In many big-data applications, finding the *total* ranking can be infeasible and / or unnecessary, and we may only be interested in the top items, bottom items, or quantiles. Consider for instance the problem of assessing the safety of neighborhoods from pairwise comparisons of Google Streetview images as is done in the Place

Pulse project (Naik et al., 2014) aimed at developing social policy (Dubey et al., 2016). Finding a complete ordering of the images in this case is impractical because many images are difficult to compare i.e., their safety scores are very close (see Section 7.2). Furthermore, a total ordering may be unnecessary from a public policy point of view: one may only be interested in the unsafe-appearing images, or the *approximate* rank of every image on the safe-unsafe spectrum.

Motivated by this applications, we model the clustered ranking problem as follows. Given K random variables, $c \geq 2$ clusters, and cluster boundaries $1 \leq \kappa_1 < \kappa_2 < \dots < \kappa_{c-1} < \kappa_c = K$, the goal is to reliably identify the κ_1 random variables with the highest means, the $\kappa_2 - \kappa_1$ random variables with the highest means among the remaining $K - \kappa_1$ random variables, and so on, by observing draws from their distributions (for a precise formulation see Section 4). The focus of this paper is on algorithms to achieve the clustering by adaptively requested draws. The clustered ranking setting applies to the scenarios above, and also subsumes many well-studied problems. The problem of finding the best item corresponds to $\kappa_1 = 1, \kappa_2 = K$. The problem of finding the top- m items corresponds to $\kappa_1 = m, \kappa_2 = K$. The problem of sorting the items into c equal-sized clusters corresponds to $\kappa_i = \text{round}(iK/c), 1 \leq i \leq c$. Finally, the complete ranking can be obtained by setting $\kappa_i = i, 1 \leq i \leq K$.

The problem of *completely sorting* items is in general hard in real-world applications, and does not exhibit gains from adaptivity. Maystre and Grossglauser (2017) who analyze the performance of Quicksort, observe in their real-world experiments:

The improvement is noticeable but modest. We notice that item parameters are close to each other on average; ... This is because there is a considerable fraction of items that have their parameters (means) very close to one another ... Figuring out the exact order of these images is therefore difficult and probably of marginal value. (1)

The fact that *adaptivity doesn't help for complete ranking* is true not just for Quicksort, but other adaptive algorithms as well - as we observe in our experiments. Adaptivity does however help for clustered ranking, and this can be explained. Consider the case when the K item means are equally separated, with a gap Δ between consecutive means. Correctly ordering any two consecutive items requires $\Omega(1/\Delta^2)$ samples, and thus *any* algorithm would require $\Omega(K/\Delta^2)$ to find a total ordering. A non-adaptive algorithm sampling the items uniformly would gather approximately equal samples from every item, and hence will find the correct ranking after roughly these many samples (up to perhaps log factors). Thus adaptivity doesn't help in this case. However, if the goal is to find only the quartiles say, an adaptive algorithm can quickly stop sampling items that are far from the quartile boundaries and gain over non-adaptive algorithms.

In this work, we make the following contributions:

- We explain why adaptive methods - and in particular Quicksort - are ineffective in high-noise regimes, and motivate the clustered ranking setting (Section 3).
- We precisely formulate the online PAC-clustered ranking problem with error tolerance ϵ and failure probability δ that can model real-valued as well as pairwise comparison feedback (Section 4).
- We propose a non-parametric PAC UCB-type algorithm **LUCBRank** to solve this problem. To the best of our knowledge, this is the first UCB-type algorithm for ranking (Section 5).
- We analyze the sample complexity of **LUCBRank** and prove an upper bound which is inversely proportional to the distance of the item to its closest cluster boundary, where the distance is measured in terms of Chernoff information (Section 6).
- We also prove a nearly matching distribution-dependent lower bound. The contribution of an item to the lower bound is inversely proportional to the distance of the item to the closest item in an adjacent cluster, with distance in this case measured using KL-divergences (Section 6.3).
- Finally, we compare the performance of our algorithm to several baselines on synthetic as well as real-world data gathered using MTurk, and observe that it performs 2 - 3x better than existing algorithms even when they have the advantage of knowing the underlying parametric model (Section 7).

1.1 Ranking using Pairwise Comparisons

As stated above and later in Section 4, our model assumes access to direct feedback for the queried item (in the form of a 5-star rating for example). However, any algorithm designed to solve the direct-feedback clustered ranking problem can also be used with pairwise comparison feedback using Borda reduction (Jamieson et al., 2015b). According to this technique, whenever the algorithm asks to draw a sample from item i , we compare item i to a randomly chosen item j , and ascribe a reward of 1 to item i if i wins the duel, and 0 otherwise. This is equivalent to the rewards being sampled from a Bernoulli distribution with means given by the Borda scores of the items. The Borda score of item i is defined as

$$p_i := \frac{1}{K-1} \sum_{j \neq i} \mathbb{P}(i > j). \quad (2)$$

Terminology: We use the term direct-feedback or pure-rewards or real-rewards to indicate a setting where we can sample directly from the item's reward distribution. Our algorithm is stated for this setting. In contrast, the pairwise-comparison or dueling setting is one where we compare two items and receive 1-bit feedback about who won the duel. Our algorithm can be translated to this setting using the technique outlined in Section 1.1.

2 Related Work

There is extensive work on ranking from noisy pairwise comparisons, we refer the reader to excellent surveys by Yan; Busa-Fekete and Hüllermeier (2014); Agarwal (2016). We discuss prominent among these next.

2.1 Ranking from Pairwise Comparisons

The pairwise comparison matrix P (where $P_{ij} = \mathbb{P}(i > j)$) and assumptions on it play a major role in the design of ranking algorithms (Agarwal, 2016). A sequence of progressively relaxed assumptions on P can be shown where ranking methods that work under restrictive assumptions fail when these are slightly relaxed (Rajkumar and Agarwal, 2014; Rajkumar et al., 2015). Spectral ranking algorithms have been proposed when comparisons are available for a fixed set of pairs (Negahban et al., 2012a,b); this corresponds to a partially observed P matrix. Braverman and Mossel (2009); Wauthier et al. (2013) propose and analyze algorithms for the noisy-permutation model; this corresponds to a P matrix which has two types of entries: $1 - p$ in the upper triangle and p in the lower triangle (assuming the true ordering of the items is $1 \dots K$). They also focus on settings where queries cannot be

repeated. Our work makes no assumptions on the P matrix and ranks items using their Borda scores. This is important given the futility of parametric models to model real-life scenarios (Shah et al., 2016).

Quicksort is another highly recommended algorithm for ranking using noisy pairwise comparisons (Yan). Maystre and Grossglauser (2017) study Quicksort under the BTL noise model, and Alonso et al. (2003) analyze Quicksort under the noisy permutation model. We comment on these in Section 3.

Jamieson and Nowak (2011) propose an algorithm for active ranking from pairwise comparisons when points can be embedded in Euclidean space. Ailon (2012) consider ranking when query responses are fixed. More recently, Agarwal et al. (2017) consider top- m item identification and ranking under limited rounds of adaptivity, Falahatgar et al. (2017) consider the problem of finding the maximum and ranking assuming strong-stochastic transitivity and the stochastic-triangle inequality. We do not need these assumptions.

Our setting is closest to the setting proposed by Heckel et al. (2016), in the context of ranking using pairwise comparisons. Our setting however applies to real-valued rewards as well as pairwise comparison feedback. Furthermore, our setting incorporates the notion of ϵ -optimality which allows the user to specify an error tolerance (Even-Dar et al., 2006). This is important in practice if the item means are very close to each other. Finally, as they note, the algorithm they propose is an elimination-style algorithm, our algorithm is UCB-style; it is known that the latter perform better in practice (Jiang et al., 2017).

2.2 Relation to Bandits

The idea of sampling items based on lower and upper confidence bounds is well-known in the bandits literature (Auer, 2002). However, these algorithms either focus on finding the best or top- m items (Audibert and Bubeck, 2010; Kalyanakrishnan et al., 2012; Kaufmann et al., 2015; Chen et al., 2017), or on minimizing regret (Bubeck et al., 2012). This is the first work to our knowledge that employs this tool for ranking.

3 Motivation

We argue that existing adaptive methods offer no gains over their non-adaptive counterparts, and clustered ranking is the panacea for many real-world applications. We provide brief theoretical justification for this claim in the discussion after quote (1), and empirically verify this behavior in Fig. 4. In this section, we focus on Quicksort, because it has been well-studied

under multiple noise models. Quicksort has optimal sample complexity when comparisons are noiseless (Sedgewick and Wayne, 2011) and is naturally appealing when comparisons are noisy (Yan; Maystre and Grossglauser, 2017). Intuitively it feels like the right thing to do - by comparing an item with the pivot and putting it left or right appropriately, Quicksort performs a binary search for the true position of an item. However it is far from optimal under two noise models as we argue next.

First, consider the noisy-permutation (NP) noise model (Feige et al., 1994) where the outcomes of pairwise comparisons are independently flipped with an error probability p . In the first stage of Quicksort, every item that is compared with the pivot and put in the wrong bucket contributes on average $\frac{K}{2}$ to the Kendall tau distance (total number of inverted pairs). Now, Kp items are put in the wrong bucket on average in the first stage of Quicksort, and hence the total number of inverted pairs is at least $\Omega(K^2p)$. Alonso et al. (2003) show that $\Theta(K^2p)$ is indeed the expected number of inversions. Surprisingly though, they also show that even when $p \simeq 1/K$, the expected number of inversions is $\Theta(K^2p)$. They conjecture that in order to obtain $\Theta(K)$ expected inversions, p needs to go down faster than $1/K$, like $\frac{1}{K \log K}$. This is far from optimal because $O(K)$ Kendall tau distance can be obtained with $K \log K$ comparisons even for constant p (Braverman and Mossel, 2009). As the above calculation shows, Alonso et al. (2003) conjecture that this is because Quicksort is extremely “brittle”:

the main contribution (to the total inversions) comes from the “first” error, in some sense

One may be able to get rid of this lack of robustness by repeating queries, but this may require prior knowledge of the error probability p and parameter tuning. Even if this were somehow possible, a good model for real-world examples is one where p increases with K , since it becomes more difficult to compare adjacent items in the true ranking as K increases. Quicksort certainly fails in this regime.

The other class of well-studied noise models are the Bradley-Terry-Luce (BTL) (Bradley and Terry, 1952) or Thurstone (Thurstone, 1927) models, which assume a K -dimensional weight vector that measures the quality of each item, and the pairwise comparison probabilities are determined via some fixed function of the qualities of pair of objects. These models are more realistic than the NP model since under these models, comparisons between items that are far apart in the true ranking are less noisy than those between nearby items. Maystre and Grossglauser (2017) analyze the expected number of inversions of Quicksort under the BTL model, and show that when the average gap be-

tween adjacent items is λ , the expected number of inversions is $O(\lambda^{-3})$. They note however that real-world datasets have extremely small λ ($\hat{\lambda}^{-1} = 376$ in their experiments) and Quicksort performs no better than random (see quote (1)). We make similar observations about the inefficacy of Quicksort (and other adaptive algorithms) in our real-world experiments (see Fig. 4).

Clustered ranking converts a high-noise problem to a low-noise one. Even if the gap between adjacent items is small, most items are far from their nearest cluster boundary, and an adaptive algorithm can stop sampling these early.

4 Setting

In this section, we precisely formulate the clustered ranking setting. For ease of reference, we use terminology from the bandits literature and refer to an item / element as arm. Also, pulling or drawing an arm is equivalent to sampling from the item's reward distribution.

Consider a multi-armed bandit with K arms. Each arm a corresponds to a Bernoulli distribution with an unknown mean p_a , denoted $\mathcal{B}(p_a)$. A draw / pull of arm a yields a reward from distribution $\mathcal{B}(p_a)$. Without loss of generality, assume the arms are numbered so that $p_1 \geq p_2 \geq \dots \geq p_K$.

Given an integer $c \geq 2$ representing the number of clusters, let $\kappa := (\kappa_1, \kappa_2, \dots, \kappa_c)$ be a collection of positive integers such that $1 \leq \kappa_1 < \kappa_2 < \dots < \kappa_c = K$. Any such collection of positive integers defines a partition of $[K]$ into c disjoint sets of the form

$$M_1^* := \{1, \dots, \kappa_1\}, M_2^* := \{\kappa_1 + 1, \dots, \kappa_2\}, \dots, \\ \dots, M_c^* := \{\kappa_{c-1} + 1, \dots, K\}. \quad (3)$$

To solve the *clustered ranking problem* given a set of cluster boundaries κ , an algorithm may sample arms of the K -armed bandit and record the result of its pulls; the algorithm is required to terminate and cluster the arms into an ordered set of disjoint sets of the form (3). We refer to this output as a *clustered ranking*.

We next define the notion of ϵ -tolerance. For some fixed tolerance $\epsilon \in [0, 1]$, let $M_{i,\epsilon}^*$ be the set of all arms that should be in cluster i upto a tolerance ϵ , i.e.

$$M_{i,\epsilon}^* := \{a : p_{\kappa_{i-1}+1} + \epsilon \geq p_a \geq p_{\kappa_i} - \epsilon\},$$

(with the convention that $p_{\kappa_0} = 0$). Note that the true set of arms in cluster i : $M_i^* := \{\kappa_{i-1} + 1, \dots, \kappa_i\}$, is a subset of $M_{i,\epsilon}^*$; the latter set contains in addition arms that are ϵ close to the boundary.

For a given mistake probability $\delta \in [0, 1]$ and a given error tolerance $\epsilon \in [0, 1]$, we call an algorithm (ϵ, δ) -PAC if, with a probability greater than $1 - \delta$, after using

a finite number of samples, it returns a rank for each arm such that the i^{th} ranked cluster according to the returned ranking is a subset of $M_{i,\epsilon}^*$ for all $1 \leq i \leq c$. Formally, if $\sigma(a)$ is the rank of arm a returned by the algorithm after using a finite number of samples, we can define the empirical cluster i as

$$\hat{M}_i := \{a : \kappa_{i-1} + 1 \leq \sigma(a) \leq \kappa_i\},$$

and we say the algorithm is (ϵ, δ) -PAC if

$$\mathbb{P}(\exists i \text{ such that } \hat{M}_i \not\subseteq M_{i,\epsilon}^*) \leq \delta. \quad (4)$$

5 Algorithm

Let $\kappa = (\kappa_1, \dots, \kappa_c = K)$ be the cluster boundaries. We describe here the **LUCBRank** algorithm using generic confidence intervals $\mathcal{I}_a = [L_a(t), U_a(t)]$, where t is the round of the algorithm. Let $N_a(t)$ be the number of draws of arm a , $S_a(t)$ be the sum of the rewards from arm a up to round t . Let $\hat{p}_a(t) = \frac{S_a(t)}{N_a(t)}$ be the corresponding empirical mean reward. Sort the arms in the decreasing order of their empirical mean rewards, and for $1 \leq i \leq c - 1$, let $J_i(t)$ denote the κ_i arms with the highest empirical mean rewards. Define

$$l_t^i := \arg \min_{a \in J_i(t)} L_a(t), \quad u_t^i := \arg \max_{a \notin J_i(t)} U_a(t) \quad (5)$$

to be the two *critical* arms from $J_i(t)$ and $J_i^c(t)$ that are likely to be misclassified (see Fig. 1).

Algorithm 1 contains the pseudocode of **LUCBRank**, which is also depicted in Fig. 1. The algorithm maintains active cluster boundaries in the set C , where a cluster boundary i is active if the overlap of confidence intervals in J_i and J_i^c is not less than ϵ . In every round, it samples both the critical arms at every active cluster boundary (lines 10-12). At the end of every round, it checks if the critical arms at any boundary are separated according to the tolerance criterion, and removes such boundaries from the active set (lines 17-21). For our experiments, we use KL-UCB (Garivier and Cappé, 2011) confidence intervals. For an exploration rate $\beta(t, \delta)$, the KL-UCB upper and lower confidence bounds for arm a are calculated as

$$U_a(t) := \max\{q \in [\hat{p}_a(t), 1] : N_a(t)d(\hat{p}_a(t), q) \leq \beta(t, \delta)\}, \quad (6)$$

$$L_a(t) := \min\{q \in [0, \hat{p}_a(t)] : N_a(t)d(\hat{p}_a(t), q) \leq \beta(t, \delta)\}. \quad (7)$$

where $d(x, y)$ is the Kullback-Leibler divergence between two Bernoulli distributions, given by $d(x, y) = x \log \frac{x}{y} + (1 - x) \log \frac{1-x}{1-y}$.

LUCBRank can also be easily modified for pairwise-comparison queries: whenever the algorithm calls for drawing an arm i , duel arm i with another arm chosen uniformly at random.

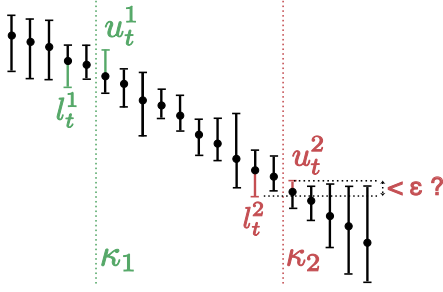


Figure 1: A visualization of LUCBRank on a bandit instance with $K = 20$ arms, $c = 3$ clusters, with boundaries at $\kappa_1 = 5, \kappa_2 = 15$. Also shown are the critical arms l_t^i, u_t^i pulled at each boundary. The algorithm stops sampling a boundary when confidence interval overlap is less than ϵ .

6 Analysis

We prove the accuracy of LUCBRank in Theorem 1, and give an upper bound on the sample complexity in Theorem 2. Our distribution-dependent lower bound for the sample complexity of any δ -PAC algorithm is stated in Theorem 3. All proofs can be found in the Appendix. Recall that $1 \leq \kappa_1 < \kappa_2 < \dots < \kappa_{c-1} < \kappa_c = K$ are the cluster boundaries.

6.1 PAC Guarantee

Theorem 1 gives choices of $\beta(t, \delta)$ such that LUCBRank is correct with probability at least δ , in the sense defined by Eq. (4).

Theorem 1. LUCBRank using $\beta(t, \delta) = \log(\frac{k_1 K t^\alpha}{\delta}) + \log \log(\frac{k_1 K t^\alpha}{\delta})$ with $\alpha > 1$ and $k_1 > (\frac{c-1}{2})^\alpha + \frac{2e}{\alpha-1} + \frac{4e}{(\alpha-1)^2}$, is correct with probability $1 - \delta$.

6.2 Sample Complexity

To state our results on sample complexity, we need the notion of Chernoff information (Cover and Thomas, 2012).

Chernoff Information: Consider two Bernoulli distributions $\mathcal{B}(x)$ and $\mathcal{B}(y)$, and let $d(x, y)$ denote the KL-divergence between these distributions. The Chernoff information $d^*(x, y)$ between these two Bernoulli distributions is defined by

$$d^*(x, y) := d(z^*, x) = d(z^*, y)$$

where z^* is the unique z such that $d(z, x) = d(z, y)$. Next we introduce some notation. For an arm a , let $g(a)$ (read group of arm a) denote the index of the cluster that arm a belongs to. Formally,

$$g(a) := \min\{1 \leq i \leq c : p_a \leq p_{\kappa_i}\}. \quad (8)$$

Let $b_i \in [p_{\kappa_i}, p_{\kappa_i+1}]$, $1 \leq i \leq c-1$ be any points in the cluster boundary gaps, and $b := (b_1, b_2, \dots, b_{c-1})$.

Algorithm 1 LUCBRank

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1: Input:  $\epsilon > 0$ , cluster boundaries  $1 \leq \kappa_1, \dots, \kappa_c = K$ 
2:  $t = 1$  (round number),  $C = \{1, \dots, c-1\}$  // active cluster boundaries
3:
4: for  $a = 1, \dots, K$  do
5:   Sample arm  $a$ , compute confidence bounds  $U_a(1)$  and  $L_a(1)$ .
6: end for
7:
8: while  $C \neq \emptyset$  do
9:   // Sample active cluster boundaries
10:  for  $i \in C$  do
11:    Sample item  $l_t^i$ .
12:    Sample item  $u_t^i$ .
13:    (If pairwise comparing, this means compare item  $l_t^i$  to a random other item, and compare item  $u_t^i$  to a random other item. See Section 1.1).
14:  end for
15:   $t = t + 1$ 
16:  Update rewards, number of plays, and confidence bounds. Compute  $l_t^i, u_t^i$  for  $i \in C$  (see Eq. (5)).
17:
18:  // Eliminate unambiguous cluster boundaries
19:  for  $i \in C$  do
20:    if  $U_{u_t^i}(t) - L_{l_t^i}(t) < \epsilon$  then
21:       $C = C \setminus i$ 
22:    end if
23:  end for
24: end while
25:
26: Return arms sorted by their empirical mean rewards.

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Define

$$\Delta_b^*(a) := \begin{cases} d^*(p_a, b_1) & a \in \{1, \dots, \kappa_1\} \\ \min(d^*(p_a, b_{g(a)-1}), d^*(p_a, b_{g(a)})) & a \in \{\kappa_1 + 1, \dots, \kappa_{c-1}\} \\ d^*(p_a, b_{c-1}) & a \in \{\kappa_{c-1} + 1, \dots, K\} \end{cases} \quad (9)$$

to be the “distance” of each arm from the closest cluster boundary. Our result on the sample complexity of LUCBRank is stated in terms of the quantity $H_{\epsilon, b}^*$, where

$$H_{\epsilon, b}^* := \sum_{a \in \{1, \dots, K\}} \frac{1}{\max(\Delta_b^*(a), \epsilon^2/2)}. \quad (10)$$

Our result on the sample complexity of LUCBRank is stated in Theorem 2.

Theorem 2. Let $b = (b_1, b_2, \dots, b_{c-1})$, where $b_i \in [p_{\kappa_i}, p_{\kappa_i+1}]$. Let $\epsilon > 0$. Let $\beta(t, \delta) = \log(\frac{k_1 K t^\alpha}{\delta}) +$

$\log \log(\frac{k_1 K t^\alpha}{\delta})$ with $k_1 > (\frac{c-1}{2})^\alpha + \frac{2e}{\alpha-1} + \frac{4e}{(\alpha-1)^2}$. Let τ be the random number of samples taken by LUCBRank before termination. If $\alpha > 1$,

$$\mathbb{P}\left(\tau \leq 2C_0(\alpha)H_{\epsilon,b}^* \log\left(\frac{k_1 K (2H_{\epsilon,b}^*)^\alpha}{\delta}\right)\right) \geq 1 - \delta$$

where $C_0(\alpha)$ is such that $C_0(\alpha) \geq (1 + \frac{1}{e})(\alpha \log(C_0(\alpha)) + 1 + \frac{\alpha}{e})$.

6.3 Distribution-Dependent Lower Bound

In this section, we state our non-asymptotic lower bound on the expected number of samples needed by any δ -PAC algorithm to cluster and rank the arms into groups of sizes $\kappa = (\kappa_1, \kappa_2 - \kappa_1, \dots, K - \kappa_{c-1})$. For simplicity, we focus on the case $\epsilon = 0$. The proof of the lower bound uses standard change of measure arguments (Kaufmann et al., 2015), which requires some continuity and well-separation assumptions. We state these next.

We consider the following class of bandit models where the clusters are unambiguously separated, i.e.

$$\mathcal{M}_\kappa = \{p = (p_1, \dots, p_K) : p_i \in \mathcal{P}, p_{\kappa_i} > p_{\kappa_i+1}, 1 \leq i < c\}, \quad (11)$$

where \mathcal{P} is a set that satisfies

$$\forall p, q \in \mathcal{P}^2, p \neq q \Rightarrow 0 < KL(p, q) < +\infty.$$

We also assume the following:

Assumption 1. For all $p, q \in \mathcal{P}^2$ such that $p \neq q$, for all $\alpha > 0$,

there exists $q_1 \in \mathcal{P}$: $KL(p, q) < KL(p, q_1) < KL(p, q) + \alpha$ and $\mathbb{E}_{X \sim q_1}[X] > \mathbb{E}_{X \sim q}[X]$,

there exists $q_2 \in \mathcal{P}$: $KL(p, q) < KL(p, q_2) < KL(p, q) + \alpha$ and $\mathbb{E}_{X \sim q_2}[X] < \mathbb{E}_{X \sim q}[X]$.

To state our lower bound, we need to define for each arm a , another “distance” from the boundary, similar to Eq. (9). Define

$$\Delta_\kappa^{\text{KL}}(a) := \begin{cases} KL(p_a, p_{\kappa_1+1}) & a \in \{1, \dots, \kappa_1\} \\ \min(KL(p_a, p_{\kappa_{g(a)}-1}), KL(p_a, p_{\kappa_{g(a)}+1})) & a \in \{\kappa_1 + 1, \dots, \kappa_{c-1}\} \\ KL(p_a, p_{\kappa_{c-1}}) & a \in \{\kappa_{c-1} + 1, \dots, K\}, \end{cases} \quad (12)$$

where $g(a)$ defined in Eq. (8) is the cluster that arm a belongs to. We highlight the differences from Eq. (9). First, the Chernoff information in Eq. (9) is replaced with KL-divergence in Eq. (12), and second, the distance is measured with the closest arm in either adjacent cluster here, as opposed to a point in the gap between the clusters in Eq. (9).

Our lower bound involves the quantity

$$\sum_{a \in \{1, \dots, K\}} \frac{1}{\Delta_\kappa^{\text{KL}}(a)} \quad (13)$$

and is as follows:

Theorem 3. Let $p \in \mathcal{M}_\kappa$, and assume that \mathcal{P} satisfies Assumption 1; any clustered ranking algorithm that is δ -PAC on \mathcal{M}_κ satisfies, for $\delta \leq 0.15$,

$$\mathbb{E}_p[\tau] \geq \left\lceil \sum_{a \in \{1, \dots, K\}} \frac{1}{\Delta_\kappa^{\text{KL}}(a)} \right\rceil \log\left(\frac{1}{2.4\delta}\right)$$

6.4 Remarks

- The tightest high-probability upper bound is obtained by setting b equal to $\arg \min_{b: b_i \in [p_{\kappa_i}, p_{\kappa_i+1}]} H_{\epsilon,b}^*$ in Theorem 2.

- Although stated for Bernoulli distributions, the results in this paper can easily be extended to rewards in the exponential family (Garivier and Cappé, 2011) by using the appropriate d function.

7 Experiments

7.1 Ranking from Direct Observations

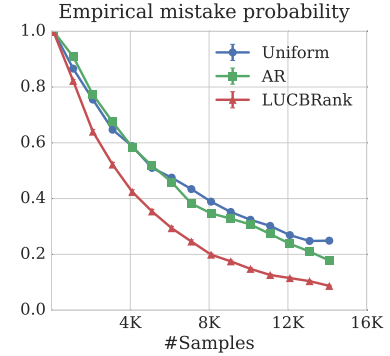


Figure 2: Exp 1 (description in text)

In our first experiment, we compare LUCBRank with uniform sampling and the Active Ranking (AR) algorithm (Heckel et al., 2016). The AR algorithm is an adaptation of the successive elimination approach to solve the clustered ranking problem. It maintains a set of unranked arms and plays every arm in this set, removing an arm from the set when it is confident of the cluster the arm belongs to. Although originally developed for pairwise comparison feedback, the AR algorithm can easily be adapted to the pure-rewards setting.

We look at the bandit instance B ($K = 15$; $p_1 = \frac{1}{2}$; $p_a = \frac{1}{2} - \frac{a}{40}$ for $a = 2, 3, \dots, K$) studied in the context of finding the best-arm (Bubeck et al., 2013). We consider the problem of finding the top-3 and the bottom-3 arms, which corresponds to $\kappa_1 = 3, \kappa_2 = 12$.

In Fig. 2, we record the probability (averaged over 1000 simulations) that the empirical clusters returned

by the algorithm do not match the true clusters. We set $\delta = 0.1$ for both LUCBRank and AR, and $\epsilon = 0$ in LUCBRank to have a fair comparison with AR. We see that the mistake probability drops faster for LUCBRank than for AR.

7.2 Ranking from Pairwise Comparisons

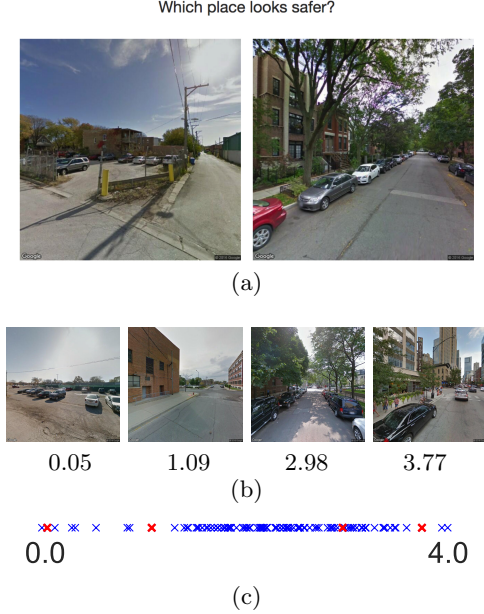


Figure 3: (a) A sample query on NEXT. (b) Four sample images and their estimated BTL scores beneath. (c) Scatter plot of all the BTL scores, with the sample image markers highlighted.

To measure the performance of our algorithm on real-world data, we selected $K = 100$ Google Streetview images in Chicago, and collected 6000 pairwise responses on MTurk using NEXT (Jamieson et al., 2015a), where we asked users to choose the safer-looking image out of two images. This experiment is similar to the Place Pulse project (Naik et al., 2014), where the objective is to assess how the appearance of a neighborhood affects its perception of safety. Fig. 3(a) shows a sample query from our experiment. We estimated the safety scores of these Streetview images from the user-responses by fitting a Bradley-Terry-Luce (BTL) model (Bradley and Terry, 1952). Given two items i and j with scores θ_i and θ_j , the BTL model estimates the probability that item i is preferred to item j as $\mathbb{P}(i > j) = \frac{e^{\theta_i}}{e^{\theta_i} + e^{\theta_j}}$. We fit a BTL model to the Streetview preferences using maximum-likelihood estimation, and used this as the ground truth to generate noisy comparisons. Fig. 3(b) shows 4 images overlayed with their estimated BTL scores (where the lowest score was set to 0), and Fig. 3(c) shows a scatter plot of the scores of all 100 images.

LUCBRank can be used in the pairwise comparison setting using Borda reduction, as described in Section 1.1.

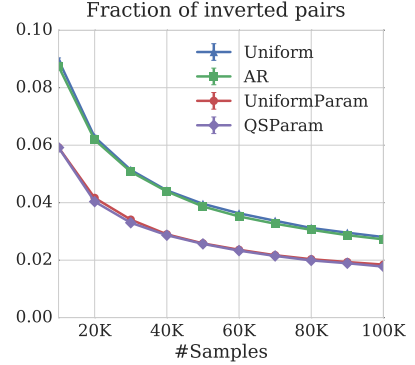


Figure 4: The futility of adaptive methods if the goal is to obtain a complete ranking. We compare uniform sampling with Active Ranking (both use non-parametric rank aggregation), and uniform sampling with quicksort (where both use parametric rank aggregation). We see that the Kendall tau distance of adaptive methods is no lower than those of their non-adaptive counterparts.

If complete ranking is the objective, adaptive methods offer no advantages when items means are close to each other as they are in this dataset. Oblivious of the generative model, a lower bound (ignoring constants and log factors) on the number of samples required by any algorithm to sort the items by their Borda scores is given by $\sum 1/\Delta_i^2$ (Jamieson et al., 2015b), where the Δ s are gaps between consecutive sorted Borda scores. For the dataset considered in this experiment, $\sum 1/\Delta_i^2 = 322$ million! We verify the futility of adaptive methods in Fig. 4, where we compare the performance of parametric as well as non-parametric adaptive methods in the literature (we describe these methods shortly) to their non-adaptive counterparts, with a goal of finding a complete ranking of the images. In the parametric algorithms (UniformParam and QSParm), we find MLE estimates of the BTL scores that best fit the pairwise responses. In the non-parametric algorithms (Uniform and AR), we estimate the scores using empirical probabilities in Eq. (2). In Fig. 4, we plot the fraction of pairs that are inverted in the empirical ranking compared to the true ranking, and see no benefits for adaptive methods. We do see gains from adaptivity in the clustered formulation (Fig. 5), as we explain shortly.

The adaptive methods in literature we compare to are AR (as in the previous section), and Quicksort (QS) (Ailon et al., 2008; Maystre and Grossglauser, 2017). The Quicksort algorithm works exactly like its non-noisy counterpart: it compares a randomly chosen pivot to all elements, and divides the elements into

two subsets - elements preferred to the pivot, and elements the pivot was preferred over. The algorithm then recurses into these two subsets. This algorithm is studied when comparisons are obtained from a BTL model, and the pairwise probabilities increase geometrically with distance between items (Maystre and Grossglauser, 2017). In this experiment, we stop the quicksort algorithm early as soon as all the subsets are inside the user-specified clusters. Continuing the algorithm further won't change the items in any cluster. This reduces the sample complexity of Quicksort.

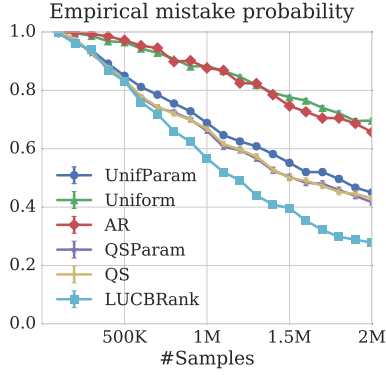


Figure 5: Probability of error in identifying the clusters: LUCBRank does better than parametric versions of other active algorithms.

We consider the problem of clustering the images into pentiles ($\kappa_i = 20i$, $1 \leq i \leq 5$). We set $\delta = 0.1$ for both LUCBRank and AR, and $\epsilon = 0$ in LUCBRank to ensure a fair comparison with AR. In Fig. 5, we record the probability (averaged over 600 simulations) that the empirical pentiles returned by the algorithm do not match the true pentiles. We find that LUCBRank has a lower mistake probability than even the parametric version of Quicksort, which assumes knowledge of the BTL model. As an aside, note that when the items are close as in this experiment, the parametric versions of Uniform and Quicksort perform similarly, and the active nature of Quicksort offers no significant advantage.

In Fig. 6(a) and (b) we plot the ratio of inter-cluster and intra-cluster inversions respectively of LUCBRank and Uniform. An inter-cluster pair is a pair of items that are in different clusters in the true ranking, while an intra-cluster pair is a pair of items from the same cluster. We see that the ratio of inter-cluster inversions goes down in Fig. 6(a), because that is the metric LUCBRank focuses on. LUCBRank does not expend effort on refining its estimate of an item's rank once its cluster has been found, and hence pays a price in the form of intra-cluster inversions (Fig. 6(b)).

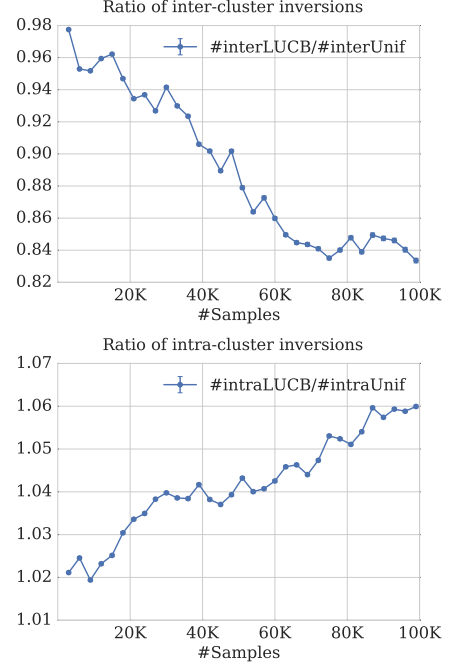


Figure 6: (a) The ratio of inter-cluster inversions of LUCBRank and Uniform. (b) The ratio of intra-cluster inversions of LUCBRank and Uniform. LUCBRank focuses on minimizing inter-cluster inversions at the cost of intra-cluster inversions.

8 Conclusion

Motivated by real-world problems in social science, we precisely formulate the clustered ranking problem. This setting is important because many real-world problems have high noise and are hard, and a complete ranking is not feasible; fortunately, it is often also not necessary. We propose a practical online algorithm for solving it, LUCBRank, and prove distribution-dependent upper and lower bounds on its sample complexity. We evaluate its performance on crowdsourced data gathered using MTurk, and observe that it performs better than existing algorithms in the literature.

We leave open several questions of interest. First, our upper bound is stated in terms of Chernoff information between the distributions, while our lower bound is stated in terms of KL-divergences, and there is a gap between the two. Second, note that the cluster boundaries need to be user-specified in our current setting. If the gap between the nearest items in adjacent clusters is small, this can adversely affect the sample complexity. Although this is partially addressed through the error-tolerance parameter ϵ , an attractive algorithm would be one which auto-tunes the positions of the cluster boundaries at the widest gaps, subject to user-specified constraints.

To the best of our knowledge, this paper presents the first bandit UCB algorithm for ranking.

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9 Appendix

9.1 PAC Guarantee

We'll use the following lemma (Kaufmann and Kalyanakrishnan, 2013) which bounds the probability of 'bad' events in round t .

Lemma 1. *Let $U_a(t)$ and $L_a(t)$ be the confidence bounds defined in Eq. (6) and Eq. (7). For any algorithm and arm a ,*

$$\mathbb{P}(U_a(t) < p_a) = \mathbb{P}(L_a(t) > p_a) \leq e(\beta(t, \delta) \log t + 1) \exp(-\beta(t, \delta))$$

We shall also need the following technical lemma, which we'll use to upper bound the probability of any bad event.

Lemma 2. *If $\beta(t, \delta) = \log(\frac{k_1 K t^\alpha}{\delta}) + \log \log(\frac{k_1 K t^\alpha}{\delta})$,*

$$\sum_{t=1}^{\infty} (\beta(t, \delta) \log t + 1) \exp(-\beta(t, \delta)) \leq \frac{\delta}{k_1 K} \left(\frac{2}{(\alpha - 1)^2} + \frac{1}{(\alpha - 1)} \right)$$

Proof. Let us consider

$$\begin{aligned} \beta(t, \delta) (\log t) e^{-\beta(t, \delta)} &= \left(\log \left(\frac{k_1 K t^\alpha}{\delta} \right) + \log \log \left(\frac{k_1 K t^\alpha}{\delta} \right) \right) (\log t) \left(\frac{\delta}{k_1 K t^\alpha} \cdot \frac{1}{\log \frac{k_1 K t^\alpha}{\delta}} \right) \\ &\leq 2 \log \left(\frac{k_1 K t^\alpha}{\delta} \right) \cdot \log t \cdot \left(\frac{\delta}{k_1 K t^\alpha} \cdot \frac{1}{\log \frac{k_1 K t^\alpha}{\delta}} \right) \\ &= 2 \log t \cdot \frac{\delta}{k_1 K t^\alpha} \end{aligned}$$

Hence

$$\begin{aligned} \sum_{t=1}^{\infty} (\beta(t, \delta) \log t + 1) \exp(-\beta(t, \delta)) &\leq \sum_{t=1}^{\infty} \left(2 \log t \cdot \frac{\delta}{k_1 K t^\alpha} + \frac{\delta}{k_1 K t^\alpha} \right) \\ &\leq \frac{\delta}{k_1 K} \left(\frac{2}{(\alpha - 1)^2} + \frac{1}{(\alpha - 1)} \right) \end{aligned}$$

□

9.1.1 Proof of Theorem 1

Theorem. *LUCBRank using $\beta(t, \delta) = \log(\frac{k_1 K t^\alpha}{\delta}) + \log \log(\frac{k_1 K t^\alpha}{\delta})$ with $\alpha > 1$ and $k_1 > 1 + \frac{2e}{\alpha - 1} + \frac{4e}{(\alpha - 1)^2}$, is correct with probability $1 - \delta$.*

Proof. Consider the event

$$W = \bigcap_{t \in \mathbb{N}} \bigcap_{a \in \{1, \dots, K\}} ((U_a(t) > p_a) \cap (L_a(t) < p_a))$$

where all arms are well-behaved i.e. their true means are inside their confidence intervals. We show that LUCBRank is correct on the event W .

Assume LUCBRank fails, which means that when it terminates, there exists a cluster i , such that arm a belongs to cluster i in the returned ranking, and $a \in M_{\epsilon, i}^{*, c}$; that is, either 1) $p_a > p_{\kappa_{i-1}+1} + \epsilon$ or 2) $p_a < p_{\kappa_i} - \epsilon$.

Consider the first case: $p_a > p_{\kappa_{i-1}+1} + \epsilon$. Consequently, there exists arm b such that $p_b \leq p_{\kappa_{i-1}+1}$, and $\tau(b) \leq \kappa_{i-1}$ in the returned ranking. Since the algorithm stopped and boundary $i-1$ was removed from the set of active boundaries C , it must be the case that $U_a(t) - L_b(t) < \epsilon$ upon stopping. Hence, the following holds:

$$\begin{aligned} & \bigcup_{t \in \mathbb{N}} (\exists a, b : p_a > p_{\kappa_{i-1}+1} + \epsilon, p_b \leq p_{\kappa_{i-1}+1}, U_a(t) - L_b(t) < \epsilon) \\ & \subseteq \bigcup_{t \in \mathbb{N}} (\exists a, b : (U_a(t) < p_b + \epsilon < p_a) \cup (L_b(t) > p_b)) \\ & \subseteq \bigcup_{t \in \mathbb{N}} \bigcup_{a \in \{1, \dots, K\}} (U_a(t) < p_a) \bigcup_{b \in \{1, \dots, K\}} (L_b(t) > p_b) \subseteq W^c \end{aligned}$$

Consider the second case: $p_a < p_{\kappa_i} - \epsilon$. Consequently, there exists an arm b such that $p_b \geq p_{\kappa_i}$, and $\tau(b) > \kappa_i$ in the returned ranking. Since the algorithm stopped and boundary i was removed from the set of active boundaries C , it must be the case that $U_b(t) - L_a(t) < \epsilon$ upon stopping. Hence, the following holds:

$$\begin{aligned} & \bigcup_{t \in \mathbb{N}} (\exists a, b : p_a < p_{\kappa_i} - \epsilon, p_b \geq p_{\kappa_i}, U_b(t) - L_a(t) < \epsilon) \\ & \subseteq \bigcup_{t \in \mathbb{N}} (\exists a, b : (U_b(t) < p_b) \cup (L_a(t) > p_b - \epsilon > p_a)) \\ & \subseteq \bigcup_{t \in \mathbb{N}} \bigcup_{b \in \{1, \dots, K\}} (U_b(t) < p_b) \bigcup_{a \in \{1, \dots, K\}} (L_a(t) > p_a) \subseteq W^c \end{aligned}$$

Hence

$$\begin{aligned} \mathbb{P}(\text{LUCBRank fails}) & \leq \mathbb{P}(W^c) \\ & \leq 2eK \sum_{t=1}^{\infty} (\beta(t, \delta) \log t + 1) \exp(-\beta(t, \delta)) && \text{(by Lemma 1)} \\ & \leq \frac{\delta}{k_1} \left(\frac{4e}{(\alpha-1)^2} + \frac{2e}{(\alpha-1)} \right) && \text{(by Lemma 2)} \\ & \leq \delta && \text{(by the constraint on } k_1) \end{aligned}$$

□

9.2 Sample Complexity

We define the event W_t which says that all arms are well-behaved in round t i.e. their true means are contained inside their confidence intervals.

$$W_t = \bigcap_{a \in \{1, 2, \dots, K\}} ((U_a(t) > p_a) \cap (L_a(t) < p_a))$$

Note that the event W defined earlier is $W = \bigcup_{t \in \mathbb{N}} W_t$.

Proposition 1 gives a sufficient condition for stopping.

Proposition 1. *Let $b_i \in [p_{\kappa_i}, p_{\kappa_i+1}]$. If $U_{u_t^i} - L_{l_t^i} > \epsilon$ and W_t holds, then either $k = l_t^i$ or $k = u_t^i$ satisfies*

$$b_i \in \mathcal{I}_k(t) \text{ and } \tilde{\beta}_k(t) > \frac{\epsilon}{2},$$

where we define $\tilde{\beta}_a(t) = \sqrt{\frac{\beta(t, \delta)}{2N_a(t)}}$

Proof. Our W_t condition is stronger than that required in the Proposition 1 in Kaufmann and Kalyanakrishnan (2013), and hence their proof applies. □

Lemma 3 is another concentration result that will be used in our sample complexity guarantee.

Lemma 3. *Let $T \geq 1$ be an integer, and $1 \leq i \leq (c-1)$ be any cluster boundary. Let $\delta > 0, \gamma > 0$ and $x \in]0, 1[$ be such that $p_a \neq x$. Then*

$$\sum_{t=1}^T \mathbb{P} \left(a = u_i^t \vee a = l_i^t, N_a(t) > \left\lceil \frac{\gamma}{d^*(p_a, x)} \right\rceil, N_a(t) d(\hat{p}_a(t), x) \leq \gamma \right) \leq \frac{\exp(-\gamma)}{d^*(p_a, x)}$$

We prove the following lemma, which states that the Chernoff information increases as the second distribution moves away from the first.

Lemma 4. *If $x < y < y'$ or $x > y > y'$, $d^*(x, y) \leq d^*(x, y')$*

Proof. We shall prove the statement for the case $x < y < y'$. The proof for $x > y > y'$ is analogous. Let z^* be the unique z such that $d(z^*, x) = d(z^*, y)$. Since $z^* < y < y'$, $d(z^*, y') \geq d(z^*, y) = d(z^*, x)$. Hence, there exists $z^{*'} \geq z^*$ such that $d^*(x, y) = d(z^*, x) \leq d(z^{*'}, x) = d(z^{*'}, y') = d^*(x, y')$. \square

Lemma 5. *Let x^* be the solution of the equation:*

$$x = \frac{1}{\gamma} \left(\log \frac{x^\alpha}{\eta} + \log \log \frac{x^\alpha}{\eta} \right)$$

Then if $\gamma < 1$ and $\eta < 1/e^e$,

$$\frac{1}{\gamma} \log \left(\frac{1}{\eta \gamma^\alpha} \right) \leq x^* \leq \frac{C_0}{\gamma} \log \left(\frac{1}{\eta \gamma^\alpha} \right)$$

where C_0 is such that $C_0 \geq (1 + \frac{1}{e}) (\alpha \log C_0 + 1 + \frac{\alpha}{e})$.

Proof. x^* is upper bounded by any x such that $\frac{1}{\gamma} \left(\log \frac{x^\alpha}{\eta} + \log \log \frac{x^\alpha}{\eta} \right) \leq x$. We look for x^* of the form $\frac{C_0}{\gamma} \log \left(\frac{1}{\eta \gamma^\alpha} \right)$.

$$\begin{aligned} \frac{1}{\gamma} \left(\log \frac{x^\alpha}{\eta} + \log \log \frac{x^\alpha}{\eta} \right) &\leq \frac{1}{\gamma} \left(1 + \frac{1}{e} \right) \log \left(\frac{x^\alpha}{\eta} \right) \\ &= \frac{1}{\gamma} \left(1 + \frac{1}{e} \right) \left(\alpha \log C_0 + \log \frac{1}{\eta \gamma^\alpha} + \alpha \log \log \frac{1}{\eta \gamma^\alpha} \right) \\ &\leq \frac{1}{\gamma} \left(1 + \frac{1}{e} \right) \left(\alpha \log C_0 + \left(1 + \frac{\alpha}{e} \right) \log \frac{1}{\eta \gamma^\alpha} \right) \\ &\leq \frac{1}{\gamma} \left(1 + \frac{1}{e} \right) \left(\alpha \log C_0 + 1 + \frac{\alpha}{e} \right) \log \frac{1}{\eta \gamma^\alpha} \end{aligned}$$

where the first and second inequalities hold because $\log x \leq \frac{x}{e}$, and the last inequality holds because $\frac{1}{\eta \gamma^\alpha} > e$. Choosing C_0 such that

$$C_0 \geq \left(1 + \frac{1}{e} \right) \left(\alpha \log C_0 + 1 + \frac{\alpha}{e} \right)$$

gives us our upper bound.

To prove the lower bound, consider the series defined by

$$\begin{aligned} x_0 &= 1 \\ x_{n+1} &= \frac{1}{\gamma} \left(\log \frac{x_n^\alpha}{\eta} + \log \log \frac{x_n^\alpha}{\eta} \right) \end{aligned}$$

First note that since $\gamma < 1$ and $\eta < 1/e^e$, the sequence is increasing. Second, note that the sequence converges to x^* . Hence

$$\begin{aligned} x^* &\geq x_2 = \frac{1}{\gamma} \left[\log \left(\frac{1}{\eta\gamma^\alpha} \left(\log \frac{1}{\eta} + \log \log \frac{1}{\eta} \right)^\alpha \right) + \log \log \left(\frac{1}{\eta\gamma^\alpha} \left(\log \frac{1}{\eta} + \log \log \frac{1}{\eta} \right)^\alpha \right) \right] \\ &= \frac{1}{\gamma} \left[\log \frac{1}{\eta\gamma^\alpha} + \alpha \log \left(\log \frac{1}{\eta} + \log \log \frac{1}{\eta} \right) + \log \log \frac{1}{\eta\gamma^\alpha} + \alpha \log \log \left(\log \frac{1}{\eta} + \log \log \frac{1}{\eta} \right) \right] \\ &\geq \frac{1}{\gamma} \log \frac{1}{\eta\gamma^\alpha} \end{aligned}$$

since $\eta < 1/e^e$. □

Corollary 1. Let $\gamma = \frac{1}{2H_{\epsilon,b}^*}$, $\eta = \frac{\delta}{k_1 K}$. Then applying Lemma 5 gives

$$2H_{\epsilon,b}^* \log \left(\frac{k_1 K (2H_{\epsilon,b}^*)^\alpha}{\delta} \right) \leq S_1^* \leq 2C_0(\alpha) H_{\epsilon,b}^* \log \left(\frac{k_1 K (2H_{\epsilon,b}^*)^\alpha}{\delta} \right)$$

9.3 Proof of Theorem 2

Theorem. Let $b = (b_1, b_2, \dots, b_{c-1})$, where $b_i \in [p_{\kappa_i}, p_{\kappa_i+1}]$. Let $\epsilon > 0$. Let $\beta(t, \delta) = \log(\frac{k_1 K t^\alpha}{\delta}) + \log \log(\frac{k_1 K t^\alpha}{\delta})$ with $k_1 > 1 + \frac{2e}{\alpha-1} + \frac{4e}{(\alpha-1)^2}$. Let τ be the random number of samples taken by LUCBRank before termination. If $\alpha > 1$,

$$\mathbb{P} \left(\tau \leq 2C_0(\alpha) H_{\epsilon,b}^* \log \left(\frac{k_1 K (2H_{\epsilon,b}^*)^\alpha}{\delta} \right) \right) \geq 1 - \delta$$

where $C_0(\alpha)$ is such that $C_0(\alpha) \geq (1 + \frac{1}{\epsilon}) (\alpha \log(C_0(\alpha)) + 1 + \frac{\alpha}{\epsilon})$.

Proof. The LUCBRank algorithm proceeds in rounds. In a round, it samples the two arms on opposite sides of an active boundary whose confidence intervals overlap the most. A boundary is active as long as this overlap is less than ϵ . Thus, the number of samples up to round T is

$$\begin{aligned} \# \text{samples}(T) &\leq 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(U_{u_t^i} - L_{l_t^i} > \epsilon)} \\ &= 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(U_{u_t^i} - L_{l_t^i} > \epsilon)} (\mathbb{1}_{W_t} + \mathbb{1}_{W_t^c}) \\ &\leq 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(U_{u_t^i} - L_{l_t^i} > \epsilon)} \mathbb{1}_{W_t} + 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{W_t^c} \\ &\leq 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \sum_{a \in \{1, 2, \dots, K\}} \mathbb{1}_{(a=l_t^i) \vee (a=u_t^i)} \mathbb{1}_{(b_i \in \mathcal{I}_a(t))} \mathbb{1}_{(\tilde{\beta}_a(t) > \frac{\epsilon}{2})} + 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{W_t^c} \\ &\quad \text{(by Proposition 1)} \end{aligned}$$

We now split the first sum into two depending on whether an arm a belongs to the set $\mathcal{A}_\epsilon = \{a \in \{1, 2, \dots, K\} :$

$$\Delta_b^* < \epsilon^2/2\}.$$

$$\begin{aligned} \#\text{samples}(T) &\leq 2 \sum_{a \in \mathcal{A}_\epsilon} \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(a=l_t^i) \vee (a=u_t^i)} \mathbb{1}_{\left(N_a(t) < \frac{\beta(t, \delta)}{\epsilon^2/2}\right)} + \\ &\quad 2 \sum_{a \in \mathcal{A}_\epsilon^c} \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(a=l_t^i) \vee (a=u_t^i)} \mathbb{1}_{(b_i \in \mathcal{I}_a(t))} + 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{W_t^c} \\ &\leq 2 \sum_{a \in \mathcal{A}_\epsilon} \frac{\beta(T, \delta)}{\epsilon^2/2} + 2 \sum_{a \in \mathcal{A}_\epsilon^c} \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(a=l_t^i) \vee (a=u_t^i)} \mathbb{1}_{\left(N_a(t) \leq \left\lceil \frac{\beta(T, \delta)}{\Delta_b^*(a)} \right\rceil\right)} + \\ &\quad \underbrace{2 \sum_{a \in \mathcal{A}_\epsilon^c} \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(a=l_t^i) \vee (a=u_t^i)} \mathbb{1}_{\left(N_a(t) > \left\lceil \frac{\beta(T, \delta)}{\Delta_b^*(a)} \right\rceil\right)} + 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{W_t^c}}_{R_T} \\ &= 2H_{\epsilon, b}^* \beta(T, \delta) + R_T \end{aligned}$$

where

$$R_T = 2 \sum_{a \in \mathcal{A}_\epsilon^c} \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{(a=l_t^i) \vee (a=u_t^i)} \mathbb{1}_{\left(N_a(t) > \left\lceil \frac{\beta(T, \delta)}{\Delta_b^*(a)} \right\rceil\right)} \mathbb{1}_{(b_i \in \mathcal{I}_a(t))} + 2 \sum_{t=1}^T \sum_{i=1}^{c-1} \mathbb{1}_{W_t^c}$$

If we define $S_1^* = \min\{x : 2H_{\epsilon, b}^* \beta(x, \delta) < x\}$, then we get that for $S > S_1^*$, the algorithm must have stopped before S samples on the event $(R_T = 0)$. Denoting the total number of samples used by the algorithm by τ , we have that, for any $S > S_1^*$, $\mathbb{P}(\tau > S) \leq \mathbb{P}(R_T \neq 0)$.

$$\begin{aligned} \mathbb{P}(\tau > S) &\leq \mathbb{P}(R_T \neq 0) \\ &\leq \mathbb{P}\left(\exists a \in \mathcal{A}_\epsilon^c, t \leq T, 1 \leq i \leq (c-1) : a = l_t^i \vee a = r_t^i, N_a(t) > \left\lceil \frac{\beta(T, \delta)}{\Delta_b^*(a)} \right\rceil, b_i \in \mathcal{I}_a(t)\right) + \mathbb{P}(W^c) \\ &\leq \mathbb{P}\left(\exists a \in \mathcal{A}_\epsilon^c, t \leq T, 1 \leq i \leq (c-1) : a = l_t^i \vee a = r_t^i, N_a(t) > \left\lceil \frac{\beta(T, \delta)}{d^*(p_a, b_i)} \right\rceil, b_i \in \mathcal{I}_a(t)\right) + \mathbb{P}(W^c) \end{aligned} \quad (14)$$

where the final inequality follows because $\Delta_b^*(a) \leq d^*(p_a, b_i) \forall 1 \leq i \leq c-1$ (by Lemma 4).

Let us look at the first term:

$$\begin{aligned} &\mathbb{P}\left(\exists a \in \mathcal{A}_\epsilon^c, t \leq T, 1 \leq i \leq (c-1) : a = l_t^i \vee a = r_t^i, N_a(t) > \left\lceil \frac{\beta(T, \delta)}{d^*(p_a, b_i)} \right\rceil, b_i \in \mathcal{I}_a(t)\right) \\ &\leq \sum_{a \in \mathcal{A}_\epsilon^c} \sum_{i=1}^{c-1} \sum_{t=1}^T \mathbb{P}\left(a = l_t^i \vee a = r_t^i, N_a(t) > \left\lceil \frac{\beta(T, \delta)}{d^*(p_a, b_i)} \right\rceil, b_i \in \mathcal{I}_a(t)\right) \\ &\leq \sum_{a \in \mathcal{A}_\epsilon^c} \sum_{i=1}^{c-1} \frac{\exp(-\beta(T, \delta))}{d^*(p_a, b_i)} \quad (\text{by Lemma 3}) \\ &\leq (c-1) \exp(-\beta(T, \delta)) \sum_{a \in \mathcal{A}_\epsilon^c} \frac{1}{d^*(p_a, b_i)} \\ &\leq (c-1) H_{\epsilon, b}^* \exp(-\beta(T, \delta)) \\ &\leq (c-1) H_{\epsilon, b}^* \exp\left(-\beta\left(\frac{S_1^*}{c-1}, \delta\right)\right) \quad (\text{if } \tau > S_1^*, T > \frac{S_1^*}{c-1}) \\ &= (c-1) H_{\epsilon, b}^* \cdot \frac{\delta(c-1)^\alpha}{k_1 K S_1^{*, \alpha}} \frac{1}{\log \frac{k_1 K S_1^{*, \alpha}}{\delta(c-1)^\alpha}} \\ &\leq (c-1)^{\alpha+1} H_{\epsilon, b}^* \cdot \frac{\delta}{k_1 K \left(2H_{\epsilon, b}^* \log\left(\frac{k_1 K (2H_{\epsilon, b}^*)^\alpha}{\delta}\right)\right)^\alpha} \frac{1}{\log \frac{k_1 K S_1^{*, \alpha}}{\delta(c-1)^\alpha}} \quad (\text{by the lower bound in Corollary 1}) \\ &\leq \frac{\delta}{k_1} \cdot \left(\frac{c-1}{2}\right)^\alpha \quad (\text{since } (c-1) \leq K \text{ and } \alpha > 1) \end{aligned}$$

For the second term, note that

$$\begin{aligned}\mathbb{P}(W^c) &\leq 2eK \sum_{t=1}^{\infty} (\beta(t, \delta) \log t + 1) \exp(-\beta(t, \delta)) \quad (\text{by Lemma 1}) \\ &\leq \frac{\delta}{k_1} \left(\frac{4e}{(\alpha-1)^2} + \frac{2e}{(\alpha-1)} \right) \quad (\text{by Lemma 2})\end{aligned}$$

Substituting in Eq. (14), we get that for $S > S_1^*$,

$$\mathbb{P}(\tau > S) \leq \frac{\delta}{k_1} \left(\left(\frac{c-1}{2} \right)^\alpha + \frac{4e}{(\alpha-1)^2} + \frac{2e}{(\alpha-1)} \right) \leq \delta$$

by the choice of k_1 . \square

9.4 Lower Bound

The proof uses standard change of measure arguments used to prove lower bounds. For bandit problems, this is succinctly expressed through Lemma 1 in Kaufmann et al. (2015) that we restate here for completeness.

Lemma 6. *Let p and p' be two bandit models with K arms such that for all a , the distributions p_a and p'_a are mutually absolutely continuous. Let σ be a stopping time with respect to (\mathcal{F}_t) and let $A \in \mathcal{F}_\sigma$. Then*

$$\sum_{a=1}^K \mathbb{E}_p[N_a(\sigma)] \text{KL}(p_a, p'_a) \geq d(\mathbb{P}_p(A), \mathbb{P}_{p'}(A))$$

where $d(x, y) = x \log(x/y) + (1-x) \log((1-x)/(1-y))$.

9.4.1 Proof of Theorem 3

Consider any arm a . By Assumption 1, there exists alternative model p' such that:

$$\text{KL}(p_a, p_{\kappa_{g(a)}+1}) < \text{KL}(p_a, p'_a) < \text{KL}(p_a, p_{\kappa_{g(a)}+1}) + \alpha \text{ and } p'_a < p_{\kappa_{g(a)}+1}$$

Note that in the model p' , arm a no longer belongs to the cluster $g(a)$. Let $\hat{M}_{g(a)}$ be the set of arms returned by an algorithm in the $g(a)^{\text{th}}$ cluster. If we define the event $A = \{a \in \hat{M}_{g(a)}\} \in \mathcal{F}_\tau$, then by definition, for any δ -PAC algorithm, $\mathbb{P}_p(A) \geq 1 - \delta$ and $\mathbb{P}_{p'}(A) \leq \delta$. Letting $N_a(\tau)$ denote the number of pulls of arm a by time τ , we have by Lemma 6 and the monotonicity of $d(x, y)$ that

$$\text{KL}(p_a, p'_a) \mathbb{E}_p[N_a(\tau)] \geq d(1 - \delta, \delta) \geq \log\left(\frac{1}{2.4\delta}\right)$$

where we use the property that for $x \in [0, 1]$, $d(x, 1-x) \geq \log \frac{1}{2.4x}$. This gives us that

$$\mathbb{E}_p[N_a(\tau)] \geq \frac{1}{\text{KL}(p_a, p_{\kappa_{g(a)}+1}) + \alpha} \log\left(\frac{1}{2.4\delta}\right)$$

Letting $\alpha \rightarrow 0$, we get

$$\mathbb{E}_p[N_a(\tau)] \geq \frac{1}{\text{KL}(p_a, p_{\kappa_{g(a)}+1})} \log\left(\frac{1}{2.4\delta}\right) \quad (15)$$

Similarly, by considering an alternative model p'' such that

$$\text{KL}(p_a, p_{\kappa_{g(a)}-1}) < \text{KL}(p_a, p''_a) < \text{KL}(p_a, p_{\kappa_{g(a)}-1}) + \alpha \text{ and } p''_a > p_{\kappa_{g(a)}-1}$$

we get

$$\mathbb{E}_p[N_a(\tau)] \geq \frac{1}{\text{KL}(p_a, p_{\kappa_{g(a)}-1})} \log\left(\frac{1}{2.4\delta}\right) \quad (16)$$

From Eq. (15), Eq. (16), and the definition of $\Delta_\kappa^{\text{KL}}(a)$ in Eq. (12), we get that

$$\mathbb{E}_p[N_a(\tau)] \geq \frac{1}{\Delta_\kappa^{\text{KL}}(a)} \log\left(\frac{1}{2.4\delta}\right)$$

Summing over all the arms yields the required bound for $\mathbb{E}_p[\tau] = \sum_{a=1}^K \mathbb{E}_p[N_a(\tau)]$.