## THE COMPUTATIONAL HARDNESS OF ESTIMATING EDIT DISTANCE\*

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Abstract. We prove the first nontrivial communication complexity lower bound for the problem of estimating the edit distance (aka Levenshtein distance) between two strings. To the best of our knowledge, this is the first computational setting in which the complexity of estimating the edit distance is provably larger than that of Hamming distance. Our lower bound exhibits a trade-off between approximation and communication, asserting, for example, that protocols with O(1) bits of communication can obtain only approximation  $\alpha \geq \Omega(\log d/\log \log d)$ , where d is the length of the input strings. This case of O(1) communication is of particular importance since it captures constantsize sketches as well as embeddings into spaces like  $l_1$  and squared- $l_2$ , two prevailing algorithmic approaches for dealing with edit distance. Indeed, the known nontrivial communication upper bounds are all derived from embeddings into  $l_1$ . By excluding low-communication protocols for edit distance, we rule out a strictly richer class of algorithms than previous results. Furthermore, our lower bound holds not only for strings over a binary alphabet but also for strings that are permutations (aka the Ulam metric). For this case, our bound nearly matches an upper bound known via embedding the Ulam metric into  $l_1$ . Our proof uses a new technique that relies on Fourier analysis in a rather elementary way.

 ${\bf Key}$  words. embedding, communication complexity, edit distance, Ulam metric, Fourier analysis, sketching

AMS subject classifications. 65W25, 68W32, 68Q17

DOI. 10.1137/080716530

1. Introduction. The *edit distance* (aka *Levenshtein distance*) between two strings is the number of insertions, deletions, and substitutions needed to transform one string into the other. This distance is of key importance in several fields, such as computational biology and text processing, and consequently computational problems involving the edit distance were studied quite extensively. The most basic problem is that of computing the edit distance between two strings of length d over alphabet  $\Sigma$ . The fastest algorithm known for the case of constant-size alphabet remains the algorithm of Masek and Paterson [25] from 1980, which runs in time  $O(d^2/\log^2 d)$ . Unfortunately, such near-quadratic time is prohibitive when working on large datasets, which is common in areas such as computational biology. A possible approach is to trade accuracy for speed and employ faster algorithms that compute the edit distance approximately (possibly as a preliminary filtering step). Currently, the best nearlinear time algorithm, due to Andoni and Onak [5], achieves an approximation factor of  $2^{\tilde{O}(\sqrt{\log d})}$ , improving the earlier results of [8, 6, 9].

Another major algorithmic challenge is to design a scheme for nearest neighbor search (NNS) under the edit distance. In this problem, we wish to design a data

<sup>\*</sup>Received by the editors February 25, 2008; accepted for publication (in revised form) October 7, 2009; published electronically April 30, 2010. A preliminary version of this paper appeared in *Proceedings of the 48th Annual IEEE Symposium on Foundations of Computer Science*, 2007.

http://www.siam.org/journals/sicomp/39-6/71653.html

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structure that preprocesses a dataset of n strings of length d each, so that when a query string is given, the query's nearest neighbor (i.e., a dataset string with the smallest edit distance to the query string) can be reported quickly. However, no efficient solutions for this problem are known, even if one allows a small approximation. All known algorithms with fast query time (polynomial in d and  $\log n$ ) either require large space or have large approximation error—Indyk [19] achieves constant approximation using  $n^{d^{\Omega(1)}}$  space, and Ostrovsky and Rabani [29] obtain  $2^{O(\sqrt{\log d \log \log d})}$  approximation using space that is polynomial in d and n.

It is thus natural to ask the following: Is it really "hard" to design algorithms for the edit distance? A natural benchmark is the Hamming distance, which is equal to the number of positions where the two strings differ. Hamming distance can be seen as edit distance where the only operation allowed is substitution. For Hamming distance, much better algorithms are known: (i) the distance between two strings can clearly be computed in O(d) time, and (ii) NNS schemes by Indyk and Motwani [20] and by Kushilevitz, Ostrovsky, and Rabani [24] achieve  $1 + \epsilon$  approximation using space that is polynomial in d and in  $n^{1/\epsilon^2}$ . Empirically, edit distance appears to be more difficult than Hamming distance, and the reason is quite clear—insertions and deletions cause portions of the string to move and create an alignment problem but there is no rigorous evidence that supports this intuition. In particular, we are not aware of a computational model in which the complexity of approximating edit distance.<sup>1</sup>

We give the first rigorous evidence for the *computational* hardness of approximating the edit distance. In fact, we show a computational model in which the complexity of estimating edit distance is significantly larger than that of Hamming distance, and this is the first setting where such a separation is known. Our results hold for two important metrics:

1. standard edit metric, i.e., edit distance on  $\{0, 1\}^d$ ;

2. the Ulam metric, which is the edit distance on permutations of length d. Here and throughout, a permutation is a string consisting of distinct characters coming from a large alphabet,  $|\Sigma| \geq d$ . This definition of permutations is nonstandard, although our results also hold under the standard one, where  $|\Sigma| = d$  (see Fact 2.4 and the discussion preceding it). Our results immediately imply lower bounds for sketching algorithms and for metric embeddings. These two algorithmic techniques received a lot of attention lately as promising approaches to many metric problems. We will discuss these implications in more detail after stating our main results.

1.1. Main results. Our main result is stated in terms of the communication complexity of the distance threshold estimation problem (DTEP) and holds for both

<sup>&</sup>lt;sup>1</sup>We are aware of only two results that come close. First, if the operations on the symbols of the strings are restricted to tests of equality, then computing edit distance between two strings over a large alphabet requires  $\Omega(d^2)$  comparisons [34]. However, this lower bound holds only for exact computation (or 1 + o(1) approximation) and for strings over a large alphabet (but not for binary strings). In fact, the lower bound breaks down even in the comparison model (when we can compare the relative order of two symbols); e.g., the algorithm of [10] runs in time  $O(d^2 \frac{\log^2 \log d}{\log^2 d})$  for computing edit distance between strings over an arbitrarily large alphabet.

Second, if we restrict our attention to *sublinear* time, i.e., algorithms that probe only a small part of the two input strings, then there exists a simple separation in terms of query complexity. Specifically, deciding whether the edit distance is  $\Omega(d)$  or  $O(d^{1-\epsilon})$  requires reading at least  $\Omega(d^{1/2-\epsilon/2})$ positions of the strings [8], while the same decision under Hamming distance is achieved easily by sampling O(1) positions. This separation has limited computational implications since it essentially shows that estimating edit distance requires reading "many" positions of the input strings.

the edit metric over  $\Sigma = \{0, 1\}$  and for the Ulam metric. In DTEP [32], for a threshold R and an approximation  $\alpha \ge 1$  fixed as parameters, we are given inputs x, y and want to decide whether ed(x, y) > R or  $ed(x, y) \le R/\alpha$ .

In the communication protocol setting, Alice and Bob, who have access to a common source of randomness, receive strings x and y, respectively, as their inputs, and their goal is to solve DTEP by exchanging messages. The communication complexity of the protocol is then defined as the minimum number of bits Alice and Bob need to exchange in order to succeed with probability at least 2/3. When x, y come from the standard edit metric, we denote the communication complexity by  $\operatorname{CC}_{\alpha,R}^{\{0,1\}^d}$ . Similarly, when x, y come from the Ulam metric, we denote the communication complexity by  $\operatorname{CC}_{\alpha,R}^{Ulam_d}$ . Our main theorem provides a lower bound on the latter, exhibiting a trade-off between communication and approximation.

THEOREM 1.1 (main theorem). There exists a constant c > 0 such that, for every string length d > 1, approximation  $\alpha > 1$ , and R satisfying  $d^{0.1} \leq R \leq d^{0.49}$ ,

$$c \cdot \operatorname{CC}_{\alpha,R}^{\operatorname{Ulam}_d} + \log(\alpha \log \alpha) \ge \log \log d.$$

We extend this result from the Ulam metric to the standard edit metric by reducing the latter to the former. The key idea, which may be of independent interest, is that substituting every alphabet symbol independently with a random bit is likely to preserve the edit distance, up to a constant factor, as stated in the following theorem.

THEOREM 1.2. Let  $P, Q \in \Sigma^d$  be two permutations, and let  $\pi : \Sigma \mapsto \{0, 1\}$  be a random function. Then

•  $ed(\pi(P), \pi(Q)) \le ed(P, Q)$  for any choice of  $\pi$ , and

•  $\Pr_{\pi}\left[\operatorname{ed}(\pi(P), \pi(Q)) \ge \Omega(1) \cdot \operatorname{ed}(P, Q)\right] \ge 1 - 2^{-\Omega(\operatorname{ed}(P, Q))}.$ 

Using our two theorems, we obtain the following.

COROLLARY 1.3. There exists a constant c > 0 such that, for every string length d > 1, approximation  $\alpha > 1$ , and R satisfying  $d^{0.1} \leq R \leq d^{0.49}$ ,

$$c \cdot \operatorname{CC}_{\alpha,R}^{\{0,1\}^d} + \log(\alpha \log \alpha) \ge \log \log d.$$

The previously known lower bounds for  $\operatorname{CC}_{\alpha,R}^{\{0,1\}^d}$  and  $\operatorname{CC}_{\alpha,R}^{\operatorname{Ulam}_d}$  are all obtained by a straightforward reduction from the same problem on the Hamming metric. These bounds assert that the communication complexity for  $\alpha = 1 + \epsilon$  is  $\Omega(1/\epsilon)$ , and in the case of sketching (aka simultaneous)<sup>2</sup> protocols is  $\Omega(1/\epsilon^2)$  [35] (see also [36, Chapter 4]), and both are clearly uninformative for (say)  $\alpha \geq 2$ . See also [30] for other related results.

The only nontrivial upper bounds currently known are (i)  $\operatorname{CC}_{\alpha,R}^{\{0,1\}^d} \leq O(1)$  for suitable  $\alpha = 2^{O(\sqrt{\log d \log \log d})}$ ; and (ii)  $\operatorname{CC}_{\alpha,R}^{\operatorname{Ulam}_d} \leq O(1)$  for suitable  $\alpha = O(\log d)$ ; and they both follow via embedding into  $\ell_1$ . See section 1.2 and Table 1.1 for more details.

Comparison with Hamming distance. The next proposition, proved (implicitly) by Kushilevitz, Ostrovsky, and Rabani [24], upper bounds the communication complexity of DTEP over the Hamming metric. Let H(x, y) be the Hamming distance between x and y.

PROPOSITION 1.4 (see [24]). Let d > 1, R > 1, and  $\epsilon > 0$ . Then there exists a communication protocol (in fact, a sketching algorithm) that given inputs  $x, y \in \Sigma^d$ 

 $<sup>^{2}</sup>$ See the formal definition of sketching in section 1.2.

distinguishes whether H(x,y) > R or  $H(x,y) \leq R/(1+\epsilon)$ , using  $O(1/\epsilon^2)$  bits of communication.

Observe that for approximation factor  $\alpha$ , which is a constant (namely, independent of d and R), the complexity of the Hamming metric is O(1), while that of the edit metric is  $\Omega(\log \log d)$ . It thus follows that edit distance is indeed provably harder to compute than Hamming distance in the context of communication protocols.

**1.2. Implications and related work.** Two promising approaches to designing algorithms for the edit metrics are via metric embeddings and via sketching, and our results preclude good approximation algorithms obtained via either of these approaches.

Embedding of edit distance into normed metrics. A current line of attack on edit distance is by embedding it into a computationally easier metric, for which efficient algorithms are known. An embedding is a mapping f from the strings into, say, an  $\ell_1$  metric such that, for all strings x, y,

$$\operatorname{ed}(x,y) \le \|f(x) - f(y)\|_1 \le D \cdot \operatorname{ed}(x,y),$$

and  $D \geq 1$  is called the embedding's *distortion* (approximation factor). An embedding with low distortion would have major consequences since it allows porting a host of existing algorithms for an  $\ell_1$  metric to the case of edit distance. For example, an (efficiently computable) embedding with distortion D gives an efficient nearest neighbor data structure for approximation (say) 2D by applying the embedding and reverting to [20, 24].

Naturally, researchers were keen to find the least distortion for an embedding into  $\ell_1$ —the problem is cited in Matoušek's list of open problems [26], as well as in Indyk's survey [18]. Table 1.1 summarizes the previously known upper and lower bounds, as well as the implications of our theorems. The reader may find more background on some variants of the edit distance in [31].

TABLE 1.1

Known bounds on distortion/approximation of embedding variants of edit distance into  $\ell_1$ , squared- $\ell_2$ , and the approximation for achieving O(1)-size sketch. Since  $\ell_1$  embeds isometrically into squared- $\ell_2$  and the latter admits O(1)-size sketch for 2-approximation, the upper bounds transfer from left to right, and the lower bounds transfer from right to left (as suggested by the arrows). n/a means no result is given (even implicitly).

Metric	Reference	Embedding into $\ell_1$	Embedding into squared-la	O(1)-size sketch
Edit distance on $\{0,1\}^d$	[29]	$2\tilde{O}(\sqrt{\log d})$	$\rightarrow$	$\rightarrow$
	[22],[23]	$\Omega(\log d)$	n/a	n/a
	[1]	$\leftarrow$	$\geq 3/2$	n/a
	This paper	$\leftarrow$	$\leftarrow$	$\Omega(\frac{\log d}{\log\log d})$
Ulam metric (edit distance on permutations)	[12]	$O(\log d)$	$\rightarrow$	$\longrightarrow$
	[13]	$\leftarrow$	$\geq 4/3$	n/a
	This paper	$\leftarrow$	$\leftarrow$	$\Omega(\frac{\log d}{\log \log d})$
Edit distances with block operations	[15], [28], [13], [14]	$\tilde{O}(\log d)$	$\rightarrow$	$\longrightarrow$

It is readily seen from the table that the only previous superconstant distortion lower bound is  $\Omega(\log d)$  for embedding edit distance into  $\ell_1$ , due to Krauthgamer and Rabani [23], building on a technique of Khot and Naor [22], who gave a bound of  $\Omega((\log d)^{1/2-o(1)})$ . Although this lower bound is important, one can potentially overcome such a lower bound by, say, embedding edit distance into a richer space, such as squared- $\ell_2$ , a real space with squared Euclidean distance, with a possibly smaller distortion—the major implications of an embedding into squared- $\ell_2$  are precisely the same as those of an embedding into  $\ell_1$ . On this front, much weaker lower bounds were known: the previous lower bound is only 3/2 [1]. To further stress how little was known, we note that one can consider even richer metrics, such as any fixed power of  $\ell_2$  (essentially equivalent to embedding a fixed root of edit distance into  $\ell_2$ ), which also has an efficient nearest neighbor data structure. For a sufficiently high (but fixed) power of  $\ell_2$ , even the 3/2 bound of [1] gets weaker and becomes arbitrarily close to 1.

Our results rule out all such embeddings indirectly by targeting a richer class of metrics—metrics for which the respective DTEP problem admits a protocol with O(1) bits of communication and O(1) approximation. (Proposition 1.4 shows that this class of metrics is indeed richer.) It follows from our communication lower bounds that every embedding of edit distance (either on 0-1 strings or on permutations) into a metric in that richer class must incur distortion  $D \ge \Omega(\frac{\log d}{\log \log d})$ , without requiring that the embedding be efficiently computable. For completeness, we state and prove this distortion lower bound explicitly for metrics which are a fixed power of  $\ell_2$ .

COROLLARY 1.5. For every fixed  $p \ge 1$ , embedding the standard edit metric or the Ulam metric into  $(\ell_2)^p$ , the pth power of  $\ell_2$ , requires distortion  $\Omega(\frac{\log d}{\log \log d})$ . The same is true also for embedding into  $\ell_1$ .

*Proof.* Suppose  $p \ge 1$  is fixed and the edit metric ed (or similarly the Ulam metric) embeds into  $(\ell_2)^p$  with distortion  $D \ge 1$ . In other words, the metric ed<sup>1/p</sup> (i.e., 1/p-power of every distance) embeds into  $\ell_2$  with distortion  $D^{1/p}$ . The DTEP problem for  $\ell_2$  metrics can be solved with (say) approximation  $1 + \frac{1}{p}$  and communication  $O(p^2)$  using Proposition 1.4 (since finite  $\ell_2$  metrics embed isometrically into Hamming space). Together, we obtain a protocol for the DTEP problem on the metric ed<sup>1/p</sup>, which achieves approximation  $D^{1/p}(1 + \frac{1}{p})$  and communication  $O(p^2)$ . Observe that the same protocol also solves DTEP on the edit metric ed, except that the threshold now is  $R^p$  instead of R, and the approximation is  $(D^{1/p}(1 + \frac{1}{p}))^p < De$ . The communication is the same  $O(p^2)$ , and thus Corollary 1.3 (or Theorem 1.1, respectively) implies that  $De \log(De) \ge 2^{-O(p^2)} \log d$ . For fixed p this completes the proof. □

For the Ulam metric, this distortion lower bound of  $\Omega(\frac{\log d}{\log \log d})$  is near-optimal, since that metric embeds into  $\ell_1$  with  $O(\log d)$  distortion [12]. The previous distortion lower bound was 4/3 [13]. Other upper and lower bounds for low-distortion embeddings appear in Table 1.1.

Sketching of edit distance. The sketch of a string x is a (randomized) mapping of x into a short "fingerprint"  $\mathbf{sk}(x)$  such that sketches of two strings,  $\mathbf{sk}(x)$  and  $\mathbf{sk}(y)$ , are sufficient to distinguish between the case where edit distance is  $\mathrm{ed}(x, y) \leq R/\alpha$ , and the case where  $\mathrm{ed}(x, y) > R$ , for fixed approximation factor  $\alpha > 1$  and parameter R > 1. The main parameter of a sketching algorithm is its sketch size, the length of  $\mathbf{sk}(x)$ .

The sketching model can also be described as a (randomized) simultaneous communication protocol as follows. Alice receives x and computes  $\mathbf{sk}(x)$ , Bob receives yand computes  $\mathbf{sk}(y)$ , and then they send their computed values to a "referee," who needs to decide whether x, y are close or far based only on the sketches. By letting either Alice or Bob play the role of the referee in this simultaneous protocol, one easily sees that the sketch size required by a sketching algorithm is always no smaller than the number of communication bits required by a (general) protocol. The following corollary thus follows immediately from our preceding communication lower bounds.

COROLLARY 1.6. For every d > 1 and  $d^{0.1} \leq R \leq d^{0.49}$ , every O(1)-size sketching algorithm of the standard edit metric or of the Ulam metric can achieve approximation of only  $\Omega(\frac{\log d}{\log \log d})$ .

Sketching with constant sketch size can be viewed as a generalization of the "embeddings approach" presented above, by using Proposition 1.4, albeit with an arbitrarily small constant factor loss in the approximation factor. An important observation is that this more general approach suffices for the purpose of designing an NNS scheme with efficient query time (assuming that computing the sketch can be done efficiently) and with polynomial storage.<sup>3</sup> Indeed, the nearest neighbor data structure for the Hamming metric of [24] could be viewed as an instantiation of the last step. In addition, sketching can be useful for the original goal of quickly estimating the distance (e.g., as a filtering step).

The sketching model is also important as a basic computational notion for massive data sets, and in recent years, an intensive research effort has led to several sketching algorithms for DTEP over different metrics. Prior to our work, there were essentially three metrics for which a sketch size's lower bounds were known:  $\ell_1$  [35] (equivalently, for  $\ell_p$ ,  $p \in (1,2]$ ),  $\ell_\infty$  [32, 7] (implying lower bounds for  $\ell_p$ , p > 2), and the Earthmover distance over  $\{0,1\}^d$  [2].

Sketching of edit distance was studied in [8, 6, 29, 12], but the only lower bound known for sketching of edit distance is trivial in the sense that it follows immediately from Hamming distance (by a straightforward reduction). This lower bound on the sketch size is  $\Omega(1/\epsilon^2)$  for approximation  $\alpha = 1 + \epsilon$  [35], which becomes uninformative for even a 2-approximation. In fact, Bar-Yossef et al. [6] write that "The state of affairs indicates that proving sketching lower bounds for edit distance may be quite hard."

**1.3.** Our techniques. Our proof of Theorem 1.1 consists of three steps. Generally speaking, we design two input distributions:  $\tilde{\mu}_0$  over "far" pairs (x, y) (i.e.,  $\operatorname{ed}(x, y) > R$ ) and  $\tilde{\mu}_1$  over "close" pairs (i.e.,  $\operatorname{ed}(x, y) \leq R/\alpha$ ). The goal then becomes to show that these distributions are indistinguishable by protocols with low communication complexity. By Yao's minimax principle, it suffices to consider deterministic protocols.

The first step reduces the problem to proving that the two distributions  $\tilde{\mu}_0, \tilde{\mu}_1$ are indistinguishable by boolean functions over  $\mathbb{Z}_p^d$ . Roughly speaking, we show that if there is a protocol using at most l bits of communication, then there exists a (deterministic) sketching protocol that uses sketch size of 1 bit and achieves an advantage of at least  $\Omega(2^{-l})$  in distinguishing between the two distributions. Let  $\mathcal{H}^A, \mathcal{H}^B : \mathbb{Z}_p^d \to \{-1, +1\}$  be the boolean functions that Alice and Bob, respectively, use as their sketch functions. We can then further restrict the sketching protocol so that the referee decides by checking whether or not  $\mathcal{H}^A(x) = \mathcal{H}^B(y)$ . This step follows the approach employed earlier in [2], with some minor technical differences.

The second step's main goal is to further characterize the advantage achieved by  $\mathcal{H}^A, \mathcal{H}^B$  in terms of a carefully crafted measure of statistical distance between the two input distributions  $\tilde{\mu}_0, \tilde{\mu}_1$ . For this approach to be effective, it is important

<sup>&</sup>lt;sup>3</sup>In particular, one can first amplify the sketching's probability of success to  $1-n^{-\Omega(1)}$ , where *n* is the number of points in the dataset, using sketch size  $O(\log n)$ . Then, the data structure preindexes all possible sketches in the amplified protocol, using only  $2^{O(\log n)} = n^{O(1)}$  space. For each possible value of the amplified sketch, the data structure stores the answer that the sketching referee would conclude from the sketch of the query and that of each dataset point. Note that, in fact, *s*-size sketches imply  $n^{O(s)}$ -size nearest neighbor data structure.

that the functions  $\mathcal{H}^A, \mathcal{H}^B$  depend only on a few coordinates of their inputs, and in order to guarantee this (indirectly), we include in  $\tilde{\mu}_0, \tilde{\mu}_1$  a noise component, which effectively destroys any dependence of  $\mathcal{H}^A, \mathcal{H}^B$  on many coordinates. Specifically, this step assumes that each distribution  $\tilde{\mu}_t$ ,  $t \in \{0, 1\}$ , has the following structure: choose  $x \in \mathbb{Z}_p^d$  uniformly at random, and then generate y from x via a sequence of two randomized operations. The first of the two is a noise operator with rate  $\rho \in (0, 1)$ ; i.e., each coordinate is modified independently with probability  $1 - \rho$  into a randomly chosen value. The second operation permutes the coordinates according to a permutation drawn from a distribution  $\mathcal{D}_t$ . Given this  $\mathcal{D}_t$ , consider the following derived distribution: take a vector  $u \in \mathbb{Z}_p^d$  with  $\lambda$  nonzero positions (called a  $\lambda$ -test) and apply a random permutation  $\pi \in \mathcal{D}_t$  to it; let  $A_u^{(t,\lambda)}$  be the resulting distribution of vectors. (Note that the support of  $A_u^{(t,\lambda)}$  contains only vectors with precisely  $\lambda$ nonzero entries.) Our measure  $\Delta_{\lambda}$ , called  $\lambda$ -test distinguishability, is the maximum, over all such  $\lambda$ -tests u, of the total variation distance between  $A_u^{(0,\lambda)}$  and  $A_u^{(1,\lambda)}$ . It pretty much captures the statistical advantage in distinguishing  $\mathcal{D}_0$  from  $\mathcal{D}_1$  (and thus  $\tilde{\mu}_0$  from  $\tilde{\mu}_1$ ) achievable by inspecting only  $\lambda$  positions of, say, y (e.g., by tracing them back to x). Altogether, our upper bound on the advantage achieved by  $\mathcal{H}^A, \mathcal{H}^B$  takes roots in the following dichotomy. If  $\mathcal{H}^B$  essentially depends on many coordinates of y (e.g., a linear function with many terms), then the advantage is bounded by  $\rho^{\lambda}$ (i.e., the noise destroys almost all the information), and if  $\mathcal{H}^B$  essentially depends on a few, say  $\lambda$ , coordinates, then the advantage is bounded by the aforementioned  $\Delta_{\lambda}$ . To prove this dichotomy, we rely on Fourier analysis which expands  $\mathcal{H}^{A}, \mathcal{H}^{B}$  into linear functions at different levels  $\lambda$ .

In the third step, we complete the description of  $\tilde{\mu}_0$ ,  $\tilde{\mu}_1$  by detailing the construction of  $\mathcal{D}_0$ ,  $\mathcal{D}_1$  and give an upper bound on the  $\lambda$ -test distinguishability  $\Delta_{\lambda}$  for these distributions. In a simplified view, each distribution  $\mathcal{D}_t$  is generated by a block rotation operation, namely, choosing a random block of length L and applying to it  $\epsilon_t L$ cyclic shifts. The difference between the two distributions is in the magnitude of the rotation (namely,  $\epsilon_t$ ).

Our use of Fourier analysis is elementary and does not involve the KKL theorem [21] or Bourgain's noise sensitivity theorem [11], which were used in the previous nonembeddability results for edit distance [22, 23]. We also note that our hard distribution is notably different from the distributions of [23] or [22], which do admit efficient communication protocols.

To prove Theorem 1.2, we give a new characterization of the Ulam distance between two strings. In particular, building on the work of [30, 17], we prove that if two strings (permutations) P, Q are at distance k = ed(P, Q), then there exist  $\Theta(k)$  pairs of characters in P, all characters at distinct positions, such that for each pair (a, b)their order in P is opposite to that in Q (if they appear in Q at all). We then exploit this characterization by a careful counting of the number of the possible low-cost alignments between P and Q, tailored to the aforementioned  $\Theta(k)$  positions.

**2. Preliminaries.** We use the notation  $[d] = \{1, 2, ..., d\}$  and  $\mathbb{Z}_p = \{0, 1, ..., p-1\}$ . For a vector  $u \in \mathbb{Z}_p^d$ , define the *weight* of u, denoted wt(u), to be the number of coordinates in u that are nonzero.

DEFINITION 2.1. For matrix  $A \in M_{n,n}(\mathbb{R})$  and  $p \in [1,\infty]$ , the p-norm of A is defined by  $||A||_p = \max\{||Av||_p : v \in \mathbb{C}^n, ||v||_p = 1\}.$ 

**2.1. Fourier analysis over**  $\mathbb{Z}_p^d$ . We review basic Fourier analysis over  $\mathbb{Z}_p^d$  for a prime  $p \geq 2$ .

The collection of functions  $f: \mathbb{Z}_p^d \to \mathbb{C}$  is a vector space of dimension  $p^d$ , equipped with an inner product given by  $\langle f,g\rangle = \mathbb{E}_{x\in\mathbb{Z}_p^d}[f(x)\cdot\overline{g(x)}]$ . For  $u\in\mathbb{Z}_p^d$ , define a character  $\chi_u(x) = e^{\frac{2\pi i}{p}(x \cdot u)}$ , where  $x \cdot u$  is the scalar product of  $x, u \in \mathbb{Z}_p^d$ . The set of characters  $\{\chi_u \mid u \in \mathbb{Z}_p^d\}$  forms an orthonormal basis, called the Fourier basis. Thus every function  $f: \mathbb{Z}_p^d \to \mathbb{C}$  admits a Fourier expansion  $f = \sum_{u \in \mathbb{Z}_p^d} \hat{f}_u \chi_u$ , where  $\hat{f}_u = \langle f, \chi_u \rangle$  is called the Fourier coefficient of f corresponding to u. Parseval's identity states that  $\mathbb{E}_{x \in \mathbb{Z}_n^d} [f(x)\overline{g(x)}] = \sum_{u \in \mathbb{Z}_n^d} \hat{f}_u \hat{\overline{g}}_u$ .

We let  $N_{\rho}$  stand for a *noise* vector over  $\mathbb{Z}_{p}^{d}$ , namely, a vector where each coordinate is set independently at random as follows: with probability  $\rho$  it is set to zero, and with probability  $1 - \rho$  it is set to a random value from  $\mathbb{Z}_p$ . We refer to  $\rho$  as the rate of the noise.

The noise operator  $T_{\rho}$  (also called Bonami–Beckner operator) operates on functions  $f : \mathbb{Z}_p^d \to \mathbb{R}$  and is defined by  $(T_{\rho}f)(x) = \mathbb{E}_{N_{\rho}}[f(x+N_{\rho})]$ . The following standard fact relates the Fourier coefficients of f with those of  $T_{\rho}f$ .

FACT 2.2. For every vector  $u \in \mathbb{Z}_p^d$ ,  $(T_\rho f)_u = \hat{f}_u \cdot \rho^{\operatorname{wt}(u)}$ . Proof. We can write  $(T_\rho f)(x) = \mathbb{E}_{N_\rho} [f(x + N_\rho)]$  as

$$\mathbb{E}_{N_{\rho}}\left[\sum_{u\in\mathbb{Z}_{p}^{d}}\hat{f}_{u}e^{\frac{2\pi i}{p}u\cdot(x+N_{\rho})}\right] = \sum_{u\in\mathbb{Z}_{p}^{d}}\hat{f}_{u}e^{\frac{2\pi i}{p}u\cdot x}\mathbb{E}_{N_{\rho}}\left[e^{\frac{2\pi i}{p}u\cdot N_{\rho}}\right] = \sum_{u\in\mathbb{Z}_{p}^{d}}\hat{f}_{u}\rho^{\mathrm{wt}(u)}\chi_{u}$$

where we used the fact that for every  $w \in \mathbb{Z}_p \setminus \{0\}$  we have  $\mathbb{E}_{v \in \mathbb{Z}_p}[e^{\frac{2\pi i}{p}wv}] = 0$ .  $\square$ Note that, for p = 2, i.e., Fourier expansion over  $\{0,1\}^d$ , this is equivalent to having  $(T_{\rho}f)_{S} = \hat{f}_{S}\rho^{|S|}$  for every  $S \subseteq [d]$ .

**2.2. Edit metric and Ulam metric.** Let  $\Sigma$  be the alphabet; we mostly consider  $\Sigma = \{0, 1\}$  or  $\Sigma = \mathbb{Z}_p = \{0, 1, \dots, p-1\}$  for  $p \in \mathbb{N}$  (we will use  $p = \Theta(d^3)$ ).

For  $x \in \Sigma^d$ , we let  $x_i$  denote the *i*th position in x whenever  $i \in [d]$  and extend the notation to  $i \notin [d]$  by defining  $x_i = x_j$ , where  $i \equiv j \pmod{d}$  and  $j \in [d]$ .

DEFINITION 2.3 (edit metrics). Let d be a positive integer. The edit metric over  $\Sigma$  is the space  $\Sigma^d$  endowed with distance function ed(x, y), which is defined as the minimum number of character substitutions/insertions/deletions to transform xinto y.

When  $|\Sigma| \geq d$ , let the Ulam metric be the space of permutations  $x \in \Sigma^d$ , where x is called a permutation if no symbol  $c \in \Sigma$  appears more than once in x. This space is endowed with the same distance function ed(x, y).

We note that allowing alphabets  $\Sigma$  bigger than [d] does not make the Ulam metric harder (at least in our communication complexity setting), and thus our main theorem carries over to the Ulam metric over permutations with alphabet  $\Sigma = [d]$  (i.e., the standard notion of permutations). In particular, one can perform the following reduction from the former problem to the latter problem.

FACT 2.4. For any string length d and alphabet  $\Sigma$ ,  $|\Sigma| \ge d$ , there is a function  $f: \Sigma^d \to \Sigma^{|\Sigma|}$  such that for every pair of permutations  $x, y \in \Sigma^d$  we have that f(x), f(y) are permutations over  $\Sigma$  and

$$\operatorname{ed}(x, y) \le \operatorname{ed}(f(x), f(y)) \le 3 \operatorname{ed}(x, y).$$

*Proof.* For given  $x \in \Sigma^d$ , construct  $f(x) \in \Sigma^{|\Sigma|}$  by appending all the alphabet symbols that are missing from x in an increasing order. Then, clearly  $ed(f(x), f(y)) \ge$ 

 $\operatorname{ed}(x, y)$ . Furthermore, we claim that  $\operatorname{ed}(f(x), f(y)) \leq 3 \operatorname{ed}(x, y)$ . Indeed, the edit distance between the starting block of length d of f(x) and of f(y) is  $\operatorname{ed}(x, y)$ . Also, if  $z \leq \operatorname{ed}(x, y)$  is the number of symbols that appear in x but not in y and vice versa, then the edit distance between the ending block of length  $|\Sigma| - d$  of f(x) and f(y) is 2z. The total edit distance between f(x) and f(y) is at most  $3 \operatorname{ed}(x, y)$ .

Note that when  $|\Sigma| = p = \Theta(d^3)$ , as is the case in our main theorem,  $\log |\Sigma| = \Theta(\log d)$ , and thus the logarithmic lower bound carries over.

We will also use the following operation on strings, which is illustrated in Figure 2.1.

DEFINITION 2.5 (rotation operations). Fix a positive integer d and an alphabet  $\Sigma$ . For  $s, L \in [d]$ , define the right rotation operation  $\overrightarrow{R}_{s,L} : \Sigma^d \to \Sigma^d$  as follows. When applied to a string x, it takes the substring of x of length L starting at position s (with wraparound) and performs on it one cyclic shift to the right (by 1 position); the rest of x remains unchanged. A left rotation  $\overleftarrow{R}_{s,L}$  is defined similarly. We call L the length of the rotation operation.



FIG. 2.1. The rotation operation  $\overrightarrow{R}_{s,L}(\cdot)$ . Here,  $\sigma$  is the substring of length L-1 starting at position s in x, and a is the character at position s + L - 1 in x.

Note that  $\overrightarrow{R}_{s,L}$  works as a permutation (and thus is a bijection on the space of strings). Also, for  $i \in [L]$ ,  $(\overrightarrow{R}_{s,L})^i$  is a rotation of the same block by *i* positions to the right. Note that a rotation operation  $\overrightarrow{R}_{s,L}$  can be simulated by at most two deletions and two insertions (and only one of each when the rotation block does not wrap around at the string's boundary). Thus,  $\operatorname{ed}(x, (\overrightarrow{R}_{s,L})^i(x)) = O(i)$  for every xand *i*.

3. Proof of main theorem. In this section we prove Theorem 1.1. Fix the values of d and R, and let us use the alphabet  $\Sigma = \mathbb{Z}_p$  for p sufficiently large so that a random string from  $\Sigma^d$  is a permutation with high probability (e.g., it suffices to set p to be the smallest prime greater than  $d^3$ ). For the rest of this section, we denote our hard distribution by  $\tilde{\mu} = \frac{\tilde{\mu}_0 + \tilde{\mu}_1}{2}$ , where  $\tilde{\mu}_0$  will be a distribution over far pairs of strings (x, y) and  $\tilde{\mu}_1$  will be a distribution over close pairs (x, y), i.e.,  $\operatorname{ed}(x, y) > R$  and  $\operatorname{ed}(x, y) \leq R/\alpha$ , respectively.

We will follow the steps outlined in section 1.3 and eventually put all the pieces together in section 3.3. Our general approach to proving the theorem uses just a few simple properties of the hard distribution, which we will specify along the way. To differentiate the underlying technique from the specifics of our hard distribution, we describe the hard distribution and prove its required properties separately in section 4.

**3.1. Reduction to boolean functions.** Our first lemma says that if there is an efficient communication protocol, then there are boolean functions with a nonnegligible advantage in distinguishing the distribution  $\tilde{\mu}_0$  from  $\tilde{\mu}_1$ . This lemma is based on the ideas from [2], although the presented proof is simpler than in [2].

LEMMA 3.1. Let  $\tilde{\mu}_0$  and  $\tilde{\mu}_1$  be distributions over far and close pairs, respectively. If  $CC_{\alpha,R}^{Ulam_d} \leq l$  for some  $l \geq 1$ , then there exist boolean functions  $\mathcal{H}^A, \mathcal{H}^B : \mathbb{Z}_p^d \rightarrow \{-1,+1\}$  such that

$$\Pr_{\tilde{\mu}_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] - \Pr_{\tilde{\mu}_1}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] \geq \frac{1}{3} \cdot 2^{-l}.$$

*Proof.* The idea is to reduce the general communication protocol to a simultaneous (i.e., sketching) protocol where Alice and Bob each send a sketch of one bit only, and the referee performs an equality test on these two bits. Then, using Yao's minimax principle, we easily obtain two deterministic boolean functions  $\mathcal{H}^A$  and  $\mathcal{H}^B$  that complete the proof.

To accomplish the reduction, consider an actual *l*-bit (randomized) protocol  $\Pi$ . We construct a one-bit sketching protocol as follows: Alice and Bob make a random guess of the entire transcript of an *l*-bit protocol using the public coins, uniform over the space of all  $2^l$  protocols (the guess is independent of the actual inputs). Each of them then checks whether the guessed transcript describes the messages they would send in the actual protocol  $\Pi$ , using the guessed transcript to simulate the other party's messages. For example, Alice starts the protocol  $\Pi$  (that depends on her input), but instead of sending the messages to Bob, she verifies that her messages are exactly the same as the ones appearing in the guessed protocol. Alice also uses the messages from the guessed protocol to simulate Bob's answers.

If at any moment Alice (or Bob) spots an inconsistency, she (or he) sends a bit chosen independently at random. Otherwise, Alice outputs 1, and Bob outputs the outcome of the guessed transcript. Observe that if the guessed transcript is not equal to the actual protocol they would have run, then at least one of the two players notices an inconsistency, and one of the bits output by Alice or Bob is random.

Thus, if x and y are such that  $\operatorname{ed}(x, y) \leq R/\alpha$  (close pair), then Alice and Bob's bits are equal with probability at least  $\frac{2}{3} \cdot 2^{-l} + (1 - 2^{-l})\frac{1}{2} = \frac{1}{2} + \frac{1}{6}2^{-l}$  (where  $\frac{2}{3}$  is the probability that the original protocol II succeeds on (x, y)). Similarly, if x and y are such that  $\operatorname{ed}(x, y) > R$  (far pair), then Alice and Bob's bits are equal with probability at most  $\frac{1}{3} \cdot 2^{-l} + (1 - 2^{-l}) \cdot \frac{1}{2} = \frac{1}{2} - \frac{1}{6}2^{-l}$ . Using Yao's minimax principle, we conclude that, for given distributions  $\tilde{\mu}_0$  and  $\tilde{\mu}_1$  over far and close pairs, respectively, there exist some fixed boolean functions  $\mathcal{H}^A, \mathcal{H}^B$  that achieve a success probability at least  $\frac{1}{2} + \frac{1}{6}2^{-l}$  on the distribution  $\tilde{\mu} = \frac{\tilde{\mu}_0 + \tilde{\mu}_1}{2}$ , or, formally,

$$\frac{1}{2} \Pr_{\tilde{\mu}_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] + \frac{1}{2} \Pr_{\tilde{\mu}_1}[\mathcal{H}^A(x) = \mathcal{H}^B(y)] \ge \frac{1}{2} + \frac{1}{6} \cdot 2^{-l}.$$

We conclude that  $\Pr_{\tilde{\mu}_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] - \Pr_{\tilde{\mu}_1}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] \geq \frac{1}{3} \cdot 2^{-l}$ . The rest of the proof of Theorem 1.1 uses these boolean functions  $\mathcal{H}^A, \mathcal{H}^B$ .

**3.2. From boolean functions to**  $\lambda$ **-tests.** Next we provide a method to lower bound the advantage achieved by the boolean functions  $\mathcal{H}^A, \mathcal{H}^B$  by relating it to a certain statistical property of the hard distribution  $\tilde{\mu}$ . Our hard distribution  $\tilde{\mu} = \frac{\tilde{\mu}_0 + \tilde{\mu}_1}{2}$  will have a specific generic construction that we describe next. For each  $t \in \{0, 1\}$ , the distribution  $\tilde{\mu}_t$  is formed via a small modification of another distribution  $\mu_t$ , which is easier to analyze (due to certain independencies) but might (rarely) produce invalid inputs. Specifically, each  $\tilde{\mu}_t$  is the distribution  $\mu_t$  conditioned on the fact that the pair  $(x, y) \in \mu_t$  is valid in the sense that x and y are both permutations and the pair (x, y) is, respectively, a far (when t = 0) or a close (when t = 1) pair. We analyze below the distributions  $\mu_0$  and  $\mu_1$  (specifically, in Lemma 3.4). For completeness, we

mention that, in the next section, we show that this analysis extends to distributions  $\tilde{\mu}_0$  and  $\tilde{\mu}_1$  using the fact that  $\tilde{\mu}_0$  and  $\tilde{\mu}_1$  are statistically very close to distributions  $\mu_0$  and  $\mu_1$ , respectively.

The distribution  $\mu_t$  consists of pairs (x, y) chosen as follows:  $x \in \mathbb{Z}_p^d$  is chosen uniformly at random, and y is constructed from x in two steps. In the first step, let  $z \triangleq x + N_\rho$ , where  $N_\rho$ , defined in the preliminaries, is noise of rate  $\rho \in (0, 1)$ , independent of t. In the second step, y is obtained from z by permuting the coordinates of z according to a distribution  $\mathcal{D}_t$ . Formally,  $\mathcal{D}_t$  is a distribution over permutation operations, where a *permutation operation* is a function  $\pi : \mathbb{Z}_p^d \to \mathbb{Z}_p^d$  for which there exists a permutation  $\hat{\pi} : [d] \to [d]$  such that  $\pi(x) \equiv (x_{\hat{\pi}(1)}, \ldots, x_{\hat{\pi}(d)})$ . We will require that  $\mathcal{D}_t$  be symmetric in the sense that, for every  $\pi$ , the permutation operations  $\pi$ and  $\pi^{-1}$  are equiprobable (in it). Notice that y has the same marginal distribution as x, i.e., uniform over  $\mathbb{Z}_p^d$ .

We now quantify the "difference" between the distributions  $\mathcal{D}_0, \mathcal{D}_1$  from the perspective of what we call  $\lambda$ -tests. For  $\lambda \in [d]$ , we define a  $\lambda$ -test to be a vector  $u \in \mathbb{Z}_p^d$ with precisely  $\lambda$  nonzero entries, i.e., wt $(u) = \lambda$ . For a distribution  $\mathcal{D}_t$  and  $\lambda \in [d]$ , let the matrix  $A^{(t,\lambda)}$  be the transition matrix of a Markov chain whose states are all the  $\lambda$ -tests and whose transitions are according to  $\mathcal{D}_t$ ; i.e., at a  $\lambda$ -test u, the process picks  $\pi \in \mathcal{D}_t$  and moves to state  $\pi(u)$  (which is also a  $\lambda$ -test). In other words, a row corresponding to u in  $A^{(t,\lambda)}$  is a vector that has, for every  $\lambda$ -test w, a coordinate of value  $\Pr_{\pi \in \mathcal{D}_t}[\pi(u) = w]$ . We denote this row by  $A_u^{(t,\lambda)}$ . Note that the matrix  $A^{(t,\lambda)}$ is symmetric (since  $\mathcal{D}_t$  is symmetric), and thus it is doubly stochastic.

DEFINITION 3.2. The  $\lambda$ -test distinguishability of  $\mathcal{D}_0, \mathcal{D}_1$ , denoted  $\Delta_{\lambda}$ , is the maximum, over all  $\lambda$ -tests u, of the total variation distance between the distributions  $A_u^{(0,\lambda)}$  and  $A_u^{(1,\lambda)}$ .

We can also write  $\Delta_{\lambda}$  using matrix norms (as per Definition 2.1) and the easy fact that  $||B||_{\infty} = \max_{i \in [n]} \sum_{j \in [n]} |B_{ij}|$  for all  $B \in M_{n,n}(\mathbb{R})$ . Later (in Fact 3.8) we shall use known inequalities because different matrix norms (in particular  $\ell_{\infty}$  and  $\ell_2$ ).

FACT 3.3.  $\Delta_{\lambda} = \|A^{(0,\lambda)} - A^{(1,\lambda)}\|_{\infty}/2.$ 

The following lemma bounds the advantage achieved by  $\mathcal{H}^A, \mathcal{H}^B$  in terms of the  $\lambda$ -test distinguishability  $\Delta_{\lambda}$  of distributions  $\mathcal{D}_0$  and  $\mathcal{D}_1$  for any pair of distributions  $\mathcal{D}_0, \mathcal{D}_1$ . Note that we have not yet specified the distributions  $\mathcal{D}_0$  and  $\mathcal{D}_1$  themselves. We will specify the distributions  $\mathcal{D}_0$  and  $\mathcal{D}_1$  in section 4, thus completing the definition of the hard distribution  $\tilde{\mu}$ .

LEMMA 3.4. Consider  $\mathcal{H}^A, \mathcal{H}^B : \mathbb{Z}_p^d \to \{-1, +1\}$  and  $\rho \in (0, 1)$ . If each  $\mu_t$ , for  $t \in \{0, 1\}$ , is defined as above from a symmetric distribution  $\mathcal{D}_t$  over permutation operations, then

$$\Pr_{\mu_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] - \Pr_{\mu_1}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] \le \max_{\lambda \in [d]} \Delta_\lambda \rho^\lambda.$$

*Proof.* For  $t \in \{0,1\}$ , define  $C^{(t)} \triangleq \mathbb{E}_{\mu_t} \left[ \mathcal{H}^A(x) \mathcal{H}^B(y) \right]$  to be the *correlation* between the two boolean functions. Note that

$$\Pr_{\mu_t}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] = \frac{1}{4} \mathbb{E}_{\mu_t} \left[ \mathcal{H}^A(x) - \mathcal{H}^B(y) \right]^2 = \frac{1}{2} - \frac{C^{(t)}}{2}$$

Thus,

$$\Pr_{\mu_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] - \Pr_{\mu_1}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] = \frac{C^{(1)} - C^{(0)}}{2}.$$

We will show that  $C^{(1)} - C^{(0)} \leq 2 \max_{\lambda \in [d]} \Delta_{\lambda} \rho^{\lambda}$ . For this purpose, it is more convenient to express each  $C^{(t)}$  in terms of the Fourier coefficients of  $\mathcal{H}^A$  and  $\mathcal{H}^B$ . Recall that  $\mu_t$  is generated by picking a random x and constructing y from x by adding to it the noise  $N_{\rho}$  and then applying a random permutation drawn from  $\mathcal{D}_t$ , namely,  $y = \pi(x + N_{\rho})$ , where  $\pi \in \mathcal{D}_t$ . Let  $\mu_t | x$  denote the distribution  $\mu_t$  conditioned on the value of x. Thus,

$$\mathbb{E}_{\mu_t} \left[ \mathcal{H}^A(x) \mathcal{H}^B(y) \right] = \mathbb{E}_{x \in \mathbb{Z}_p^d} \left[ \mathcal{H}^A(x) \cdot \mathbb{E}_{\mu_t \mid x} \left[ \mathcal{H}^B(y) \right] \right].$$

Define  $f^{(t)}(x) \triangleq \mathbb{E}_{\mu_t|x} \left[ \mathcal{H}^B(y) \right]$ . Then

$$f^{(t)}(x) = \mathbb{E}_{N_{\rho}} \left[ \mathbb{E}_{\pi \in \mathcal{D}_{t}} \left[ \mathcal{H}^{B}(\pi(x + N_{\rho})) \right] \right]$$

Since  $C^{(t)} = \mathbb{E}_x \left[ \mathcal{H}^A(x) f^{(t)}(x) \right]$ , we can switch to the Fourier basis by applying Parseval's identity and get

(3.1) 
$$C^{(t)} = \sum_{u \in \mathbb{Z}_p^d} \widehat{(\mathcal{H}^A)}_u \widehat{(f^{(t)})}_u,$$

where  $(\widehat{\mathcal{H}^A})_u$  and  $\widehat{(f^{(t)})}_u$  are the Fourier coefficients of  $\mathcal{H}^A$  and  $f^{(t)}$ , respectively. The next proposition, which we shall prove shortly, expresses the level  $\lambda$  Fourier

The next proposition, which we shall prove shortly, expresses the level  $\lambda$  Fourier coefficients of  $f^{(t)}$  in terms of those of  $\mathcal{H}^B$ . Let  $(\widehat{(f^{(t)})}_u)_{u:\mathrm{wt}(u)=\lambda}$  be the vector of the Fourier coefficients of  $f^{(t)}$  indexed by u's of weight  $\mathrm{wt}(u) = \lambda$ . Define  $(\widehat{(\mathcal{H}^B)}_u)_{u:\mathrm{wt}(u)=\lambda}$  similarly.

PROPOSITION 3.5. For all  $\lambda \in [d]$  and  $\mathcal{H}^B : \mathbb{Z}_p^d \to \mathbb{C}$ ,

(3.2) 
$$\left(\widehat{(f^{(t)})}_{u}\right)_{u:\mathrm{wt}(u)=\lambda} = \rho^{\lambda} A^{(t,\lambda)} \cdot \left(\widehat{(\mathcal{H}^{B})}_{u}\right)_{u:\mathrm{wt}(u)=\lambda}.$$

This proposition naturally leads us to break each  $C^{(t)}$  into the terms corresponding to each Fourier level  $\lambda$ . Define the  $\lambda$ th *correlation* to be

(3.3) 
$$C_{\lambda}^{(t)} \triangleq \sum_{u \in \mathbb{Z}_p^d: \mathrm{wt}(u) = \lambda} \widehat{(\mathcal{H}^A)}_u \overline{\widehat{(f^{(t)})}}_u$$

Then,  $C^{(1)} - C^{(0)} = \sum_{\lambda=0}^{d} (C_{\lambda}^{(1)} - C_{\lambda}^{(0)})$ . We can now bound each  $C_{\lambda}^{(1)} - C_{\lambda}^{(0)}$  in terms of  $\Delta_{\lambda}$  and  $\rho$ .

Let  $\omega_{\lambda}^{A} = \|(\widehat{(\mathcal{H}^{A})}_{u})_{u:wt(u)=\lambda}\|_{2}$  be the  $\ell_{2}$ -weight of the level  $\lambda$  Fourier coefficients of  $\mathcal{H}^{A}$ , and define  $\omega_{\lambda}^{B}$  similarly. By Parseval's identity,  $\sum_{\lambda=0}^{d} (\omega_{\lambda}^{A})^{2} = \mathbb{E}_{x}[\mathcal{H}^{A}(x) \cdot \overline{\mathcal{H}^{A}(x)}] = 1$ , and similarly  $\sum_{\lambda=0}^{d} (\omega_{\lambda}^{B})^{2} = 1$ .

PROPOSITION 3.6. For all  $\lambda \in [d]$ ,

$$C_{\lambda}^{(1)} - C_{\lambda}^{(0)} \le 2\Delta_{\lambda}\rho^{\lambda} \cdot \omega_{\lambda}^{A}\omega_{\lambda}^{B}.$$

We will prove the proposition shortly by a straightforward calculation. In addition,  $C_0^{(1)} = C_0^{(0)}$  because the 0th level Fourier coefficient of  $f^{(t)}$  equals  $\mathbb{E}_{x \in \mathbb{Z}_p^d} \left[ f^{(t)}(x) \right]$  $= \mathbb{E}_{y \in \mathbb{Z}_p^d} \left[ \mathcal{H}^B(y) \right]$ , which does not depend on  $t \in \{0, 1\}$ . Given the above proposition,

we thus have

$$C^{(1)} - C^{(0)} = \sum_{\lambda=0}^{d} \left( C_{\lambda}^{(1)} - C_{\lambda}^{(0)} \right) \le \sum_{\lambda=1}^{d} 2\Delta_{\lambda}\rho^{\lambda} \cdot \omega_{\lambda}^{A}\omega_{\lambda}^{B}$$
$$\le \sum_{\lambda=1}^{d} 2\Delta_{\lambda}\rho^{\lambda} \cdot \frac{\left(\omega_{\lambda}^{A}\right)^{2} + \left(\omega_{\lambda}^{B}\right)^{2}}{2} \le 2\max_{\lambda\in[d]} \Delta_{\lambda}\rho^{\lambda},$$

where we used the geometric-arithmetic mean inequality. This finishes the proof of Lemma 3.4. П

It remains to prove Propositions 3.5 and 3.6.

Proof of Proposition 3.5. Define a new function  $g^{(t)}: \mathbb{Z}_p^d \to \mathbb{R}$  as

$$g^{(t)}(z) \triangleq \mathbb{E}_{\pi \in \mathcal{D}_t} \left[ \mathcal{H}^B(\pi(z)) \right]$$

Then  $f^{(t)} = T_{\rho}g^{(t)}$ , and thus  $\widehat{(f^{(t)})}_{u} = \widehat{(g^{(t)})}_{u} \cdot \rho^{\operatorname{wt}(u)}$  for all  $u \in \mathbb{Z}_{p}^{d}$  (by Fact 2.2). It remains to prove that

(3.4) 
$$\left(\widehat{(g^{(t)})}_{u}\right)_{u:\mathrm{wt}(u)=\lambda} = A^{(t,\lambda)} \cdot \left(\widehat{(\mathcal{H}^{B})}_{u}\right)_{u:\mathrm{wt}(u)=\lambda}$$

Similarly to the operator  $T_{\rho}$ , we define the operator  $\mathcal{O}_t$  as

$$(\mathcal{O}_t \mathcal{H}^B)(x) \triangleq \mathbb{E}_{\pi \in \mathcal{D}_t} \left[ \mathcal{H}^B(\pi(x)) \right]$$

Since  $g^{(t)} = \mathcal{O}_t \mathcal{H}^B$ , we proceed to analyze how the operator  $\mathcal{O}_t$  works on the Fourier coefficients of a function  $\mathcal{H}^B$ .

FACT 3.7. For a permutation operation  $\pi$ , define  $\mathcal{P}_{\pi}$  to be an operator on functions  $\psi : \mathbb{Z}_p^d \to \mathbb{R}$ , given by  $(\mathcal{P}_{\pi}\psi)(x) \triangleq \psi(\pi(x))$ . Then,  $(\mathcal{P}_{\pi}\psi)_u = \hat{\psi}_{\pi(u)}$ . Now, the operator  $\mathcal{O}_t$  defined earlier is simply a convex combination of several

 $\mathcal{P}_{\pi}$ , where  $\pi$  is drawn from  $\mathcal{D}_t$ . Thus, with the above fact, for every  $u \in \mathbb{Z}_p^d$ ,

(3.5) 
$$\widehat{(g^{(t)})}_u = (\widehat{\mathcal{O}_t \mathcal{H}^B})_u = \mathbb{E}_{\pi \in \mathcal{D}_t} \left[ (\widehat{\mathcal{H}^B})_{\pi(u)} \right].$$

Consequently, the vector of level  $\lambda$  Fourier coefficients of  $g^{(t)}$  can be written as a product of the matrix  $A^{(t,\lambda)}$  and the vector of the (same) level  $\lambda$  Fourier coefficients of  $\mathcal{H}^B$ , which proves Proposition 3.5. 

We will need the following fact for the proof of Proposition 3.6. Recall that  $||A||_p$ denotes the *p*-norm of such a matrix A, as per Definition 2.1.

FACT 3.8. Let  $B \in M_{n,n}(\mathbb{R})$  be a symmetric matrix. Then,  $||B||_2 \leq ||B||_{\infty}$ . *Proof.* It is known that

$$||B||_1 = \max_{j \in [n]} \sum_{i \in n} |B_{ij}|$$
 and  $||B||_{\infty} = \max_{i \in [n]} \sum_{j \in [n]} |B_{ij}|$ .

and since B is symmetric, these two norms are equal. By the Riesz-Thorin interpolation theorem,  $||B||_2 \le \max\{||B||_1, ||B||_\infty\} = ||B||_\infty$ . (The Riesz-Thorin interpolation theorem states that for every 1  $\leq$  p < q < r  $\leq$   $\infty$  and a real matrix A we have  $||A||_q \le \max\{||A||_p, ||A||_r\}.)$ 

*Proof of Proposition* 3.6. For every  $\lambda$ , the matrix  $A^{(t,\lambda)}$  is symmetric, and so is  $A^{(1,\lambda)} - A^{(0,\lambda)}$ . Thus,

$$\begin{split} C_{\lambda}^{(1)} - C_{\lambda}^{(0)} &= \sum_{u \in \mathbb{Z}_{p}^{d}: \mathrm{wt}(u) = \lambda} \widehat{(\mathcal{H}^{A})}_{u} \cdot \left(\overline{(f^{(1)})}_{u} - \overline{(f^{(0)})}_{u}\right) \\ &\leq \left\| \left( \widehat{(\mathcal{H}^{A})}_{u} \right)_{u:\mathrm{wt}(u) = \lambda} \right\|_{2} \cdot \left\| \left( \overline{(f^{(1)})}_{u} - \overline{(f^{(0)})}_{u} \right)_{u:\mathrm{wt}(u) = \lambda} \right\|_{2} \\ &= \omega_{\lambda}^{A} \cdot \left\| \rho^{\lambda} \left( A^{(1,\lambda)} - A^{(0,\lambda)} \right) \left( \overline{(\mathcal{H}^{B})}_{u} \right)_{u:\mathrm{wt}(u) = \lambda} \right\|_{2} \\ &\leq \rho^{\lambda} \cdot \omega_{\lambda}^{A} \cdot \left\| A^{(1,\lambda)} - A^{(0,\lambda)} \right\|_{2} \left\| \left( \widehat{(\mathcal{H}^{B})}_{u} \right)_{u:\mathrm{wt}(u) = \lambda} \right\|_{2} \\ &\leq \rho^{\lambda} \cdot \omega_{\lambda}^{A} \omega_{\lambda}^{B} \cdot \left\| A^{(1,\lambda)} - A^{(0,\lambda)} \right\|_{\infty} \\ &= 2\Delta_{\lambda} \cdot \rho^{\lambda} \cdot \omega_{\lambda}^{A} \omega_{\lambda}^{B}, \end{split}$$

where we used (3.3), Cauchy–Schwarz, Proposition 3.5, Definition 2.1, Fact 3.8, and Fact 3.3, respectively.

**3.3.** Putting it all together. We proceed to proving Theorem 1.1, using the machinery just developed in sections 3.1 and 3.2. Recall that we still need to exhibit a suitable hard distribution. We outlined the construction of our hard distribution in section 3.2; the construction relies on two distributions  $\mathcal{D}_0$  and  $\mathcal{D}_1$  which were not specified. The next lemma asserts that the desired hard distribution exists. More precisely, it asserts that it can be constructed to satisfy the required properties, such as a small  $\lambda$ -test distinguishability.

LEMMA 3.9 (hard distribution). There exist constants  $\theta, c_1, c_2, d_0 > 0$  such that, for all  $d > d_0$ ,  $p > d^3$ ,  $1 < \alpha \leq O(\frac{\log d}{\log \log d})$ , and  $d^{0.1} \leq R \leq d^{0.49}$ , there exist symmetric distributions  $\mathcal{D}_0^*$  and  $\mathcal{D}_1^*$  over permutation operations on  $\mathbb{Z}_p^d$  (as defined in section 3.2) with the following guarantees.

- (a) For all  $\lambda \geq 1$ , the  $\lambda$ -test distinguishability of  $\mathcal{D}_0^*$  and  $\mathcal{D}_1^*$  is  $\Delta_{\lambda} \leq c_1 \cdot \lambda \frac{\log \alpha}{\log d} \cdot \frac{R}{d}$ .
- (b) Define each distribution μ<sub>t</sub> from D<sup>\*</sup><sub>t</sub> as described in section 3.2, setting ρ = 1 − θ R/α/d. Define the distribution μ̃<sub>t</sub> to be the restriction (i.e., conditioning) of μ<sub>t</sub> to the event that the sampled pair (x, y) ∈ μ<sub>t</sub> be legal, in the sense that x, y ∈ Z<sup>d</sup><sub>p</sub> are permutations and are, respectively, a far pair (for t = 0) or a close pair (for t = 1). Then for each t ∈ {0,1}, the total variation distance between μ̃<sub>t</sub> and μ<sub>t</sub> is at most d<sup>-c<sub>2</sub></sup>.

We prove this lemma separately in section 4, where we include a full description of  $\mathcal{D}_0^*$  and  $\mathcal{D}_1^*$ . Here, we use the lemma to complete the proof of the main theorem.

Proof of Theorem 1.1. First, consider the hard distribution given by Lemma 3.9. Next, by Lemma 3.1, there must exist functions  $\mathcal{H}^A, \mathcal{H}^B$  such that

$$\Pr_{\tilde{\mu}_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] - \Pr_{\tilde{\mu}_1}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] \ge \frac{1}{3} \cdot 2^{-\mathrm{CC}_{\alpha,R}^{\mathrm{Ulam}_d}}$$

Applying Lemma 3.4 to the distributions  $\mu_0, \mu_1$  and using the fact that  $\tilde{\mu}_0$  and  $\tilde{\mu}_1$ , respectively, are statistically close to  $\mu_0$  and  $\mu_1$  (Lemma 3.9(b)), we deduce that

$$\Pr_{\tilde{\mu}_0}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] - \Pr_{\tilde{\mu}_1}[\mathcal{H}^A(x) \neq \mathcal{H}^B(y)] \le \max_{\lambda \in [d]} \Delta_\lambda \rho^\lambda + d^{-c_2}.$$

Combining the two inequalities above and plugging in the upper bound on  $\Delta_{\lambda}$  and the value of  $\rho$  from Lemma 3.9(a), we have

$$\frac{1}{3} \cdot 2^{-\mathrm{CC}_{\alpha,R}^{\mathrm{Ulam}_d}} \leq \max_{\lambda \in [d]} \left[ c_1 \cdot \lambda \frac{\log \alpha}{\log d} \cdot \frac{R}{d} \cdot \left( 1 - \theta \frac{R/\alpha}{d} \right)^{\lambda} \right] + d^{-c_2}$$
$$\leq \frac{c_1}{\theta} \cdot \alpha \cdot \frac{\log \alpha}{\log d} \cdot \max_{x \ge 0} x \cdot e^{-x} + d^{-c_2}$$
$$= O\left(\frac{\alpha \log \alpha}{\log d}\right),$$

which concludes the proof of Theorem 1.1. 

4. Construction of the hard distribution. In this section we prove Lemma 3.9. We start by giving the detailed construction of our hard distribution  $\tilde{\mu} = \frac{\tilde{\mu}_0 + \tilde{\mu}_1}{2}$ . Then, in sections 4.1 and 4.2 we prove, respectively,  $\lambda$ -test indistinguishability (part (a)) and statistical closeness (part (b)) properties of the hard distribution.

The hard distribution construction follows the outline given in section 3.2. We first specify the distributions  $\mathcal{D}_0^*, \mathcal{D}_1^*$  over permutation operators, which form the bulk of the construction. Once these distributions are specified, we obtain the intermediary distributions  $\mu_0$  and  $\mu_1$  as already described in section 3.2. We finalize the description by constructing  $\tilde{\mu}_t$  from  $\mu_t$ , for each  $t \in \{0, 1\}$ , by conditioning on the pair  $(x, y) \in \mu_t$ being a legal pair, namely, that  $x, y \in \mathbb{Z}_p^d$  are permutations and are, respectively, a far pair (for t = 0) or a close pair (for t = 1).

Fix  $\epsilon_0 \triangleq 1/2$ , and select  $\epsilon_1 = \Theta(\frac{1}{\alpha})$  as follows. Let  $\beta \triangleq \frac{1-\epsilon_1}{1-\epsilon_0} = 2(1-\epsilon_1)$ , and let  $\xi_1 \triangleq \lceil \log_2(C_1 \alpha) \rceil$ , for a sufficiently large constant  $C_1 > 0$  (in particular  $C_1 = 805$  will suffice). Let  $\epsilon_1$  be the solution to the equation  $(1 - \epsilon_1) = \epsilon_1 \beta^{\xi_1}$  satisfying  $\epsilon_1 \leq \frac{2}{C_1 \alpha}$ . The existence of  $\epsilon_1$  follows from the following claim, whose proof is deferred to the end of the construction.

CLAIM 4.1. Let  $\alpha > 1$  and  $C_1 > 1$  be sufficiently large. Then there exists  $\epsilon_1$  with  $\frac{1}{3C_1\alpha} < \epsilon_1 \leq \frac{2}{C_1\alpha} \text{ such that } (1-\epsilon_1) = \epsilon_1(2(1-\epsilon_1))^{\xi_1}, \text{ where } \xi_1 = \lceil \log_2(C_1\alpha) \rceil.$ We thus have, by construction,

(4.1) 
$$\epsilon_0 = (1 - \epsilon_0) = (1 - \epsilon_1)\beta^{-1} = \epsilon_1 \beta^{\xi_1 - 1}.$$

For each  $t \in \{0, 1\}$ , we define the distribution  $\mu_t$  over (x, y) such that ed(x, y) is almost surely  $\Theta(\epsilon_t R)$ . Choose  $x \in \Sigma^d = \mathbb{Z}_p^d$  uniformly at random. Then set  $z \triangleq x + N_{\rho}$ , where  $N_{\rho} \in \mathbb{Z}_p^d$  is a noise of rate  $\rho \triangleq 1 - \epsilon_1 R/d$  (i.e., each position is randomized with probability  $1 - \rho = \epsilon_1 R/d$ ). We shall obtain y from z by applying a number of random rotation operations, each picked independently from a specific distribution. We use the following notation:

- $m \triangleq 0.01 \cdot \log_{\beta} d = \Theta(\log d)$  is the number of possible lengths of a rotation operation;
- $L_{\min} \triangleq d^{0.01}$  determines the minimum length of a rotation operation (modulo
- a factor of β);
  w ≜ C<sub>2</sub> · R/m·L<sub>min</sub> is the number of rotation operations that we apply for a sufficiently large constant C<sub>2</sub> > 0 to be determined later (in section 4.2).

Generate a sequence  $(r_1, r_2, \ldots, r_w)$  of w rotations by picking each  $r_i$  independent and identically distributed (i.i.d.) according to the following distribution  $\mathcal{D}_t^{\text{rot}}$ :

1. Pick  $l_i \in [m]$  randomly so that  $\Pr[l_i = l] = \frac{\beta^{-l}}{\zeta}$  for each  $l \in [m]$ , where  $\zeta = \sum_{l=1}^{m} \beta^{-l}$  is the normalization constant.

2. Pick a starting position  $s_i \in [d]$  uniformly at random, and rotate the block that starts at position  $s_i$  and has length (with wraparound)  $L_i = \beta^{l_i} L_{\min}$  by  $\epsilon_t L_i$  positions, either to the right or to the left, at random. We choose  $r_i$  at random from the set  $\{(\tilde{R}_{s,L_i})^{\epsilon_t L_i} \mid s \in [d], \tilde{R} \in \{\vec{R}, \overline{R}\}\}$ .

We note that  $(\tilde{R}_{s,L_i})^{\epsilon_t L_i}$  is not well defined when  $\epsilon_t L_i$  or  $L_i$  are not integers. Overloading the notation, we define  $(\vec{R}_{s,L_i})^{\epsilon_t L_i}$  for noninteger  $\epsilon_t L_i, L_i$  as follows. Let  $B_1$  be the block that starts at position  $s_i$  and has length  $\lfloor (1 - \epsilon_t)L_i \rfloor$ , and let  $B_2$  be the block immediately following  $B_1$  of length  $\lfloor \epsilon_t L_i \rfloor$ , i.e.,

$$B_1 = [s:s + \lfloor (1 - \epsilon_t)L_i \rfloor - 1], \quad B_2 = [s + \lfloor (1 - \epsilon_t)L_i \rfloor : s + \lfloor (1 - \epsilon_t)L_i \rfloor + \lfloor \epsilon_t L_i \rfloor - 1].$$

Then,  $(\overrightarrow{R}_{s,L_i})^{\epsilon_t L_i}$  swaps blocks  $B_1$  and  $B_2$ . We define  $(\overleftarrow{R}_{s,L_i})^{\epsilon_t L_i}$  similarly. To obtain y, we apply to  $z = x + N_{\rho}$  the sequence of rotations  $r_1, \ldots, r_w$ , i.e.,

$$y \triangleq r_w(r_{w-1}(\ldots r_1(z)\ldots)) = (r_w \circ \cdots \circ r_2 \circ r_1)(x+N_\rho).$$

In the language of section 3.2, the distribution  $\mathcal{D}_t^*$  of permutation operations is simply the distribution of  $\pi = r_w \circ r_{w-1} \circ \cdots \circ r_1$ , where  $r_1, \ldots, r_w$  are drawn independently from  $\mathcal{D}_t^{\text{rot}}$ .

Intuitively, each rotation operation  $r_i$ , or more precisely its distribution  $\mathcal{D}_t^{\text{rot}}$ , is designed to achieve the following goal. Consider a position  $j \in [d]$ , and assume for simplicity  $j \in [0.1d, 0.9d]$ . Let the random variable  $Z_{t,j} \in \mathbb{Z}$  be the displacement (change in position) of position j under a (random) rotation operation  $r_i \in \mathcal{D}_t^{\text{rot}}$ ; i.e.,  $Z_{t,j} \in \mathbb{Z}$  is the unique value such that  $r_i(e_j) = e_{j+Z_{t,j}}$ , where  $e_k$  denotes the kth standard basis vector. By construction,  $Z_{t,j}$  is symmetric around 0, i.e.,  $\Pr[Z_{t,j}] =$  $k = \Pr[Z_{t,j} = -k]$ , and its distribution does not depend on j; i.e.,  $Z_{t,j}$  and  $Z_{t,j'}$  have the same distribution (but they are correlated). Moreover, its support, i.e., values k > 0 with probability  $\Pr[Z_{t,j} = k] > 0$ , forms a geometric sequence (because the block length L has a geometric distribution). Let us now condition on the event that position j be included in the rotation block, i.e.,  $Z_{t,j} \neq 0$ . Then the distribution of  $Z_{t,j}$  is almost uniform over the support—this follows from the distribution of L and of s and by (4.1). Furthermore, the distributions of  $Z_{0,j}$  and  $Z_{1,j}$  (when we condition on them being nonzero) are almost identical, because their supports differ only at the boundaries, i.e., at the smallest and largest displacements, again due to (4.1), and they are both almost uniform. We repeat the rotation operation many times in order to obtain a high concentration in the distance between y and z.

To finalize the construction, it remains to define  $\tilde{\mu}_t$  for  $t \in \{0, 1\}$ . We note that we cannot set  $\tilde{\mu}_t$  to be exactly  $\mu_t$  because the latter may sometimes generate pairs (x, y) that are not far or close, respectively, or are not even permutations altogether.  $(x \text{ and } y \text{ are not always permutations since each of the two strings is uniformly at$ random and may have a multiple occurrence of the same symbol.) We thus define $<math>\tilde{\mu}_0$  to be the distribution  $\mu_0$  restricted to (i.e., conditioned on) pairs of permutations (x, y) with  $\operatorname{ed}(x, y) > R$ , and similarly  $\tilde{\mu}_1$  is the distribution  $\mu_1$  restricted to pairs of permutations with  $\operatorname{ed}(x, y) \leq R/\alpha$ .

It remains only to prove Claim 4.1, namely, that the desired  $\epsilon_1$  exists.

Proof of Claim 4.1. Define function  $f(x) : [0,1] \to \mathbb{R}$  as  $f(x) = x \cdot (1-x)^{\xi_1 - 1} 2^{\xi_1 - 1}$ . Note that  $\epsilon_1$  is the solution to the equation f(x) = 0. For  $x = 1/(3C_1\alpha)$ ,  $f(x) \le \frac{1}{3C_1\alpha}(1 - \frac{1}{3C_1\alpha})^{\xi_1 - 1} \cdot 2^{\log_2(C_1\alpha) + 1} - 1 < 0$ . Similarly, for  $x = \frac{2}{C_1\alpha}$ ,  $f(x) \ge \frac{2}{C_1\alpha}(1 - \frac{2(\xi_1 - 1)}{C_1\alpha}) \cdot 2^{\log_2(C_1\alpha)} - 1 \ge 2(1 - \frac{2(\log_2(C_1\alpha) - 1)}{C_1\alpha}) - 1 > 0$ , provided  $C_1$  is a sufficiently large constant. By the continuity of f(x), there exists some  $x \in \left[\frac{1}{3C_{1}\alpha}, \frac{2}{C_{1}\alpha}\right]$  satisfying f(x) = 0.

In the rest of this section we prove the two properties required from our hard distribution, stated in Lemma 3.9: that  $\mathcal{D}_0^*$  and  $\mathcal{D}_1^*$  have small  $\lambda$ -test distinguishability (Lemma 3.9(a)), and that each  $\tilde{\mu}_t$  is very close to  $\mu_t$ , for both  $t \in \{0, 1\}$ (Lemma 3.9(b)).

Here and throughout the big  $O(\cdot)$  notation may hide dependence on constants used in the construction of the hard distribution, namely  $C_1$  and  $C_2$ . Furthermore, although the parameters  $\beta$  and  $\zeta$  are not constants (they depend on  $\alpha$ ), we can bound  $1.5 < \beta < 2$ , which guarantees that  $\frac{1}{1-\beta^{-1}} \leq O(1)$  and  $\frac{1}{\zeta} \leq \beta \leq O(1)$ .

4.1.  $\lambda$ -test indistinguishability. We prove Lemma 3.9(a) via the following lemma.

LEMMA 4.2. Let  $\Delta_{\lambda}$  be the  $\lambda$ -test distinguishability of  $\mathcal{D}_{0}^{*}$  and  $\mathcal{D}_{1}^{*}$ . Then for all  $\lambda \geq 1$  we have  $\Delta_{\lambda} \leq O(\lambda \frac{\log \alpha}{\log d} \cdot \frac{R}{d})$ . *Proof.* Fix a  $\lambda$ -test  $u \in \mathbb{Z}_{p}^{d}$ , and let

$$\delta_{\lambda}(u) = \max_{T \subseteq \mathbb{Z}_p^d} \left| \Pr[r^{(0)}(u) \in T] - \Pr[r^{(1)}(u) \in T] \right|$$

be the total variation distance between the distributions  $r^{(0)}(u)$  and  $r^{(1)}(u)$ , where  $r^{(t)} \in \mathcal{D}_t^{\text{rot}}$  for  $t \in \{0, 1\}$ . The heart of this lemma is the following bound, which we shall prove below:

(4.2) 
$$\delta_{\lambda}(u) \le O\left(\lambda \log \alpha \cdot \frac{L_{\min}}{d}\right).$$

We shall also prove shortly the claim that  $\Delta_{\lambda} \leq w \cdot \max_{u} \delta_{\lambda}(u)$ . The lemma then follows immediately from (4.2), and this claim, by plugging the former into the latter and recalling  $w = C_2 \cdot \frac{R}{m \cdot L_{\min}}$ , is the number of rotation operations. Since  $\lambda \frac{\log \alpha}{\log d} \cdot \frac{R}{d} > 1$  for  $\lambda \ge d^{0.95}$ , it actually suffices to prove (4.2) only for  $\lambda < d^{0.95}$ .

We now prove the above claim, that  $\Delta_{\lambda} \leq w \cdot \max_{u} \delta_{\lambda}(u)$ , by induction. Let  $v_i^t = r_i^{(t)}(r_{i-1}^{(t)}(\dots r_1^{(t)}(u)\dots))$  for  $t \in \{0,1\}$  and  $i \in [w]$ . We prove that, for any  $T \subseteq \mathbb{Z}_p^d$ , we have  $|\Pr[v_i^0 \in T] - \Pr[v_i^1 \in T]| \le i \cdot \max_v \delta_\lambda(v)$ . The base case i = 1 holds by the definition of  $\delta_{\lambda}$ , and so we turn to the inductive step:

$$\begin{aligned} \Pr[v_i^0 \in T] &= \sum_v \Pr[v_{i-1}^0 = v] \Pr[r_i^{(0)}(v) \in T] \\ &\leq \sum_v \Pr[v_{i-1}^0 = v] \left( \Pr[r_i^{(1)}(v) \in T] + \delta_\lambda(v) \right) \\ &\leq \max_v \delta_\lambda(v) + \sum_r \Pr[r_i^{(1)} = r] \Pr[r(v_{i-1}^0) \in T] \\ &\leq \max_v \delta_\lambda(v) + \sum_r \Pr[r_i^{(1)} = r] \left( \Pr[r(v_{i-1}^1) \in T] + (i-1) \cdot \max_v \delta_\lambda(v) \right) \\ &= i \cdot \max_v \delta_\lambda(v) + \Pr[v_i^1 \in T]. \end{aligned}$$

Proving the same inequality with the roles of t = 0 and t = 1 reversed, we obtain that  $\Delta_{\lambda} = \max_{T \subset \mathbb{Z}_{x}^{d}} |\Pr[v_{w}^{0} \in T] - \Pr[v_{w}^{1} \in T]| \le w \cdot \max_{u} \delta_{\lambda}(u).$ 

In the rest of the proof of Lemma 4.2, we prove the bound (4.2). The proof consists of two parts. The first part proves the bound for  $\lambda = 1$  and contains the

main intuition why our distribution is hard. The second part builds on the first one to show the bound for general  $\lambda$ .

Part 1:  $\lambda = 1$ . We prove that  $\delta_1(u) \leq O(\log \alpha \cdot \frac{L_{\min}}{d})$  next. In this part, we shall assume that L and  $\epsilon_t L$  are integers, deferring the full treatment of this technicality to the second part.

Since  $\lambda = 1$ , we have only one nonzero entry in u, say, at position j. For  $t \in \{0, 1\}$ , let  $j_t$  be the random variable denoting the position of the symbol  $u_j$  in the vector  $r^{(t)}(u)$  obtained by applying the random rotation  $r^{(t)} \in \mathcal{D}_t^{\text{rot}}$  on u. Also, let  $Z_t$  be the displacement of  $j_t$  with respect to j on the cycle  $\mathbb{Z}_p$ , and namely  $Z_t = (j_t - j_t)$  $j + d/2 \pmod{d} - d/2$  (where the addition/subtraction of d/2 is for the purpose of accounting for string boundaries). It is not hard to see that the distribution of  $Z_t$ does not depend on the value of j.

The total variation distance between the distributions of  $r^{(0)}(u)$  and of  $r^{(1)}(u)$ equals to the total variation distance between  $Z_0$  and  $Z_1$ . We compute the latter via its complement, i.e., the probability mass that is "common" to the two distributions, which is, formally,  $\sum_{z \in [-d,d]} \min_{t \in \{0,1\}} \Pr_{r^{(t)}}[Z_t = z].$ 

First, we can compute the probability that  $Z_t = 0$ , i.e., the symbol  $u_j$  remains at position j, as follows:

$$\Pr[Z_t = 0] = \frac{d - \mathbb{E}[L]}{d} = 1 - m \cdot \frac{L_{\min}}{\zeta d},$$

irrespective of the value of  $t \in \{0, 1\}$ .

Next, consider the case when  $Z_t \neq 0$ , and note that  $\Pr_{r^{(0)}}[Z_0 \neq 0] = \Pr_{r^{(1)}}[Z_1 \neq 0]$  $[0] = m \cdot \frac{L_{\min}}{d}$ . We show that, conditioned on  $Z_t \neq 0$ , the variable  $Z_t$  is uniform over most of its support, denoted  $S_t$ . Moreover  $S_0$  and  $S_1$  have almost the same size and almost completely overlap. Formally, we prove the following claim.

CLAIM 4.3. There exists a set  $S \subset \mathbb{Z} \setminus \{0\}$  satisfying the following:

- There is  $\nu > 0$  such that for each  $t \in \{0,1\}$  and  $z \in S$  we have  $\Pr_{r^{(t)}}[Z_t =$  $z] = \nu.$
- For each  $t \in \{0,1\}$  we have  $\Pr_{r^{(t)}}[Z_t \in S] \ge \frac{m-\xi_1}{m} \cdot \Pr_{r^{(t)}}[Z_t \neq 0].$

We first show how Claim 4.3 lets us prove that  $\delta_1(u) \leq O(\log \alpha \cdot \frac{L_{\min}}{d})$ . Indeed, one can observe that  $\delta_1(u)$  is bounded by the probability that  $\Pr_{r^{(0)}}[Z_0 \notin S \cup \{0\}] =$  $\Pr_{r^{(1)}}[Z_1 \notin S \cup \{0\}],$  which we can bound as

$$\delta_1(u) \le 1 - \Pr_{r^{(t)}}[Z_t = 0] - \Pr_{r^{(t)}}[Z_t \in S] \le \frac{\xi_1}{m} \cdot \Pr_{r^{(t)}}[Z_t \neq 0] = O(\log \alpha) \cdot \frac{L_{\min}}{\zeta d}.$$

*Proof of Claim* 4.3. We show the claim for

$$S = \{ \pm (1 - \epsilon_1) \beta^l L_{\min} \mid l = 1, \dots, m - \xi_1 \}$$

and  $\nu = \frac{1}{2} \cdot \frac{L_{\min}}{\zeta d}$ . Let us consider the case that  $Z_t \neq 0$ . Then, the magnitude of the displacement,  $|Z_t|$ , must be either  $\epsilon_t L$  or  $(1-\epsilon_t)L$ , where  $L = \beta^l L_{\min}$  for some  $l \in [m]$ . In particular,  $Z_t \neq 0$  iff the position j falls inside the rotation block of the operation  $r^{(t)}$ , and either (i) j falls into the bigger part of size  $(1 - \epsilon_t)L$  (that does not wrap around) and hence  $|Z_t| = \epsilon_t L$ ; or (ii) j falls into the smaller part of size  $\epsilon_t L$  (that does wrap around), and hence  $|Z_t| = L - \epsilon_t L = (1 - \epsilon_t)L$ . Moreover, conditioned on the magnitude of  $Z_t$ , the sign of  $Z_t$  is equiprobable to be either positive or negative (depending on whether the rotation block rotates to the right or left).

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When t = 0, we can compute the probability that  $|Z_0| = \frac{1}{2}L = \frac{1}{2}\beta^l L_{\min}$  for some  $l \in [m]$  as follows. We have  $Z_0 = L/2$  when we choose block length  $L = \beta^l L_{\min}$ , which happens with probability  $\beta^{-l}/\zeta$ , and additionally either (i) position j is inside the "bigger" part of the block, of size  $(1 - \epsilon_0)L = L/2$ , and the block moves to right, or (ii) position j is inside the "smaller" part of the block, of size  $\epsilon_0 L = L/2$ , and the block moves to left. Formally,

$$\Pr_{r^{(0)}}\left[Z_{0} = \frac{L}{2}\right] = \Pr_{r^{(0)}}\left[Z_{0} = -\frac{L}{2}\right] = \frac{\beta^{-l}}{\zeta} \cdot \frac{(1-\epsilon_{0})\beta^{l}L_{\min}}{d} \cdot \frac{1}{2} + \frac{\beta^{-l}}{\zeta} \cdot \frac{\epsilon_{0}\beta^{l}L_{\min}}{d} \cdot \frac{1}{2} = \frac{L_{\min}}{\zeta d} \cdot \frac{1}{2} = \nu.$$

Note that  $z = \frac{1}{2}\beta^l L_{\min}$  may be written as  $z = (1 - \epsilon_1)\beta^{l-1}L_{\min}$  (using (4.1)), and thus  $z \in S$  whenever  $l \in \{2, \ldots, m - \xi_1 + 1\}$ . Now let t = 1. When  $|Z_1| = \epsilon_1\beta^{l+\xi_1}L_{\min} = (1 - \epsilon_1) \cdot \beta^l L_{\min} \in S$  for  $l \in S$ 

 $\{1, \ldots, m - \xi_1\}$  (the equality here is by (4.1)), we again have that

$$\Pr_{r^{(1)}}[Z_1 = \epsilon_1 \beta^{l+\xi_1}] = \frac{\beta^{-l-\xi_1}}{\zeta} \cdot \frac{(1-\epsilon_1)\beta^{l+\xi_1}L_{\min}}{d} \cdot \frac{1}{2} + \frac{\beta^{-l}}{\zeta} \cdot \frac{\epsilon_1 \beta^l L_{\min}}{d} \cdot \frac{1}{2} = \frac{L_{\min}}{\zeta d} \cdot \frac{1}{2} = \nu.$$

Finally, note that  $\Pr_{r^{(t)}}[Z_t \in S] = \sum_{z \in S} \Pr_{r^{(t)}}[Z_t = z] = 2(m - \xi_1) \cdot \nu = \frac{m - \xi_1}{m} \cdot \Pr_{r^{(t)}}[Z_t \neq 0]$ . This concludes the proof of Claim 4.3.

Part 2 :  $\lambda \geq 2$ . When we have  $\lambda \geq 2$  nonzero entries in u, the intuition is to group these nonzero entries into one or more "atomic intervals" and then reduce to the case  $\lambda = 1$  with the role of "symbol  $u_i$ " being replaced by an atomic interval. For example, when there are  $\lambda = 2$  nonzero entries in u, most of the block lengths L fall into two categories:

- L is much larger than the distance between the positions of the two nonzero entries—in this case, the two nonzero symbols from u move jointly (atomically) most of the time, and thus the interval connecting the two symbols behaves roughly as the "symbol  $u_i$ " in the  $\lambda = 1$  scenario;
- L is much smaller than the distance between the two positions—in this case, each of the two nonzero entries can be treated independently as in  $\lambda = 1$  case, and we lose only a factor of  $\lambda$  (by "union bound").

Furthermore, we can bound the number of values of L that do not satisfy one of the above properties. A relatively straightforward bound is  $O(\lambda^2)$  (all pairwise distances between the nonzero entries) times  $O(\xi_1)$  (the same extra factor as in the  $\lambda = 1$  case). This analysis would give a bound of  $\delta_{\lambda}(u) \leq O(\lambda^3 \log \alpha \cdot \frac{L_{\min}}{d})$ . In what follows we obtain a stronger bound, with only a linear dependence on  $\lambda$ , using a more careful analysis. (For the impact of a weaker bound see the calculation in section 3.3.)

More generally, we partition the nonzero entries of u such that each part consists of "nearby" entries, while the parts are "far" amongst themselves. We then view each part as a contiguous A-interval (stands for atomic interval). Once we manage such an approximation, we have several A-intervals (at most  $\lambda$ ), and we expect each one to move atomically: all nonzero entries from the same A-interval will move in the same direction by the same displacement most of the time. The main challenge lies in the fact that the notion of nearby entries depends on the length L of the rotation block, and we say two nonzero entries are nearby if their positions differ by at most L. Thus, for each possible block length L, we have a possibly different partition of entries into A-intervals (partitions are progressively coarser with bigger L). The main technical work is to analyze the structure of these A-intervals over all lengths L.

We proceed with a complete proof below. For a block length  $L = \beta^l L_{\min}$ , we define the graph  $G_L$  as follows.  $G_L$  is an undirected graph on  $\lambda$  vertices, where each vertex corresponds to a nonzero entry in u. For convenience, we use the term "entry" when we refer to the position of a nonzero entry of u and equivalently a vertex of  $G_L$  (in contrast, we will use the term "node" for another graph structure defined later). We connect two entries  $i, j \in [d]$  if  $|i - j|^* \leq L$ , where  $|i - j|^* =$  $\min\{|i - j|, d - |i - j|\}$  computes distance on the d-cycle. For a graph  $G_L$ , we focus on its connected components, which may be viewed as intervals in  $\mathbb{Z}_d$ . Specifically, to each connected component  $C \subset V$  we assign the interval I(C), an interval defined as the minimal interval (with wraparound) on  $\mathbb{Z}_d$  that contains all entries in C. Overloading the notation, we write an interval I(C) = [i, j] to mean that  $I(C) = \{i, i + 1, \ldots, j\}$ if  $i \leq j$  and  $I(C) = \{i, i + 1, \ldots, d, 1, 2, \ldots, j\}$  if j < i. The length of interval I = [i, j]is  $\operatorname{len}(I) = |I| = (j - i + 1) \pmod{d}$ . Note that, for every connected component C, every two consecutive entries in I(C) are at distance at most L; thus, the length of any interval I(C) can be at most  $L \cdot \lambda < d^{0.99}$ ; also if I(C) = [i, j], then both i and jare nonzero entries of u.

An A-interval is then an interval I(C) that corresponds to some connected component C. Each block length L induces a potentially different graph  $G_L$ , which in turn induces a different set of A-intervals. The following observation relates A-intervals induced by different  $G_L$ 's.

OBSERVATION 4.4. If two entries are in the same A-interval (equivalently, connected component) in  $G_L$  for some L, then they are also in the same A-interval in  $G_{L'}$  for any  $L' \geq L$ .

We use this observation to define a forest on all the A-intervals as follows. The forest consists of m levels, where nodes at level  $l \in [m]$  correspond to the A-intervals for  $L = \beta^l L_{\min}$  (i.e., the connected components in  $G_L$ ). For a forest node v at level l we write I(v) for the corresponding A-interval. The edges in the forest are defined as follows: for two forest nodes  $v_1, v_2$  on two consecutive levels, l and l + 1, respectively, we connect  $v_1$  to  $v_2$  iff  $I(v_1) \subseteq I(v_2)$ . This construction is well defined due to Observation 4.4. Nodes at level 1 will be called *leaves*. Notice that every forest node at level l > 1 indeed has at least one edge to a node at level l - 1; i.e., nonleaf nodes have at least one child. Let  $n_l \in [\lambda]$  be the number of nodes at level l.

We now wish to bound the error incurred by considering an A-interval to be an atomic object. Specifically, a too long A-interval is likely to move not atomically, in the sense that the interval is "cut" by the rotation block. We bound the error of our "approximation" using the probability that a random position  $s \in [d]$  (one of the two block boundaries) falls inside these A-intervals at a random level l. The latter probability is proportional to the expected sum of lengths of the A-intervals of  $G_L$  when we choose the block length L randomly according to the distribution  $\mathcal{D}_t^{\text{rot}}$ .

CLAIM 4.5. Let  $s \in [d]$  be chosen uniformly at random, and let  $l \in [m]$  be chosen randomly with probability  $\beta^{-l}/\zeta$ . Then,

$$\Pr_{s,l}[s \text{ is inside one of the A-intervals at level } l] \le O\left(\lambda \frac{L_{\min}}{d}\right).$$

*Proof.* Consider any two *consecutive* nonzero entries of u, and let J be the interval between them (with wraparound), including one of the endpoints, say the left one. We compute next the probability that s is contained in this interval J and interval J is contained in an A-interval I(v) for a forest node v at level l. Note that summing this probability over all  $\lambda$  intervals J gives the final quantity we want.

By definition, an interval J is inside an A-interval at level l iff  $|J| \leq \beta^l L_{\min}$ . Thus, for a fixed J, the probability that both  $s \in J$  and J is contained in an A-interval at

level l is at most

$$\frac{|J|}{d} \cdot \sum_{l \in [m]: \ |J| \le \beta^l L_{\min}} \frac{\beta^{-l}}{\zeta} \le \frac{|J|}{d} \cdot \frac{L_{\min}}{\zeta \cdot |J|} \cdot \frac{1}{1 - \beta^{-1}} \le O\left(\frac{L_{\min}}{d}\right).$$

We have exactly  $\lambda$  such intervals J; thus the total contribution is  $O(\lambda \frac{L_{\min}}{d})$ . We now continue with computing the total variation distance  $\delta_{\lambda}(u)$  between  $r^{(0)}(u)$  and  $r^{(1)}(u)$ , where  $r^{(0)} \in \mathcal{D}_0^{\text{rot}}$  and  $r^{(1)} \in \mathcal{D}_1^{\text{rot}}$ . As in Part 1 ( $\lambda = 1$ ), we will bound the total variation distance between them by estimating the probability mass "common" to the two distributions.

First we compute the probability that all nonzero entries of u stay put (as in Part 1).

CLAIM 4.6. For each  $t \in \{0, 1\}$ , we have that

$$\Pr_{r^{(t)}}[r^{(t)}(u) = u] \ge 1 - O\left(\lambda \frac{L_{\min}}{d}\right) - \sum_{l=1}^{m} n_l \cdot \frac{L_{\min}}{\zeta d}.$$

*Proof.* The complement event is that at least one nonzero entry of u is displaced. Whenever it occurs, at least one of the following holds:

- the left or right endpoint of the rotation block belongs to an A-interval induced by  $G_L$ ; or else
- the rotation block contains inside it an entire A-interval induced by  $G_L$ .

The probability of the first event is bounded by, using Claim 4.5,

$$2\Pr_{s,L}[s \text{ is inside one of the A-intervals at level } l] \le O\left(\lambda \frac{L_{\min}}{d}\right).$$

The probability of the second event can be bounded by the probability that the rotation block includes the leftmost endpoint of some A-interval at level l:

$$\sum_{l=1}^{m} \frac{\beta^{-l}}{\zeta} \sum_{v \text{ at level } l} \Pr_{s} \left[ \text{left endpoint of } I(v) \text{ is inside } [s, s+L-1] \right]$$
$$\leq \sum_{l=1}^{m} \frac{\beta^{-l}}{\zeta} \cdot \frac{n_{l} \cdot L}{d} = \sum_{l=1}^{m} n_{l} \frac{L_{\min}}{\zeta d}.$$

The claim follows from the last two inequalities by applying a union and then considering the complement event. 

We now prove a claim that should be seen as the analogue of Claim 4.3 from Part 1, which characterizes the common weight of the distributions of  $r^{(0)}(u)$  and  $r^{(0)}(u)$  when some entries (more precisely, A-intervals) move. In contrast to Part 1, here we have to also consider the case when an A-interval does not behave atomically, i.e., when the rotation block intersects the A-interval of some node v at a level  $l \in [m]$ . This will contribute some additional error term that depends on the length of the interval I(v) and which we will bound using Claim 4.5.

Let us define the random variable  $Z_t(I)$ , for an interval I = I(v) corresponding to a forest node v, and  $t \in \{0, 1\}$ .  $Z_t(I)$  denotes the (position) displacement of the entries from the interval I under rotation  $r^{(t)} \in \mathcal{D}_1^{\text{rot}}$  when the interval I moves atomically and no entry outside I moves. We set  $Z_t(I) = \bot$  if the interval I does not move atomically and/or some other entry outside I moves as well under  $r^{(t)}$ .

CLAIM 4.7. There exists a set  $S \subset \mathbb{Z} \setminus \{0\}$  satisfying the following:

• For each interval I = I(v) corresponding to a forest node v at level  $l^* \in \{\xi_1 + 1, \ldots, m\}$ , and for each  $t \in \{0, 1\}$  and  $z \in S$ ,

$$\Pr_{r^{(t)}} [Z_t(I) = z] \ge \frac{1}{2} \cdot \frac{L_{\min}}{\zeta d} - \beta^{-(l^* - \xi_1)} \cdot \frac{2 \operatorname{len}(I)}{\zeta d}$$

• Call two intervals I(v) and I(v') distinct if they have at least one distinct endpoint; then

$$\sum_{z \in S} \sum_{\text{distinct } I=I(v)} \min_{t \in \{0,1\}} \left\{ \Pr_{r^{(t)}} \left[ Z_t(I) = z \right] \right\} \ge \sum_{l=\xi_1+1}^m n_l \frac{L_{\min}}{\zeta d} - O\left(\lambda \frac{L_{\min}}{d}\right).$$

*Proof.* We show the claim for  $S = \{ \pm \lfloor (1 - \epsilon_1)\beta^l L_{\min} \rfloor \mid l = 1, \dots, m - \xi_1 \}.$ 

Fix an interval I = I(v) for a node at level  $l^* \ge \xi_1 + 1$ . Consider the displacement  $z = \lfloor \epsilon_1 \beta^{l^*} L_{\min} \rfloor = \lfloor (1 - \epsilon_1) \beta^{l^* - \xi_1} L_{\min} \rfloor \in S$  (the equality again is by (4.1)). We now bound  $\Pr[Z_1(I) = z]$ , namely the probability that all the entries in I(v) are moved (atomically) z positions to the right (and all the other entries stay put), under the distribution  $\mathcal{D}_1^{\text{rot}}$ . We have  $\Pr[Z_1(I) = z]$  when either (i)  $l = l^*$ , interval I is completely inside the "bigger" part of the block, of size  $\lfloor (1 - \epsilon_1)\beta^l L_{\min} \rfloor$ , and the block moves to right, or (ii)  $l = l^* - \xi_1$ , interval I is completely inside the "smaller" part of the block, of size  $\lfloor \epsilon_1 \beta^{l - \xi_1} L_{\min} \rfloor$ , and the block moves to left. Note that in both cases all entries outside I stay put as they are at (position) distance at least  $\beta^{l^*} L_{\min} + 1$  from I and thus cannot be inside the rotation block. Formally,

$$\Pr_{r^{(1)}} [Z_1(I) = z] = \Pr_{r^{(1)} = (\tilde{R}_{s,L})^{\epsilon_1 L}, \ L = \beta^l L_{\min}} \left[ Z_1(I) = z, \ l = l^*, \ \tilde{R} = \overrightarrow{R} \right] + \Pr_{r^{(1)} = (\tilde{R}_{s,L})^{\epsilon_1 L}, \ L = \beta^l L_{\min}} \left[ Z_1(I) = z, \ l = l^* - \xi_1, \ \tilde{R} = \overleftarrow{R} \right] \geq \frac{\beta^{-l^*}}{\zeta} \cdot \frac{1}{2} \cdot \frac{(1 - \epsilon_1)\beta^{l^*} L_{\min} - 1 - \operatorname{len}(I)}{d} + \frac{\beta^{-(l^* - \xi_1)}}{\zeta} \cdot \frac{1}{2} \cdot \frac{\epsilon_1 \beta^{l^* - \xi_1} L_{\min} - 1 - \operatorname{len}(I)}{d} (4.3) \geq \frac{1}{2} \frac{L_{\min}}{\zeta d} - \frac{\beta^{-(l^* - \xi_1)}}{\zeta} \cdot \frac{2 \operatorname{len}(I)}{d}.$$

Similarly, we can give the exact same lower bound for each of the following four events:  $Z_1(I) = \pm z$  and  $Z_0(I) = \pm z$ .

We can now bound the probability mass that is common to the two distributions  $r^{(0)}(u)$  and  $r^{(1)}(u)$  for the event that there is a distinct interval I such that  $Z_t(I) = z$  for some  $z \in S$ :

(4.4) 
$$\sum_{z \in S} \sum_{\text{distinct } I = I(v)} \min_{t \in \{0,1\}} \left\{ \Pr_{r^{(t)}} \left[ Z_t(I) = z \right] \right\}$$
$$\geq \sum_{l=\xi_1+1}^m \sum_{v \text{ at level } l} \Pr\left[ Z_t(I(v)) \in \left\{ \pm \left\lfloor \epsilon_1 \beta^l L_{\min} \right\rfloor \right\} \right]$$

because, for each node v at level  $l^* \ge \xi_1 + 1$ , we can consider the interval I = I(v)and the displacement of  $z = z(v) = \lfloor \epsilon_1 \beta^{l^*} L_{\min} \rfloor \in S$ . Then all the events  $Z_t(I(v)) =$   $\pm z(v)$  are mutually exclusive (over the choice of such v), and hence we obtain the sum from (4.4). Furthermore, using (4.3), we obtain

(4.5)  

$$\sum_{z \in S} \sum_{\text{distinct } I = I(v)} \min_{t \in \{0,1\}} \left\{ \Pr_{r^{(t)}} \left[ Z_t(I) = z \right] \right\}$$

$$\geq \sum_{l^* = \xi_1 + 1}^m 2 \sum_{v \text{ at level } l^*} \left( \frac{1}{2} \cdot \frac{L_{\min}}{\zeta d} - \frac{\beta^{-(l^* - \xi_1)}}{\zeta} \cdot \frac{2 \operatorname{len}(I(v))}{d} \right)$$

$$\geq \sum_{l^* = \xi_1 + 1}^m n_{l^*} \frac{L_{\min}}{\zeta d} - 4 \sum_{l^* = 1}^m \sum_{v \text{ at level } l^*} \frac{\beta^{-l^*}}{\zeta} \cdot \frac{\operatorname{len}(I(v))}{d},$$

where we recall that  $n_l$  is the number of nodes at level  $l^*$ . The last inequality follows from the fact that, for each interval I = I(v) of a node v at level  $l^* \ge \xi_1 + 1$ , we can charge len(I(v)) to lengths of the intervals of the descendants of v at level  $l^* - \xi_1$ .

Finally, observe that the last term in (4.5), namely  $\sum_{l^*} \sum_{v} \frac{\beta^{-l^*}}{\zeta} \cdot \frac{\operatorname{len}(I(v))}{d}$ , is equal precisely to the probability that a random position *s* falls into an A-interval at level  $l^*$ , where  $l^*$  is chosen at random according to the distribution  $l^* = l$  with probability  $\beta^{-l}/\zeta$ . Thus we can use Claim 4.5 to bound it from above,

$$\sum_{l^*=1}^m \sum_{v \text{ at level } l^*} \frac{\beta^{-l^*}}{\zeta} \cdot \frac{\operatorname{len}(I(v))}{d} \le O(\lambda) \cdot \frac{L_{\min}}{\zeta d},$$

which together with (4.5) completes the proof of Claim 4.7.

To summarize, the total probability mass that we accounted to be common for t = 0 and t = 1 is the sum of (our lower bounds on) the probability that all entries stay put, plus the probability that exactly one distinct interval I = I(v) is displaced by precisely  $z \in S$  positions. Combining Claims 4.6 and 4.7 and using the trivial bound of  $n_l \leq \lambda$  for all  $l \in [m]$ , we obtain

$$1 - \delta_{\lambda}(u) \ge 1 - O\left(\lambda \cdot \frac{L_{\min}}{d}\right) - \sum_{l=1}^{m} n_l \cdot \frac{L_{\min}}{\zeta d} + \sum_{l=\xi_1+1}^{m} n_l \frac{L_{\min}}{\zeta d} - O\left(\lambda \frac{L_{\min}}{d}\right)$$
$$\ge 1 - O(\lambda\xi_1) \cdot \frac{L_{\min}}{\zeta d}.$$

Finally, using the fact that  $\xi_1 = O(\log \alpha)$ , we conclude (4.2), which completes the proof of Lemma 4.2.  $\Box$ 

4.2. Statistical closeness of the distributions  $\tilde{\mu}_t$  and  $\mu_t$ . We prove Lemma 3.9(b) via the following lemma.

LEMMA 4.8. For every  $t \in \{0, 1\}$ , the total variation distance between  $\tilde{\mu}_t$  and  $\mu_t$  is at most  $d^{-\Omega(1)}$ .

*Proof.* First we recall that  $\tilde{\mu}_t$  is equal to the distribution  $\mu$  conditioned on the fact that the generated pair  $(x, y) \in \mu$  be legal; i.e., both x and y are permutations, and (x, y) are far or close for t = 0 or t = 1, respectively. Since both x and y are random from  $\mathbb{Z}_p^d$ , and  $p > d^3$ , we obtain that x and y are both permutations with probability at least 1 - O(1/d).

Thus, the total variation distance between  $\tilde{\mu}_0$  and  $\mu_0$  is at most  $\Pr_{\mu_0}[\operatorname{ed}(x, y) \leq R] + O(1/d)$ . Similarly, the total variation distance between  $\tilde{\mu}_1$  and  $\mu_1$  is at most  $\Pr_{\mu_1}[\operatorname{ed}(x, y) > R/\alpha] + O(1/d)$ . Thus, it suffices to prove that  $\Pr_{\mu_0}[\operatorname{ed}(x, y) \leq R] \leq C$ 

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 $d^{-\Omega(1)}$  and  $\Pr_{\mu_1}[\operatorname{ed}(x, y) > R/\alpha] \leq d^{-\Omega(1)}$ . Remember that x is chosen at random, then  $z = x + N_{\rho}$ , and finally y is obtained from z via a sequence of rotation operations.

We choose the constant  $C_2 = 20\zeta/\epsilon_0 = 40\zeta$  and condition on the event that x, y, and z all be permutations, which happens with probability  $\geq 1 - O(1/d)$ . We can also describe the distribution  $\mu_t$  as follows. Start with a permutation z, and let xbe the permutation obtained by modifying every coordinate in z to a new symbol independently with probability  $1 - \rho$ . We may assume, without loss of generality (by renaming symbols), that z is the identity permutation of length d, i.e., for all  $i \in [d]$  we have z(i) = i, and furthermore with probability  $\rho$  we have x(i) = z(i) and x(i) = i + d otherwise. Next, let y be the permutation obtained from z by applying w random rotation operations chosen from  $\mathcal{D}_t^{\text{rot}}$ .

It will then suffice to prove the following two claims.

CLAIM 4.9. For both  $t \in \{0, 1\}$ ,

$$\Pr_{\mu_t} \left[ \operatorname{ed}(x, z) \le 2\epsilon_1 R \right] \ge 1 - e^{-d^{\Omega(1)}}.$$

CLAIM 4.10. For both  $t \in \{0, 1\}$ ,

$$\Pr_{\mu_t} \left[ 0.1 \le \frac{\operatorname{ed}(z, y)}{R \cdot C_2 \epsilon_t / \zeta} \le 10 \right] \ge 1 - d^{-\Omega(1)}.$$

We can now obtain the lemma statement from the above two claims, using a union bound and applying the triangle inequality  $|\operatorname{ed}(x,y) - \operatorname{ed}(z,y)| \leq \operatorname{ed}(x,z)$  (see also Figure 4.1). Indeed, we obtain (i) that for the distribution  $\mu_0$ , with high probability,  $\operatorname{ed}(x,y) \geq (0.1C_2\epsilon_0/\zeta - 2\epsilon_1)R = (2-2\epsilon_1)R > R$ ; and (ii) that for the distribution  $\mu_1$ , with high probability,  $\operatorname{ed}(x,y) \leq (10C_2\epsilon_1/\zeta + 2\epsilon_1)R = 402\epsilon_1R \leq \frac{804}{C_1\alpha} \cdot R < R/\alpha$ .



FIG. 4.1. The relative positions of x, y, z under the distributions  $\mu_0$  and  $\mu_1$ , respectively.

It remains to prove Claims 4.9 and 4.10.

Proof of Claim 4.9. One can verify that ed(x, z) is upper bounded by the number of substitutions performed when constructing x from z. This number of substitutions may be bounded using a straightforward Chernoff bound.

THEOREM 4.11 (Chernoff bound; cf. [27]). Let  $X_i$ , i = 1..., d, be i.i.d. random Poisson trials with  $\mathbb{E}[X_i] = q$  for some  $q \in (0, 1)$ . Then  $\Pr[|\sum X_i - qd| > \frac{1}{2}qd] \leq 2e^{-qd/10}$ .

In our case probability of substitution is  $q = (1 - \rho)(1 - 1/p)$ , where the second factor is the probability that the substituted symbol is different from the original symbol. Since  $\rho = 1 - \epsilon_1 R/d$ , we get

$$\Pr_{\mu_t}\left[\operatorname{ed}(x,z) \leq 2\epsilon_1 R\right] \geq 1 - e^{-\Omega(\epsilon_1 R)} \geq 1 - e^{-d^{\Omega(1)}}. \quad \Box$$

*Proof of Claim* 4.10. We first show an upper bound on ed(z, y) by analyzing the sum of magnitudes of all the rotation operations. Recall that there are w rotation

operations; a single rotation operation works on a block of (random) length  $L = \beta^l L_{\min}$  and incurs edit distance at most (in fact, exactly)  $2\lfloor \epsilon_l L \rfloor$ . For  $l \in [m]$ , let the random variable  $Z_l$  denote the number of rotation operations in which the block length equals  $\beta^l L_{\min}$ . Observe that  $Z_l$  has Binomial distribution  $B(w, \frac{\beta^{-l}}{\zeta})$  and its expectation is  $\mathbb{E}[Z_l] = w \cdot \frac{\beta^{-l}}{\zeta} \geq \frac{C_2 R}{m L_{\min}} \cdot \frac{d^{-0.01}}{\zeta} \geq d^{\Omega(1)}$ . By a straightforward Chernoff bound (Theorem 4.11),

$$\Pr\left[Z_l \ge 2 \mathbb{E}[Z_l]\right] \le e^{-\Omega(\mathbb{E}[Z_l])} \le e^{-d^{\Omega(1)}}$$

Taking a union bound over these events for l = 1, ..., m, we conclude that with high probability

$$\operatorname{ed}(z,y) \leq \sum_{l=1}^{m} \left( 2w \frac{\beta^{-l}}{\zeta} \cdot 2\epsilon_t \beta^l L_{\min} \right) = \frac{4C_2 \epsilon_t R}{\zeta}.$$

We proceed to show a lower bound on  $\operatorname{ed}(z, y)$  by counting inversions, i.e., pairs of symbols  $(a_1, b_1), \ldots, (a_k, b_k)$ , such that each  $a_j$  appears before  $b_j$  in z, but  $a_j$  appears after  $b_j$  in y. It is easy to verify that if the inversions are disjoint, in the sense that the symbols  $a_1, b_1, \ldots, a_k, b_k$  are all distinct, then  $\operatorname{ed}(z, y) \geq k$  (because in every alignment of z with y, for each  $j = 1, \ldots, k$ , at least one of  $a_j, b_j$  must incur an edit operation). For each of the w rotation operations we take  $\lfloor \epsilon_t L \rfloor$  pairs—simply take the  $\lfloor \epsilon_t L \rfloor$  symbols that were at the beginning of the block and match them to the  $\lfloor \epsilon_t L \rfloor$  symbols that were at the end of the block. It follows, using Chernoff bounds as above, that with probability at least  $1 - e^{-d^{\Omega(1)}}$  this process picks at least  $\frac{1}{2} \cdot \frac{C_2 \epsilon_t R}{\zeta}$  pairs of symbols, but this count might include repetitions. Furthermore, a pair "inverted" in one rotation operation may be inverted back by another rotation. To mitigate this concern, fix a pair (a, b) taken at some jth rotation operation. The probability that symbol a was inside a rotated block in at least one other rotation is at most (using the independence between rotations and a union bound)

$$(w-1)\sum_{l=1}^{m} \left(\frac{\beta^{-l}}{\zeta} \cdot \frac{\beta^{l}L_{\min}}{d}\right) < \frac{wmL_{\min}}{\zeta d} = \frac{C_{2}R}{\zeta d}.$$

A similar argument applies to symbol b, and clearly if both a and b were not inside a rotated block of any of the other w - 1 rotations, then either (a, b) or (b, a) is an inversion between z and y. It remains to apply a union bound over the  $\frac{C_2 \epsilon_t R}{2\zeta}$  pairs of symbols the above process produces, and indeed the probability that at least one of them fails is at most

$$2 \cdot \frac{C_2 \epsilon_t R}{2\zeta} \cdot \frac{C_2 R}{\zeta d} \le O\left(\frac{R^2}{d}\right) \le d^{-\Omega(1)}.$$

We conclude that with probability at least  $1 - d^{-\Omega(1)}$  the above process produces  $\frac{C_2 \epsilon_t R}{2\zeta}$  disjoint inversions, and thus  $\operatorname{ed}(y, z) \geq \frac{C_2 \epsilon_t R}{2\zeta}$ . This completes the proof of Claim 4.10.

We thus have finalized the proof of Lemma 4.8.  $\Box$ 

5. Reducing Ulam to edit distance on 0-1 strings. In this section, we prove Theorem 1.2. We make no attempt to optimize the constants.

The basic intuition behind this proof is quite simple. The first part (the upper bound on  $ed(\pi(P), \pi(Q))$ ) is immediate, and the main challenge is to prove the lower

bound on  $\operatorname{ed}(\pi(P), \pi(Q))$ . To prove the lower bound, we proceed by ruling out all "potential certificates" that  $\operatorname{ed}(\pi(P), \pi(Q))$  is small. Specifically, a "potential certificate" is a potential fixed alignment between  $\pi(P)$  and  $\pi(Q)$  of low cost, i.e., a fixed monotone mapping that matches monotonically all but at most  $\frac{1}{100} \operatorname{ed}(P, Q)$  of the positions in  $\pi(P)$  and  $\pi(Q)$ . We then analyze the probability that such an alignment is "successful," in the sense that every pair of positions that is matched under the potential alignment has equal symbols. Indeed, we show this probability is exponentially small because many of the pairs matched are independent coin tosses. We then apply a union bound over all potential alignment of small cost. Although a direct union bound is not sufficient (there are too many potential alignments to consider), we reduce the number of potential low-cost alignments by partitioning the set of all such alignments into a smaller number of groups of "equivalent alignments."

We proceed to set up some basic terminology and notation and to provide two lemmas that will be used in the proof of the theorem.

For two permutations P, Q, we say that an index (position)  $i \in [d]$  in P is missing (from Q) if the symbol P(i) does not appear inside Q.<sup>4</sup> We say that a pair of indices  $\{i, j\} \subseteq [d]$  is an inversion (in P with respect to Q) if the two characters P(i), P(j) appear in Q but in the opposite relative order than in P, formally given by  $(i-j)(Q^{-1}(P(i)) - Q^{-1}(P(j))) < 0$ . We also say that index j is inverted with respect to i.

An alignment of two strings  $x, y \in \Sigma^d$  is a mapping  $A : [d] \mapsto [d] \cup \{\bot\}$  that is monotonically increasing on  $A^{-1}([d]) = \{i \in [d] \mid A(i) \in [d]\}$ . Intuitively, A models a candidate longest common subsequence between x and y, and thus it maps indices in x to their respective indices in y and takes the value  $\bot$  when there is no respective index in y (i.e., the respective position of x is not in the candidate subsequence). A disagreement in the alignment A is an index  $i \in [d]$  for which  $A(i) \neq \bot$  and  $x(i) \neq$ y(A(i)). The alignment is called successful if it has no disagreements. The cost of an alignment is the number of positions in x (equivalently, in y) that are not mapped to a respective index in the other string, namely  $|A^{-1}(\bot)| = d - |A^{-1}([d])| = d - |A([d])|$ , where  $A([d]) = \{A(i) \mid i \in [d]\}$ . It is easy to verify that, for all x, y,

(5.1) 
$$\frac{1}{2}\operatorname{ed}(x,y) \le \min_{A} \operatorname{cost}(A) \le \operatorname{ed}(x,y),$$

where the minimum is taken over all successful alignments A.

In the following lemma, we present a property of strings P and Q that will let us prove that, for a fixed potential alignment between  $\pi(P)$  and  $\pi(Q)$ , the probability of the alignment being successful is very small.

LEMMA 5.1. Let P, Q be two permutations of length d that contain the same symbols, i.e., P([d]) = Q([d]). Then there exists a collection of  $m \ge \operatorname{ed}(P,Q)/4$  inversions  $\{i_1, j_1\}, \ldots, \{i_m, j_m\}$  such that  $i_1, j_1, \ldots, i_m, j_m$  are all distinct.

*Proof.* Fix P, Q. Define an (undirected) graph G with vertex set [d] and an edge  $\{i, j\}$  whenever  $\{i, j\}$  is an inversion. Let  $E^* \subseteq E(G)$  be a matching in G (i.e., no two edges in  $E^*$  share an endpoint) that is maximal with respect to containment. Observe that  $E^*$  is a collection of inversions whose indices are all distinct (as desired), and it remains only to bound  $m = |E^*|$  from below. Following [30], we achieve the latter using the well-known relation between maximal matching and vertex-cover.<sup>5</sup>

 $<sup>{}^{4}\</sup>mathrm{Remember}$  that we have defined a permutation P as a string with a large alphabet where every symbol appears at most once.

<sup>&</sup>lt;sup>5</sup>Another proof may be obtained using the O(1)-approximation in [17, Theorem 3.3].

Let  $V^*$  be the set of vertices incident to any edge in  $E^*$ ; thus  $|V^*| = 2|E^*|$ . Clearly,  $V^*$  is a vertex-cover of G; namely, every edge (inversion) must have at least one endpoint in  $V^*$ . It follows that  $V \setminus V^*$  contains no edges (inversions) and thus immediately yields a successful alignment A between P and Q. Formally, the subsequence of P obtained by removing the positions  $V^*$  is also a subsequence of Q, and Ais the monotone map matching them. Thus,  $A(i) = \bot$  iff  $i \in V^*$  and  $cost(A) = 2|E^*|$ . Finally, using (5.1) we get  $m = |E^*| = \frac{1}{2}cost(A) \ge \frac{1}{4}ed(P,Q)$ .

We now give a lemma that essentially lets us partition all potential alignments into a small number of groups of equivalent alignments.

LEMMA 5.2. Let P, Q be two permutations of length d. Fix  $0 < \gamma < 1/2$  and a subset  $S \subseteq [d]$ . For an alignment A of P and Q (not necessarily successful), let  $A_{|S}: S \to [d] \cup \{\bot\}$  be a function that is equal to the function A on the domain S. Define

 $F = \{A_{|S} \mid A \text{ is an alignment of } P \text{ and } Q \text{ with } \operatorname{cost}(A) \leq \gamma |S|\}.$ 

Then  $|F| \leq (3e/\gamma)^{2\gamma|S|}$ .

*Proof.* Let us denote s = |S|. An alignment of P with Q of cost at most  $\gamma s$  can be described as deleting exactly  $\gamma s$  symbols from P and exactly  $\gamma s$  symbols from Q. (We assume here for simplicity that  $\gamma s$  is an integer; otherwise, we round it up and change constants accordingly.) Clearly, we can bound |F| by the number of such alignments between P and Q, namely  $|F| \leq {d \choose \gamma s} {d \choose \gamma s}$ , but we aim to get a bound that depends on s = |S| and not on d by more carefully counting restrictions  $A_{|S}$ .

An alignment A of P with Q of cost at most  $\gamma s$  can be described as first deleting exactly  $\gamma s$  characters from P and then inserting into the resulting string exactly  $\gamma s$ characters. Observe that  $A_{|S}$  is completely determined from the following information: (a) which positions in S are deleted; (b) how many characters are deleted between every two successive indices in S; and (c) how many characters are inserted between every two successive indices in S. (When we say two successive indices in S, it should be interpreted to also include 0 and d+1 as indices in S, and in particular (b) describes also how many characters before the first index in S are deleted from P.) Indeed, for each  $i \in S$ , data (a) determines whether  $A(i) = \bot$ . If  $A(i) \neq \bot$ , then  $A(i) = i - d_i + a_i$ , where  $d_i$  is the total number of deletions among indices  $1, \ldots, i - 1$ , which can be determined from data (a) and (b), and  $a_i$  is the total number of insertions before position i, which can be determined from data (c).

It remains to upper bound the number of possible outcomes to data (a)–(c). Clearly, the outcomes for (a) and (b) together can be upper bounded by the number of outcomes of throwing  $\gamma s$  indistinguishable balls into 2s + 2 bins (a bin per element in S which may get at most one ball, a bin per each interval between elements in S, and one extra bin to account for the case when the cost is strictly less than  $\gamma s$ ). This upper bound is equal to  $\binom{2s+2+\gamma s}{\gamma s}$  possible outcomes. The outcomes of data (c) can be similarly upper bounded by  $\binom{s+1+\gamma s}{\gamma s}$ . Together, we obtain that

$$|F| \le \binom{2s+2+\gamma s}{\gamma s} \binom{s+1+\gamma s}{\gamma s} \le \left(\frac{e(2+2\gamma)}{\gamma}\right)^{\gamma s} \left(\frac{e(1+2\gamma)}{\gamma}\right)^{\gamma s} \le \left(\frac{3e}{\gamma}\right)^{2\gamma s},$$

which proves the lemma.  $\hfill \Box$ 

Having established the two lemmas, we proceed to prove the theorem, which states that with high probability  $\Omega(\operatorname{ed}(P,Q)) \leq \operatorname{ed}(\pi(P),\pi(Q)) \leq \operatorname{ed}(P,Q)$ .

Proof of Theorem 1.2. Fix two permutations P and Q of length d. The inequality  $ed(\pi(P), \pi(Q)) \leq ed(P, Q)$  follows immediately from the observation that every sequence of edit operations to transform P into Q can also be applied to transform  $\pi(P)$  and  $\pi(Q)$ . It thus remains to prove the other direction. Assume for now that P and Q use the same symbols, i.e., P([d]) = Q([d]). We will later explain how the general case follows using a similar argument.

Apply Lemma 5.1 to P, Q, and extract  $m \ge \operatorname{ed}(P, Q)/4$  inversions  $\{i_1, j_1\}, \ldots, \{i_m, j_m\}$  such that  $i_1, j_1, \ldots, i_m, j_m$  are all distinct. Define  $S = \{i_1, j_1, \ldots, i_m, j_m\}$ ; hence |S| = 2m. Fix  $\gamma = 1/100$ , and let F be defined as in Lemma 5.2 (with respect to our P, Q, S, and  $\gamma$ ). By that lemma,  $|F| \le (3e/\gamma)^{2\gamma|S|} = (3e/\gamma)^{4\gamma m}$ . Note that F does not depend on  $\pi$ .

For every  $f \in F$ , let  $\mathcal{E}_f$  be the event that all  $i \in S$  with  $f(i) \neq \bot$  satisfy  $\pi(P(i)) = \pi(Q(f(i)))$ . That is,

$$\mathcal{E}_f = \bigwedge_{i \in S \setminus f^{-1}(\bot)} \{ \pi(P(i)) = \pi(Q(f(i))) \}.$$

We claim that

(5.2) 
$$\Pr\left[\operatorname{ed}(\pi(P), \pi(Q)) < \frac{1}{2}\gamma \cdot \operatorname{ed}(P, Q)\right] \leq \Pr\left[\bigcup_{f \in F} \mathcal{E}_f\right].$$

To prove the claim we show that  $\operatorname{ed}(\pi(P), \pi(Q)) < \frac{1}{2}\gamma \cdot \operatorname{ed}(P,Q)$  implies that at least one of the events  $\mathcal{E}_f$  happens. Indeed, suppose there is a successful alignment A between  $\pi(P)$  and  $\pi(Q)$  that has  $\operatorname{cost} \frac{1}{2}\gamma \cdot \operatorname{ed}(P,Q) \leq 2\gamma m = \gamma |S|$ . Since A is successful, for all  $i \in S \setminus A^{-1}(\bot)$ , we must have  $\pi(P(i)) = \pi(Q(A(i)))$ . Furthermore, we can think of A as an alignment between P and Q, and then by definition its restriction  $A_{|S|}$  must be in F.

We now bound  $\Pr[\mathcal{E}_f]$  for any fixed  $f \in F$ ; i.e.,  $f = A_{|S}$  for some alignment A of cost at most  $\gamma|S| = 2\gamma m$ . Since S is the union of m inversions  $\{i_t, j_t\}$  with distinct indices, for at least  $(1 - 2\gamma)m$  of these inversions, we have that  $f(i_t), f(j_t) \neq \bot$ . For every such inversion  $\{i_t, j_t\}$ , it cannot be that both  $P(i_t) = Q(f(i_t))$  and  $P(j_t) = Q(f(j_t))$  (as that would contradict the fact that the alignment A is increasing). Let  $a_t \neq b_t$  denote these two differing symbols (i.e., either  $a_t = P(i_t), b_t = Q(f(i_t))$  or  $a_t = P(j_t), b_t = Q(f(j_t)))$ ; the event  $\mathcal{E}_f$  can occur only if  $\pi(a_t) = \pi(b_t)$ . We thus obtain  $(1 - 2\gamma)m$  requirements of the form  $\pi(a_t) = \pi(b_t)$ . These requirements have distinct symbols  $a_t$  in their left-hand sides (since they come from distinct positions in P), and similarly, the right-hand sides contain distinct symbols  $b_t$ . Altogether, every symbol in  $\Sigma$  may appear in at most two requirements, and thus we can extract (say, greedily) a subcollection containing at least one half of these requirements, namely, at least  $(1-2\gamma)m/2 \ge m/4$  requirements, such that every symbol appears in at most one requirement. Since  $\pi$  is a random function, the probability that all these requirements are satisfied is at most  $2^{-m/4}$ , and we conclude that  $\Pr[\mathcal{E}_f] \le 2^{-m/4}$ .

To complete the proof of the theorem, we plug the last bound into (5.2) and use a union bound and Lemma 5.2, which altogether gives

$$\Pr\left[\operatorname{ed}(\pi(P), \pi(Q)) < \frac{1}{2}\gamma \cdot \operatorname{ed}(P, Q)\right] \le \left(\frac{3e}{\gamma}\right)^{4\gamma m} \cdot 2^{-m/4} \le 2^{-m/8}.$$

Finally, we extend the proof to the case where P and Q differ on some symbols; i.e., there is at least a symbol in P that is not in Q (and vice versa). Define  $\Sigma' =$ 

 $P([d]) \cap Q([d])$  to be the set of symbols that appears in both P and Q. Let P' be the string obtained by deleting from P the symbols not in  $\Sigma'$ , and let Q' be obtained similarly from Q. Clearly, P' and Q' are permutations, they have the same length  $d' = |\Sigma'|$ , and they use exactly the same symbols. Furthermore,  $ed(P,Q) = ed(P',Q') + \Theta(d-d')$ . Applying Lemma 5.1 to P',Q', we get  $m \ge ed(P',Q')/4$  inversions  $\{i'_1,j'_1\},\ldots,\{i'_m,j'_m\}$  such that  $i'_1,j'_1,\ldots,i'_m,j'_m$  are all distinct. Translating these positions to P yields m inversions  $\{i_1,j_1\},\ldots,\{i_m,j_m\}$  between P and Q such that  $i_1,j_1,\ldots,i_m,j_m$  are all distinct. We then let S contain the indices in these inversions and also the d-d' indices in P containing the symbols not in  $\Sigma'$ . It is not difficult to see that we will still get  $|S| \ge \Omega(ed(P,Q))$ . Inversions will give rise to requirements of the form  $\pi(a) = \pi(b)$  as before, and each index i, where  $P(i) \notin \Sigma'$ , gives rise to a requirement  $\pi(P(i)) = \pi(Q(f(i)))$ . Altogether, after removing indices i such that  $f(i) = \bot$ , we still get at least |S|/8 requirements whose variables  $\pi(a), \pi(b)$  are all distinct.  $\Box$ 

### 6. Concluding remarks.

Poincaré inequality. Our communication complexity lower bounds imply that embedding the edit and Ulam metrics into  $\ell_1$ , and into powers thereof, requires distortion  $\Omega(\frac{\log d}{\log \log d})$ . But our proof also yields a Poincaré-type inequality as follows. Indeed, using (i) a variant of Lemma 3.4, where  $\Pr[\mathcal{H}^A(x) \neq \mathcal{H}^B(x)]$  is replaced with  $\mathbb{E}[\mathcal{H}^A(x) - \mathcal{H}^B(x)]^2$  and  $\mathcal{H}^A, \mathcal{H}^B$  are real (rather than boolean) functions with  $\mathbb{E}_x \left[\mathcal{H}^A(x)\right]^2 = \mathbb{E}_x \left[\mathcal{H}^B(x)\right]^2 = 1$ , together with (ii) Lemma 4.2 for suitable parameters  $R = d^{1/4}$  and  $\alpha = \Theta(\frac{\log d}{\log \log d})$ , we get that for all  $f: \mathbb{Z}_p^d \to \mathbb{R}$  (and thus all  $f: \mathbb{Z}_p^d \to \ell_2$ )

(6.1) 
$$\mathbb{E}_{(x,y)\in\mu_0} \left[f(x) - f(y)\right]^2 - \mathbb{E}_{(x,y)\in\mu_1} \left[f(x) - f(y)\right]^2 \le \frac{1}{10} \mathbb{E}_{x,y\in\mathbb{Z}_p^d} \left[f(x) - f(y)\right]^2.$$

In fact, the aforementioned nonembeddability into  $\ell_1$  (actually into the bigger space squared- $\ell_2$ ) can be proved *directly* from the Poincaré inequality (6.1) as follows. Consider a *D*-distortion embedding into squared- $\ell_2$ ; namely, let  $\phi : \mathbb{Z}_p^d \to \ell_2$  be such that, for all permutations  $x, y \in \mathbb{Z}_p^d$ ,

$$ed(x,y)/D \le \|\phi(x) - \phi(y)\|_2^2 \le ed(x,y)$$

Schoenberg [33] proved (see, e.g., [16, Theorem 9.1.1]) that, for every  $\lambda > 0$ , applying the transform  $x \mapsto 1 - e^{-\lambda x}$  on the distances of a squared- $\ell_2$  metric always results with a squared- $\ell_2$  metric. Thus, there exists a mapping  $\psi : \mathbb{Z}_p^d \to \ell_2$  satisfying

$$\|\psi(x) - \psi(y)\|_2^2 = 1 - e^{-\|\phi(x) - \phi(y)\|_2^2 \cdot \alpha/R}.$$

We thus get, using Lemma 4.8, that

$$\mathbb{E}_{\mu_0} \|\psi(x) - \psi(y)\|_2^2 - \mathbb{E}_{\mu_1} \|\psi(x) - \psi(y)\|_2^2 - \frac{1}{10} \cdot \mathbb{E}_{x,y \in \mathbb{Z}_p^d} \|\psi(x) - \psi(y)\|_2^2 \\ \geq \frac{1}{e} - \frac{1}{e^{\alpha/p}} - \frac{1}{10} - d^{-\Omega(1)}.$$

Combining this inequality with (6.1) implies that  $D \ge \Omega(\alpha) = \Omega(\frac{\log d}{\log \log d})$ .

Overcoming nonembeddability into  $\ell_1$ . One moral of our nonembeddability lower bound is that some of the usual approaches for designing algorithms for Ulam and edit metrics cannot give approximation better than  $\tilde{\Omega}(\log d)$ . In particular, to design

for these metrics a nearest neighbor algorithm that achieves a sublogarithmic (in d) approximation factor, we must depart from embedding into  $\ell_1$  or into other spaces that admit O(1)-size sketches.

This challenge was accomplished for the Ulam metric in a recent paper [3], which designs a nearest neighbor algorithm achieving  $O(\log \log d)$  approximation using polynomial (in n + d) space and sublinear (in n) query time. These guarantees bypass the  $\ell_1$  nonembeddability barrier proved in the current paper by relying on a constant-distortion embedding of the Ulam metric into an alternative, richer host space (namely, iterated product of simple spaces such as  $\ell_1$ ) which, despite the richer structure, turns out to have reasonably good algorithms.

Sketching complexity. After the preliminary version of this paper appeared in 2007, there has been further progress on the communication complexity of estimating the Ulam and edit distances. First, [3] showed that the Ulam metric can be sketched in  $(\log d)^{O(1)}$  space with constant approximation. Second, [4] showed that the Ulam metric requires  $\Omega(\frac{\log d/\log \log d}{\alpha})$  communication complexity, for approximation  $\alpha > 1$ , which improves our lower bound of  $\Omega(\log(\frac{\log d}{\alpha \log \alpha}))$  from Theorem 1.1. In particular, for a constant approximation, this significantly improves the communication lower bound from  $\Omega(\log \log d)$  to  $\Omega(\frac{\log d}{\log \log d})$ . We note that the lower bound of [4] is nonetheless based on Theorem 1.1 in the setting where approximation  $\alpha = \Theta(\frac{\log d}{\log \log d})$ . This lower bound of [4] for the Ulam metric extends to edit distance over binary strings by applying Theorem 1.2.

The two aforementioned results together imply that a (sufficiently large) constant approximation of the Ulam distance has sketching complexity  $(\log d)^{\Theta(1)}$ .

Acknowledgments. We thank Parikshit Gopalan, Piotr Indyk, T.S. Jayram, Ravi Kumar, Ilan Newman, and Yuri Rabinovich for numerous early discussions on nonembeddability of the Ulam and edit metrics, which undoubtedly were a precursor of the current work. We also thank James Lee and Assaf Naor for enlightening discussions about the Poincaré inequality. Finally, we thank the reviewers for careful reading and suggestions that helped improve the clarity and exposition of the article.

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