Statistical Inference for Entropy

Karina Marks

March 26, 2017

Contents

Chapter 1

Introduction

1.1 Entropy

Entropy $H(S)$, can be thought of as a representation of the average information content of an observation; sometimes referred to as a measure of unpredictability or disorder.

"$H(S)$ is the quantity of surprise you should feel upon reading the result of a measurement" (Fraser and Swinney, 1986) [10]. Thus the "entropy of S can be seen as the uncertainty of S" [15].

1.1.1 Shannon Entropy

The Shannon entropy of a random vector $X \in \mathbb{R}^d$, from a continuous distribution, with density function f is given by;

$$H = - \int_{x \in \mathbb{R}^d} f(x)log(f(x))dx \tag{1.1}$$

(This is often referred to as the differential entropy.)

If a random vector X has a discrete distribution with probability mass function $f(x)$, $x \in \mathbb{Z}^d$, then we have the Shannon entropy ;

$$H = - \sum_{x \in \mathbb{Z}^d} f(x)log(f(x))$$

1.1.2 Rényi and Tsallis Entropy

The Rényi and Tsallis entropies are for the order $q \neq 1$ and the construction of them relies upon the generalisation of the Shannon entropy, equation 1.1. For a random vector $X \in \mathbb{R}^d$, from a continuous distribution with density function f, we define;

Rényi entropy

$$H_q^* = \frac{1}{1-q} log \left(\int_{\mathbb{R}^d} f^q(x) dx \right) \qquad (q \neq 1) \tag{1.2}$$

Tsallis entropy

$$H_q = \frac{1}{q-1} \left(1 - \int_{\mathbb{R}^d} f^q(x) dx \right) \qquad (q \neq 1) \tag{1.3}$$

Moreover, if we consider a random vector X from a discrete distribution, with probability mass function $f(x)$ for $x \in \mathbb{Z}^d$, then for $q \neq 1$, the Rényi and Tsallis entropies are defined as follows;

$$H_q^* = \frac{1}{1-q} log \left(\sum_{x \in \mathbb{Z}^d} f^q(x) \right)$$

$$H_q = \frac{1}{q-1} \left(1 - \sum_{x \in \mathbb{Z}^d} f^q(x) \right)$$

When the order of the entropy $q \to 1$, both the Rényi, (1.2), and Tsallis, (1.3), entropies tend to the Shannon entropy, (1.1), as this is a special case for when $q = 1$.

1.2 Background

1.2.1 Properties of Entropy

I will begin by exploring properties specific to the Shannon entropy; and then progress to those for other types of entropy. Kapur and Kesavan's book on *Entropy Optimization Principles with Applications* [18], gives an account of some properties of the discrete Shannon entropy H. Whilst Johnson [17], discusses properties for the continuous Shannon entropy - the differential entropy, and

compares them to the discrete properties in his book on *Information Theory and The Central Limit Theorem.*

The Shannon entropy of a discrete random variable is always non-negative $H \geq 0$, and is strictly positive unless if f is any of the N degenerate distributions (X is deterministic). That is, the only case for which $H = 0$ is when $f(x_i) = 1$ if $i = k$, $k \in [1, N]$ and $f(x_i) = 0$ otherwise, for a discrete distribution. However, for the continuous Shannon entropy, H can be either positive of negative. This is obvious to see if we consider a random variable Y with the continuous uniform distribution on $[0, c]$, then;

$$H = -\int_0^c \frac{1}{c} \log\left(\frac{1}{c}\right) dx = \log c$$

thus, depending on the value of c we can obviously have either a positive or negative entropy.

Moreover, for a discrete random variable X, the Shannon entropy is both shift and scale invariant, which means for all $a \neq 0$ and $b \in \mathbb{R}$ we have;

$$H(aX + b) = H(X) \tag{1.4}$$

On the contrary, for a continuous random variable Y, only the shift invariance holds, thus for all $a \neq 0$ and $b \in \mathbb{R}$ we have the following relationship;

$$H(aY + b) = H(Y) + \log a \tag{1.5}$$

The maximum value of the continuous entropy H is attained with different distributions, dependent on a number of factors. The maximising distribution f depends on how it is supported, and what conditions hold on the moments of X. For example, when $supp\{f\} = \mathbb{R}$, for a fixed variance σ^2, the maximum of H is attained when f is from the Normal/Gaussian distribution, in comparison to any other distribution with the same σ^2 [18].

Additionally, for the Shannon entropy, there is an interesting result on an upper bound of H, one which depends on the distribution of X, which can be used to show if the entropy of a sample X is finite. Consider a random variable X, then for $s > 0$ we have;

$$H(X) \leq \frac{1}{s} \log\left(\frac{2^s e \Gamma^s(\frac{1}{s}) \mathbb{E}(|X|^s)}{s^{s-1}}\right) \tag{1.6}$$

This indicates that if the s-th moment of X if finite, $\mathbb{E}(|X|^s) < \infty$, for some $s > 0$ then the Shannon entropy of X will also be finite, $H(X) < \infty$ [**?**].

Shannon entropy, as mentioned earlier, is a special case of the Rényi and Tsallis entropies as $q \to 1$. Thus, for a general q-entropy, other properties hold. Namely, H_q is concave when $q > 0$ (and convex when $q < 0$), implying for the Shannon entropy, when $q = 1$, that H is concave.

There are also other special cases that have certain properties; for example the Rényi entropy with $q = 2$ is known as the quadratic Rényi entropy, which is defined as follows for a continuous random variable X with density function f;

$$H_2^* = -log \left(\int_{\mathbb{R}^d} f^2(x) dx \right) \tag{1.7}$$

Moreover, another special case is considering the Rényi entropy as $q \to \infty$, if the limit exists, is defined as the minimum entropy, since it's the smallest possible value of H_q^*;

$$H_\infty^* = -\log \sup_{x \in \mathbb{R}^d} f(x)$$

Furthermore, there are some interesting relationships between the different specific types of entropy, for example Leonenko and Seleznjev [24] show the following relationship between H_2^* and H_∞^*;

$$H_\infty^* \leq H_2^* \leq 2H_\infty^* \tag{1.8}$$

Additionally, they show an approximate relationship between the Shannon entropy, H, and the quadratic Rényi entropy, H_2^* ;

$$H_2^* \leq H \leq \log(d) + \frac{1}{d} - e^{-H_2^*}$$

where d is the dimension of the distribution.

1.2.2   Applications of Entropy

Entropy began as a concept in thermodynamics, about the idea of that within any irreversible system, a small amount of heat energy is aways lost. Entropy has more recently found application in the field of information theory, where it describes a similar loss, this time of missing information or data in systems of information transmission. Thus, entropy has many applications across both these areas.

I will be concentrating on Shannon entropy - also mentioning Rényi and Tsallis entropies - which concern information theory; therefore I will consider

applications accordingly. I will give a short overview of some of its applications; however, this is not an exhaustive list, since the application of entropy are extensive.

The estimation of Shannon entropy is useful in various science/engineering applications, such as independent component analysis, image analysis, genetic analysis, speech recognition, manifold learning, and time delay estimation.

Independent component analysis (ICA), in signal processing, is a computational method for decomposing large, often very complex, multivariate data to find underlying/hidden factors or components. The computation of ICA depends on knowing the entropy of the sample; and in most cases this must be estimated, as an exact entropy is not always known. Kraskov, Stögbauer and Grassberger [19] discussed how estimating the mutual information (MI) using entropy estimators is useful for assessing the independence of components from ICA. Learned-Miller and Fisher [21] also presented another example of how to use estimation of entropy to obtain a new algorithm for the ICA problem.

Image analysis is the investigation of an image and the extraction of useful information. Hero and Michel [14] first discuss the applications of Rényi entropy in image processing, then Neemuchwala, Hero and Carson [25] discuss how in image analysis an important task is that of image retrieval, which uses entropy estimation to compute entropic similarities that are used to match a reference image to another image. Moreover Du, Wang, Guo and Thouin, [8] considered the importance of entropy-based image thresholding; using both Shannon and relative entropy.

Genetic analysis is the study and research of genes and molecules to find information on biological systems. Statistical analysis of specific cells can help us understand how genomic entropy can help diagnose diseases and cancers. Wieringen and Vaart, [29] discuss how chromosomal disorganisation increases as cancer progresses, they mention how the K-L estimator can be used to help find this disorganisation/entropy; thus finding that "as cancer evolves, and the genomic entropy increases, the transcriptomic entropy is also expected to surge".

Speech recognition (SR) is the inter-disciplinary sub-field of computational linguistics that develops methodologies and technologies that enables the recognition and translation of spoken language into text by computers. Shen, Hung and Lee [27] discuss how an entropy based algorithm can conduct accurate SR in noisy environments. Moreover, Kuo and Gao [20] focus on a method where

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the probability of a state or word sequence given an observation sequence is computed directly from the maximum entropy direct model.

It is also important to note the statistical applications of entropy; there are some tests on goodness-of-fit established by the estimation of entropy. Vasicek explored the test for normality; that its entropy exceeds that of any other distributions with the same variance [30]. Dudewicz and van der Meulen [9] discussed the property mentioned in section 1.2.1, that the uniform distribution maximises the entropy. Moreover, others have explored different distributions and their entropic properties; see [11, 5]

1.2.3   Other Estimators of Entropy

There are several estimation methods for the nonparametric estimation of the Shannon entropy of a continuous random sample. The paper *Nonparametric Entropy Estimation: An Overview* (J.Beirlant, E.Dudewicz, L.Gyorfi, E.van der Muelen, 2001) [3], gives an overview of the properties of these various methods. Also, the paper *Causality detection based on information-theoretic approaches in time series analysis* (K.Hlaváčková, M.Paluš, M.Vejmelka, and J.Bhattacharya, 2007) gives a more detailed look into these different types of estimators. I will outline a summary below to the types of estimators, which will lead us to understand why we choose the Kozachenko-Leonenko estimator for entropy.

First, I must set out the types of consistency, so we can see more obviously how it compares to the K-L estimator, for $X_1, ..., X_N$ a i.i.d sample from the distribution $f(X)$, where H_N is the estimator of $H(f)$. Then we have (as $N \to \infty$);

- Asymptotic Unbias

$$\mathbb{E}(H_N) - H(f) \to 0 \tag{1.9}$$

- Weak Consistency

$$H_N \xrightarrow{p} H(f) \tag{1.10}$$

- Mean Square Consistency

$$\mathbb{E}\{(H_N - H(f))^2\} \to 0 \tag{1.11}$$

- Strong Consistency (Asymptotic Normality)

$$\sqrt{N}(\hat{H}_{N,k} - H) \xrightarrow{d} N(0, \sigma^2) \qquad (1.12)$$

$$\mathbb{E}(\hat{H}_{N,k} - H)^2 = \frac{\sigma^2}{N}\left(1 + O\left(\frac{1}{N}\right)\right) \qquad (1.13)$$

This is the consistency shown with the K-L estimator in Theorem 2.

The types of nonparametric estimators can be split into 3 categories; plug-in estimates, estimates based on sample-spacings and estimates based on nearest neighbour distances. The latter is the Kozachenko-Leonenko estimator, which is the main focus of this paper and will be explored in more detail in section **??**.

The plug-in estimates [3], [15] are based upon a consistent density estimate f_N, of density f, which depends on the sample $X_1, ..., X_N$, I will consider two of these; the most obvious estimator of this type if the integral estimate of entropy. Given by;

$$H_N = -\int_{A_N} f_N(x)log(f_N(x))dx \qquad (1.14)$$

where the set A_N excludes the tail values of f_N. When the sample is from a 1-dimensional distribution, Dmitriev and Tarasenko, [7] for $A_N = [-b_N, b_N]$ and f_N the kernel density estimator; proved a strong consistency for this estimator. However, if f_N is not estimated in this form, due to the numeric integration, for dimensions $d \geq 2$, Joe [16] points out that this estimator is not practical and thus proposed the next plug-in estimator for entropy - the resubstitution estimator.

The resubstitution estimate is of the form;

$$H_N = -\frac{1}{N}\sum_{i=1}^{N} log(f_N(X_i))dx \qquad (1.15)$$

which was first proposed in 1976, by Ahmad and Lin [2] who showed the mean-square consistency of this estimator, where f_N is a kernel density estimate. Joe [16] then went on to obtain the asymptotic bias and variance, and whilst satisfying certain conditions reduced the mean square error. Moreover, Hall and Morton [13] went on to say that under more restrictive conditions we have strong consistency for 1-dimensional distributions; however, when $d = 2$ the root-n consistent estimator will have significant bias.

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There are also two more plug-in estimates discussed in this paper; the splitting data and cross-validation estimates. Where in the first estimator, strong consistency is shown for a general dimension d, under some conditions on f. And in the latter estimator, strong consistency holds for a kernel estimate of f and for other estimates of f under some conditions we have root-n consistency when $1 \leq d \leq 3$.

Hence, so far the estimates for entropy looked at are only consistent whilst under strong conditions on f and f_N and mostly for a 1-dimensional distribution. So it is important to look at the next category of estimates - estimates of entropy based on sample-spacings; namely the m-spacing estimate. Sometimes it in not practical to estimate f_N, so this estimate is found based on spacings between the sample observations.

This estimator is only defined for samples of 1-dimension, where we assume $X_1, ..., X_N$ are an i.i.d sample, and let $X_{N,1} \leq X_{N,2} \leq ... \leq X_{N,N}$ be the corresponding ordered sample, then $X_{N,i+m} - X_{N,m}$ is the m-spacing.

Firstly we look at this estimator of the form, with fixed m;

$$H_{m,N} = \frac{1}{N} \sum_{i=1}^{N-m} log \left(\frac{N}{m} (X_{N,i+m} - X_{n,i}) \right) - \Psi(m) + log(m) \qquad (1.16)$$

where $\Psi(x)$ is the digamma function - more detailed explanation in section ??. For a sample from a uniform distribution this estimator has been shown to be consistent; proved by Tarasenko [28]. Under some conditions on f, on its boundedness, the weak consistency and asymptotic normality was shown by Hall [12].

To decrease the asymptotic variance of the estimator, we consider the estimator when $m_N \to \infty$, which is defined slightly differently;

$$H_{m,N} = \frac{1}{N} \sum_{i=1}^{N-m_N} log \left(\frac{N}{m_N} (X_{N,i+m_N} - X_{n,i}) \right) \qquad (1.17)$$

for this estimator the weak and strong consistencies are proved under the assumption that as $N \to \infty$, $m_N \to \infty$ and $\frac{m_N}{N} \to 0$, for densities with bounded support.

The last category of estimators discussed by Beirlant, Dudewicz, Gyorfi and Muelen are those based on nearest neighbour distances. The main focus of my paper is on the Kozachenko-Leonenko estimator for entropy; which is the estimator covered in this section of their paper. I will not go into detail for

this estimator now; however, I will mention that strong consistency holds for dimension $d \leq 3$, but higher dimensions can cause problems. Henceforth, it is important to note that recently a new estimator has been proposed by Berrerrt, Samworth and Yuan [4], formed as a weighted average of k-nearest neighbour estimators for different values of k. This estimator has shown promising results in higher dimensions, where under the same assumptions as for the K-L estimator, the strong consistency condition holds.

Chapter 2

Kozachenko-Leonenko Estimator

2.1 History

This estimator was first introduced by L.Kozachenko and N.Leonenko, in 1987, where they first published the article *Sample Estimate of the Entropy of a Random Vector*, in the paper *Problems of Information Transmission*. Using the nearest neighbour method, they created a simple estimator for the Shannon entropy of an absolutely continuous random vector from a independent sample of observations, to then establish conditions under which we have asymptotic unbiasedness and consistency.

Since then, there has been major developments in the estimator; firstly in 2007, N.Leonenko, L.Pronzato, V.Savani, proposed a similar alternative to this estimator in their paper *a Class of Renyi Information Estimators for Mulitdimensional densities*, this time using the k-nearest neighbour method, to consider estimators for the Rényi and Tsallis entropies. Then as the order of these entropies $q \to 1$, they defined the k-nearest neighbour estimator for the Shannon entropy, where k is fixed, and these estimators (under less rigorous conditions) are both consistent and asymptotically unbiased.

Moreover, in 2016, a new idea was proposed by T.Berrett, R.Samsworth and M.Yuan, written in *Efficient Mulitvariate Entropy Estimation via k-Nearest Neighbour Distances*; that the value chosen for k, depends upon the sample size

N. Also, this idea is then extended to a new estimator; "formed as a weighted average of Kozachenko-Leonenko estimators for different values of k". I will not be exploring this new estimator in depth; however, the understanding of the value of k depending on N will be examined in detail. Additionally, for $d = 1, 2, 3$, under some conditions on k, we are shown that the bias of the estimator acts in terms of $N^{-\frac{2}{d}}$; something which will also later be explored.

Lastly, also in 2016, S.Delattre and N.Fournier wrote the paper; *On the Kozachenko-Leonenko Entropy Estimator*, where they studied in detail the bias and variance of this estimator considering all 3 proposed values of k - $k = 1$, k fixed or k depends on N. The also provided a development for the bias of this estimator when $k = 1$, in dimensions $d = 1, 2, 3$, in terms of $O(N^{-\frac{1}{2}})$, and in higher dimensions, in terms of powers of $N^{\frac{-2}{d}}$. This is an idea that will be considered in the focus of this paper; for $d = 1, 2$ to show how the bias acts for large N when $k = 1$.

2.1.1 Estimator with k=1

Firstly, I considered an article *On Statistical Estimation of Entropy of Random Vector* (N.Leonenko and L.Kozachenko, 1987) [22], which considers estimating the Shannon entropy of an absolutely continuous random sample of independent observations, with unknown probability density $f(x), x \in \mathbb{R}^d$. As $f(x)$ is unknown this is not easily estimated accurately for a random sample, and by just estimating the density $\hat{f}(x)$ to replace the actual density $f(x)$ in the formula for the entropy we get highly restrictive consistency conditions.

Therefore, the following estimator was proposed for the Shannon entropy of a random sample $X_1, X_2, ..., X_N$ of d-dimensional observations;

$$H_N = d\log(\bar{\rho}) + \log(c(d)) + \log(\gamma) + \log(N-1) \tag{2.1}$$

where $c(d) = \frac{\pi^{\frac{d}{2}}}{\Gamma(\frac{d}{2}+1)}$ is the volume of the d-dimensional unit ball, the Euler constant is $\log(\gamma) = \exp\left[-\int_0^\infty e^{-t}\log(t)dt\right] = -\Psi(1)$ and $\bar{\rho} = \left[\prod_{i=1}^N \rho_i\right]^{\frac{1}{N}}$, with ρ_i the nearest neighbour distance from X_i to another member of the sample $X_j, i \neq j$.

It is important to note that one can write the Euler constant $-\Psi(1) = \log(\exp(-\Psi(1))) = \log(\frac{1}{\exp(\Psi(1))})$, this notation is what is used in the latter papers, so it is useful to introduce it here. $\Psi(x)$ is the Digamma function, and

when $x = 1$, this is just the negative Euler constant. Thus this estimator can be written if the form;

$$H_N = \log(\bar{\rho}^d) + \log(c(d)) - \Psi(1) + \log(N-1)$$

$$= \log\left(\left[\prod_{i=1}^{N} \rho_i\right]^{\frac{d}{N}}\right) \log(c(d)(N-1)) + \log\left(\frac{1}{\exp(\Psi(1))}\right)$$

$$= \frac{1}{N}\sum_{i=1}^{N}\log(\rho_i^d) + \log\left(\frac{c(d)(N-1)}{\exp(\Psi(1))}\right)$$

$$= \frac{1}{N}\sum_{i=1}^{N}\log(\rho_i^d) + \frac{1}{N}\sum_{i=1}^{N}\log\left(\frac{c(d)(N-1)}{\exp(\Psi(1))}\right)$$

$$= \frac{1}{N}\sum_{i=1}^{N}\log\left(\frac{\rho_i^d c(d)(N-1)}{\exp(\Psi(1))}\right) \tag{2.2}$$

Under some conditions on the density function, this estimator is asymptotically unbiased and under stronger conditions it is also a consistent estimator for the Shannon entropy.

The estimator here is in a simple form, which is later developed into something more sophisticated, using the nearest neighbour method, but considering larger values of k (here $k = 1$). This estimator is developed so that the consistency and asymptotic unbias of the estimator holds under less constrained conditions.

2.1.2 Estimator with k fixed

The next paper I am exploring on estimation is *a Class of Renyi Information Estimators for Multidimensional densities* (N.Leonenko, L.Pronzato, V.Savani, 2007) [23], which looks at estimating the Rényi (H_q^*) and Tsallis (H_q) entropies, when $q \neq 1$, and the Shannon ($\hat{H}_{N,k,1}$) entropy. Where these are taken for a random vector $X \in \mathbb{R}^d$ with density function $f(x)$, by using the kth nearest neighbour method, with a fixed values of k.

For the Rényi and Tsallis entropies, this is achieved by considering the integral $I_q = \int_{\mathbb{R}^d} f^q(x)dx$, and generating its estimator, which is defined as $\hat{I}_{N,k,q} = \frac{1}{N}\sum_{i=1}^{N}(\zeta_{N,k,q})^{1-q}$. Where, $\zeta_{N,k,q} = (N-1)C_k V_d(\rho_{k,N-1}^{(i)})^d$, $V_d = \frac{\pi^{\frac{d}{2}}}{\Gamma(\frac{d}{2}+1)}$ is the volume of d-dimensional unit ball, $C_k = \left[\frac{\Gamma(k)}{\Gamma(k+1-q)}\right]^{\frac{1}{1-q}}$ and $\rho_{k,N-1}^{(i)}$ is the kth nearest neighbour distance from the observation X_i to some other X_j.

The estimator $\hat{I}_{N,k,q}$, provided $q > 1$ and I_q exists - and for any $q \in (1, k+1)$ if f is bounded - is thus found to be an asymptotically unbiased estimator for I_q. Also, provided $q > 1$ and I_{2q-1} exists - and for any $q \in (1, \frac{k+1}{2})$, when $k \geq 2$ if f is bounded - $\hat{I}_{N,k,q}$ is thus a consistent estimator for I_q.

Moreover, by simple formulas both the Rényi and Tsallis entropies can be written in terms of this estimated value;

$$\hat{H}_q^* = \frac{1}{1-q} log(\hat{I}_{N,k,q}) \tag{2.3}$$

$$\hat{H}_q = \frac{1}{q-1}(1 - \hat{I}_{N,k,q}) \tag{2.4}$$

thus, under the latter conditions, provide consistent estimates of these entropies as $N \to \infty$ for $q > 1$.

Furthermore, this paper goes on to discuss an estimator for the Shannon entropy, H_1 by taking the limit of the estimator for the Tsallis entropy, $\hat{H}_{N,k,q}$ as $q \to 1$, again with a fixed value of k. This estimator is similar to that proposed in 1987, equation 2.2; however, it is now extended from the nearest neighbour to the kth nearest neighbour;

$$\hat{H}_{N,k,1} = \frac{1}{N} \sum_{i=1}^{N} log(\xi_{N,i,k}) \tag{2.5}$$

where $\xi_{N,i,k} = (N-1)\exp[-\Psi(k)]V_d(\rho_{k,N-1}^{(i)})^d$, with V_d and $\rho_{k,N-1}^{(i)}$ defined as in the estimation of I_q and the digamma function $\Psi(z) = \frac{\Gamma'(z)}{\Gamma(z)}$. The digamma function at $k = 1$ is given by $\Psi(1) = -log(\gamma)$, the Euler constant, which was used for the $k = 1$ version of this estimator. Under the following less restrictive conditions; f is bounded and I_{q_1} exists for some $q_1 > 1$; then H_1 exists and the estimator $\hat{H}_{N,k,1}$ is a consistent estimator for the Shannon entropy. This means that for large N, we have $\hat{H}_{N,k,1} \overset{L_2}{\to} H$; which implies that as $N \to \infty$, both $N^{\frac{1}{2}}(\hat{H}_{N,k,1} - H) \overset{d}{\to} N(0, \sigma^2)$ - it's asymptotically efficient - and $\mathbb{E}(\hat{H}_{N,k,1}) \to H$ - it's asymptotically unbiased.

2.1.3 Estimator with k dependent on N

The last main paper, whose results I will be exploring is *Efficient Multivariate Entropy Estimation via k-Nearest Neighbour Distances* (T.Berrett, R.Samworth, M.Yuan, 2016) [4], which initially studies the K-L estimator, and the conditions under which it is efficient and asymptotically unbiased (for a value of k depending on the sample size N).

Considering dimensions $d \leq 3$, and a sample size N from distribution with density $f(x)$, they defined the k-nearest neighbour estimator of entropy - just as in section 2.1.2 - to be;

$$\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^{N} log \left[\frac{\rho_{(k),i}^{d} V_d (N-1)}{e^{\Psi(k)}} \right] \tag{2.6}$$

where $\rho_{(k),i}$, V_d and $\Psi(k)$ are all defined as in the 2007 paper. However, the difference here is in the conditions under which the estimator is consistent and asymptotically unbiased.

Here, some conditions on the finiteness of the α moment of f and the continuity and differentaibility of f are proposed, with $k \in \{1, ..., O(N^{1-\epsilon})\}$, for some $\epsilon > 0$, we have asymptotic unbias of the estimator; where the bias can be expressed as;

$$\mathbb{E}(\hat{H}_N) - H = O\left(max \left\{ \frac{k^{\frac{\alpha}{\alpha+d}-\epsilon}}{N^{\frac{\alpha}{\alpha+d}-\epsilon}}, \frac{k^{\frac{\beta}{d}}}{N^{\frac{\beta}{d}}} \right\} \right) \qquad N \to \infty \tag{2.7}$$

Also, they considered the asymptotic normality of the estimator, given the α moment of f is finite (for $\alpha > d$), and some conditions on the continuity and differentiability of f hold and with $k \in \{k_0, ..., k_1\}$. Then the variance of the estimator is given by;

$$Var(\hat{H}_{N,k}) = \frac{\sigma^2}{N} + o(\frac{1}{N}) \tag{2.8}$$

as $N \to \infty$, where $\sigma^2 = Var(log(f(x)))$, and we define k_0, k_1 such that $\frac{k_0}{log^5(N)} \to \infty$ and $k_1 = O(N^\tau)$, where $\tau < min \left\{ \frac{2\alpha}{5\alpha+3d}, \frac{\alpha-d}{2\alpha}, \frac{4}{4+3d} \right\}$.

Moreover, T.Berrett, R.Samsworth and M.Yuan also go on to show that a consequence of the variance, given the dimension of the sample $d \leq 3$, with the same conditions, we have the asymptotic normality;

$$\sqrt{N}(\hat{H}_{N,k} - H) \xrightarrow{d} N(0, \sigma^2) \tag{2.9}$$

and

$$N\mathbb{E}(\hat{H}_{N,k} - H)^2 \to \sigma^2 \tag{2.10}$$

where the estimator is asymptotically efficient and the asymptotic variance here is the best possible.

It is important to note that for higher dimensions $(d > 3)$, these results do not necessarily hold; since I am just considering the specific dimensions $d = 1$

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and $d = 2$, there is no need to detail this. However, they do then go on to discuss a more appropriate estimator for higher dimensions, given sufficient smoothness, which is efficient in arbitrary dimensions, which was previously mentioned in section 1.2.3 - Other Estimators of Entropy.

2.2 Focus of this Paper

I now wish to more explicitly introduce the Kozachenko-Leonenko estimator of the entropy H, in the form that I will be considering. Let $X_1, X_2, ..., X_N, N \geq 1$ be independent and identically distributed random vectors in $\mathbb{R}^d$, and denote $\|.\|$ the Euclidean norm on $\mathbb{R}^d$.

- For $i = 1, 2, ..., N$, let $X_{(1),i}, X_{(2),i}, .., X_{(N-1),i}$ denote an order of the X_k for $k = \{1, 2, ..., N\} \setminus \{i\}$, such that $\|X_{(1),i} - X_i\| \leq \cdots \leq \|X_{(N-1),i} - X_i\|$. Let the metric ρ, defined as;

$$\rho_{(k),i} = \|X_{(k),i} - X_i\| \tag{2.11}$$

  denote the kth nearest neighbour or X_i.

- For dimension d, the volume of the unit d-dimensional Euclidean ball is defined as;

$$V_d = \frac{\pi^{\frac{d}{2}}}{\Gamma(1 + \frac{d}{2})} \tag{2.12}$$

  where the Γ function is defined as;

$$\Gamma(m) = \int_0^\infty x^{m-1} e^{-x} dx \tag{2.13}$$

- For the kth nearest neighbour, the digamma function is defined as;

$$\Psi(k) = -\gamma + \sum_{j=1}^{k-1} \frac{1}{j} \tag{2.14}$$

  where $\gamma = 0.577216$ is the Euler-Mascheroni constant (where the digamma function is chosen so that $\frac{e^{\Psi(k)}}{k} \to 1$ as $k \to \infty$).

Then the Kozachenko-Leonenko estimator for entropy, H, is given by;

$$\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^N log \left[\frac{\rho_{(k),i}^d V_d (N-1)}{e^{\Psi(k)}} \right] \tag{2.15}$$

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where, $\rho^d_{(k),i}$ is defined in (2.11), V_d is defined in (2.12) and $\Psi(k)$ is defined in (2.14).

This paper focuses only on distributions for $d \leq 3$, more specifically, I will first be considering samples from 1-dimensional distributions, $d = 1$. Therefore, the volume of the 1-dimensional Euclidean ball is given by $V_1 = \frac{\pi^{\frac{1}{2}}}{\Gamma(\frac{3}{2})} = \frac{\sqrt{\pi}}{\frac{\sqrt{\pi}}{2}} = 2$. Hence the Kozachenko-Leonenko estimator is of the form;

$$\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^{N} log \left[\frac{2\rho_{(k),i}(N-1)}{e^{\Psi(k)}} \right] \tag{2.16}$$

Later, I will be considering samples from 2-dimensional distributions; thus, $d = 2$ and the volume of the 2-dimensional Euclidean ball is given by $V_2 = \frac{\pi^{\frac{2}{2}}}{\Gamma(2)} = \frac{\pi}{1} = \pi$. Hence, the estimator takes the form;

$$\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^{N} log \left[\frac{\pi \rho^2_{(k),i}(N-1)}{e^{\Psi(k)}} \right] \tag{2.17}$$

I will be looking at the asymptotic bias and variance of the estimator for different values of k, the main theorems I will be working by are those from section 2.1.3, where we have the conditions 1, 2 and 3, which imply the results stated by Theorems 1 and 2.

(NB: these conditions and theorems have been tweaked slightly to only explicitly consider distributions of dimension $d = 1, 2$, since the only distributions being considered in this paper are of dimension 1 or 2)

Condition 1 *(β) For density f bounded, denoting $m := \lfloor \beta \rfloor$ and $\eta := \beta - m$, we have that f is m times continuously differentiable and there exists $r_* > 0$ and a Borel measurable function g_* such that for each $t = 1, 2, ..., m$ and $\|y - x\| \leq r_*$, we have;*

$$\|f^{(t)}(x)\| \leq g_*(x)f(x)$$

,

$$\|f^{(m)}(y) - f^{(m)}(x)\| \leq g_*(x)f(x)\|y - x\|^{\eta}$$

and $sup_{x:f(x) \geq \delta} g_*(x) = o(\delta^{-\epsilon})$ as $\delta \downarrow 0$, for each $\epsilon > 0$.

Condition 2 *(α) For density $f(x)$ and dimension d, we have;*

$$\int_{\mathbb{R}^d} \|x\|^{\alpha} f(x)dx < \infty$$

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Condition 3 *Assume that condition 1 holds for $\beta = 2$ and condition 2 holds for some $\alpha > d$. Let $k_0^* = k_{0,N}^*$ and $k_1^* = k_{1,N}^*$ denote two deterministic sequences of positive integers with $k_0^* \leq k_1^*$, with $\frac{k_0^*}{\log^5 N} \to \infty$ and with $k_1^* = O(N^\tau)$, where*

$$\tau < \min\left\{\frac{2\alpha}{5\alpha + 3d}, \frac{\alpha - d}{2\alpha}, \frac{4}{4 + 3d}\right\}$$

Theorem 1 (Asymptotic Unbiasedness) *Assume that conditions 1 and 2 hold for some $\beta, \alpha > 0$. Let $k^* = k_N^*$ denote a deterministic sequence of positive integers with $k^* = O(N^{1-\epsilon})$ as $N \to \infty$ for some $\epsilon > 0$. Then, for $d \leq 2$ (or $d \geq 3$) with $\beta \leq 2$ (or $\alpha \in (0, \frac{2d}{d-2})$), then for every $\epsilon > 0$ we have;*

$$\mathbb{E}(\hat{H}_N) - H = O\left(max\left\{\frac{k^{\frac{\alpha}{\alpha+d} - \epsilon}}{N^{\frac{\alpha}{\alpha+d} - \epsilon}}, \frac{k^{\frac{\beta}{d}}}{N^{\frac{\beta}{d}}}\right\}\right) \tag{2.18}$$

uniformly for $k \in \{1, ..., k^*\}$, as $N \to \infty$.

Theorem 2 (Efficiency and Consistency) *Assume that $d \leq 3$ and that condition 1 holds for $\beta = 2$ and condition 2 holds for some $\alpha > d$, then by condition 3 (where extra assumptions are made for $d = 3$), for the estimator $\hat{H}_{N,k}$ we have;*

$$\sqrt{N}(\hat{H}_{N,k} - H) \overset{d}{\to} N(0, \sigma^2) \tag{2.19}$$

and

$$N\mathbb{E}(\hat{H}_{N,k} - H)^2 \to \sigma^2 \tag{2.20}$$

as $N \to \infty$ uniformly for $k \in \{k_0^*, ..., k_1^*\}$, where $\sigma^2 = Var(log(f(x))$, for density function $f(x)$. Thus, the estimator is asymptotically efficient and its asymptotic variance is the best attainable.

By the above, we can now say that $\hat{H}_{N,k}$ is an consistent and asymptotically unbiased estimator of exact entropy H; thus is a consistent estimator. This is due to using the central limit theorem, on the estimator for entropy $\hat{H}_{N,k}$, which states that;

$$\frac{\hat{H}_{N,k} - \mathbb{E}\hat{H}_{N,k}}{\sqrt{Var(\hat{H}_{N,k})}} \overset{d}{\to} N(0, 1)$$

By section 2.1.3, we can assume that $Var(\hat{H}_{N,k}) = \frac{Var(\log f(x))}{N} + O(\frac{1}{N}) \approx \frac{\sigma^2}{N}$.

Accordingly, the left side of the central limit theorem above can be written as;

$$\frac{\hat{H}_{N,k} - \mathbb{E}\hat{H}_{N,k}}{\sqrt{Var(\hat{H}_{N,k})}} = \frac{\sqrt{N}(\hat{H}_{N,k} - \mathbb{E}\hat{H}_{N,k})}{\sigma}$$

$$= \frac{\sqrt{N}}{\sigma}[(\hat{H}_{N,k} - H) - (\mathbb{E}\hat{H}_{N,k} - H)]$$

$$= \frac{\sqrt{N}(\hat{H}_{N,k} - H)}{\sigma} - \frac{N(\mathbb{E}\hat{H}_{N,k} - H)}{\sigma\sqrt{N}}$$

So we can see that from Theorem 2; $\sqrt{N}(\hat{H}_{N,k} - H) \xrightarrow{d} N(0, \sigma^2)$ as $N \to \infty$. Whilst from Theorem 1 we have $\mathbb{E}\hat{H}_{N,k} - H \to 0$ as $N \to \infty$. Thus as $N \to \infty$ this tends to the standard normal distribution, $N(0, 1)$, and the central limit theorem holds.

I will be exploring the bias is more detail later to see which one of the two ideas show to be more true for the behaviors of the bias for a large value of N, in dimension $d = 1$ or $d = 2$.

- With a fixed k, by [6], for $\beta \in (0, 2] \cap (0, d]$, we choose $a \in (0, \frac{\beta}{d}]$, then;

$$|Bias(\hat{H}_{N,k})| = O\left(\frac{1}{N^a}\right) \tag{2.21}$$

- With k depending on N, by [4], for $\beta \in (0, 2]$, we again choose $a \in (0, \frac{\beta}{d}]$, then;

$$|Bias(\hat{H}_{N,k})| = O\left(\left(\frac{k}{N}\right)^a\right) \tag{2.22}$$

Moreover, I will be exploring the optimal value of k, for reducing the bias, for each of these distributions according to the simulations and regression analysis. This value of k may depend on the sample size N; something which has been explored theoretically in section 2.1 but will analytically be shown later in this paper.

Chapter 3

Monte-Carlo Simulations

In this chapter I will explore simulations of the bias of estimator (2.15) in comparison to the size of the sample estimated from, with respect to different values of k; by exploring 1-dimensional distributions and then progressing onto 2-dimensional. Firstly, the distributions considered will be analysed to determine if they satisfy the conditions 1, 2 and 3 stated for Theorems 1 and 2 to hold. Then, I will explore the estimator of entropy for simulations of samples from certain distributions, for different values of k.

The motivation for these simulations is to explore the consistency of this estimator for different values of k; the relationship between the size of the bias of the estimator $\hat{H}_{N,k}$, $Bias(\hat{H}_{N,k})$, and the sample size, N. Throughout this analysis we will be considering the absolute value of this bias, since when taking its logarithm, we need a positive value. Using Theorem 1, we can write that the bias of the estimator approaches 0 as $N \to \infty$. This is because we can write $Bias(\hat{H}_{N,k}) = \mathbb{E}(\hat{H}_{N,k}) - H$, which in equation (2.18) implies $Bias(\hat{H}_{N,k}) \to 0$ as $N \to \infty$. Thus, there must be a type of inverse relationship between the modulus of the bias of the estimator, $|Bias(\hat{H}_{N,k})|$, and N. We believe this relationship is of the form;

$$|Bias(\hat{H}_{N,k})| = \frac{c}{N^a} \tag{3.1}$$

for $a, c > 0$ [6, 4]. By taking the logarithm of this, we can generate a linear

relationship, which is easier to analyse, and is given by;

$$log|Bias(\hat{H}_{N,k})| \approx log(c) - a[log(N)] + \epsilon$$
$$\approx \zeta - a[log(N)] \qquad (3.2)$$

where $\epsilon > 0$ is some small error term. I will investigate the consistency of this estimator for a sample from a specified distribution, dependent on the value of k, this mean finding the optimum value of k for which $|Bias(\hat{H}_{N,k})| \to 0$ for $N \to \infty$. For the relationship in equation (3.1), this will happen for larger values of a and relatively small c, as $N \to \infty$. As previously mentioned, there is evidence supporting that the bias becomes either of order $(\frac{1}{N})^a$ (equation (2.21)) or $(\frac{k}{N})^a$ (equation (2.22)). This leads to also examining the dependence of c / ζ on the value of k.

As I wish to consider the difference in accuracy of the estimator when using different values of k, let us denote the approximate values for a and c dependent on k as a_k and c_k.

I will conduct a range of analysis, for each distribution, to consider how this estimator acts in reality, the process of analysis will be as follows;

1. Create a summary table of the mean absolute value of the bias of the estimator for $N = 100, 25000$ and 50000 for all values of k that satisfy Condition 3. I could also consider the variance of the bias at the values of N stated above, for all applicable values of k. However, we will find that the $Var|Bias(\hat{H}_{50000,k})| \to 0$ for $k \to 10$, by the definition of the estimator using the nearest neighbour method. Taking a larger k in the nearest neighbour method will produce less varied results, this is because more smoothing takes place for a larger k, eventually - if k is made large enough - the output will be constant and the variance negligible regardless of the inputted values. Thus, considering the variance of the bias of the estimator in comparison to k is not necessarily informative.

2. Graphical representations of the linear relationship shown in equation 3.2, of $log(N)$ against $log|Bias(\hat{H}_{N,k})|$ for sample sizes $N = 100, 200, 300, ..., 50000$ (which are taken 500 times and averaged), for each value of k.

3. Tabulate the results from the regression analysis; I will first discuss the coefficient of determination (R^2), this is a measure of how well the regression model describes the observed data [26]. Next I will consider the

standard error/deviation of the model (σ), this is a measure of accuracy of predictions. Lastly, I will go onto consider the values of a_k and c_k from relationship shown in equation 3.1, for each k, which is the regression line that minimizes the sum of squared deviations (σ) of prediction.

4. Graphically compare the values of a_k and c_k for each k.

3.1   1-dimensional Gaussian/Normal Distribution

I will begin by exploring entropy of samples from the normal distribution $N(0, \sigma^2)$, where without loss of generality we can use the mean $\mu = 0$ and change the variance σ^2 as needed. The normal distribution has an exact formula to work out the entropy, given the variance σ^2. Using equation (1.1) and the density function for the normal distribution $f(x) = \frac{1}{\sqrt{(2\pi)}\sigma} \exp\left(\frac{-x^2}{2\sigma^2}\right)$ for $x \in \mathbb{R}$, given $\mu = 0$. We can write the exact entropy for the normal distribution, using equation (1.1);

$$
\begin{aligned}
H &= -\int_{x\in\mathbb{R}^d} f(x) log(f(x)) dx \\
&= -\int_{\mathbb{R}} \frac{1}{\sqrt{(2\pi)}\sigma} \exp\left(\frac{-x^2}{2\sigma^2}\right) log\left[\frac{1}{\sqrt{(2\pi)}\sigma} \exp\left(\frac{-x^2}{2\sigma^2}\right)\right] dx \\
&= \int_{\mathbb{R}} \frac{1}{\sqrt{(2\pi)}\sigma} \exp\left(\frac{-x^2}{2\sigma^2}\right) \left(log(\sqrt{(2\pi)}\sigma) + \frac{x^2}{2\sigma^2}\right) \\
&= \frac{log(\sqrt{(2\pi)}\sigma)}{\sqrt{(2\pi)}\sigma} \int_{\mathbb{R}} \exp\left(\frac{-x^2}{2\sigma^2}\right) dx + \frac{1}{2\sqrt{(2\pi)}\sigma} \int_{\mathbb{R}} \frac{x^2}{2\sigma^2} \exp\left(\frac{-x^2}{\sigma^2}\right) dx \\
&= log(\sqrt{(2\pi)}\sigma) + \frac{1}{2}
\end{aligned}
$$

Thus the exact entropy for the normal distribution is given by;

$$
H = log(\sqrt{(2\pi e)}\sigma) \tag{3.3}
$$

I will first explore samples from 1-dimensional standard normal distribution with mean $\mu = 0$ and variance $\sigma^2 = 1$, $N(0, 1)$, to consider the behavior of the Kozachenko-Leonenko estimator. The exact entropy of this distribution is given by equation (3.3), with $\sigma^2 = 1$;

$$
H = log(\sqrt{(2\pi e)}) \approx 1.418939 \tag{3.4}
$$

Since, I am first considering the 1-dimensional normal distribution, the estimator takes the form in equation (2.16), which is given by;

$$\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^{N} log \left[\frac{2\rho_{(k),i}(N-1)}{e^{\Psi(k)}} \right]$$

3.1.1 Estimator Conditions

The density of the normal distribution satisfies Conditions 1, 2 and 3, due to the below analysis. Firstly, to satisfy Condition 1, for density function $f(x) = \frac{1}{\sqrt{(2\pi)}} \exp\left(\frac{-x^2}{2}\right)$ for $x \in \mathbb{R}$, given $\mu = 0$ and $\sigma^2 = 1$, it must be such that;

- f is bounded - obvious, since for any probability distribution we always have $f(x) \geq 0$, additionally for the normal distribution we have that $f(x) = \frac{1}{\sqrt{(2\pi)}} \exp\left(\frac{-x^2}{2}\right) < 0.4$, $\forall x \in \mathbb{R}$. Hence, f is bounded above and below; so bounded.

- f is m-times differentiable - using Hermite polynomials, defined as;

$$H_m(x) = (-1)^m e^{\frac{x^2}{2}} \frac{d^m}{dx^m} \left(e^{\frac{-x^2}{2}} \right)$$

multiplying this by the coefficient in the distribution of $f(x)$, $\frac{1}{\sqrt{(2\pi)}}$, we then get;

$$\frac{d^m}{dx^m} f(x) = \frac{H_m(x)}{(-1)^m} \frac{1}{\sqrt{2\pi}} e^{\frac{-x^2}{2}}$$
$$= \frac{H_m(x)}{(-1)^m} f(x)$$

where $\frac{H_m(x)}{(-1)^m}$ is a polynomial; thus f is m-times differentiable.

- $\exists r_* > 0$ and a Borel measurable function g_*, with $\|y - x\| \leq r_*$ so that $\|f^{(t)}(x)\| \leq g_*(x)f(x)$ and $\|f^{(m)}(x) - f^{(m)}(x)\| \leq g_*(x)f(x)\|y-x\|^\eta$, for some g_* such that $\sup_{\{x:f(x)<\delta\}} g_*(x) = O(\delta^{-\epsilon})$ as $\delta \searrow 0$ for some $\epsilon > 0$.

  Since we are considering a 1-dimensional distribution, we can write the norms $\|\cdot\|$ as $|\cdot|$. Moreover, considering that for Theorems 1 and 2, we have the value of $\beta \geq 2$; thus choosing $\beta = 2$, and since $m = \lfloor\beta\rfloor = \lfloor 2 \rfloor = 2 = \beta$ and $\eta = \beta - m$, we have that $\eta = 0$. Thus we need $|f^{(t)}(x)| \leq g_*(x)f(x)$, which is obvious by above, in view of writing $|\frac{d^t}{dx^t}f(x)| = g_*(x)f(x)$, where we choose $g_*(x) = |\frac{H_t(x)}{(-1)^t}| = |H_t(x)|$, for $t = 1, 2, ..., m$, and $|f(x)| = f(x)$,

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since $f(x) > 0$. Also, g_* is a polynomial and is hence Borel measurable over $\mathbb{R}$, and for any polynomial we obviously have $\sup_{\{x:f(x)<\delta\}} g_*(x) = O(\delta^{-\epsilon})$ as $\delta \searrow 0$ for some $\epsilon > 0$. Additionally, we need $|f^{(m)}(x) - f^{(m)}(x)| \leq g_*(x)f(x)|y - x|^0 = g_*(x)f(x)$. We currently have;

$$
\begin{aligned}
|f^{(m)}(x) - f^{(m)}(x)| &= \left| \frac{H_m(x)}{(-1)^m} f(x) - \frac{H_m(y)}{(-1)^m} f(y) \right| \\
&\leq \left| \frac{H_m(x)}{(-1)^m} f(x) \right| + \left| \frac{H_m(y)}{(-1)^m} f(y) \right| \\
&= g_*(x)f(x) + g_*(y)f(y) \\
&\leq g_*(x)f(x)
\end{aligned}
$$

since we know that $f(x) > 0$ for all $x \in \mathbb{R}$, and $g_*(x) = |H_m(x)| > 0$, which is similar to the g_* before; thus satisfies the conditions for it.

Next, to satisfy Condition 2, for the density function f of the normal distribution, must fulfill that;

- The α-moment of f must be finite, so $\int_{\mathbb{R}^d} \|x\|^\alpha f(x)dx < \infty$ - this is always true for the normal distribution, all of its moments are finite, since they are defined with respect to σ^n, for some n, and $\sigma < \infty$.

Lastly, to satisfy Condition 3, we must find the values of k for which the estimator provides a uniform convergence for Theorems **??** and **??**. To do this we must have, for some $\alpha > d = 1$, let k_0^* and k_1^* denote two deterministic sequences of positive integers with $k_0^* \leq k_1^*$. Taking $\alpha := 2$, we must have;

- $k_1^* = O(N^\tau)$, where $\tau < \min\left\{ \frac{2\alpha}{5\alpha+3d}, \frac{\alpha-d}{2\alpha}, \frac{4}{4+3d} \right\} = \min\left\{ \frac{4}{13}, \frac{1}{4}, \frac{4}{7} \right\} = \frac{1}{4}$, so we can choose $\tau := \frac{2}{9} < \frac{1}{4}$ so that we have $k_1^* = O(N^{\frac{2}{9}})$

- $\frac{k_0^*}{\log^5 N} \to \infty$ - for this to be true we need to choose $k_0^* := N^A$ for some $A > 0$. Considering that $k_0^* \leq k_1^*$ and $k_1^* = O(N^{\frac{2}{9}})$, thus $A \in (0, \frac{2}{9})$. So we can choose $A := \frac{1}{\eta}$ for some large η, which gives that $k_0^* = O(N^{\frac{1}{\eta}}) \approx 1$.

Thus, on account of the values of N being considered in the simulations; $N = 100, 200, ..., 50000$, we have that for the smallest $N = 100$, the values of k for which Theorem 1 and 2 both hold, are $k \in \{k_0^*, ..., k_1^*\} = \{1, ..., 100^{\frac{2}{9}}\} = \{1, ..., 2.782\} \approx \{1, 2\}$. Also, for the middle value $N = 25,000$, we have the values of k to be in $\{k_0^*, ..., k_1^*\}$, where $k_1^* \approx 25000^{\frac{2}{9}} = 9.491 \approx 9$, thus $k \in \{1, ..., 9\}$.

Moreover, for the largest $N = 50,000$, we must consider $k \in \{1, ..., k_1^*\} = \{1, ..., 50000^{\frac{2}{9}}\} = \{1, ..., 11.072\} \approx \{1, 2, ..., 11\}$.

Overall, due to Conditions 1, 2 and 3 being met, we can say that for the normal distribution, Theorems 1 and 2 hold; henceforth, we can say that the Kozachenko-Leonenko estimator, of a sample from the 1-dimensional normal distribution is an asymptotically unbiased and consistent estimator for entropy, for some values of $k \in \{1, 2, ..., 11\}$, depending on the sample size N.

3.1.2   Simulation Results

I will now conduct some simulations to consider this for each value of k separately, each time considering 500 samples of size N from this distribution, finding the estimator in each case and take the average of these estimators to find our entropy estimator. I will then consider the relationship show in equation (3.2) for each sample and work out the average for the values of a and c, for each $k \in \{1, 2, ..., 11\}$.

For $N = 100$, $N = 25,000$ and $N = 50,000$, using the results from **??**, we can create a table to compare the mean values of the bias of the estimator for the different values of k considered.

The results shown in table 3.1 show that for a larger N, the modulus of the bias of the estimator is smaller, this is true for all values of k except when $k = 2, 3, 7, 8$, for which the bias is smaller when $N = 25,000$ in comparison to the larger value of N. There are a number of reasons why this could be; however, it is first important to notice that when finding the values of k that satisfy condition3, we found that for $N = 100$, we must have $k \in \{1, 2\}$, for $N = 25,000$ we have $k \in \{1, 2, ..., 9\}$ and for $N = 50,000$ we have $k \in \{1, 2, ..., 11\}$.

For the smallest values of $N = 100$, we expect the best value of k to be either 1 or 2; and the table agrees with this showing that the smallest bias occurs at $k = 1$ for a small sample size.

When $N = 25,000$ we have that for $k \in \{2, ..., 8\}$ that the bias is very small, especially for the values of $k = 3, 4, 7, 8$ with the smallest bias appearing when $k = 3$; which fits with the previous analysis that the best value of k will lie within 1 and 9.

Now considering the largest sample size $N = 50,000$, the bias when $k = 5$ sticks out since it is $\approx 10^{-3}$ smaller than the other bias values in the table. However, for all other values of k the bias is still extremely small in comparison

Table 3.1: *1-dimensional normal distribution, comparison of k*

k	$|Bias(\hat{H}_{100,k})|$	$|Bias(\hat{H}_{25000,k})|$	$|Bias(\hat{H}_{50000,k})|$
1	0.0031912	0.0006312	0.0004428
2	0.0195347	0.0000092	0.0003632
3	0.0167902	0.0000056	0.0002278
4	0.0264708	0.0001657	0.0001196
5	0.0238265	0.0002138	0.0000003
6	0.0311576	0.0001546	0.0001471
7	0.0356302	0.0000217	0.0003024
8	0.0396299	0.0000984	0.0001021
9	0.0460706	0.0003620	0.0002070
10	0.0458648	0.0002752	0.0002611
11	0.0387339	0.0003332	0.0002458

This table is comparing the values of $|Bias(\hat{H}_{N,k})|$ for the values of k with $N = 100$, $N = 25,000$ and $N = 50,000$, when the estimator is taken over 500 samples

Table 3.2: *1-dimensional normal distribution, $k = 5$ for large N*

N	$|Bias(\hat{H}_{N,5})|$
49100	0.0000639
49200	0.0001463
49300	0.0001700
49400	0.0001037
49500	0.0000711
49600	0.0003221
49700	0.0001047
49800	0.0000644
49900	0.0001240
50000	0.0000003

This table is comparing the values of $Var|Bias(\hat{H}_{N,5})|$ for the large values of N.

to the bias for $N = 100$ and even in comparison to $N = 25,000$ in some places. This extreme difference could be an outlier in my data; thus in table 3.2 I have shown the values for the modulus of the bias, when $k = 5$, for different, also large values of N. This table does indeed show that $|Bias(\hat{H}_{50000,5})| \approx 0.0000003$ is an anomaly in the data, and that $k = 5$ is not necessarily the best value of k for $N = 50,000$. Thus, we cannot yet draw any major conclusions about the best value of k for the estimator of a sample this size.

I now wish to consider the equation 3.2 and plot the simulated data, to fit a regression line for each value of k separately, these are shown in Figures 3.1 and 3.2. All of these graphs agree with the relationship previously stated between the sample size and the bias of the estimator; they all show that the logarithm of this equations gives a negative linear relationship - with relatively small error bars.

Moreover, I would like to consider the coefficient of determination (R^2) for each of the above regression lines, this value provides an estimate of the strength of the relationship between the model and the response variable. Also, I would
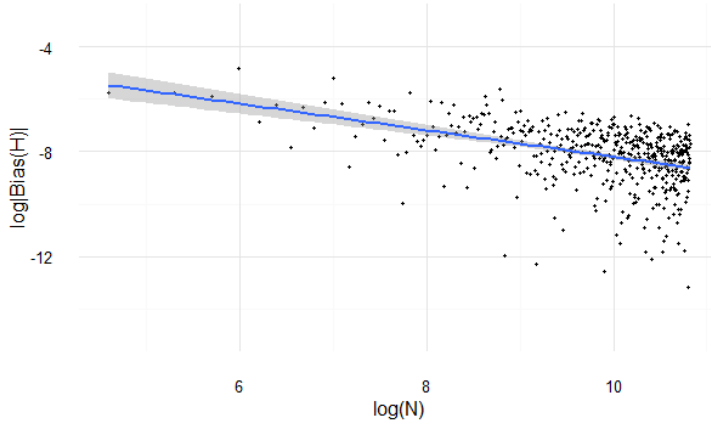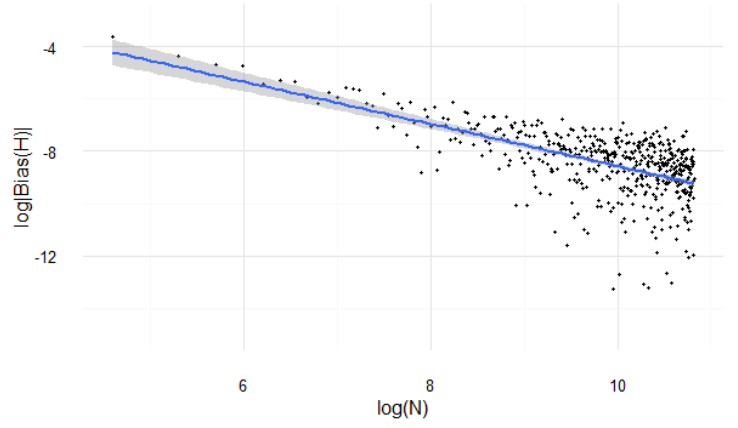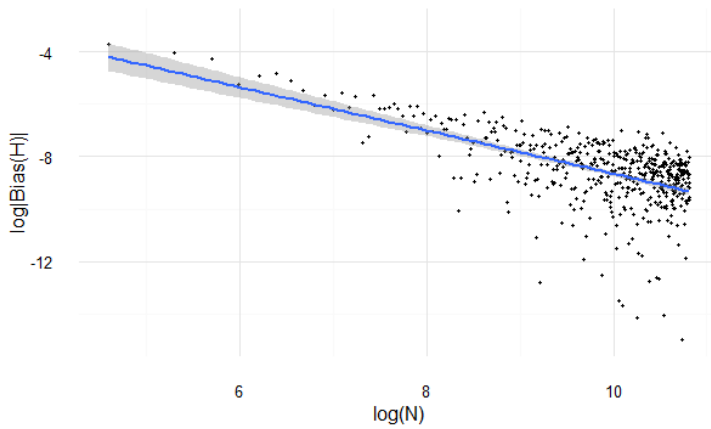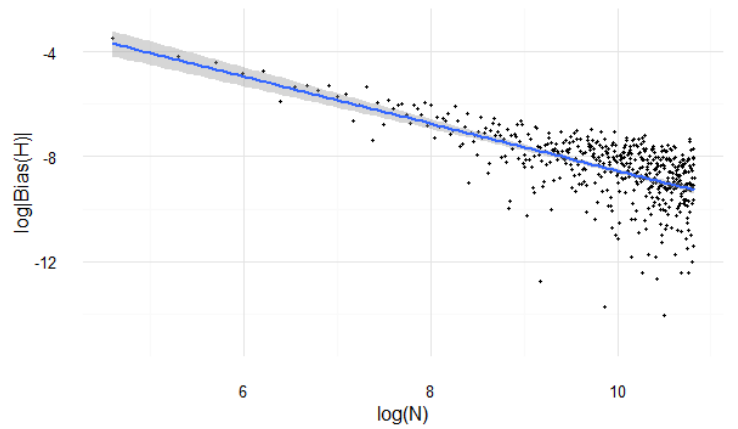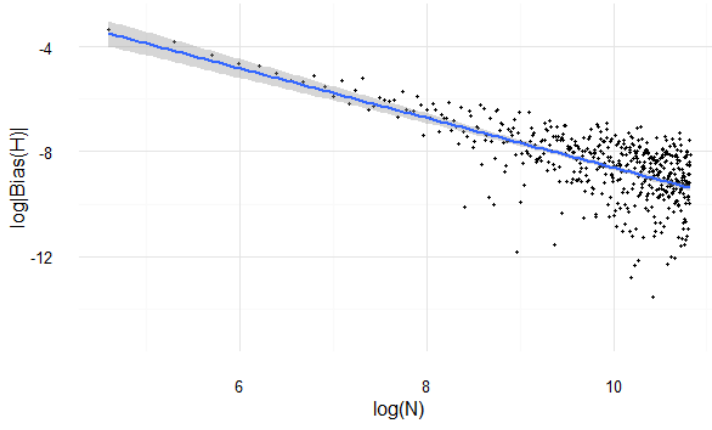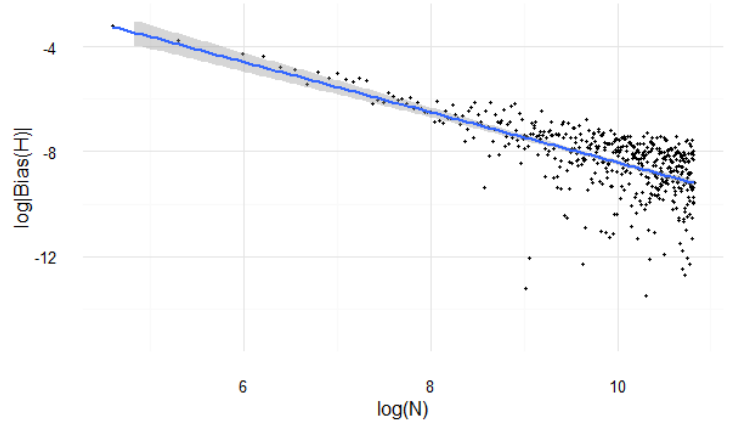
(a) *k=1*
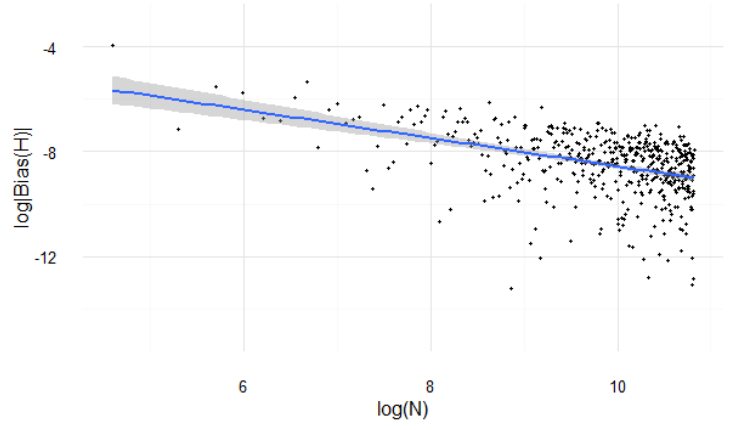
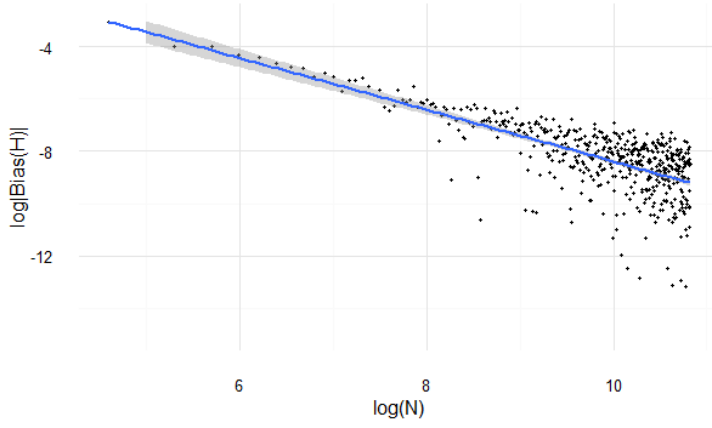(b) *k=2*
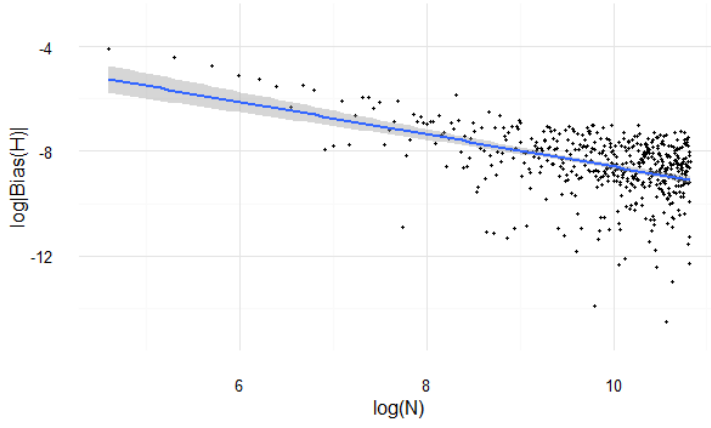
(c) *k=3*

(d) *k=4*

(e) *k=5*

(f) *k=6*

Figure 3.1: *1-dimensional normal distribution with different k = 1, ..., 6*
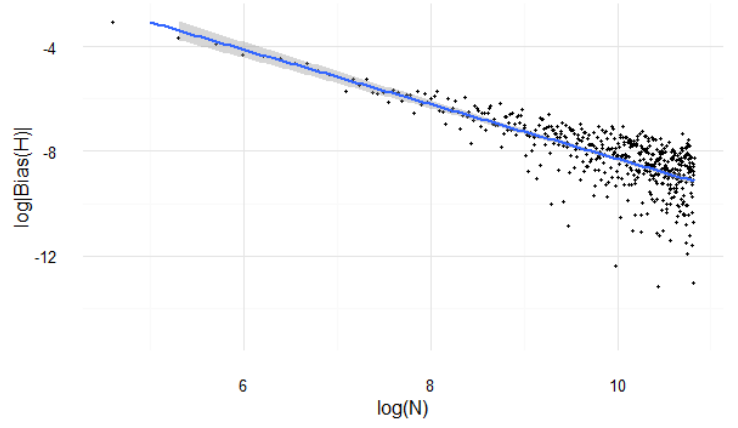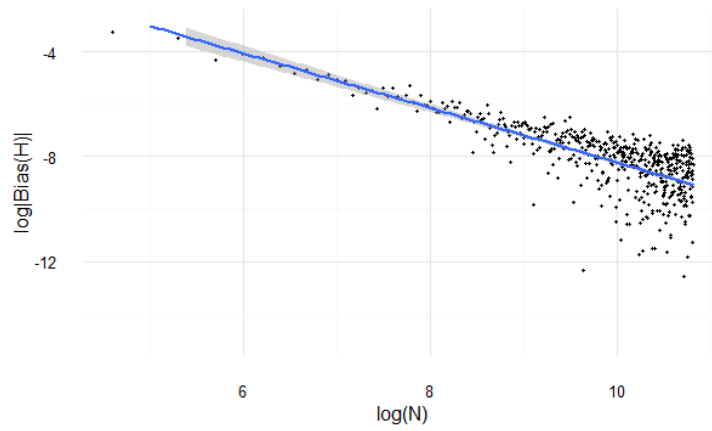
(a) *k=7*

(b) *k=8*

(c) *k=9*

(d) *k=10*

(e) *k=11*

Figure 3.2: *1-dimensional normal distribution with different* $k = 7, ..., 11$

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Table 3.3: *Comparison of the coefficient of determination and the standard deviations of the regression for each value of k for the 1-dimensional normal distribution*

k	R^2	σ
1	0.1766	1.0661
2	0.1793	1.1477
3	0.2292	1.1053
4	0.3556	1.0759
5	0.3322	1.1752
6	0.4260	1.0180
7	0.4532	1.0155
8	0.4623	1.0088
9	0.4962	0.9730
10	0.5227	0.9759
11	0.5839	0.8566

like to consider the standard error/deviation (σ), for each of the different graphs, which shows a measure of the predictions' accuracy. These are all depicted for each value of k in table 3.3.

Both columns of this table essentially point to the same conclusion; the larger the value of k, the more accurate the linear model is to fitting the data. This is shown by the R^2 value increasing towards 1 and the σ values decreasing positively.

The R^2 is very small for $k \leq 3$, which points towards the line being a poor fir to the data; however, due the the standard deviation being $\sigma^2 \approx 1.1$, we cannot say that these lines are poorly fitting; since the majoring of the data is within a very small range of the line.

The most important information found from the regression analysis is shown in table 3.4; where the values of a_k and c_k are given for each value of k.

As k runs from $1 \rightarrow 11$, we have that a_k and c_k both increase, with smooth values of a_k and a large jump, in the value of c_k, between $k = 3$ and 4, and $k = 9$ and 10. The higher the value of a_k, the stronger the negative relationship

Table 3.4: *Comparison of coefficients of regression a_k and c_k from equation 3.1, for 1-dimensional normal distribution*

k	a_k	c_k
1	0.5054	0.0433
2	0.5490	0.0459
3	0.6169	0.0894
4	0.8181	0.6690
5	0.8486	0.8235
6	0.8976	1.5514
7	0.9464	2.3576
8	0.9574	3.2021
9	0.9883	4.4558
10	1.0454	8.5402
11	1.0386	8.7457

is between the two variables in question, so for a larger values of a_k, we have that $|Bias(\hat{H}_{N,k})| \to 0$ for large N faster than smaller values of a_k. This is due to the relationship between $|Bias(\hat{H}_{N,k})|$ and a_k shown in equation (3.1)

Recall, from section 2.2 we have that the bias acts in one of two ways (equations 2.21 and 2.22); it is either of $O\left(\frac{1}{N^a}\right)$ or $O\left(\left(\frac{k}{N}\right)^a\right)$. Thus we have $|Bias(\hat{H}_{N,k})| \approx \frac{c_k}{N^{a_k}}$ where either c_k is constant or it depends on k and a_k - more specifically is $O(k^{a_k})$. There is evidence here to support the latter claim. If we consider the jump between $k = 3$ and 4 shown in the value of c_k, and consider the results in table 3.5.

This shows that the proportional behaviour between k^{a_k} and c_k also has a large jump when k goes from $3 \to 4$. This agrees with the claim of c_k depending on k in this fashion; however, in table **??** we mentioned another jump between $k = 9$ and $k = 10$, and the evidence here does not show a large jump in the same area. We cannot yet make any conclusions about the dependence of c_k on k; this motivates a graphical representation of the value of c_k against k to see if there is any relation, Figure 3.3.

Interestingly, plot 3.3 (a) shows an almost exponential relationship between

Table 3.5: *Considering the dependence of k on c_k*

k	k^{a_k}	c_k	$\frac{k^{a_k}}{c_k}$
1	1	0.0433	23.095
2	1.4631	0.0459	31.875
3	1.9694	0.0894	22.029
4	3.1085	0.6690	4.646
5	3.9187	0.8235	4.759
6	4.9942	1.5514	3.219
7	6.3067	2.3576	2.675
8	7.3218	3.2021	2.287
9	8.7716	4.4558	1.969
10	11.1020	8.5402	1.300
11	12.0668	8.7457	1.380
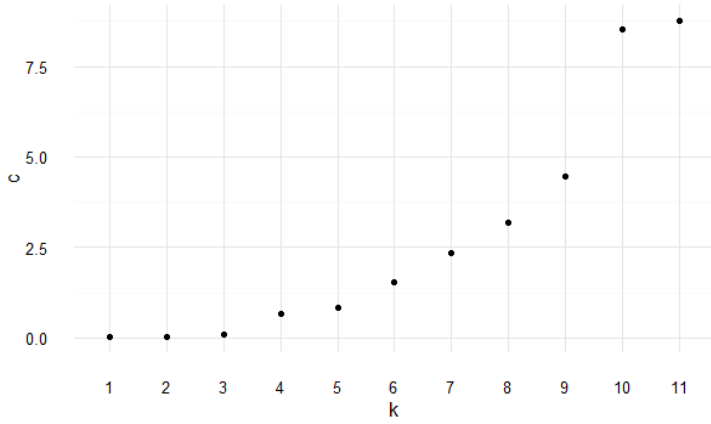
(a) *The values of k against the values of c_k*
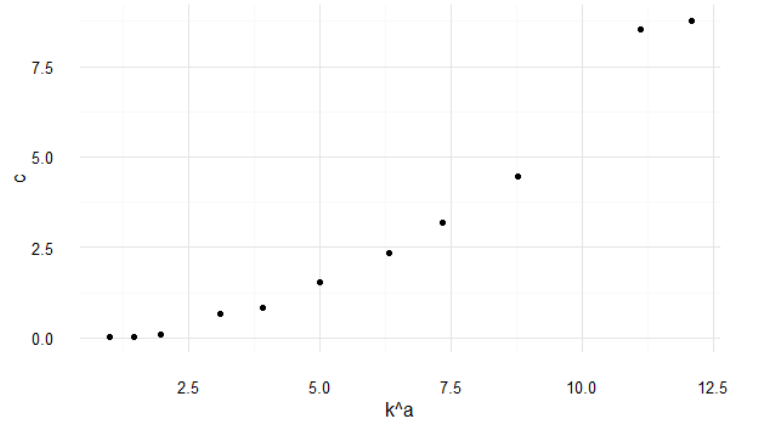
(b) *The values of k^{a_k} against the corresponding values of c_k*

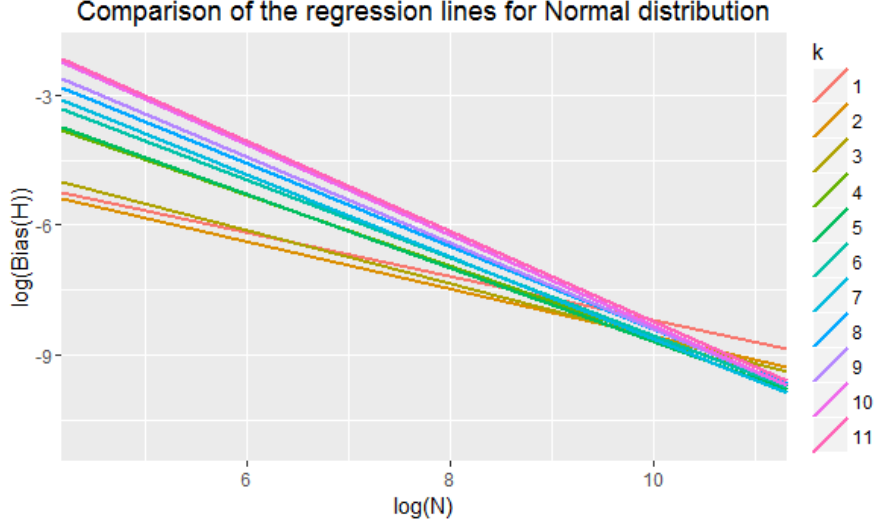Figure 3.3: *Graphically representing the relationship between c_k and k*

Figure 3.4: *Plot of regression lines for* $\log|Bias(\hat{H}_{N,k})|$ *against* $\log(N)$, *for* $k = 1, 2, ..., 11$, *for samples from the normal distribution*

the values of c_k and the values of k. This leads me to believe that there is some kind of relationship between the two variables, and looking at plot 3.3(b) this shows that there's a strong possibility that the relationship is of the form stated in equation **??**.

To better study the linear relationship between the logarithm of the bias and the logarithm of the sample size, I have generated a comparison plot, shown in Figure 3.4.

From this we can see obviously that for smaller values of N, smaller values of $log(N)$, the smallest bias occurs when $k = 2$, since this line is the lowest for the data up until $log(N) \approx 9$ - i.e. $N \approx 13,000$. For a larger sample size, we cannot accurately see in this graph which line is the best. This motivates us to look at a section of the graph when $9 \leq log(N) \leq 11$ - i.e. $8,000 \leq N \leq 50,000$, which is shown in Figure 3.5.

From this graph we can obviously discount $k = 1$ for large N, since this is the most gradual descent; thus the bias will be largest for this k. Also, both the lines for $k = 2$ and $k = 3$ are more gradual in their descent at larger N, so are probably not the best to choose. Even though, for $k = 9, 10$ and 11, the slope is the steepest - a_k is largest - the intercept is larger so around the biggest sample
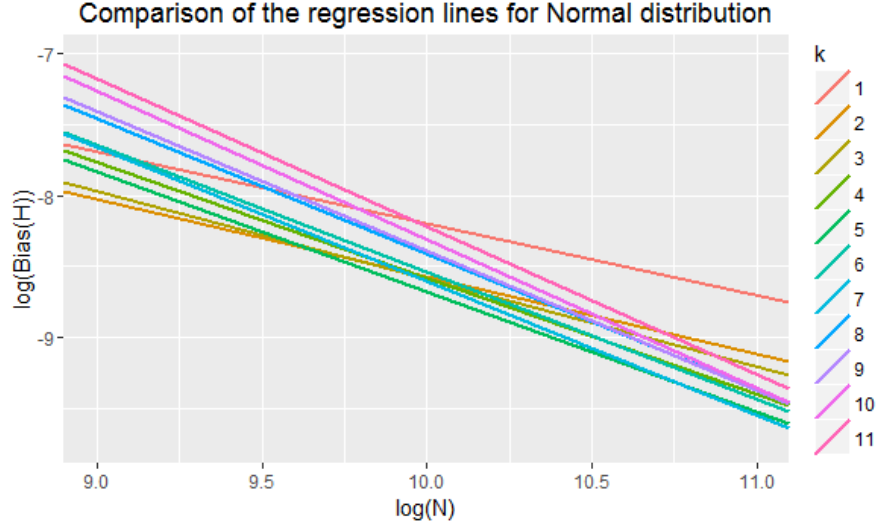
Figure 3.5: *Figure 3.4 zoomed in around large N*

size considered $N = 50,000$, $log(N) \approx 10.8$, there is not the smallest bias. Actually, for large values of $N \leq 50,000$ we can see from this graph that the best lines appear to be those which are blue/green; $k = 4, 5, 6, 7, 8$. Where the lowest lines around the maximal sample size are those for $k = 5$ and $k = 7$; thus these values of k could possible be the best nearest neighbour value to choose, when looking at a sample of size $N \approx 50,000$ from the normal distribution.

3.2   1-dimensional Uniform Distribution

I will now explore the entropy of samples from the 1-dimensional uniform distribution, $U[a, b]$. This distribution also has an exact formula to work out the entropy for. We can find this formula by considering the density function, f, from the uniform distribution, which is given by;

$$f(x) = \begin{cases} \frac{1}{b-a} & a \leq x \leq b \\ 0 & otherwise \end{cases}$$

Using the definition of Shannon entropy given in equation (1.1), we can find the exact entropy for the uniform distribution;

$$
\begin{aligned}
H &= -\int_{x \in \mathbb{R}^d} f(x) log(f(x)) dx \\
&= -\int_a^b \frac{1}{b-a} log\left[\frac{1}{b-a}\right] dx \\
&= -\frac{1}{b-a} log\left[\frac{1}{b-a}\right] \int_a^b dx \\
&= -log\left[\frac{1}{b-a}\right]
\end{aligned}
$$

Thus, the actual value of entropy for the uniform distribution is given by;

$$
H = log[b-a] \tag{3.5}
$$

Similarly to the 1-dimensional normal distribution, we have for the 1-dimensional uniform distribution that $d = 1$ so $V_1 = 2$, thus our estimator takes the form of equation (2.16);

$$
\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^N log\left[\frac{2\rho_{(k),i}(N-1)}{e^{\Psi(k)}}\right]
$$

Moreover, the samples considered will not be from the standard uniform, but from the uniform distribution $U[0, 100]$. This is because, using the standard uniform, $U[0,1]$, would fail since taking $N = 50,000$ samples between 0 and 1 would generate problems as the density function would be $f(x) = 1, \quad 0 \leq x \leq 1$, which would incur working on a very small scale; i.e taking a points with distance between them. Thus, I will be using the density function $f(x) = 0.01, \quad 0 \leq x \leq 100$, which is from the $U[0, 100]$ distribution and gives the exact entropy to be;

$$
H = log(100) \approx 4.605170 \tag{3.6}
$$

3.2.1   Estimator Conditions

For Theorems 1 and 2 to be satisfied by the estimators generated by samples from the uniform distribution, this density function must meet the Conditions 1, 2 and 3. Firstly, to satisfy Condition 1, for the density function $f(x) = 0.01$ for $0 \leq x \leq 100$, it must be such that;

- f is bounded - obviously, since the density function for the uniform distribution is constant for $x \in [a, b]$ and 0 otherwise; hence is bounded.

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- f is m-times differentiable - as f is constant this holds

- $\exists r_* > 0$ and a Borel measurable function g_*, with $\|y - x\| \leq r_*$ so that $\|f^{(t)}(x)\| \leq g_*(x)f(x)$ and $\|f^{(m)}(x) - f^{(m)}(x)\| \leq g_*(x)f(x)\|y - x\|^\eta$, for some g_* such that $\sup_{\{x:f(x)<\delta\}} g_*(x) = O(\delta^{-\epsilon})$ as $\delta \searrow 0$ for some $\epsilon > 0$. Since we are considering a 1-dimensional distribution, we can write the norms $\|\cdot\|$ as $|\cdot|$. Moreover, considering that for Theorems 1 and 2, we have the value of $\beta \geq 2$; thus choosing $\beta = 2$, and since $m = \lfloor \beta \rfloor = \lfloor 2 \rfloor = 2 = \beta$ and $\eta = \beta - m$, we have that $\eta = 0$. Thus we need $|f^{(t)}(x)| \leq g_*(x)f(x)$, which is obvious since f is constant for $f^{(t)}(x) = 0$ for all $t \geq 1$ TODO - sort this out

Next, to satisfy Condition 2, for the density function f of the uniform distribution, must fulfill that;

- The α-moment of f must be finite, so $\int_{\mathbb{R}^d} \|x\|^\alpha f(x)dx < \infty$ - this is true, since for the 1-dimensional uniform distribution, $f(x)$ is constant; thus we would be integrating a polynomial $|x|^\alpha$, over a finite interval $a \leq x \leq b$, which is always finite.

Lastly, to satisfy Condition 3, we must find the values of k for which the estimator provides a uniform convergence for Theorems 1 and 2. These values are independent of the distribution that the sample is from, and only depends on the size of the sample, the dimension of the distribution that sample is taken from and the value chosen for α, where we have chosen $\alpha = 2$. The values of k found in section **??** are $\{1, 2\}$ for $N = 100$, $\{1, 2, ..., 9\}$ for $N = 25,000$ and then $\{1, 2, ..., 11\}$ for $N = 50,000$.

Thus, due to the above conditions for Theorems 1 and 2 being met, we can say that the Kozachenko-Leonenko estimator, of a sample from the uniform distribution is an asymptotically unbiased and consistent estimator for entropy, for specific $k \in \{1, 2, ..., 11\}$, depending on the sample size N.

3.2.2 Simulation Results

I will now conduct simulations, in a similar manner as for the normal distribution, where for each value of k separately, I will consider 500 samples of size N from the uniform distribution, finding the estimator in each case and take the average of these estimators to find our entropy estimator. I will then consider

Table 3.6: *1-dimensional uniform distribution, comparison of k*

k	$\lvert Bias(\hat{H}_{100,k})\rvert$	$\lvert Bias(\hat{H}_{25000,k})\rvert$	$\lvert Bias(\hat{H}_{50000,k})\rvert$
1	0.0005189	Inf	Inf
2	0.0047466	0.0001745	0.0001163
3	0.0083912	0.0001776	0.0000899
4	0.0169364	0.0001177	0.0000490
5	0.0152168	0.0000588	0.0000509
6	0.0148205	0.0000817	0.0000538
7	0.0218339	0.0002918	0.0000663
8	0.0250401	0.0001884	0.0000487
9	0.0297655	0.0001406	0.0001184
10	0.0337164	0.0001337	0.0000949
11	0.0381473	0.0001693	0.0000417

This table is comparing the values of $\lvert Bias(\hat{H}_{N,k})\rvert$ for the values of k with $N = 100$, $N = 25,000$ and $N = 50,000$, when the estimator is taken over 500 samples

the relationship show in equation (3.2) for each sample and work out the average for the values of a and c, for each $k \in \{1, 2, ..., 11\}$.

For $N = 100$, $N = 25,000$ and $N = 50,000$, using the results from **??**, we can create a table to compare the mean values of the bias of the estimator for the different values of k considered.

Looking at the results shown in table 3.6, for $N = 100$, we can see that the smallest bias, quite obviously, occurs when the estimator is taken with $k = 1$, and that for other values of k there is a significant difference in the size of the bias.

When looking at the larger values of N, it is first important to note that `Inf` $= \infty$ is present in the table as for a large value of N, it is difficult to work out the estimator when $k = 1$. This is because there are only extremely small distances between the closest samples of this distribution, and since any program used to compute these numbers will fail with extremely small numbers

and default them to - `Inf`. We only need there to be two samples with an infinitely small distance to make the whole estimator become - `Inf`, which is `Inf` when looking at the modulus. This was mentioned earlier and was the motivation for taking the uniform distribution over $[0, 100]$; however, because of this, from now onwards I will not be considering $k = 1$ in the estimator for the uniform distribution.

Thus, looking at the results shown in table 3.6, ignoring $k = 1$ and considering $N = 25,000$, it appears to show smallest bias for values of $k \in \{4, 5, 6\}$; more specifically $k = 5$ appears to have the smallest bias. This fits with the previous analysis done on the value of k for different sample sizes, which stated that we must have $k \in \{1, 2, ..., 9\}$. Next, examining the table for $N = 50,000$, we can see that the smallest bias looks to be when $k \in \{4, 5, 6, 7, 8, 11\}$, however, for all other values of k we still have a small bias < 0.000012. Thus, we cannot yet draw any conclusions about the optimal value of k for large N and this distribution.

Next I wish to examine the graphs showing the relationship in equation 3.2, by plotting the simulated data and fitting a regression line for each value of k separately, in figures 3.6 and 3.7. Thus showing a negative linear relationship between the logarithm of the bias of the estimator and the logarithm of the sample size N. Also, looking more closely, the regression lines fitted to the data appear to become more steep for higher values of k, whilst the standard error bars appear to become smaller.

Another important thing to consider, before looking at the equations of the regression lines, is to see how well these lines actually fit the data. To do this I have, as before, examined the coefficient of determination and the standard deviation of the lines, Table 3.7.

This table is very similar to that for the normal distribution in that both columns point towards the same conclusion; the larger that k is, the more accurate the linear model is to fitting the data. This is shown by the R^2 value generally increasing towards 1 and the σ values decreasing positively - the deviation about the line is decreasing for higher k. There is slight fluctuation in the middle values of k, where for $k = \{6, 7, 8\}$ there is not an exact direction that the relationship is going. However, there is nothing to say that this isn't normal behaviour, since we do not yet know the relationship between the bias and k.
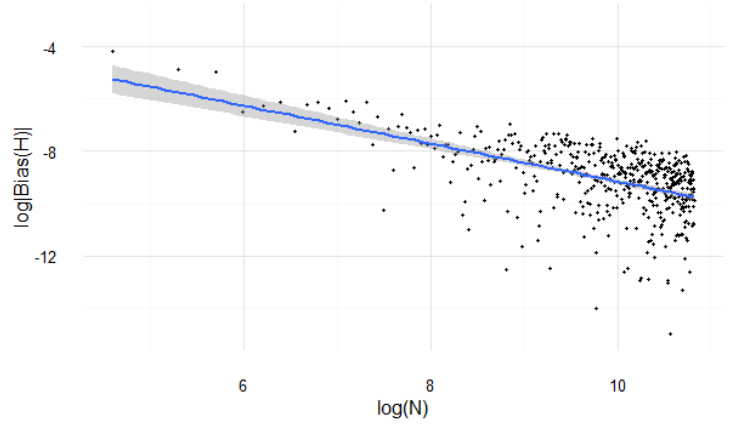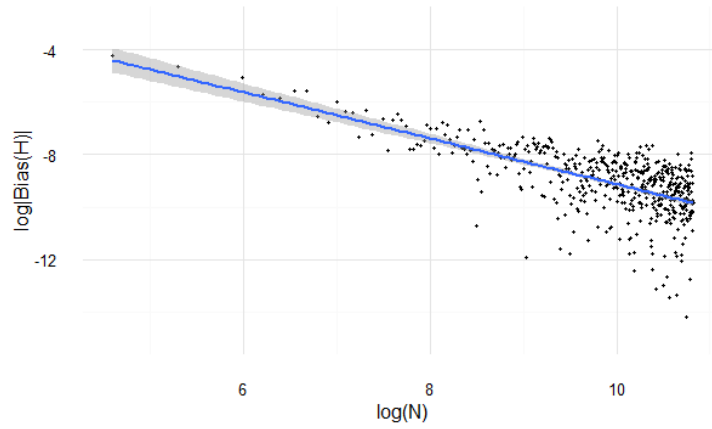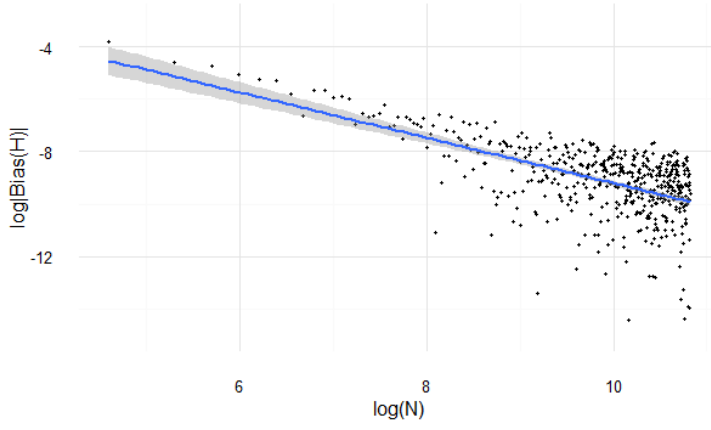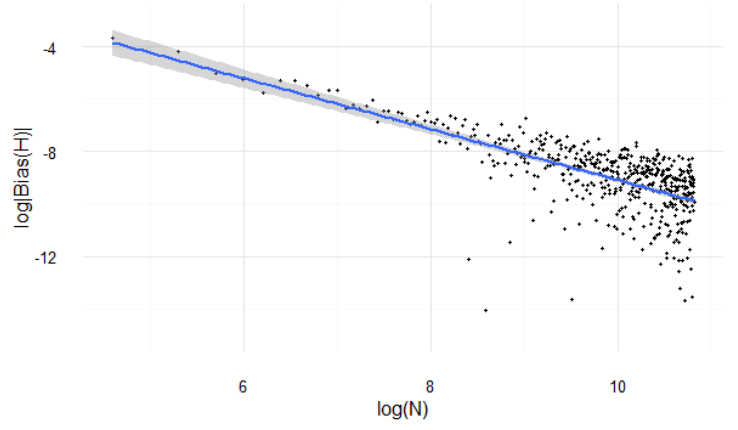
(a) *k=2*

(b) *k=3*
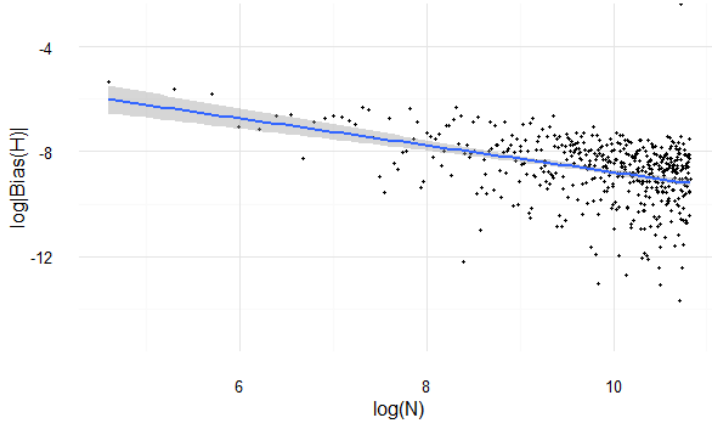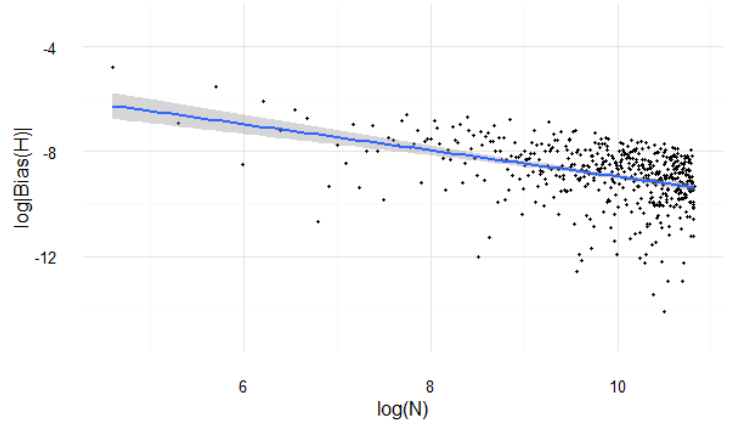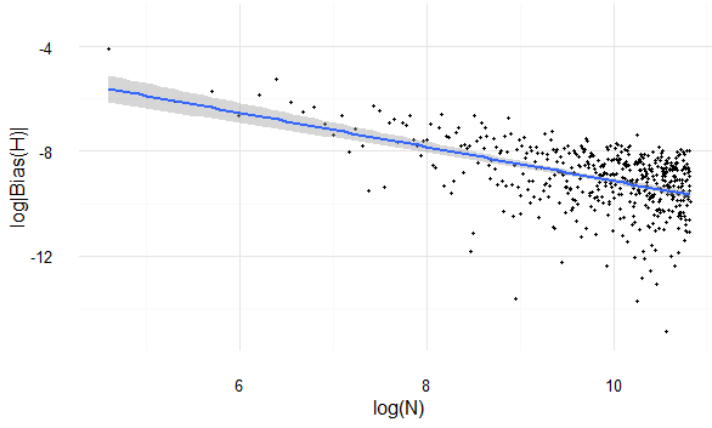
(c) *k=4*
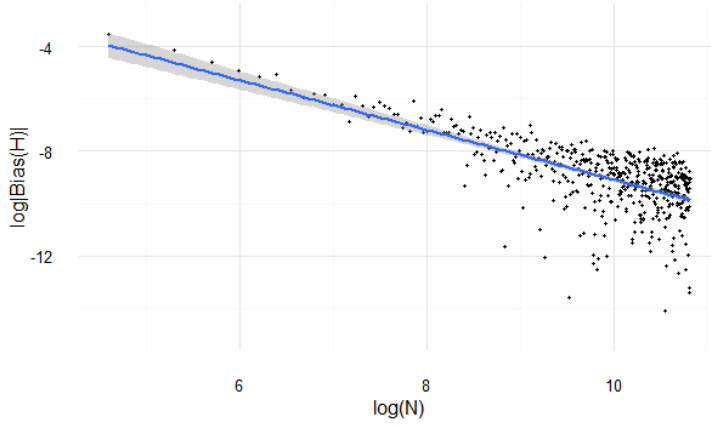
(d) *k=5*

(e) *k=6*

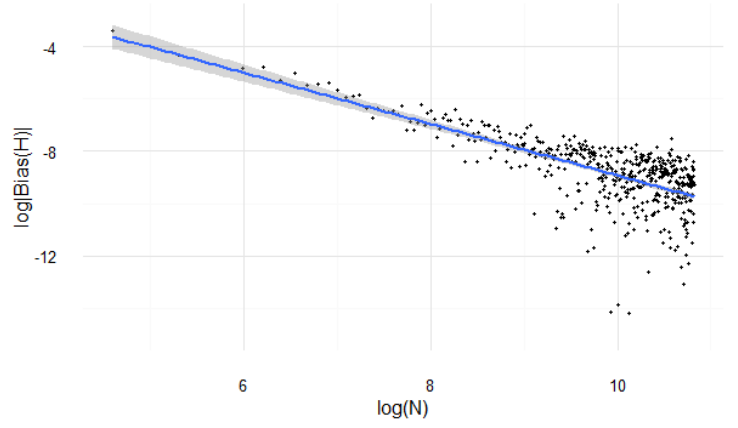Figure 3.6: *1-dimensional Uniform distribution with different $k = 2, ..., 6$*

(a) *k=7*

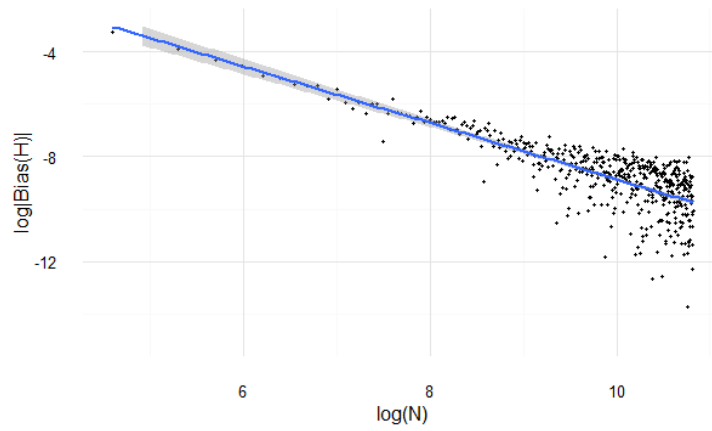(b) *k=8*

(c) *k=9*

(d) *k=10*

(e) *k=11*

Figure 3.7: *1-dimensional Uniform distribution with different k = 7, ..., 11*

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Table 3.7: *Comparison of the coefficient of determination and the standard deviations of the regression for each value of k for the 1-dimensional uniform distribution*

k	R^2	σ
2	0.1721	1.0982
3	0.1576	1.1402
4	0.2485	1.1200
5	0.2784	1.1459
6	0.4027	1.0286
7	0.3605	1.1252
8	0.4689	1.0073
9	0.4653	0.9940
10	0.5088	0.9408
11	0.6215	0.8208

Moreover, R^2 is very small for $k \leq 4$, which points towards the line being a poor fir to the data; however, due the the standard deviation being $\sigma \approx 1.1$, we cannot discard the importance of these lines; since most of the data is in a very small range about the line. Additionally, for $k = 11$ we have the strongest relationship seen so far - the largest R^2 and smallest σ values appear - this could indicate that for this distribution, when $k = 11$ the linear relationship between the logarithm of the bias and the logarithm of the sample size is most likely to be true.

Possibly the most important information found from the regression analysis, is shown in Table 3.8; where the values of a_k and c_k are given for each value of $k = \{2, 3, ..., 11\}$.

We wish to have a value of $a_k \geq 0.5$ to show the relationship desired, and this is true for all values of k in this table. As k increases, we have that both a_k and c_k increase, expect for $k = 8$, which has larger values that $k = 9$. However, this slight change around $k = 8$ does not necessarily imply anything dramatic since the overall trend seems to fit with that found for the normal distribution.

To better see if the relationship of the bias of the estimator and the sample

Table 3.8: *Comparison of coefficients of regression a_k and c_k from equation 3.1, for 1-dimensional uniform distribution*

k	a_k	c_k
2	0.5125	0.0258
3	0.5048	0.0199
4	0.6593	0.0779
5	0.7286	0.1531
6	0.8645	0.6090
7	0.8648	0.5758
8	0.9688	1.8377
9	0.9492	1.5095
10	0.9801	2.4041
11	1.0765	6.6932

size is of the form of equations 2.21 and 2.22; either $O\left(\frac{1}{N^a}\right)$ or $O\left(\left(\frac{k}{N}\right)^a\right)$. Thus I will consider the value of c_k for each k, to see if it is dependent on k_k^a or not, shown in table 3.9.

This shows that the proportional behaviour between k^{a_k} and c_k does not imply that $c_k = O(k_k^a)$, since there is no exact trend in the numbers, other than a slight decrease as k increases. However, this increase is not uniform so we can make little assumptions about the behaviour of c_k, perhaps a better way to show this relationship is through a graphical representation of c_k against k to see what form it takes.

Figure 3.8, for the uniform distribution tells a slightly different story to that of the normal. It does seem to show an almost exponential relationship, as before, however, there are a larger number of outlier values that do not fit within a smooth line in these graphs than in those for the normal distribution, Figure 3.3. However, the relationship that is shown leads me to believe that there is something important between the two variables, which could possibly be of the form stated in equation **??**.

To better study the linear relationship between the logarithm of the bias and the logarithm of the sample size, I have generated a comparison plot, shown in

Table 3.9: *Considering the dependence of k on c_k*

k	k^{a_k}	c_k	$\frac{k^{a_k}}{c_k}$
2	1.4265	0.0258	55.291
3	1.7412	0.0199	87.498
4	2.4942	0.0779	32.019
5	3.2305	0.1531	21.101
6	4.7066	0.6090	7.729
7	5.3807	0.5758	9.345
8	7.4975	1.8377	4.080
9	8.0495	1.5095	5.332
10	9.5521	2.4041	3.973
11	13.2148	6.6932	1.974
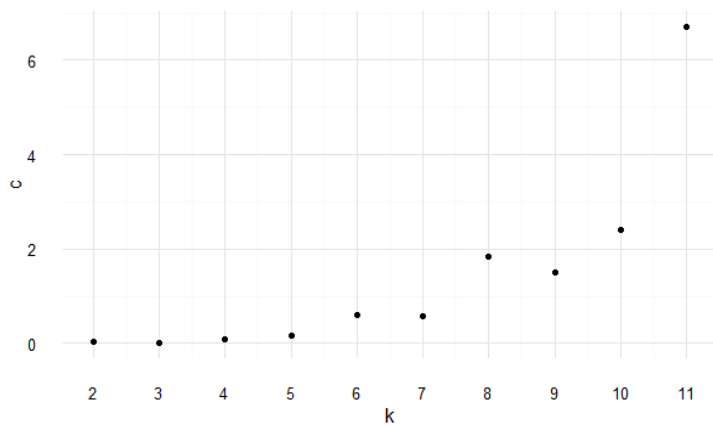
(a) *The values of k against the values of c_k*
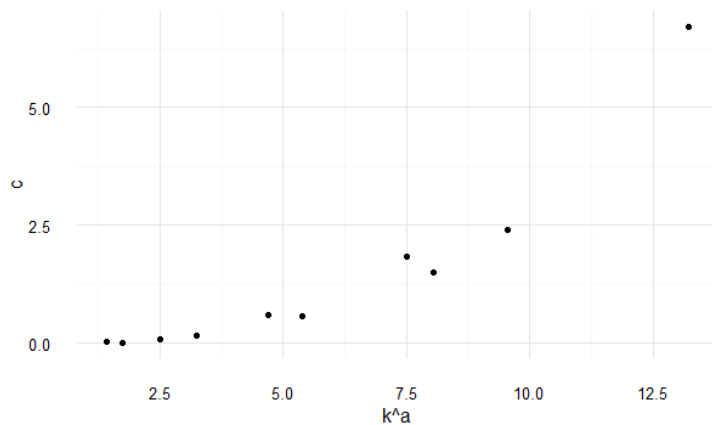
(b) *The values of k^{a_k} against the corresponding values of c_k*

Figure 3.8: *Graphically representing the relationship between c_k and k for the uniform distribution*
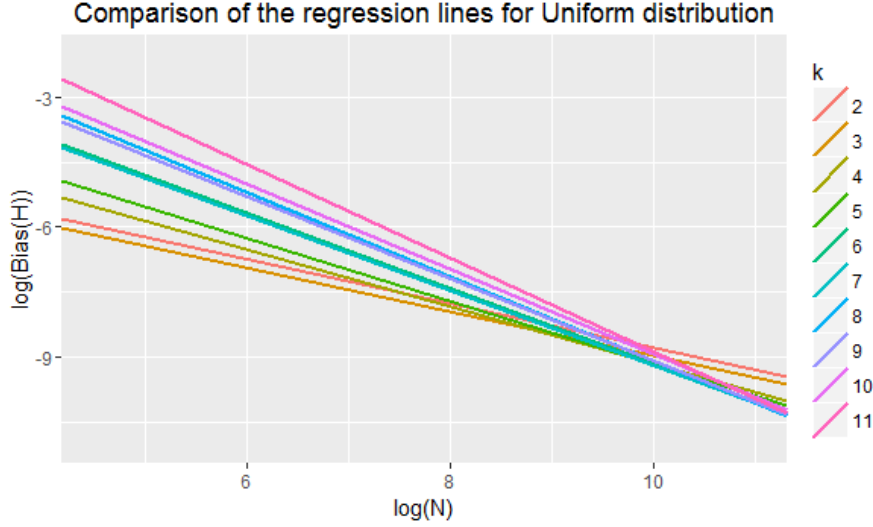
Figure 3.9: *Plot of regression lines for* $\log|Bias(\hat{H}_{N,k})|$ *against* $\log(N)$, *for* $k = 2, 3, ..., 11$, *for samples from the uniform distribution*

Figure 3.4.

For the smaller values of N, when $N \approx 100$, we can see obviously from this plot that the lowest line occurs from $k = 3$, and this value of k stays the best for the estimator until $log(N) \approx 8.5$, so the sample size $N \approx 5,000$. Above this value, it appears to be that the smallest bias occurs from when $k \in \{7, 8, 9, 11\}$, to see this more accurately consider an enlarged version of this graph for $5,000 < N < 50,000$, Figure **??**.

This graph shows that for large $N \leq 50,000$, the lowest line on the graph is when the estimator is found using $k = 7$; thus there is a possibility that this value of k could be the best nearest neighbour value to choose, when considering sample size $N \approx 50,000$ from the uniform distribution.

3.3   1-dimensional Exponential Distribution

I will now be looking at the entropy of samples from the exponential distribution $exp(\lambda)$, where $\lambda > 0$ is the rate or inverse scale parameter. In a similar fashion to the previous distributions, the exponential also has an exact formula for

Figure 3.10: *Figure 3.10 zoomed in around large N.*

the entropy, given the rate parameter λ. Using equation (1.1) and the density function for the exponential distribution $f(x) = \lambda e^{-\lambda x}$ for $x \in [0, \infty)$ and for $\lambda > 0$, we can write the exact entropy;

$$
\begin{aligned}
H &= -\int_{x \in \mathbb{R}^d} f(x) log(f(x)) dx \\
&= -\int_0^\infty \lambda e^{-\lambda x} log[\lambda e^{-\lambda x}] dx \\
&= -\lambda \int_0^\infty \lambda e^{-\lambda x} [log(\lambda) - \lambda x] dx \\
&= \lambda \int_0^\infty \lambda e^{-\lambda x} - log(\lambda) e^{-\lambda x} dx \\
&= -\lambda \left[x e^{-\lambda x} \right]_0^\infty + \int_0^\infty \lambda e^{-\lambda x} dx + log(\lambda) \left[e^{-\lambda x} \right]_0^\infty \\
&= 0 + (log(\lambda) - 1) \left[e^{-\lambda x} \right]_0^\infty \\
&= -(log(\lambda) - 1)
\end{aligned}
$$

Thus we have the the exact value of entropy for the exponential distribution, given the rate parameter $\lambda > 0$, is;

$$
H = 1 - log(\lambda) \tag{3.7}
$$

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Moreover, I am again considering a 1-dimensional distribution; thus $V_d = V_1 = 2$, and;

$$\hat{H}_{N,k} = \frac{1}{N} \sum_{i=1}^{N} log \left[\frac{2\rho_{(k),i}(N-1)}{e^{\Psi(k)}} \right]$$

is the form of the Kozachenko-Leonenko estimator that I will be considering here, equation (2.16).

I have decided to choose to explore the exponential distribution with rate parameter $\lambda = 0.5$, this is because, we know that for the exponential distribution we must have the rate parameter $\lambda > 0$ and if I choose $\lambda > e \approx 2.7183$ we get a negative values of entropy, $H < 0$. This will introduce problems when considering the modulus of the bias; hence, for this analysis it will be more beneficial to consider a positive value of entropy. Also, for $\lambda \geq 1$, we have a very small value of entropy, $0 \leq H \leq 1$, which would cause problems when calculating large samples and their entropy. Therefore, I have chosen the rate parameter such that $\lambda \in (0, 1)$, so $\lambda = 0.5$, the exact entropy is given by;

$$H = 1 - log(0.5) \approx 1.693147 \tag{3.8}$$

3.3.1 Estimator Conditions

Samples from the exponential distribution must satisfy the conditions of Theorem 2 and 1, to be an asymptotically unbiased and consistent estimator. For these theorems to hold, this distribution must satisfy the Conditions 1, 2 and 3.

Firstly, to satisfy Condition 1, the density function $f(x) = \frac{e^{\frac{-x}{2}}}{2}$ for $x \in [0, \infty)$, where we have chosen $\lambda = 0.5$, must be such that;

- f is bounded - this is true, since for any probability distribution we have $f(x) \geq 0$, also for the exponential distribution we always have for $x \in [0, \infty)$ that $f(x) \leq 1$, so f is a bounded function.

- f is m-times differentiable - this is obvious, since if we consider the mth derivative of the density function f we get;

$$\frac{d^m}{dx^m}(f(x)) = (-1)^m \lambda^{m+1} e^{\frac{-x}{\lambda}}$$
$$= (-1)^m \lambda^m f(x)$$

where $(-1)^m \lambda^m$ is a polynomial, which exists for all $x \in [0, \infty)$, thus f is m-times differentiable.

- $\exists r_* > 0$ and a Borel measurable function g_*, with $\|y - x\| \le r_*$ so that $\|f^{(t)}(x)\| \le g_*(x)f(x)$ and $\|f^{(m)}(x) - f^{(m)}(x)\| \le g_*(x)f(x)\|y - x\|^\eta$, for some g_* such that $\sup_{\{x:f(x)<\delta\}} g_*(x) = O(\delta^{-\epsilon})$ as $\delta \searrow 0$ for some $\epsilon > 0$

  Since we are considering a 1-dimensional distribution, the norm $\|\cdot\|$ can be written as $|.|$. Moreover, considering that for Theorems 1 and 2, we have the value of $\beta \ge 2$; thus choosing $\beta = 2$, and since $m = \lfloor \beta \rfloor = \lfloor 2 \rfloor = 2 = \beta$ and $\eta = \beta - m$, we have that $\eta = 0$, just as previously. Thus we need $|f^{(t)}(x)| \le g_*(x)f(x)$, which is obvious by above, in view of writing $|\frac{d^t}{dx^t}f(x)| = g_*(x)f(x)$, where we choose $g_*(x) = |(-1)^m \lambda^m| = |\lambda^m| = \lambda^m$, for $t = 1, 2, ..., m$, and $|f(x)| = f(x)$, since $f(x) > 0$. Also, g_* is a polynomial and is hence Borel measurable over $\mathbb{R}$, and for any polynomial we obviously have $\sup_{\{x:f(x)<\delta\}} g_*(x) = O(\delta^{-\epsilon})$ as $\delta \searrow 0$ for some $\epsilon > 0$. Additionally, we need $|f^{(m)}(x) - f^{(m)}(y)| \le g_*(x)f(x)|y-x|^0 = g_*(x)f(x)$. We currently have;

$$|f^{(m)}(x) - f^{(m)}(y)| = |(-1)^m \lambda^m f(x) - (-1)^m \lambda^m f(y)|$$
$$\le \lambda^m |f(x) + f(y)|$$
$$\le g_*(x)(|f(x)| + |f(y)|)$$
$$\le g_*(x)f(x)$$

  since we know that $f(x) > 0$ for all $x \in \mathbb{R}$, and $g_*(x) = \lambda^m > 0$, which is the g_* before; thus satisfies the conditions for it.

Next, to satisfy Condition 2, for the density function f of the exponential distribution, must fulfill that;

- The α-moment of f must be finite, so $\int_{\mathbb{R}^d} \|x\|^\alpha f(x)dx < \infty$ - this is true since the moments of the exponential distribution are given by;

$$\int_{\mathbb{R}^d} \|x\|^\alpha f(x)dx = \int_0^\infty |x|^\alpha \lambda e^{-\lambda x}dx$$
$$= \frac{\alpha!}{\lambda^\alpha} < \infty$$

  for all $\alpha \in \mathbb{N}$, which is obviously finite.

Lastly, to satisfy Condition 3, we must find the values of k for which the estimator provides a uniform convergence for Theorems **??** and **??**. As previously, these values are independent of the distribution that the sample is from,

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and only depends on the size of the sample, the dimension of the distribution that sample is taken from and the value chosen for α, where we have chosen $\alpha = 2$. Thus, the values of k found in section 3.1.1 are $\{1, 2, ..., 11\}$.

Due to the above conditions for Theorems 1 and 2 being met, we can say that the Kozachenko-Leonenko estimator, of a sample from the exponential distribution is an asymptotically unbiased and consistent estimator for entropy.

3.3.2   Simulation Results

I will be exploring the simulation results using the same process as for the previous two distributions, which begins with examining the values of the bias of the estimator at certain values of N for all different $k \in \{1, 2, ..., 11\}$.

In a similar manner to before, using $N = 100$, $N = 25,000$ and $N = 50,000$, from **??**, we can create a table to compare the mean values of the bias of the estimator for the different values of k considered. Recall that the information is taken over 500 sample of size N and the mean of these estimators is considered, then the bias is found by taking the modulus of this value minus the exact entropy, equation 3.8.

For the sample size $N = 100$, Table 3.10 shows that the optimal value of k to choose for the estimator is obviously given by $k = 1$, since the value of the bias is $\approx 10^{-3}$ smaller than that for all the other values of k. Considering $k = 1$ for larger sample sizes, we have ∞ - an error, similar to that for uniform distribution. This happens due to limitations in computer programs; since, we cannot have continuous data, it must be discretised, so for infinitely small distances between two points, we get a $-\infty$ value for the distance to the values closest neighbour, which only has to happen between two points for the whole estimator to become infinite. Because of this, I will not be examining the value of the estimator for $k = 1$, when considering large sample sizes.

Looking at the magnitude of the bias for sample sizes $N = 25,000$, when $k \in \{2, 3, ..., 11\}$, we can see that the smallest bias occurs at $k = 6$; but it is also quite small for the estimator found with $k = 4, 5$ and 11. Because of the closeness of the size of the bias for all the values of k just mentioned, we cannot currently draw up any conclusions about which value of k is definitly better to use - in terms of reducing the bias.

Considering the information found from samples of size $N = 50,000$, it appears to show that the estimator for $k = 4$ has a significantly smaller bias

Table 3.10: *1-dimensional exponential distribution, comparison of k*

k	$|Bias(\hat{H}_{100,k})|$	$|Bias(\hat{H}_{25000,k})|$	$|Bias(\hat{H}_{50000,k})|$
1	0.0008253	Inf	Inf
2	0.0210589	0.0008092	0.0000643
3	0.0173133	0.0001295	0.0004322
4	0.0146824	0.0000769	0.0000023
5	0.0155819	0.0000654	0.0000749
6	0.0194462	0.0000117	0.0001435
7	0.0127644	0.0005499	0.0000490
8	0.0174169	0.0000440	0.0004291
9	0.0216625	0.0003314	0.0000560
10	0.0177220	0.0004974	0.0000544
11	0.0162163	0.0000472	0.0006244

This table is comparing the values of $|Bias(\hat{H}_{N,k})|$ for the values of k with $N = 100$, $N = 25,000$ and $N = 50,000$, when the estimator is taken over 500 samples

than that for the other values of k. The next smallest bias is $\approx 10^1$ larger than that for $k = 4$, and occurs when $k = 2, 5, 7, 9$ and 10. Since we are only considering the bias at this specific sample size, and not looking at the information about the value of the bias found for samples of a similar size, we cannot yet draw any conclusive decisions from this table.

Considering the relationship shown in equation 3.2, we can plot the simulated data to see if the data does infact show this linear relationship proposed earlier. I have plotted the data seperately for each value of k, and have fitted a linear regression line to each plot to indicate the general trend of the plot. I have not considered $k = 1$ since a large proportion of the simulated values are infinite; thus, the graph does not make much sense. These plots are shown in Figures 3.11 and 3.12.

These graphs all show what we expected, and what has been confirmed with the previous two distributions, that the logarithm of the sample size against the logarithm of the bias of the estimator does indeed show a linear relationship. In these graphs there seems to be little difference between the values of k, in both the gradient of the line and in how close the data actually fits the line. This motivates us to consider both the coefficient of determination and the standard deviations of the regression for each value of k, which is given in Table 3.11.
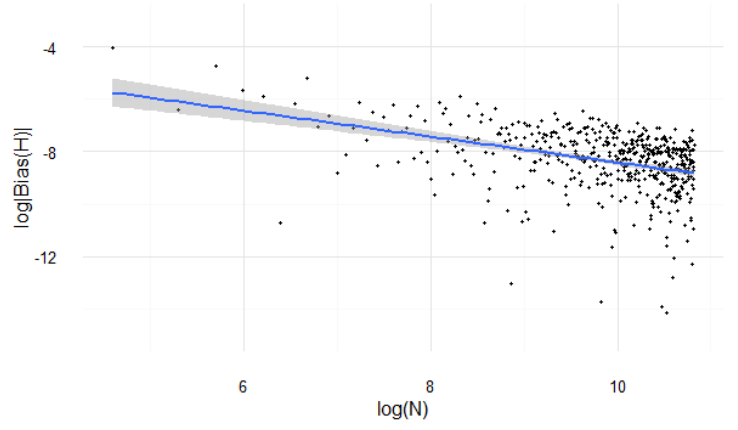
The R^2 value (coefficient of determination) is ≈ 0.207, with very little variation, for all values of k. This is an indication that the regression line could be a poor fit to the data, since a value of 1 means a perfect fit. However, the standard error of the regression line is small, ≈ 1.15 for all k, again with little variation. The standard error for the lines here is similar to that for both the uniform and normal distributions. Thus, this doesn't indicate a poor fit, just that .. TODO wtf does this show???

The most important information found from these graphs is the equation of the regression line. I have collated all the data about the coefficients a_k and c_k and displayed them in Table 3.12.

This distribution shows interesting results, that are somewhat different to those previously shown in the other two distributions. In a similar fashion to previously there is a general increase in a_k with k; however, in comparison to before where c_k acted similarly, here c_k does not increase uniformly with k. Both the values of a_k and c_k are unvaried between each k, with a_k ranging from 0.4940 to 0.6606 and c_k ranging from 0.0266 to 0.1324. Interestingly, the
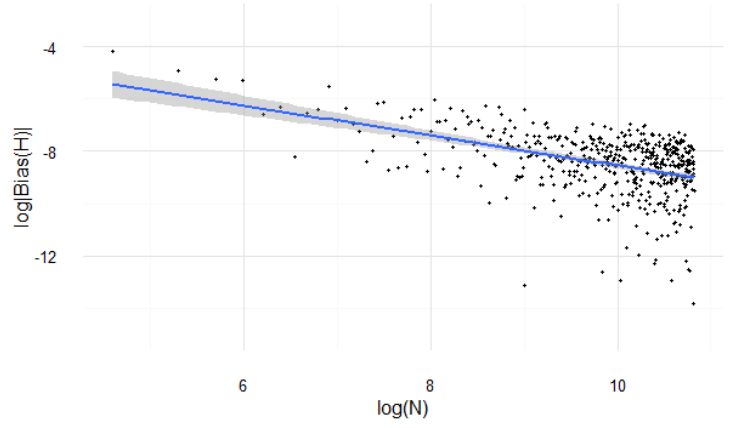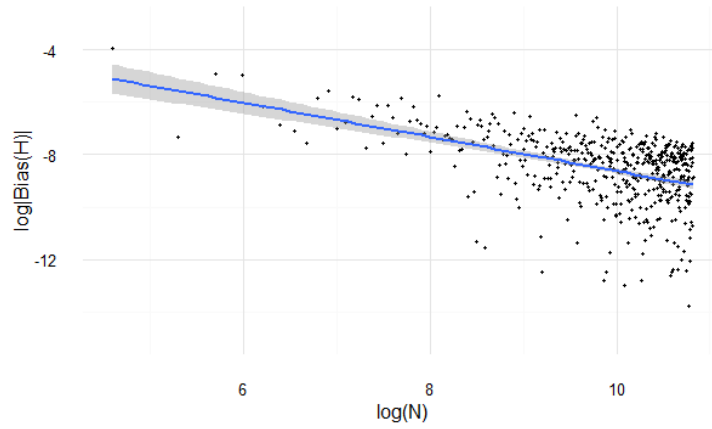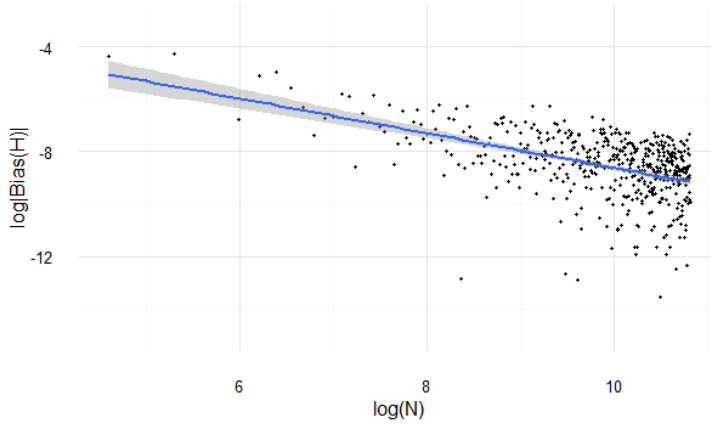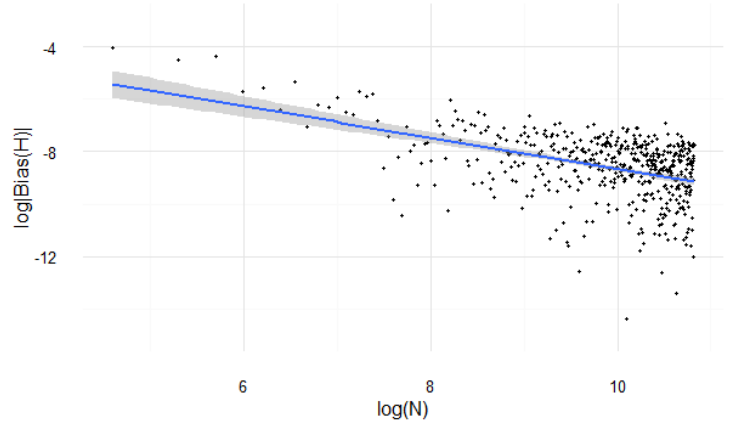
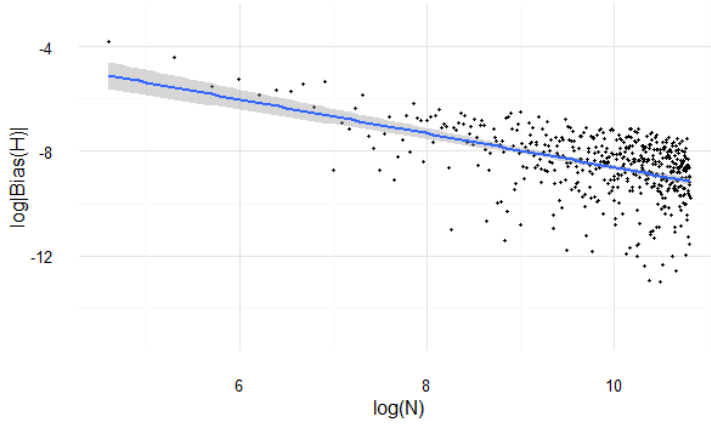(a) *k=2*

(b) *k=3*
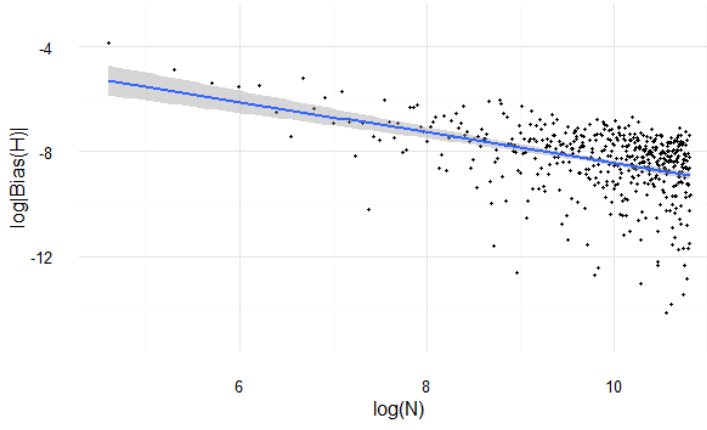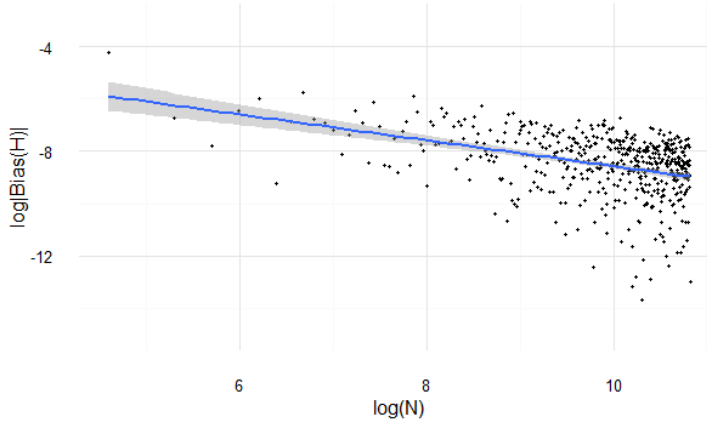
(c) *k=4*

(d) *k=5*

(e) *k=6*

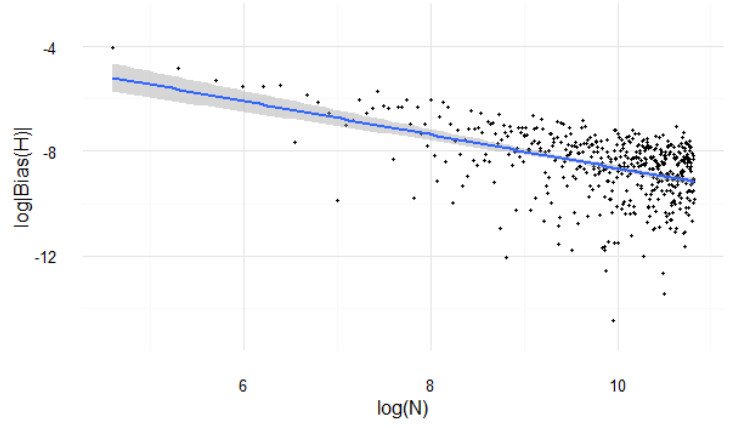Figure 3.11: *1-dimensional Exponential distribution with different $k = 2, ..., 6$*

(a) *k=7*

(b) *k=8*

(c) *k=9*

(d) *k=10*
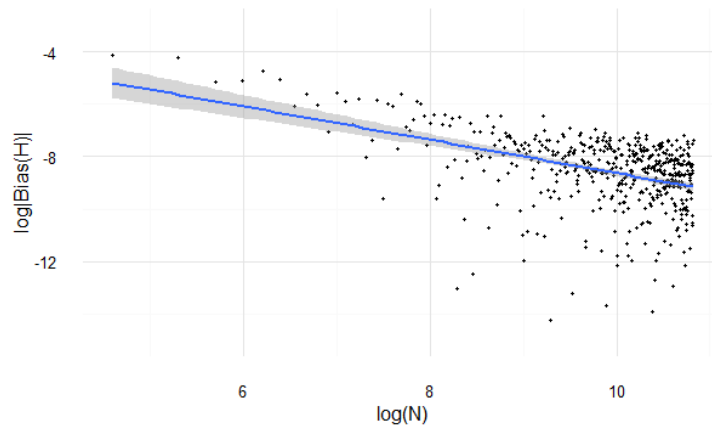
(e) *k=11*

Figure 3.12: *1-dimensional Exponential distribution with different* $k = 7, ..., 11$

Table 3.11: *Comparison of the coefficient of determination and the standard deviations of the regression for each value of k for the 1-dimensional exponential distribution*

k	R^2	σ
2	0.1843	1.1972
3	0.1596	1.1073
4	0.1538	1.1321
5	0.2151	1.0687
6	0.2277	1.1694
7	0.2474	1.1253
8	0.2107	1.1473
9	0.2509	1.0939
10	0.2200	1.2150
11	0.2017	1.2369

Table 3.12: *Comparison of coefficients of regression a_k and c_k from equation 3.1, for 1-dimensional exponential distribution*

k	a_k	c_k
2	0.5824	0.0739
3	0.4941	0.0310
4	0.4940	0.0266
5	0.5727	0.0602
6	0.6500	0.1199
7	0.6605	0.1324
8	0.6067	0.0739
9	0.6480	0.1189
10	0.6606	0.1267
11	0.6365	0.1042

Table 3.13: *Considering the dependence of k on c_k*

| k	k^{a_k}	c_k	$\frac{k^{a_k}}{c_k}$
2	1.0000	0.0739	13.532
3	1.4084	0.0310	45.434
4	1.7207	0.0266	64.687
5	2.2121	0.0602	36.745
6	2.8466	0.1199	23.742
7	3.2656	0.1324	24.665
8	3.2563	0.0739	44.063
9	3.8477	0.1189	32.361
10	4.2695	0.1267	33.697
11	4.3301	0.1042	41.556

smallest values of a_k and c_k, occur at $k = 4$ then at $k = 3$. Additionally, the largest two values occur at $k = 7$ and $k = 10$. The optimal value of k will hopefully become more apparent when plotting all the regression lines against one and other in Figure ??.

Firstly, I wish to look at the relationship between k and c_k in the exponential distribution. To do this, I will look at the values of c_k for each k and see if it depends on k^{a_k} or not. This should help us to decipher if the bias is of $O\left(\frac{1}{N^a}\right)$ or $O\left(\left(\frac{k}{N}\right)^a\right)$. This information is shown in table 3.13.

For the previous distributions there has been an obvious increase when looking at the values of $\frac{k^{a_k}}{c_k}$ for increasing k. However, for the exponential distribution this is not the case, the values seem to be all over the place, and plotting these in Figure 3.13 confirms this.

These graphs do not show the almost exponential relationship seen before, they don't appear to show any relationship between k and c_k. This could imply that for the exponential distribution we have $Bias|\hat{H}_{N,k}| = O\left(\frac{1}{N^a}\right)$, in comparison the normal and uniform which showed results in favour of the opposing idea; $Bias|\hat{H}_{N,k}| = O\left(\left(\frac{k}{N}\right)^a\right)$.

To find the optimal value of k for this distribution, I have generated a comparison plot of all of the regression lines, shown in Figure 3.14. Hopefully this
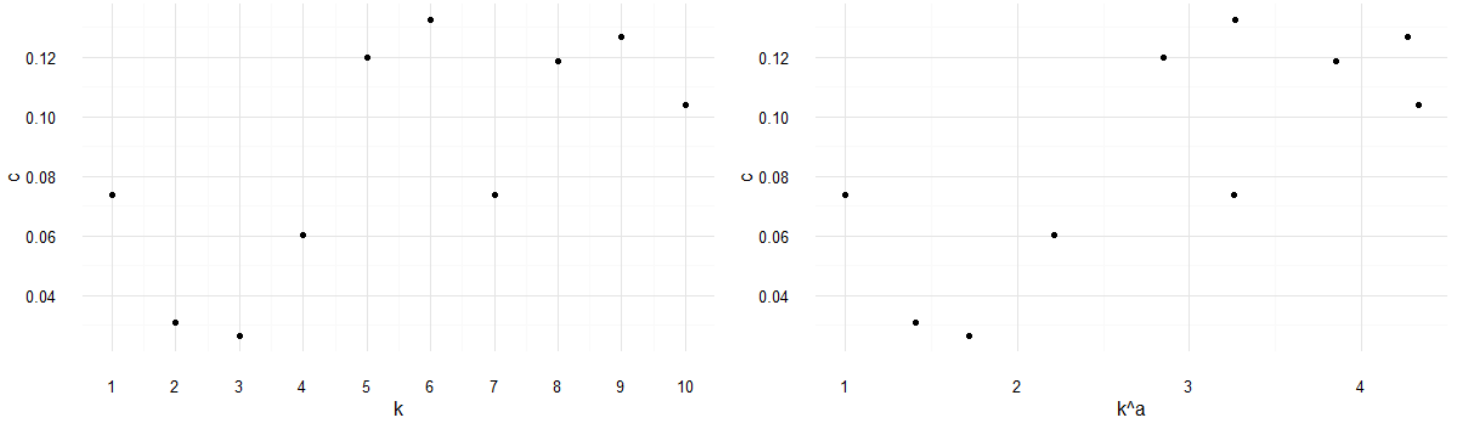
55

(a) *The values of k against the values of c_k*

(b) *The values of k^{a_k} against the corresponding values of c_k*

Figure 3.13: *Graphically representing the relationship between c_k and k for the exponential distribution*

can shed some light on which value of k appears to be the best for use for the estimator, depending on the sample size N.

At smaller sample size $N \leq 5,000$, the graph appears to show that the estimator with $k = 4$ has the lowest line; hence, the smallest bias and the best one to use for the estimation. However, for larger N this is not true, to better visualise what is happening for large N, I have created an enlarged version in Figure 3.15.

This graph shows that for large sample sizes, $5,000 \leq N \leq 50,000$ that the estimator with $k = 10$ has the smallest bias; thus appears to be the best one to use for estimating entropy from this distribution.
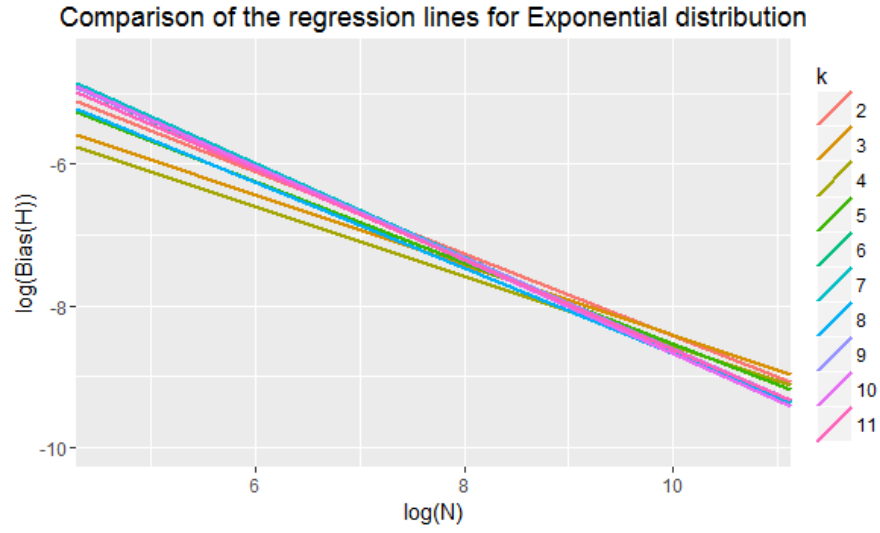
Figure 3.14: *Plot of regression lines for $\log |Bias(\hat{H}_{N,k})|$ against $\log(N)$, for $k = 2, 3, ..., 11$, for samples from the exponential distribution*

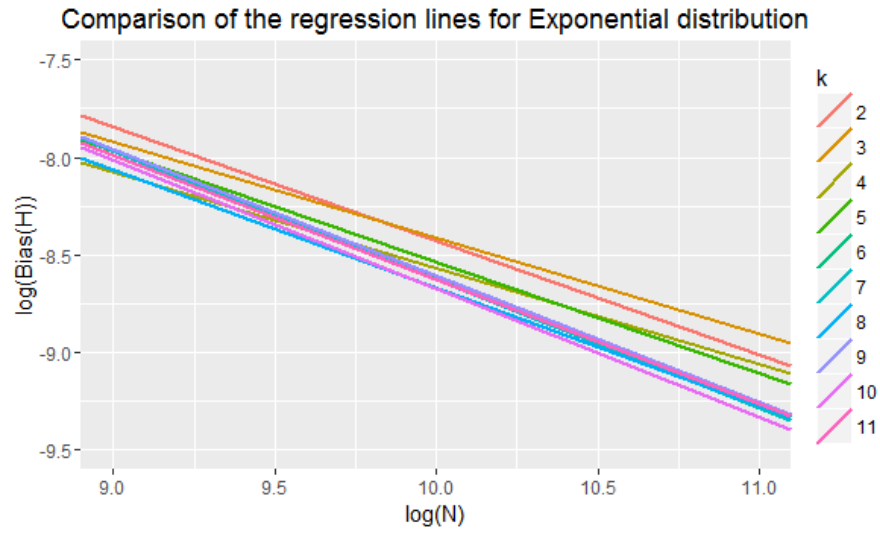Figure 3.15: *Figure 3.14 zoomed in around large N.*

Chapter 4

Conclusion

Appendix A

Data

Appendix B

Code

For the simulations in this project, I used R; "a language and environment for statistical computing and graphics" [1]. I created a package called `EntropyEst`, with the functions needed to created the Kozachenko-Leonenko entropy estimator (KLEE). The functions exported form this package are; `KLEE`, `GammaFun`, `Rho`, `VolD`, `NormalEnt`, `UniformEnt`, `ExpoEnt` and `EntBias`. The functions in this package I then used to run simulations on samples from different statistical distributions to create the results in this paper.

To create my package `Entropy-Estimators`, I used two of Hadley Wickham's [31] packages; `devtools` and `roxygen2`. `Entropy-Estimators` also has only 1 dependent package (alongside the base R packages); `FNN`, which was used for the kth nearest neighbour function.

When running simulations and creating the graphical information in this paper, I also used 4 more packages; `ggplot2` for the graphical representation of the data, `Rcpp` to create a C++ for loop for faster computation, `dplyr` for the manipulation of data, and `readr` to read and write the csv files of data.

I will outline the important code used for the simulations in this appendix; however, the full package and a complete account of the code used can be found on my GitHub page *https://github.com/KarinaMarks/Entropy-Estimators*.

B.1 The Estimator

Using the definition of the Kozachenko-Leonenko estimator, found in Section
??, I created the function `KLEE`. To do this I first had to make the following
functions; `GammaFun`, `Rho`, `VolD`. For the purpose of this paper, I have only
created the function to work out the estimator for a 1-dimensional sample.

Firstly, I created the function `Rho`, defined in equation 2.11, which takes
three arguments;

- `X`; a vector of a sample to work out the distance between a value and its
  kth nearest neighbour

- `k`; the order or nearest neighbour to be used

- `d`; the dimension of the sample, default value 1 - since currently can only
  work for 1-dimensional samples

and returns a vector of distances, where its first entry is the distance from
the first value of `X` to its kth nearest neighbour. The function is given by the
following code;

```
Rho <- function(X, k, d=1) {
  if (d == 1){
    # find the length of the sample
    n <- length(X)
    # check that k is not larger than the length of the
        vector
    stopifnot(n > k)

    # creating the matrix of kth nn distances for X
    NNdist <- FNN::knn.dist(data=X, k=k)

    # return the kth column of the matrix
    NNdist[,k]
  } else {
    return("Dimension is too high for this estimator")
  }
}
```

Next, I created the function `GammaFun`, defined in equation 2.13, which takes one argument `m` and returns a numeric value, using the following code;

```
GammaFun <- function(m) {
  # check that m > 0
  stopifnot(m > 0)
  # writing the function for the integrand
  integrand <- function(x) {x^(m-1)*exp(-x)}
  # integrating the integrand from 0 to infinity
  res <- integrate(integrand, 0, Inf)
  # selecting the result of the integral from the
      integrate class, which is a list structure
  as.numeric(res$value)
}
```

Then, using the function above for Γ, we can define the function `VolD`, equation **??**, which is the volume of the d-dimensional unit euclidean ball. This function takes one argument `d` which is the dimension and returns a numeric value using the following code;

```
VolD <- function(d) {
  # the formula to find the d-dimensional euclidean unit
      ball
  (pi^(d/2))/GammaFun(1 + (d/2))
}
```

The only other function needed for the estimator is the digamma function which is defined in base R; thus, we can now define the estimator `KLEE`. This takes three arguments;

- `X`; a vector of a sample to estimate the entropy of

- `k`; the order or nearest neighbour to be used

- `d`; the dimension of the sample, default value 1 as before for `Rho`

and returns a numeric value which is the estimator of entropy for this distribution. The code written to define this function is as follows;

```
KLEE <- function(X, k, d=1) {
  if (d==1){
```

```
  # length of the sample
  n <- length(X)
  # check that k is smaller than the length of the
    sample
  stopifnot(k < n)
  # define the vector Roe of nearest neighbour
    distances
  NN <- Rho(X, k)
  # find the volume of the unit ball
  V1 <- VolD(1)
  # return the estimator
  (1/n)*sum(log((NN*V1*(n-1))/exp(digamma(k))))
} else {
  # this would be changed to include higher dimensions
  return("Dimension must be 1")
}
}
```

B.1.1 Example

To show that the KLEE function works, I am giving the example below where I can work out the estimation of entropy by hand, and check that my function works for this.

Consider a sample $X = \{3, 6, 1, 7, 2\}$, from this we can easily find the estimator of entropy for $k = 1$ and $k = 2$. For both estimators the sample szie $n = 5$, the volume of the d-dimensional euclidean ball is $V_d = V_1 = 2$. Then for $k = 1$, the vector of nearest neighbour distances is given by $NN_1 = \{1, 1, 1, 1, 1\}$ and the digamma function is given by $\Psi(1) = -\gamma \approx -0.57722$. Thus we have;

$$\hat{H}_{5,1}(X) = \frac{1}{5} \sum_{i=1}^{5} \frac{\log(NN_{1,i} \cdot 2 \cdot (5-1))}{\exp(-\gamma)}$$

$$= \frac{1}{5\exp(-\gamma)} \sum_{i=1}^{5} \log(8)$$

$$= \frac{1}{\exp(-\gamma)} \log(8)$$

$$\approx somethingwrong$$

For $k = 2$ we have the vector of nearest neighbour distances is given by $NN_2 = \{2, 3, 2, 4, 1\}$ and the digamma function is given by $\Psi(2) = -\gamma + 1 \approx 0.42278$. Thus we have;

$$\hat{H}_{5,2}(X) = \frac{1}{5} \sum_{i=1}^{5} \frac{\log(NN_{2,i} \cdot 2 \cdot (5-1))}{\exp(-\gamma+1)}$$

$$= \frac{1}{5\exp(-\gamma+1)} \sum_{i=1}^{5} \log(8NN_{2,i})$$

$$= \frac{1}{5\exp(-\gamma+1)} \left(\log(16) + \log(24) + \log(16) + \log(32) + \log(8)\right)$$

$$\approx something?wrong$$

TODO- look at both above?

We can also find this using the function KLEE, by the following code;

```
> x <- c(3, 6, 1, 7, 2)
> KLEE(x, k=1)
[1] 2.656658
> KLEE(x, k=2)
[1] 2.430898
```

B.2   Exact Entropies

To consider the bias of the estimator, I had to find the exact value of entropy from a 1-dimensional normal, uniform and exponential distribution. The function written to return this for the normal distribution is NormalEnt with

parameter **sd**, the standard deviation of the sample, we do not need the mean value for finding the entropy of the normal distribution. The function is defined as follows;

NormalEnt <− **function**(**sd**){
  (**log**(**sqrt**(2∗pi∗**exp**(1))∗**sd**))
}

With **sd =1**, as is true in the samples considered here, we find the entropy to be given by;

> NormalEnt(**sd**=1)
[1] 1.418939

The function for the uniform distribution is **UniformEnt**, with parameters **min** and **max**, is defined as;

UniformEnt <− **function**(**min**, **max**){
  log(**max** − **min**)
}

Here we use **min=0** and **max=100** in the samples considered; thus we find the exact entropy to be given by;

> UniformEnt(**min** = 0, **max** = 100)
[1] 4.60517

Lastly, for the exponential distribution we have the function **ExpoEnt**, with only one parameter **rate**, defined below;

ExpoEnt <− **function**(rate){
  1 − **log**(rate)
}

In this paper we are using the exponential distribution with parameter **rate=1.5**, thus;

> ExpoEnt(rate = 1.5)
[1] 0.5945349

B.3   Simulations

*In this section I used the packages **readr** to save the data, **dplyr** for the manipulation of data and **Rcpp** for creating a fast loop over hundreds of iterations.*

I created functions `normalloop`, `uniformloop` and `expoloop`, in C++ which, for each sample size N creates M samples of that size, finds the estimator for sample and puts the result in a vector of length M. These functions are as follows;

```
cppFunction('
          NumericVector normalloop(int M, int N, int k){
            NumericVector est(M);
            NumericVector x(N);
            for (int i = 0; i < M; i++) {
            int sd=1;
            Function KLEE("KLEE");
            Function rnorm("rnorm");
            x=rnorm(N, sd=sd);
            est[i]=as<double>(KLEE(x ,k=k));
            }
            return Rcpp::wrap(est);
            }
            ')
```

for the normal distribution, and for the uniform distribution;

```
cppFunction('
          NumericVector uniformloop(int M, int N, int k,
             int min, int max){
            NumericVector est(M);
            NumericVector x(N);
            for (int i = 0; i < M; i++) {
            Function KLEE("KLEE");
            Function runif("runif");
            x=runif(N, min=min, max=max);
            est[i]=as<double>(KLEE(x ,k=k));
            }
            return Rcpp::wrap(est);
```

```
            }
            ')
```

Lastly for the exponential distribution;

```
cppFunction('
          NumericVector expoloop(int M, int N, int k,
             float rate){
          NumericVector est(M);
          NumericVector x(N);
          for (int i = 0; i < M; i++) {
          Function KLEE("KLEE");
          Function rexp("rexp");
          x=rexp(N, rate=rate);
          est[i]=as<double>(KLEE(x ,k=k));
          }
          return Rcpp::wrap(est);
          }
          ')
```

Using these functions I created each column of the tables, where each table is a different distribution, each column is a different value of $k \in \{1, 2, ..., 11\}$ and each row is a different sample size $N \in \{100, 200, 300, ..., 50000\}$. Below is how the column with $k = 1$ for the normal distribution was created, all other columns were done similarly;

```
# initalise the data frame with all sample sizes n
data.frame(n = seq(100, 50000, 100)) %>%
  # group by n to use summarise on each n
  dplyr::group_by(n) %>%
  # for each n the mean of the normalloop function is
      found, taken over 500 samples of size n
  summarise(Ent = mean(normalloop(M=500, N=n, k=1, rate
      =0.5), na.rm=TRUE))
```

B.4   Analysis

In this section I use the packages ggplot2 for the graphs, dplyr for the data manipulation and readr to read in my csv data files.

Once I obtained all the simulated data, I found the modulus of the bias for each sample size N, each k and each distribution. I then selected the information for all k with $N = 100, 25000$ and 50000, to display in Tables 3.1, 3.6 and **??**, this involved taking `Data` and subtracting either `NormalEnt(sd=1)`, `UniformEnt(min=0, max=100)` or `ExpoEnt(rate=0.5)` from the estimators, depending on the distribution.

Next, I plotted graphs for each k of the logarithm of the bias of the estimator $\hat{H}_{N,k}$ against the logarithm of the sample size N, shown in Figures 3.1, 3.2, 3.6, 3.7, 3.11 and 3.12. I used the following code to do this, changing the `y` value to either `k1, k2, ..., k11` depending on which value of k I was plotting. Also the `data` would be read in from a different file for each distribution, the code below shows plotting the simulations from the normal distribution with $k = 1$.

```
# read in the data as a data frame
data <- as.data.frame(read_csv("./Data/data_normal.csv"))

# find the modulus of the bias for all n and k
data[-1] <- abs(data[-1] - NormalEnt(1))

# take the logarithm of everything
logdata <- log(data)

# the max and min x values
xmin <- min(logdata$n)
xmax <- max(logdata$n)

# the min and max y values
ymin <- -15 # this is because there are only 5 values
    smaller than -15
ymax <- ceiling(max(logdata[-1]))

# plot the graph for each k - here k=1
```

```
# defining the data
ggplot(data=logdata, aes(x=n, y=k1)) +
  # plotting the points
  geom_point(size=0.8) +
  # adding a linear regression line
  geom_smooth(method="lm") +
  # labelling the axis
  xlab("log(N)") +
  ylab("log|Bias(H)|") +
  # setting the axis limits
  xlim(c(xmin, xmax)) +
  ylim(c(ymin, ymax)) +
  # choosing the graph theme
  theme_minimal()
```

Additionally, I created a summary table of the useful information needed, containing the coefficients of the intercept ζ and the gradient $-a_k$ from the regression analysis, the coefficient of determination R^2 and the standard error σ also from the regression analysis. I also modified ζ and $-a_k$ to find both a_k and c_k. The code below shows how I did this for the normal distribution, and it is similar for the other two distributions, just changing the data inputted and the exact value of entropy used to find the bias.

```
# read in the data as a data frame
data <- as.data.frame(read_csv("./Data/data_normal.csv"))

# find the modulus of the bias for all n and k - removing
    the 1st column, n
data[-1] <- abs(data[-1] - NormalEnt(1))

# take the logarithm of everything
logdata <- log(data)

# initalise and empty df with everything in
Info <- data.frame(k = 1:11, ak = rep(0, 11),
                    zeta = rep(0, 11), powera = rep(0, 11),
                    ck = rep(0, 11), rsquared = rep(0, 11),
```

$$\text{sigma} = \textbf{rep}(0, 11))$$

```r
# fill in data frame
for (k in 1:11){
  # find linear relationship of logarithm of bias against
      logrithm of n
  reg <- lm(logdata[[k+1]] ~ logdata$n)

  # the coeffs of log(bias)
  zeta <- round(reg$coefficients[["(Intercept)"]], 4)
  ak <- round(reg$coefficients[["logdata$n"]], 4)

  # the coeffs of normal bias
  ck <- round(exp(reg$coefficients[["(Intercept)"]]), 4)
  powera <- -ak

  # find the R squared value
  rsquared <- summary(reg)$r.squared

  # find the standard error
  sigma <- summary(reg)$sigma

  # fill in the each row for k=k
  Info[k,] <- c(k, ak, zeta, powera, ck, rsquared, sigma)
}

# save the Info data to a csv file
write_csv(Info, "../Data/normal_info.csv")
```

These tables are shown in Appendix **??**, and from these I found the information in Tables 3.3, 3.4, 3.7, 3.8, 3.11 and 3.12. Then to create Tables 3.5, 3.9 and 3.13, I just had to modify the summary tables found above to include two extra columns with the k^{a_k} and $\frac{k^{a_k}}{c_k}$, which was done by the following;

```r
# read in the summary data as a data frame
Info <- as.data.frame(read_csv("./Data/normal_info.csv"))
```

```
# make sure k is an integer not a factor for the
    following computation
Info$k <- as.integer(Info$k)

# create a new data frame, Info2 with c, k^a and (k^a)/c
Info2 <- Info %>%
    mutate("k^a" = k^-ak, "(k^a)/c" = ((k^-ak)/ck)) %>%
    select('k^a', ck, '(k^a)/ck')
```

From this table I than created the graphs shown in Figures 3.3, 3.8 and 3.13, using the below code.

```
# Graph (a) k against c
ggplot(data=Info2, aes(x=k, y=ck)) +
    # plotting the points
    geom_point() +
    # x axis labels
    scale_x_continuous(breaks = c(2:11), labels = c(2:11))
        +
    theme_minimal()

# Graph (b) k^a against c
ggplot(data=Info2, aes(x='k^a', y=ck)) +
    # plotting the points
    geom_point() +
    theme_minimal()
```

The last part of analysis conducted, was plotting all the regression lines of the logarithm of N against the logarithm of the bias, for each k on the same graph. To do this I used the summary data, read in as Info, and the xmin, xmax, ymin and ymax found when plotting the graphs for each k separately. The following code was then used to create the graphs in Figures 3.4, 3.10 and 3.14;

```
# make k a factor
Info$k <- as.factor(Info$k)

# plot graph of comparison for each k
ggplot()+
```

```
# add the lines for each k
geom_abline(aes(intercept=zeta, slope=a, colour=k),
    data=Info, size=1) +
# set the axis limits
ylim(c(ymin, ymax)) +
xlim(c(xmin, xmax))+
# set the axis labels
xlab("log(N)") +
ylab("log(Bias(H))") +
# set the graph title
ggtitle("Comparison of the regression lines for Normal
    distribution")


# plot graph of comparison for each k − enlarged
ggplot()+
  # add the ines for each k
  geom_abline(aes(intercept=zeta, slope=a, colour=k),
      data=Info, size=1) +
  # set tha axis limits − smaller this time for the
      enlarged plot
  ylim(c(−9.5, −7.5)) +
  xlim(c(9, 11))+
  # set the axis labels
  xlab("log(N)") +
  ylab("log(Bias(H))") +
  # set the graph title
  ggtitle("Comparison of the regression lines for Normal
      distribution")
```

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