PHENOMENOLOGY OF DOUBLE DESCENT IN FINITE-WIDTH NEURAL NETWORKS

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ABSTRACT

'Double descent' delineates the generalization behaviour of models depending on the regime they belong to: under- or over-parameterized. The current theoretical understanding behind the occurrence of this phenomenon is primarily based on linear and kernel regression models — with informal parallels to neural networks via the Neural Tangent Kernel. Therefore such analyses do not adequately capture the mechanisms behind double descent in finite-width neural networks, as well as, disregard crucial components — such as the choice of the loss function. We address these shortcomings by leveraging influence functions in order to derive suitable expressions of the population loss and its lower bound, while imposing minimal assumptions on the form of the parametric model. Our derived bounds bear an intimate connection with the spectrum of the Hessian at the optimum, and importantly, exhibit a double descent behaviour at the interpolation threshold. Building on our analysis, we further investigate how the loss function affects double descent — and thus uncover interesting properties of neural networks and their Hessian spectra near the interpolation threshold.

1 INTRODUCTION

Double-descent (DD) (Belkin et al., 2019) refers to the phenomenon of population (or test) loss exhibiting a second descent when the model is over-parameterized beyond a certain threshold (dubbed as the interpolation threshold). While such a behaviour has been originally observed in multiple contexts before (Loog et al., 2020), this terminology has been made commonplace by the influential work of (Belkin et al., 2019), who posited DD as a way to reconcile the traditional statistical wisdom with modern-day usage in machine learning. In particular, when models such as deep neural networks are employed in the over-parameterized regime, this second descent in test loss is seen as an extension of the usual U-shaped bias-variance curve — thus providing a basis to reason about the amazing generalization abilities of neural networks despite their extreme surplus of parameters.

As a result, there has been a vast number of studies that investigate double-descent further and others that seek to explain the underlying mechanisms. The former category of works (Nakkiran et al., 2019) has served to make this phenomenon conspicuous even in the context of frequently employed deep networksand shown the existence of such double-descent like behaviour based on other axes such as the amount of data or number of epochs (Nakkiran, 2019; Nakkiran et al., 2019); while the latter category of studies (Hastie et al., 2019; Mei & Montanari, 2019; Advani et al., 2020; Bartlett et al., 2020) has provided the theoretical grounds for the occurrence of such a phenomenon, almost always, in linear or kernel regression. And while the connection of (infinite-width) neural networks to Neural Tangent Kernel (Jacot et al., 2018) provides interesting parallels, but ultimately its applicability to finite-width neural networks — i.e., models of practical significance — is unclear.

Hence, our aim is to go beyond such informal correspondences, and instead, thoroughly understand and characterize the phenomenology of double descent *in finite-width neural networks*. In other words, we strive to develop a theoretical analysis that can explain the source of double descent in the class of models that actually brought this phenomenon to the limelight, as well as can cater to the idiosyncrasies that arise with the usage of neural networks, e.g., the choice of the loss function.

To this end, we derive an expression for the population risk by considering the change in the loss at a particular sample — when excluded from the training set and when included. We do so by utilizing the tool of influence functions from robust statistics, which provides us with a closed-form estimate

of the change in parameters. Subsequently, we establish a lower bound to the population risk and leverage results from Random Matrix Theory to show that it diverges at the interpolation threshold — thereby capturing double descent.

The use of influence functions (Hampel et al., 1986; Efron & Stein, 1981) makes our analysis applicable to a broad class of parametric estimators such as maximum-likelihood-type — which includes the class of neural networks but also applies to other settings such as linear and kernel regression. As a result, this approach allows us to understand the effect of the loss functions used to train neural networks — which, as we will see, starkly influences the nature of double-descent — and reveals novel insights into the properties of neural networks near the interpolation threshold.

Contributions

- Section 3: We derive a generalization expression based on a 'add-one-in' (akin to 'leave-one-out') procedure that requires minimal assumptions, and in principle, allows us to analyze the population risk of any neural network or linear/kernel regression model.
- Section 4: We show how this yields a suitable asymptotic lower-bound, which diverges at the interpolation threshold and helps explain double descent for neural networks via the Hessian spectra at the optimum. We additionally support the theoretical arguments by an empirical investigation of the involved quantities, which also suggests its applicability in the non-asymptotic setting.
- Section 5: We thoroughly study the effect of the loss function by exploring in detail the double descent behaviour for cross-entropy loss, such as the location of the interpolation threshold, Hessian spectra near interpolation, as well as discuss novel insights implied by these observations.
- Section 6: Lastly, as a by-product, we derive a generalized closed-form expression to estimate leave-one-out loss using influence functions, as well as, via that highlight an important aspect about its utility in deep learning.

2 Setup and Background

Let us consider the general setting of statistical estimation. Assume that the input samples $\boldsymbol{z} \in \mathcal{Z}$ are drawn i.i.d. from some (unknown) distribution $\mathcal{D}$. A parametric model can be defined as a family of probability distributions $\mathcal{D}_{\boldsymbol{\theta}}$ on the sample space $\mathcal{Z}$, where $\boldsymbol{\theta} \in \boldsymbol{\Theta}$. The main task is to find an estimate $\widehat{\boldsymbol{\theta}}$ of the parameters such that "$\mathcal{D}_{\widehat{\boldsymbol{\theta}}} \approx \mathcal{D}$", but given a finite set of samples $S = \{z_1, \cdots, z_n\}$ drawn i.i.d from $\mathcal{D}$. The parameter estimator, $\widehat{\boldsymbol{\theta}}$, is considered to be provided by some statistic $T(\boldsymbol{z}_1, \cdots, \boldsymbol{z}_n)$. Alternatively, if we denote the empirical distribution of the dataset S by $\mathcal{D}_n$, then the parameter estimator can be written as a functional, $\widehat{\boldsymbol{\theta}}_{\mathcal{D}} := \widehat{\boldsymbol{\theta}}(\mathcal{D}_n) = T(\mathcal{D}_n)$.

2.1 Influence Functions: a quick primer

Quite often we are interested in analyzing how a slight contamination in the distribution affects, say, the estimated parameters $\widehat{\boldsymbol{\theta}}$, or some other statistic T. Specifically, when this contamination can be expressed in the form of a Dirac distribution $\delta_{\boldsymbol{z}}$, the (standardized) change in the statistic T can be obtained via the concept of influence functions.

Definition 1. *(Hampel et al. (1986)) The influence function* IF *of a statistic T based on some distribution $\mathcal{D}$, is given by,* $\mathrm{IF}(\boldsymbol{z}; T, \mathcal{D}) = \lim_{\epsilon \to 0} \dfrac{T\left((1-\epsilon)\mathcal{D} + \epsilon\delta_{\boldsymbol{z}}\right) - T(\mathcal{D})}{\epsilon}$, *evaluated on a point $\boldsymbol{z} \in \mathcal{Z}$ where such a limit exists.*

Essentially, the influence function (IF) involves a directional derivative of T at $\mathcal{D}$ along the direction of $\delta_{\boldsymbol{z}}$. More generally, one can view influence functions from the perspective of a Taylor series expansion (or what is referred to in this context as the first-order von Mises expansion) of T at $\mathcal{D}$, evaluated on some distribution $\widetilde{\mathcal{D}}$ close to $\mathcal{D}$: $T(\widetilde{\mathcal{D}}) = T(\mathcal{D}) + \int \mathrm{IF}(\boldsymbol{z}; T, \mathcal{D})\, d(\widetilde{\mathcal{D}} - \mathcal{D})(\boldsymbol{z}) + \mathcal{O}(\epsilon^2)$. So, as evident from this, it is also possible to utilize higher-order influence functions as considered in Debruyne et al. (2008). However, the first-order term is usually the dominating term and to ensure tractability when we later consider neural networks — we will restrict our attention to first-order influence functions hereafter. Influence functions have a long history as an important tool, particularly in robust statistics (Hampel et al., 1986), but lately also in deep learning (Koh & Liang,

2017). Not only they have useful properties (see Appendix B.1 for a detailed background), but they form a natural tool to analyze changes in distribution (e.g., leave-one-out estimation) as illustrated in the upcoming sections.

2.2 Influence function for Maximum likelihood type Estimators

In general, we do not explicitly have the analytic form of the estimator $\widehat{\boldsymbol{\theta}}$ — but implicitly as the solution to an optimization problem (e.g., parameters of neural networks obtained via training). So, let us consider influence functions for the class of 'maximum likelihood type' estimators (or M-estimators), i.e. $\widehat{\boldsymbol{\theta}}$ satisfying the implicit equation, $\underset{\boldsymbol{z}\sim\mathcal{D}}{\mathbb{E}}\left[\boldsymbol{\psi}(\boldsymbol{z};\widehat{\boldsymbol{\theta}})\right] = 0$.

Let us assume that the function $\boldsymbol{\psi}(\boldsymbol{z};\boldsymbol{\theta}) := \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})$, where the (loss) function $\boldsymbol{\ell} : \mathcal{Z}\times\boldsymbol{\Theta}\mapsto\mathbb{R}$ is twice-differentiable in the parameters $\boldsymbol{\theta}$. When the loss $\boldsymbol{\ell}$ is the negative log-likelihood, we recover the usual maximum likelihood estimator, $\widehat{\boldsymbol{\theta}} = \arg\min_{\boldsymbol{\theta}\in\boldsymbol{\Theta}} \mathbb{E}_{\boldsymbol{z}\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})\right]$, assuming regularity conditions to exchange the order of expectation and derivatives. The following proposition describes the influence function for such an estimator $\widehat{\boldsymbol{\theta}}$ (the proof of which can be found in Appendix A).

Proposition 2. *The (first-order) influence function IF of the M-estimator $\widehat{\boldsymbol{\theta}}_{\mathcal{D}}$ based on the distribution $\mathcal{D}$, evaluated at point $\boldsymbol{z}$, takes the following form:* $\mathrm{IF}(\boldsymbol{z};\widehat{\boldsymbol{\theta}}_{\mathcal{D}},\mathcal{D}) = -\left[\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}})\right]^{-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\mathcal{D}})$, *where, the Hessian matrix $\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}}) := \nabla^2_{\boldsymbol{\theta}}\mathcal{L}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}})$ contains the second derivative of the loss $\mathcal{L}(\boldsymbol{\theta}) := \underset{\boldsymbol{z}\sim\mathcal{D}}{\mathbb{E}}\left[\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})\right]$ with respect to the parameters $\boldsymbol{\theta}$.*

Remark. The Hessian in neural networks is typically rank deficient (Sagun et al., 2017; Singh et al., 2021), so for the sake of our analysis we will consider a regularization term $\lambda\mathbf{I}$, $\lambda > 0$, along with the $\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}})$ in the inverse and call this modification IF_{λ}. Later, we will take the limit $\lambda \to 0$.

3 Expression of the population risk

As we lack access to the true distribution, we usually take the route of empirical risk minimization and consider $\widehat{\boldsymbol{\theta}}_S := \arg\min_{\boldsymbol{\theta}\in\boldsymbol{\Theta}} \mathcal{L}_S(\boldsymbol{\theta})$ with $\mathcal{L}_S(\boldsymbol{\theta}) := \frac{1}{n}\sum_{i=1}^{n}\boldsymbol{\ell}(\boldsymbol{z}_i,\boldsymbol{\theta})$ and recall, $S = \{\boldsymbol{z}_i\}_{i=1}^{n}$ is the training set of size n. However, in regards to performance, our actual concern is the population risk $\mathcal{L}(\boldsymbol{\theta}) := \mathbb{E}_{\boldsymbol{z}\sim D}\left[\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})\right]$, which is measured empirically via the loss $\mathcal{L}_{S'}(\boldsymbol{\theta})$ on some unseen (test) set S'. To analyze double descent (Belkin et al., 2019) — i.e., with increasing model capacity, the population risk exhibits a peak before descending again (besides the first descent as usual) — we first derive a suitable expression of the population risk that will also apply to neural networks.

'Add-one-in' procedure . Let us consider how the parameter estimate $\widehat{\boldsymbol{\theta}}_S$ changes when an additional example $\boldsymbol{z}' \sim D$ is included in the training set S. We can cast this as a contamination of the corresponding empirical distribution $\mathcal{D}_n$ to yield an altered distribution $\mathcal{D}'_{n+1} = (1-\epsilon)\mathcal{D}_n + \epsilon\,\delta_{\boldsymbol{z}'}$, with the contamination amount $\epsilon = \frac{1}{n+1}$. Thanks to the influence function of the parameter estimate (Eq. 4 for $\mathcal{D} = \mathcal{D}_n$), we do not have to retrain and explicitly measure the change in final parameters, but we have an analytic expression given as follows, $\mathrm{IF}_{\lambda}(\boldsymbol{z}';\widehat{\boldsymbol{\theta}}_S,\mathcal{D}_n) = -\left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_S)$, where the superscript S in $\mathbf{H}^S_{\mathcal{L}}$ denotes the computation of the Hessian on the set S. Now using the chain rule of influence functions, we can write the influence on $\boldsymbol{\ell}_{\boldsymbol{z}'} := \boldsymbol{\ell}(\boldsymbol{z}',\boldsymbol{\theta})$ as, $\mathrm{IF}_{\lambda}(\boldsymbol{\ell}_{\boldsymbol{z}'};\widehat{\boldsymbol{\theta}}_S,\mathcal{D}_n) = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_S\big)^{\top}\left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_S\big)$. When $|S| = n$ is large enough to ignore $\mathcal{O}(n^{-2})$ terms, the (infinitesimal) definition of influence function is equivalent to using the finite-difference form. This allows us to directly express the change in loss and which we leverage to derive an expression of the test loss, in Theorem 3 below (the proof is located in Appendix A).

Theorem 3. *Consider the parameter estimator $\widehat{\boldsymbol{\theta}}_S$ based on the set of input samples S of $|S| = n$. Then the population risk $\mathcal{L}(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}\sim D}\left[\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_S)\right])$ takes the following form,*

$$\mathcal{L}(\widehat{\boldsymbol{\theta}}_S) = \widetilde{\mathcal{L}}_S(\widehat{\boldsymbol{\theta}}_S) + \frac{1}{n+1}\,Tr\left(\left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1}\mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S)\right) + \mathcal{O}\left(\frac{1}{n^2}\right), \tag{1}$$

where $\widetilde{\mathcal{L}}_S(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}\left(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}}\right)\right]$ *is the expectation of 'one-sample training loss' and* $\mathbf{C}_{\mathcal{L}}^{\mathcal{D}}(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S)\,\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S)^\top\right]$ *is the (uncentered) covariance of loss gradients.*

Remark. Note that the first term in the right-hand side of eq. 1 is not exactly the training loss, but can be interpreted as a closely related quantity (as discussed in Section A.1.1). However, this quantity is inconsequential in regards to our analysis of double descent, as the training loss on individual samples is close to zero near interpolation.

Related work. The more interesting quantity in eq. 1 is the second term on the right, which is reminiscent of the Takeuchi Information Criterion (Takeuchi, 1976; Stone, 1977) and a similar term appears in several works on neural networks (Murata et al., 1994; Thomas et al., 2019) as well as in the analyses of least-squares regression (Flammarion & Bach, 2015; Défossez & Bach, 2015; Pillaud-Vivien et al., 2018). However, an important difference is that in our expression the Hessian $\mathbf{H}_{\mathcal{L}}^S$ evaluated on the training set appears, whereas it is based on the entire true data distribution in prior works. This seemingly minor difference is in fact crucial, since studying double descent necessarily involves analyzing the relation between the training set size and the model capacity (such as the number of parameters). Also, directly taking the results from previous works and merely approximating the involved quantities based on the training set, such as the Hessian and the covariance $\mathbf{C}_{\mathcal{L}}^{\mathcal{D}}$, does not work either. Since in such a scenario, when the limit of the regularization strength $\lambda \to 0$, results from prior works reduce to something ineffective for further analysis. E.g., for mean-squared loss, to the rank of Hessian at the optimum (see Section A.2.1 for more details).

4 Lower bound on the population risk

As a brief outline, our strategy to theoretically illustrate the double descent behaviour will be to show that the lower bound of the population risk diverges around a certain threshold. Before proceeding further, we would like to emphasize that so far *we have not employed any assumptions on the structural form of the model or the neural network*, as well as neither on the data distribution.

So let us now introduce some relevant notations to describe the precise setting for our upcoming result. We assume that the samples $\boldsymbol{z}$ are tuples $(\boldsymbol{x}, \boldsymbol{y})$, where the input $\boldsymbol{x} \in \mathbb{R}^d$ has dimension d and the targets $\boldsymbol{y} \in \mathbb{R}^K$ are of dimension K. Let us consider the neural network function is $\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}) := \boldsymbol{f}(\boldsymbol{x}, \boldsymbol{\theta}) : \mathbb{R}^d \times \mathbb{R}^p \mapsto \mathbb{R}^K$, where we have taken $\boldsymbol{\theta} \in \mathbb{R}^p$. The parameter estimator, $\widehat{\boldsymbol{\theta}}_S$, in this context refers to the parameters obtained by training to convergence using the training set S, and which we will henceforth denote by $\boldsymbol{\theta}^\star := \widehat{\boldsymbol{\theta}}_S$ (to emphasize the fact that we are at the local optimum). Also, the loss function $\boldsymbol{\ell}(\boldsymbol{z}, \boldsymbol{\theta})$, with a slight abuse of notation, refers to $\boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}\big)$ in this setting. Let us additionally define a shorthand $\boldsymbol{\ell}_i := \boldsymbol{\ell}(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_i), \boldsymbol{y}_i)$. Next, let us discuss in more detail the two matrices that appear in Theorem 3, before we introduce the assumptions we require.

First, we have the *Hessian matrix of the loss*, which in general can be decomposed as a sum of two other matrices: outer-product Hessian $\mathbf{H}_o^S(\boldsymbol{\theta})$ and functional Hessian $\mathbf{H}_f^S(\boldsymbol{\theta})$, i.e.,

$$\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}) = \mathbf{H}_o^S(\boldsymbol{\theta}) + \mathbf{H}_f^S(\boldsymbol{\theta}) = \frac{1}{n}\sum_{i=1}^n \nabla_{\boldsymbol{\theta}}\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_i)\big[\nabla_{\boldsymbol{f}}^2\,\boldsymbol{\ell}_i\big]\,\nabla_{\boldsymbol{\theta}}\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_i)^\top \;+\; \frac{1}{n}\sum_{i=1}^n \boldsymbol{\ell}_i\,\nabla_{\boldsymbol{\theta}}^2\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_i)\,, \quad (2)$$

where, $\nabla_{\boldsymbol{\theta}}\boldsymbol{f}_{\boldsymbol{\theta}} \in \mathbb{R}^{p\times K}$ is the Jacobian of the function and $\nabla_{\boldsymbol{f}}^2\,\boldsymbol{\ell} \in \mathbb{R}^{K\times K}$ is the Hessian of the loss with respect to the function outputs. Next, the other matrix that appears in eq. 1 is the (uncentered) *covariance of loss gradients*, $\mathbf{C}_{\mathcal{L}}^{\mathcal{D}}(\boldsymbol{\theta}) = \mathbb{E}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}}\left[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}\big)\,\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}\big)^\top\right]$, with $\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}\big) = \nabla_{\boldsymbol{\theta}}\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x})\,\nabla_{\boldsymbol{f}}\boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}\big)$ where $\nabla_{\boldsymbol{f}}\boldsymbol{\ell} \in \mathbb{R}^K$ is the gradient of the loss with respect to the function output. Similar to the covariance of loss gradients, let us define the covariance of function Jacobians, which we will denote by $\mathbf{C}_{\boldsymbol{f}}^S$ when computed over the samples in set S. We can express it concisely as $\mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}) = \frac{1}{|S|}\mathbf{Z}_S(\boldsymbol{\theta})\mathbf{Z}_S(\boldsymbol{\theta})^\top$, with $\mathbf{Z}_S := [\nabla_{\boldsymbol{\theta}}\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_1) \cdots \nabla_{\boldsymbol{\theta}}\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_{|S|})]^\top \in \mathbb{R}^{p\times K|S|}$.

4.1 Main result

We employ the following assumptions to obtain an appropriate lower bound of the population risk:

Assumption A1. *The loss function $\boldsymbol{\ell}\left(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\right)$ is convex in the first argument, i.e., $\nabla^2_{\boldsymbol{f}}\, \boldsymbol{\ell} \succeq \mathbf{0}$.*

Assumption A2. *The functional Hessian at the optimum $\boldsymbol{\theta}^\star$ is zero, i.e., $\mathbf{H}^S_f(\boldsymbol{\theta}^\star) = \mathbf{0}$.*

Assumption A3. *There exists a sample $(\boldsymbol{x}, \boldsymbol{y}) \sim \mathcal{D}$ such that $\sigma^2_{(\boldsymbol{x},\boldsymbol{y})} := \|\nabla_{\boldsymbol{f}} \boldsymbol{\ell}(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y})\|^2 > 0$.*

Assumption A4. *The columns of $\mathbf{Z}_S(\boldsymbol{\theta}^0)$ are isotropic vectors in $\mathbb{R}^p$ at initialization $\boldsymbol{\theta}^0$.*

Assumption A5. *The minimum non-zero eigenvalue λ_{min} of covariance of function Jacobians at optimum $\boldsymbol{\theta}^\star$ is bounded above and below by the corresponding one at initialization $\boldsymbol{\theta}^0$, i.e., $A_{\boldsymbol{\theta}^0}\, \lambda_{min}(\mathbf{C}^S_{\boldsymbol{f}}(\boldsymbol{\theta}^0)) \leq \lambda_{min}(\mathbf{C}^S_{\boldsymbol{f}}(\boldsymbol{\theta}^\star)) \leq B_{\boldsymbol{\theta}^0}\, \lambda_{min}(\mathbf{C}^S_{\boldsymbol{f}}(\boldsymbol{\theta}^0))$, with constants $0 < A_{\boldsymbol{\theta}^0},\, B_{\boldsymbol{\theta}^0} < \infty$.*

Finally, we are in a position to state our main theorem (all of our proofs can be found in Appendix A).

Theorem 4. *Under the assumptions, A1, A2, A3, A4, A5 and under the limit of external regularization $\lambda \to 0$, we obtain the following lower bound on the population risk stated in eq. 1,*

$$\mathcal{L}(\boldsymbol{\theta}^\star) \;\geq\; \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) \;+\; \frac{1}{n+1} \cdot \frac{\sigma^2_{min}\; A_{\boldsymbol{\theta}^0}}{\lambda_r\left(\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star)\right)} \;. \tag{3}$$

where we consider the convention $\lambda_1 \geq \cdots \geq \lambda_r$ for the eigenvalues with $r := \mathrm{rank}(\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star))$, and σ^2_{min} denotes the minimum non-zero $\sigma^2_{(\boldsymbol{x},\boldsymbol{y})}$ over $(\boldsymbol{x}, \boldsymbol{y}) \sim \mathcal{D}$.

The key takeaway of this theorem is that the lower bound on population risk is inversely proportional to the minimum non-zero eigenvalue λ_r of the Hessian at the optimum $\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star)$, which — as we will see shortly — largely characterizes the double descent like the behaviour of the population risk.

Note on the assumptions. The assumptions we considered are not restrictive and actually hold in a wide variety of scenarios. Assumption A1 only asks the loss to be convex with respect to the predicted output $\hat{\boldsymbol{y}} = \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x})$ and only at the optimum $\boldsymbol{\theta}^\star$, and most importantly, *not with respect to the parameters θ*. E.g., this holds for mean-squared loss, cross-entropy, and hinge loss, to name a few. Next, assumption A2 is known from prior works (Sagun et al., 2017; Singh et al., 2021) to hold empirically. Also, in the setting of double descent, near the interpolation threshold the individual losses are themselves close to zero, thereby making the functional Hessian vanish at the optimum. Together A1 and A2 guarantee that the Hessian $\mathbf{H}^S_{\mathcal{L}}$ at the optimum is positive semi-definite.

Further the assumption A3 just requires the existence of such a point from the true distribution $\mathcal{D}$. Notice, otherwise, we would have zero population risk, which is obviously of no interest. The benefit of this assumption is that we can analyze $\mathbf{C}^{\mathcal{D}}_{\boldsymbol{f}}(\boldsymbol{\theta}^\star)$ instead of $\mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\boldsymbol{\theta}^\star)$ by taking the minimum non-zero $\sigma^2_{\min}$ outside of the corresponding expression. The isotropic assumption A4 at initialization is just a simplifying assumption that makes the results succinct, but otherwise one can just apply commensurate transformations at the initialization or else get a factor of $\lambda_{\min}(\boldsymbol{\Sigma})$ in the numerator (where $\boldsymbol{\Sigma}$ is the corresponding covariance in such a case). The last assumption A5 essentially guarantees that, at the optimum, the minimum non-zero eigenvalue of the covariance of function gradients does not change much relative to that at initialization. This can be interpreted as smoothness in the spectrum of $\nabla^2_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x})$. In Figure. 2a, we empirically justify this assumption, but also in the adjoining Figure 2b validate our lower bound to the population risk throughout the double descent curve (additional details of which can be found in Appendix C.5).

4.2 Double descent behaviour

Now that we have established this lower bound, we will recover the existence of the double descent behaviour as a corollary. Let us consider the case of mean-squared error (MSE) loss, $\boldsymbol{\ell}(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}) = \frac{1}{2}\|\boldsymbol{y} - \boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x})\|^2$. In this case, the Hessian of the loss with respect to the function output is just the identity, i.e., $\nabla^2_{\boldsymbol{f}} \boldsymbol{\ell} = \mathbf{I}$. Further, by assumption A2, the Hessian at the optimum simply reduces to, $\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star) = \mathbf{C}^S_{\boldsymbol{f}}(\boldsymbol{\theta}^\star)$. Next, we will employ the following additional assumption,

Assumption A6. *The columns of $\mathbf{Z}_S(\boldsymbol{\theta}^0)$ are sub-Gaussian independent random vectors at initialization $\boldsymbol{\theta}^0$.*

This assumption may seem more demanding but is still mild in comparison to that in literature (Pennington & Bahri, 2017), which even consider all the entries of the covariance of function gradients at the optimum to be independent and identically distributed. Whereas, our assumption just requires

sub-gaussianity *only at initialization*. Similar sub-gaussianity assumptions are also common in the regression-based analyses of double descent (Muthukumar et al., 2019; Bartlett et al., 2020).

Nevertheless, the benefit of such an assumption is that it allows us to precisely characterize the behaviour of the minimum eigenvalue that appears in Theorem 4 using results from Random Matrix Theory (Vershynin, 2010), and thereby that of double descent, as described in the Corollary 5.

Corollary 5. *For MSE loss, under the setting of Theorem 4 and the additional assumption A6, the population risk takes the following form and diverges almost surely to ∞ at $p = \frac{1}{\sqrt{c}}Kn$,*

$$\mathcal{L}(\boldsymbol{\theta}^\star) \geq \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) + \frac{\sigma_{min}^2 \; A_{\boldsymbol{\theta}^0} \; K}{B_{\boldsymbol{\theta}^0} \left(\sqrt{Kn} - c\sqrt{p}\right)^2},$$

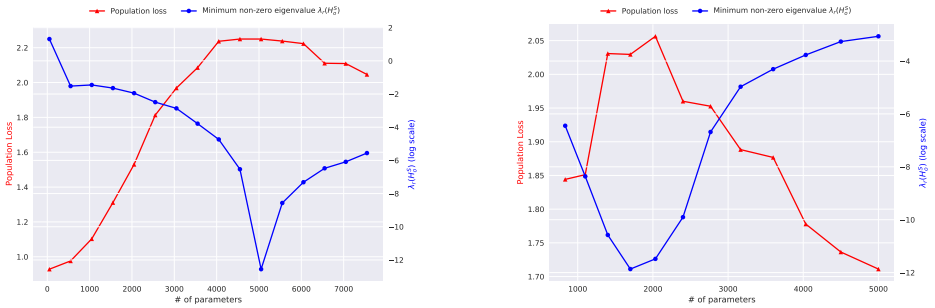
in the asymptotic regime of $p, n \to \infty$ but their ratio is a fixed constant, and where $c > 0$ is a constant that depends only on the sub-Gaussian norm of columns of $\mathbf{Z}_S(\boldsymbol{\theta}^0)$.

Takeaways. There are some important things to take note of. (a) Typically, the point where the second descent occurs, i.e, the interpolation threshold, is portrayed as $p \approx n$. This is despite the fact that empirically, for MSE loss with K targets, the more accurate location is $p \approx Kn$ — only briefly alluded to in prior works (Belkin et al., 2019; Greydanus, 2020). The above corollary provides a theoretical basis for the location of interpolation threshold. (b) Next, not only do we obtain a divergence at the interpolation threshold, but also illustrates that the complexity term will get smaller as its denominator increases for $p >> n$, — thereby also capturing the trend of population risk. (c) Lastly, while the above result holds in the asymptotic setting, we empirically show that such a behaviour also takes place for p and n as small as a few thousands, as discussed in Section 4.3 ahead.

Interpretation of the interpolation threshold location. Intuitively, based on the above results, double descent occurs at the transition when the Hessian rank (at the optimum) starts being dictated by the # of samples (i.e., when over-parameterized) rather than being governed by the # of parameters (i.e., when under-parameterized), e.g., for MSE when $p \approx Kn$. In fact, in practice, there might be redundancies due to either duplicate or linearly dependent samples as well as parameters. To illustrate this, in Figure 10 we show a simple example of linear regression with redundancies in the design matrix, and where the interpolation threshold can be changed arbitrarily by changing the extent of redundancy. But thinking in terms of the Hessian rank can help avoid such inconsistencies.

Other facets of double descent. (i) Our analysis additionally explains the empirical observation of *label noise* accentuating the peak (Nakkiran et al., 2019), since the complexity term contains a multiplicative factor of σ_{min}^2 which increases proportionately with label noise. (ii) The exact term that appears in our proof, before we take the limit of regularization $\lambda \to 0$, is $\left(\lambda_r(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star)) + \lambda\right)^{-1}$. This reveals why, when using a *regularization* of a suitable magnitude, double descent is not prominent or disappears, as also noted in (Nakkiran et al., 2019; 2020), — since this term can no longer explode.

4.3 Empirical Verification



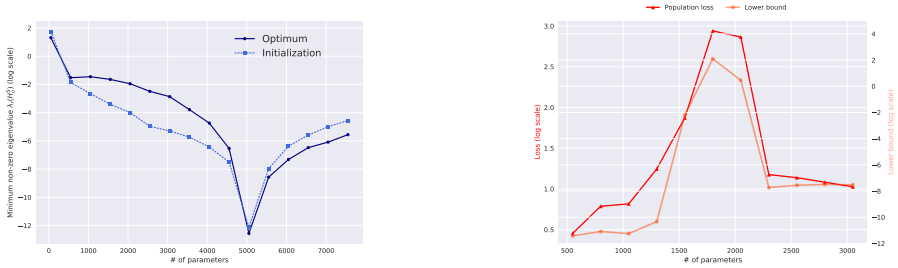
(a) $L = 2$, $n = 500$, $K = 10$ **(b)** $L = 3$, $n = 200$, $K = 10$

Figure 1: MSE Loss: Behaviour of the population (test) loss and minimum non-zero eigenvalues for the setting of $L = 2$ and $L = 3$ layer networks on downscaled MNIST (Greydanus, 2020).

To empirically demonstrate the validity of our theoretical results, we carry out the entire procedure of obtaining double descent for neural networks. Namely, this involves training a large set of neu-

ral networks — with layer widths sampled in regular intervals — for sufficiently many number of epochs. However, in our case, there is another factor which makes this whole process even more arduous — as the lower bound depends on the minimum non-zero eigenvalue of the Hessian. As a result, we cannot resort to efficient Hessian approximations based on, say, Hessian-vector products (Pearlmutter, 1994), but rather we need to compute entire Hessian spectrum — which has $\mathcal{O}(p^3)$ computational and memory costs. Hence, we cannot but restrict our empirical investigation to smaller network sizes. Nevertheless, a positive outcome is that this way we can be thoroughly assure of the accuracy of our empirical investigations — since we compute the exact Hessian and its spectrum, and that too in FLOAT64 precision.

In terms of the dataset, we primarily utilize MNIST1D (Greydanus, 2020), which is a downscaled version of MNIST yet designed to be significantly harder than the usual version. However, we also present results on CIFAR10 and the usual (easier) MNIST, which alongside other empirical details, can be found in Appendix C. Figure 1 shows the results of running the double descent experiments for the settings of two and three layer fully-connected networks with ReLU activation trained for 5K epochs via SGD. Alongside the population loss (empirically measured on a test set) — which peaks at the interpolation threshold of $p \approx KN$ — we plot the trend of the minimum non-zero eigenvalue. As predicted by our theory, this eigenvalue indeed tends to zero (note the log scale), in both the settings, around the precise neighborhood where the population loss takes its maximum value.



(a) Minimum non-zero eigenvalue stays close. **(b)** Lower bound captures the trend of population loss.

Figure 2: Empirical validity of assumption A5 & the lower bound of Theorem 4 in Double Descent.

5 THE CASE OF CROSS-ENTROPY

Greydanus (2020) notes that unlike MSE loss, in case of cross-entropy the population loss peaks at $p \approx n$ (in contrast to $p \approx Kn$ for MSE). To better understand this aspect, let us start by checking the form of the Hessian at the optimum for cross-entropy loss, which by our assumption A2, will be given by the outer-product term. as in eq. 2. Now, the main difference is that the Hessian of the loss with respect to function outputs (named 'output-Hessian'), instead of being identity like in MSE, is given by $\nabla^2_{\boldsymbol{f}}\ell(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}) = \text{diag}(\boldsymbol{p}) - \boldsymbol{p}\boldsymbol{p}^\top$, where $\boldsymbol{p} = \text{softmax}(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}))$ denotes the predicted class-probabilities obtained from applying the softmax operation. In general, this output-Hessian matrix of size $K \times K$, is rank-deficient with rank $K-1$. Thus, the $\text{rank}(\mathbf{H}^S_o(\boldsymbol{\theta}) = \min(p, (K-1)\,n)$ for any $\boldsymbol{\theta}$, and the initial surmise would be that the interpolation threshold is at $p \approx (K-1)\,n$.

But, this is not inline with the stated observation from Greydanus (2020). Hence, let us take another closer look at the Hessian, and in particular, $\nabla^2_{\boldsymbol{f}}\ell$. Notice that near interpolation, when the loss on individual training samples tends to zero, the predicated class-probability $\boldsymbol{p}$ will tend to 1 for the correct class (as per the training label) and 0 elsewhere. This suggests that Hessian matrix collapses to $\mathbf{0}$ since $\nabla^2_{\boldsymbol{f}}\ell(\boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}), \boldsymbol{y}) = \text{diag}(\boldsymbol{p}) - \boldsymbol{p}\boldsymbol{p}^\top \to \mathbf{0}$. And indeed, this is true based on our empirical results, displayed in Figure 3 given the network is trained sufficiently long. We clearly observe that the population loss diverges at $p \approx n$, however the test error, although not as conspicuous, still shows a slight peak (similar to the curves in (Nakkiran et al., 2019) without label-noise). Next, from the right sub-figure (plotted in log-scale), it also becomes evident that the entire Hessian spectrum — from the maximum to the minimum eigenvalue — collapses to zero near the interpolation threshold.

Fact 6. *Thus, we have that tor cross-entropy loss, near the interpolation threshold of $p \approx n$, the entire Hessian matrix vanishes at the optimum, i.e.,* $\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star) = \mathbf{0}$.

Importantly, the above fact (and our empirical results) report such a behaviour only around $p \approx n$, and not well into the over-parameterized regime $p >> n$. Therefore, a straightforward consequence

Figure 3: (Left) Double descent curve for CE loss and **(right)** log-scale plots of the test and train loss, alongside Hessian eigenvalues, for a one-hidden layer neural network trained for 40K epochs.

of the above fact and empirical observations is that $\lambda_r(\mathbf{H}_{\mathcal{L}}^{S}(\boldsymbol{\theta}^\star)) = 0$ near this interpolation threshold, implying that the population loss in case of cross-entropy loss diverges at $p \approx n$.

Vanishing Hessian hypothesis. Figure 3 right, seems to suggest, rather surprisingly, that even the training loss has the same trend as these Hessian eigenvalue statistics. Note, this is not the case that networks were not trained long enough — rather, we run them for $40,000$ epochs. In fact, we even trained models, right to the interpolation threshold up to $80,000$ epochs. Yet, the training loss only drops by factor of 2 and is still $> 1e-5$, compared to exact 0 at machine precision in half the epochs for networks lying in the region where the test loss diverges. This leads us to posit the following hypothesis: *around $p \approx n$, SGD finds critical points with zero Hessian for neural networks trained with CE loss and over-parameterization beyond the interpolation threshold helps completely avoid or significantly decelerate convergence to such critical points.* Further investigation into this hypothesis is beyond the current scope, but forms an exciting direction for future work.

6 Leave-one-out Estimate via Influence Functions

The principle behind leave-one-out (LOO) is to leave behind one sample, optimize the model from scratch on the remaining samples, then evaluate on the left-behind sample, and finally average over the choice of the left-out sample. LOO is known to provide a reasonable estimate of the population loss (Pontil, 2002). In our setting for double descent, we prefer the add-one-in procedure since it directly gives the population loss itself. However, LOO can still be useful from a practical perspective as it relies only on the training set — assuming we can analytically estimate it via some closed-form expression which avoids the need to retrain training set size many models.

Similar to the add-one-in procedure from before, leave-one-out can be cast as a slight contamination of the distribution. We can express the new distribution $\mathcal{D}_{n-1}^{\setminus i}$ with the i-th sample removed as follows, $\mathcal{D}_{n-1}^{\setminus i} = (1-\epsilon)\mathcal{D}_n + \epsilon\,\delta_{\boldsymbol{z}_i}$ with $\epsilon = \frac{-1}{n-1}$, where $\mathcal{D}_n$ refers to the original empirical distribution over the training set. Now, we can carry out similar steps like for add-one-in, and derive the change in the parameter estimate as well as the change in loss over the left-out sample. But, here we additionally analyze the effect of *incorporating the second-order influence function*, apart from the usual first-order influences. This provides us with estimates $\mathrm{LOO}^{(1)}$ and $\mathrm{LOO}^{(2)}$, the expressions of which can be found in Appendix A.5. As a quick test-bed, we investigate the fidelity of these two approaches relative to the exact formula $\mathrm{LOO}^{\mathrm{LS}}$ that exists in the case of least-squares.

Theorem 7. *Consider the particular case with the training inputs gathered into the data matrix $\mathbf{X} \in \mathbb{R}^{n\times d}$ and the targets collected in the vector $\boldsymbol{y} \in \mathbb{R}^n$. Under the assumption that the number of samples n is large enough such that $n \approx n-1$, we have that $\mathrm{LOO}^{(2)} = \mathrm{LOO}^{LS} = \frac{1}{n}\sum_{i=1}^{n}\left(\frac{y_i - \boldsymbol{\theta}^\top \boldsymbol{x}_i}{1-\mathbf{A}_{ii}}\right)^2$, and $\mathrm{LOO}^{(1)} = \frac{1}{n}\sum_{i=1}^{n}\left(y_i - \boldsymbol{\theta}^\top \boldsymbol{x}_i\right)^2 \frac{\mathbf{A}_{ii}}{1-\mathbf{A}_{ii}}$. where, $\mathbf{A}_{ii} = \boldsymbol{x}_i^\top(\mathbf{X}^\top\mathbf{X})^{-1}\boldsymbol{x}_i$ denotes the i-th diagonal entry of the matrix $\mathbf{A} = \mathbf{X}(\mathbf{X}^\top\mathbf{X})^{-1}\mathbf{X}^\top$ (i.e., the so-called 'hat-matrix') and $\boldsymbol{\theta}$ denotes the usual solution of $\boldsymbol{\theta} = (\mathbf{X}^\top\mathbf{X})^{-1}\mathbf{X}^\top\boldsymbol{y}$ obtained via ordinary least-squares. Moreover, it can be shown that $\mathrm{LOO}^{(2)} \geq 4\,\mathrm{LOO}^{(1)}$ for $\mathbf{A}_{ii} \in [0,1]$.*

This result is reassuring since it shows that LOO expressions from influence-based analysis are indeed accurate, and *we recover the least-squares formula as a special case through* $\mathrm{LOO}^{(2)}$. Further,

Corollary 8. *For any finite-width neural network, the first and second-order influence function give similar formulas for LOO like that in Theorem 7, but with* $\mathbf{A} = \mathbf{Z}_S(\boldsymbol{\theta}^\star)^\top \left(\mathbf{Z}_S(\boldsymbol{\theta}^\star)\mathbf{Z}_S(\boldsymbol{\theta}^\star)^\top\right)^{-1} \mathbf{Z}_S(\boldsymbol{\theta}^\star)$, *where* $\mathbf{Z}_S(\boldsymbol{\theta}^\star) := [\nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}_1), \cdots, \nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}_n)]$ *and* $\theta^\star$ *are the parameters at convergence for MSE loss.*

Finally, the above result, in particular $\mathrm{LOO}^{(2)} \geq 4\,\mathrm{LOO}^{(1)}$ raises an important aspect about the current practice of using first-order influence functions in deep learning — which underestimates LOO even for the simplest setting of least squares regression. While this is not necessarily a significant concern for a theoretical analysis, say that of double descent, such constant factors can be relevant from a practical viewpoint. This also seems to provide a theoretical justification for why Basu et al. (2020) find a suitable usage of first-order influence functions needs a small change in the parameters. However, an empirical investigation on this front remains outside the scope of our work.

7 Discussion

Summary. We derived an expression of the population risk via influence functions and obtained a lower bound to it — with fairly minimal assumptions — that applies to any finite-width neural network trained with commonly used loss functions. The lower bound is inversely related to the smallest non-zero eigenvalue of the Hessian of the loss at the optimum. When specialized to the MSE loss, this provably exhibits a double descent behaviour in the asymptotic regime and we empirically demonstrated that this holds even in much smaller non-asymptotic settings. We also analyzed the intriguing phenomenology of double descent across different losses — through our Hessian-based framework — which explained existing empirical observations as well as uncovered novel aspects of neural networks near interpolation. Finally, as a by-product, we presented theoretical results for leave-one-out estimation using influence functions in the case of neural networks.

Related theoretical work on Double Descent. We carve out a niche in the growing set of studies on double descent by focusing primarily on — *finite-width neural networks*. For the linear/kernel regression setting (or lately, the nearly equivalent two-layer network with frozen hidden-layer), there is a plethora of existing work (Advani et al., 2020; Bartlett et al., 2020; Mei & Montanari, 2019; Muthukumar et al., 2019; Geiger et al., 2020; Ba et al., 2020), that analyzes double descent. Thus, our aim is not to make these existing analyses tighter, but rather to take a step towards developing analyses that directly hold for finite-width neural networks. Therefore, unlike the above works, we do not impose any restrictive assumptions on the structure of neural network, like two-layer networks, or the optimization methods used to train them, like gradient flow. Yet, our work also bears a natural connection between the matrix whose spectrum comes to be of concern in the prior works — the input-covariance or kernel matrix in linear or kernel regression — while that of the outer-product Hessian in our work for MSE loss (which has the same spectrum as the 'empirical' NTK). But our strategy also makes our work applicable to cross-entropy, e.g., where we bring to light the interesting observations near the interpolation threshold.

A closely related work, Kuzborskij et al. (2021), links the population risk *for least-squares* to the minimum non-zero eigenvalue of the input covariance matrix — but although via an upper-bound, which is insufficient for illustrating double descent. Nevertheless, using an analogy to their regression result, they study the minimum eigenvalue of the covariance matrix consisting of penultimate-layer features and conjecture this as a possible extension to neural networks. However, the input-covariance matrix in least-squares is also the Hessian, and as we have thoroughly established the Hessian (at the optimum) is indeed the relevant object — thus contradicting their conjecture.

Limitations and directions for future work. There are many important aspects surrounding double descent that remain unanswered, in the context of neural networks: (a) Analogous to (Hastie et al., 2019) for linear regression, what are the conditions for the global optimum to lie in the over-parameterized regime instead of under-parameterized? (b) Given the vanishing Hessian hypothesis, is there a qualification to the regime where the flat-minima generalizes better hypothesis (Hochreiter & Schmidhuber, 1997; Keskar et al., 2016) holds — since the Hessian is the flattest possible here. (d) On the technical side: better characterization of the mentioned cross-entropy phenomenon as well as non-asymptotic results. Overall, we hope that our work will encourage foray into further studies of double descent, that are specifically built for finite-width neural networks.

REPRODUCIBILITY STATEMENT

- All the omitted proofs to the theoretical results can be found in the Appendix A.
- In regards to empirical results, we provide all the relevant details and additional results in the Appendix C.
- The corresponding code for the experiments will also be released shortly.

REFERENCES

Madhu S Advani, Andrew M Saxe, and Haim Sompolinsky. High-dimensional dynamics of generalization error in neural networks. *Neural Networks*, 132:428–446, 2020.

Jimmy Ba, Murat Erdogdu, Taiji Suzuki, Denny Wu, and Tianzong Zhang. Generalization of two-layer neural networks: An asymptotic viewpoint. In *International Conference on Learning Representations*, 2020. URL https://openreview.net/forum?id=H1gBsgBYwH.

Peter L Bartlett, Philip M Long, Gábor Lugosi, and Alexander Tsigler. Benign overfitting in linear regression. *Proceedings of the National Academy of Sciences*, 117(48):30063–30070, 2020.

Samyadeep Basu, Philip Pope, and Soheil Feizi. Influence functions in deep learning are fragile, 2020.

Mikhail Belkin, Daniel Hsu, Siyuan Ma, and Soumik Mandal. Reconciling modern machine-learning practice and the classical bias–variance trade-off. *Proceedings of the National Academy of Sciences*, 116(32):15849–15854, 2019. ISSN 0027-8424. doi: 10.1073/pnas.1903070116. URL https://www.pnas.org/content/116/32/15849.

Michiel Debruyne, Mia Hubert, and Johan A.K. Suykens. Model selection in kernel based regression using the influence function. *Journal of Machine Learning Research*, 9(78):2377–2400, 2008. URL http://jmlr.org/papers/v9/debruyne08a.html.

Alexandre Défossez and Francis Bach. Averaged least-mean-squares: Bias-variance trade-offs and optimal sampling distributions. In *Artificial Intelligence and Statistics*, pp. 205–213. PMLR, 2015.

Bradley Efron and Charles Stein. The jackknife estimate of variance. *The Annals of Statistics*, pp. 586–596, 1981.

Yuguang Fang, Kenneth A Loparo, and Xiangbo Feng. Inequalities for the trace of matrix product. *IEEE Transactions on Automatic Control*, 39(12):2489–2490, 1994.

Nicolas Flammarion and Francis Bach. From averaging to acceleration, there is only a step-size, 2015.

Mario Geiger, Arthur Jacot, Stefano Spigler, Franck Gabriel, Levent Sagun, Stéphane d'Ascoli, Giulio Biroli, Clément Hongler, and Matthieu Wyart. Scaling description of generalization with number of parameters in deep learning. *Journal of Statistical Mechanics: Theory and Experiment*, 2020(2):023401, 2020.

Sam Greydanus. Scaling down deep learning, 2020.

Frank R Hampel, Elvezio M Ronchetti, Peter J Rousseeuw, and Werner A Stahel. *Robust statistics: the approach based on influence functions*, volume 196. John Wiley & Sons, 1986.

Trevor Hastie, Andrea Montanari, Saharon Rosset, and Ryan J Tibshirani. Surprises in high-dimensional ridgeless least squares interpolation. *arXiv preprint arXiv:1903.08560*, 2019.

Sepp Hochreiter and Jürgen Schmidhuber. Flat minima. *Neural computation*, 9(1):1–42, 1997.

Peter J Huber. *Robust statistics*, volume 523. John Wiley & Sons, 2004.

Arthur Jacot, Franck Gabriel, and Clément Hongler. Neural tangent kernel: Convergence and generalization in neural networks, 2018.

Nitish Shirish Keskar, Dheevatsa Mudigere, Jorge Nocedal, Mikhail Smelyanskiy, and Ping Tak Peter Tang. On large-batch training for deep learning: Generalization gap and sharp minima. *arXiv preprint arXiv:1609.04836*, 2016.

Pang Wei Koh and Percy Liang. Understanding black-box predictions via influence functions, 2017.

Ilja Kuzborskij, Csaba Szepesvári, Omar Rivasplata, Amal Rannen-Triki, and Razvan Pascanu. On the role of optimization in double descent: A least squares study. *arXiv preprint arXiv:2107.12685*, 2021.

Marco Loog, Tom Viering, Alexander Mey, Jesse H. Krijthe, and David M. J. Tax. A brief prehistory of double descent. *Proceedings of the National Academy of Sciences*, 117(20):10625–10626, May 2020. ISSN 1091-6490. doi: 10.1073/pnas.2001875117. URL http://dx.doi.org/10.1073/pnas.2001875117.

Song Mei and Andrea Montanari. The generalization error of random features regression: Precise asymptotics and the double descent curve. *Communications on Pure and Applied Mathematics*, 2019.

N. Murata, S. Yoshizawa, and S. Amari. Network information criterion-determining the number of hidden units for an artificial neural network model. *IEEE Transactions on Neural Networks*, 5(6): 865–872, 1994. doi: 10.1109/72.329683.

Vidya Muthukumar, Kailas Vodrahalli, and Anant Sahai. Harmless interpolation of noisy data in regression. *2019 IEEE International Symposium on Information Theory (ISIT)*, Jul 2019. doi: 10.1109/isit.2019.8849614. URL http://dx.doi.org/10.1109/ISIT.2019.8849614.

Preetum Nakkiran. More data can hurt for linear regression: Sample-wise double descent. *arXiv preprint arXiv:1912.07242*, 2019.

Preetum Nakkiran, Gal Kaplun, Yamini Bansal, Tristan Yang, Boaz Barak, and Ilya Sutskever. Deep double descent: Where bigger models and more data hurt, 2019.

Preetum Nakkiran, Prayaag Venkat, Sham Kakade, and Tengyu Ma. Optimal regularization can mitigate double descent, 2020.

Barak A Pearlmutter. Fast exact multiplication by the hessian. *Neural computation*, 6(1):147–160, 1994.

Jeffrey Pennington and Yasaman Bahri. Geometry of neural network loss surfaces via random matrix theory. In *International Conference on Machine Learning*, pp. 2798–2806. PMLR, 2017.

Loucas Pillaud-Vivien, Alessandro Rudi, and Francis Bach. Exponential convergence of testing error for stochastic gradient methods. In *Conference on Learning Theory*, pp. 250–296. PMLR, 2018.

Massimiliano Pontil. Leave-one-out error and stability of learning algorithms with applications. *International Journal of Systems Science*, 2002.

Levent Sagun, Leon Bottou, and Yann LeCun. Eigenvalues of the hessian in deep learning: Singularity and beyond, 2017.

Sidak Pal Singh, Gregor Bachmann, and Thomas Hofmann. Analytic insights into structure and rank of neural network hessian maps, 2021.

Mervyn Stone. An asymptotic equivalence of choice of model by cross-validation and akaike's criterion. *Journal of the Royal Statistical Society: Series B (Methodological)*, 39(1):44–47, 1977.

Kei Takeuchi. The distribution of information statistics and the criterion of goodness of fit of models. *Mathematical Science*, 153:12–18, 1976.

Valentin Thomas, Fabian Pedregosa, Bart van Merriënboer, Pierre-Antoine Mangazol, Yoshua Bengio, and Nicolas Le Roux. Information matrices and generalization. *arXiv preprint arXiv:1906.07774*, 2019.

Roman Vershynin. Introduction to the non-asymptotic analysis of random matrices. *arXiv preprint arXiv:1011.3027*, 2010.

A OMITTED PROOFS

Proposition 9. *The influence function IF of the M-estimator $\widehat{\boldsymbol{\theta}}_{\mathcal{D}}$ based on the distribution $\mathcal{D}$, evaluated at point $\boldsymbol{z}$, takes the following form:*

$$\mathrm{IF}(\boldsymbol{z};\widehat{\boldsymbol{\theta}}_{\mathcal{D}},\mathcal{D}) = -\left[\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}})\right]^{-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{D})\,, \tag{4}$$

where, the Hessian matrix $\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}}) := \nabla^2_{\boldsymbol{\theta}}\,\mathcal{L}(\widehat{\boldsymbol{\theta}})$ is the matrix of second-derivatives of the loss $\mathcal{L}(\boldsymbol{\theta}) := \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})]$ with respect to the parameters $\boldsymbol{\theta}$.

Proof. $\widehat{\boldsymbol{\theta}}_{\mathcal{D}} = \arg\min_{\boldsymbol{\theta}\in\boldsymbol{\Theta}} \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}} \boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})$. But, instead of this formulation, we can also define the estimator as one that satisfies the following implicit equation (assuming the derivative can be moved inside expectation),

$$\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] = 0\,, \tag{5}$$

which is nothing but the first-order stationary point condition. Now, consider a contaminated distribution $\widetilde{\mathcal{D}} := (1-\epsilon)\mathcal{D} + \epsilon\,\delta_{\boldsymbol{z}}$. The corresponding implicit equation, to be satisfied by the estimator $\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}$ corresponding to $\widetilde{\widehat{\boldsymbol{\theta}}}$, can be written as follows:

$$\begin{aligned} \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\widetilde{\widehat{\boldsymbol{\theta}}}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] &= 0 \\ (1-\epsilon)\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] + \epsilon\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta}) &= 0 \end{aligned} \tag{6}$$

Take the derivative of the above expression with respect to ϵ (plus interchanging derivative and expectation), we get:

$$\begin{aligned} \frac{d}{d\epsilon}(1-\epsilon)\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] &= -\frac{d}{d\epsilon}\epsilon\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta}) \\ -\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] + (1-\epsilon)\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})]\frac{d\boldsymbol{\theta}}{d\epsilon} &= -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta}) - \epsilon\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})\frac{d\boldsymbol{\theta}}{d\epsilon} \end{aligned} \tag{7}$$

Since, $\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}$ satisfies the above eq. 7, let us substitute it in place of $\boldsymbol{\theta}$ and analyze the case for $\epsilon \to 0$. We are left with the following (after removing terms multiplied with ϵ):

$$-\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}})] + \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}})]\frac{d\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}}{d\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}) \tag{8}$$

Now, the first term goes to zero as $\epsilon \to 0$, because eq. 6 holds for $\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}$, as shown below :

$$\begin{aligned} -\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}})] &= -\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}})] + \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\widetilde{\mathcal{D}}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}})] \\ &= \epsilon\left(\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_0,\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}) - \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}})]\right)\,. \end{aligned}$$

Further, as $\epsilon \to 0$, one can replace $\widehat{\boldsymbol{\theta}}_{\widetilde{\mathcal{D}}}$ by $\widehat{\boldsymbol{\theta}}_{\mathcal{D}}$ in the expressions of the gradient and Hessian of $\boldsymbol{\ell}$. Then, assuming that the Hessian $\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}}) := \nabla^2_{\boldsymbol{\theta}}\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] = \mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}\left[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\mathcal{D}})\right]$ is invertible, this yields:

$$\mathrm{IF}(\boldsymbol{z};\widehat{\boldsymbol{\theta}}_{\mathcal{D}},\mathcal{D}) = -\left[\mathbf{H}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_{\mathcal{D}})\right]^{-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{D})\,, \tag{9}$$

□

A.1 EXPRESSION OF THE POPULATION RISK

Theorem 10. *Consider the parameter estimator $\widehat{\boldsymbol{\theta}}_S$ based on the set of input samples S of $|S| = n$. Then the population risk $\mathcal{L}(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}\sim D}\left[\boldsymbol{\ell}(\boldsymbol{z}, \widehat{\boldsymbol{\theta}}_S)\right])$ takes the following form,*

$$\mathcal{L}(\widehat{\boldsymbol{\theta}}_S) \;=\; \widetilde{\mathcal{L}}_S(\widehat{\boldsymbol{\theta}}_S) \;+\; \frac{1}{n+1}\, Tr\left(\left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda \mathbf{I}\right]^{-1} \mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S)\right), \tag{10}$$

where $\widetilde{\mathcal{L}}_S(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}\left(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}}\right)\right]$ denotes the expectation of 'one-sample training loss' and $\mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S)\, \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S)^\top\right]$ is the (uncentered) covariance of loss gradients.

Proof. Let us recall the expression of influence function for the loss on the new sample $\boldsymbol{z}'$,

$$\mathrm{IF}_\lambda(\boldsymbol{\ell}_{\boldsymbol{z}'}; \widehat{\boldsymbol{\theta}}_S, \mathcal{D}_n) \;=\; -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)^\top \left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1} \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)\,. \tag{11}$$

When $|S| = n$ is large enough to ignore $\mathcal{O}(n^{-2})$ terms, the (infinitesimal) definition of influence function is equivalent to using the finite-difference form. Then the change in loss can be expressed as,

$$\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}}\big) - \boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big) \;=\; \frac{-1}{n+1}\, \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)^\top \left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1} \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big), \tag{12}$$

where we have multiplied both sides by $\Delta\epsilon = \epsilon - 0 = \frac{1}{n+1}$. We leverage this relation to derive an expression of the test loss as follows. Starting from eq. 12, we average out over the choice of an additional sample $\boldsymbol{z}' \sim \mathcal{D}$:

$$\mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] \;-\; \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)\right] \;=$$
$$\frac{-1}{(n+1)} \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)^\top \left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1} \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)\right].$$

Using properties of the trace and moving expectation inside, we get

$$\mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}\big(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_S\big)\right] \;=\; \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] \;+\; \frac{1}{n+1}\,\mathrm{Tr}\left(\left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) + \lambda\mathbf{I}\right]^{-1} \mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S)\right), \tag{13}$$

where, $\mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\boldsymbol{\theta}) := \mathbb{E}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}', \boldsymbol{\theta})\, \nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}', \boldsymbol{\theta})^\top\right]$.

Now the term on the left-hand side is nothing but the population risk $\mathcal{L}(\widehat{\boldsymbol{\theta}}_S) := \mathbb{E}_{\boldsymbol{z}\sim D}\left[\boldsymbol{\ell}(\boldsymbol{z}, \widehat{\boldsymbol{\theta}}_S)\right])$, and we coin the first term on the right-hand side as the expectation of 'one-sample training loss'.

□

A.1.1 INTERPRETATION OF THE 'ONE-SAMPLE TRAINING LOSS'

First, note that the first term on the RHS of eq. 13 is not exactly the training loss but rather some related quantity. To see this better, let us rewrite as follows:

$$\mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] = \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\underbrace{\left[\boldsymbol{\ell}(\boldsymbol{z}', \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}}) - \frac{1}{|S|}\boldsymbol{\ell}(S, \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right]}_{\Delta\mathrm{TR}(\boldsymbol{z}', S\cup\{\boldsymbol{z}'\})} + \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\frac{1}{|S|}\boldsymbol{\ell}(S, \widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right],$$

where, $\boldsymbol{\ell}(S, \boldsymbol{\theta}) = \sum_{i=1}^{|S|} \boldsymbol{\ell}(\boldsymbol{z}_i, \boldsymbol{\theta})$ is the sum of the loss over the samples in S. Further, the expression $\Delta\mathrm{TR}(\boldsymbol{z}', S \cup \{\boldsymbol{z}'\})$ refers to the deviation between the loss of a training sample (here, $\boldsymbol{z}'$) relative

to the average loss on rest of the training samples. The 'TR' in this symbol refers to this deviation being computed on the given *training* set.

$$\begin{aligned}\mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] &= \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}[\Delta\text{TR}(\boldsymbol{z}',S\cup\{\boldsymbol{z}'\})] + \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\frac{1}{|S|}\boldsymbol{\ell}(S,\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] \\ &= \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}[\Delta\text{TR}(\boldsymbol{z}',S\cup\{\boldsymbol{z}'\})] + \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\frac{1}{|S|}\left(\boldsymbol{\ell}(S\cup\{\boldsymbol{z}'\},\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}}) - \boldsymbol{\ell}(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right)\right]\end{aligned}$$

Notice, the last term in the expression above is the same as the term on the left hand side — albeit with an additional scaling factor of $\frac{1}{|S|}$ and negative sign in front. Rearranging this results in the following equation:

$$\frac{|S|+1}{|S|}\mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] = \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}[\Delta\text{TR}(\boldsymbol{z}',S\cup\{\boldsymbol{z}'\})] + \frac{|S|+1}{|S|}\mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\frac{1}{|S\cup\{\boldsymbol{z}'\}|}\boldsymbol{\ell}(S\cup\{\boldsymbol{z}'\},\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right]$$

The only extra thing we have done is to multiply and divide by $|S\cup\{\boldsymbol{z}'\}| = |S|+1$ in the last term in the right hand side. Notice the rightmost term is an expectation (over $\boldsymbol{z}'$) of the average training loss of the training set $S\cup\{\boldsymbol{z}'\}$ and to which we assign the shorthand $\text{TR}(S\cup\{\boldsymbol{z}'\})$. Also, it is evident from here that the 'one-sample training loss' is a quantity very much related to the training loss. Lastly, considering the large $|S| = n$ limit, we have $\lim_{n\to\infty}\frac{n+1}{n} = 1$, and which thereby yields:

$$\widetilde{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S) := \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}\left[\boldsymbol{\ell}(\boldsymbol{z}',\widehat{\boldsymbol{\theta}}_{S\cup\{\boldsymbol{z}'\}})\right] = \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}[\Delta\text{TR}(\boldsymbol{z}',S\cup\{\boldsymbol{z}'\})] + \mathop{\mathbb{E}}_{\boldsymbol{z}'\sim\mathcal{D}}[\text{TR}(S\cup\{\boldsymbol{z}'\})]$$

A.2 Lower bound to the population risk

Lemma 11. *Consider two matrices $\mathbf{A}\in\mathbb{R}^{m\times m}$ and $\mathbf{B}\in\mathbb{R}^{m\times m}$, where $\mathbf{A}$ is symmetric and $\mathbf{B}$ is symmetric and positive semi-definite. Then the following holds,*

$$\lambda_{min}(\mathbf{A})Tr(\mathbf{B}) \;\le\; Tr(\mathbf{AB}) \;\le\; \lambda_{max}(\mathbf{A})Tr(\mathbf{B})$$

Proof. See Fang et al. (1994). □

Theorem 12. *Under the assumptions, A1, A2, A3, A4, A5 and under the limit of external regularization $\lambda\to 0$, we obtain the following lower bound on the population risk stated in eq. 1,*

$$\mathcal{L}(\boldsymbol{\theta}^\star) \;\ge\; \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) \;+\; \frac{1}{n+1}\cdot\frac{\sigma^2_{min}\;A_{\boldsymbol{\theta}^0}}{\lambda_r\left(\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star)\right)}\;. \tag{14}$$

where we consider the convention $\lambda_1\ge\cdots\ge\lambda_r$ for the eigenvalues with $r := \text{rank}(\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star))$, and σ^2_{min} denotes the minimum non-zero $\sigma^2_{(\boldsymbol{x},\boldsymbol{y})}$ over $(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}$.

Proof. Let us recall the expression for population risk that we proved in Theorem 3, for $\widehat{\boldsymbol{\theta}}_S = \boldsymbol{\theta}^\star$,

$$\mathcal{L}(\boldsymbol{\theta}^\star) \;=\; \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) \;+\; \frac{1}{n+1}\,\text{Tr}\left(\left[\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star)+\lambda\mathbf{I}\right]^{-1}\mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\boldsymbol{\theta}^\star)\right) \;+\; \mathcal{O}\left(\frac{1}{n^2}\right)\;, \tag{15}$$

In particular, we would like to analyze the complexity term $T := \text{Tr}\left(\left[\mathbf{H}^S_{\mathcal{L}}(\widehat{\boldsymbol{\theta}}_S)+\lambda\mathbf{I}\right]^{-1}\mathbf{C}^{\mathcal{D}}_{\mathcal{L}}(\boldsymbol{\theta}^\star)\right)$ on the right-hand side and lower bound it. The full expression of this term T is given by,

$$T = \mathop{\mathbb{E}}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}}\left[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}),\boldsymbol{y})^\top\left[\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star)+\lambda\mathbf{I}\right]^{-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}),\boldsymbol{y})\right] \tag{16}$$

Now using the chain rule, we have that $\nabla_{\boldsymbol{\theta}} \ell(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}) = \nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}) \nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)$. Then the above equation is equivalent to,

$$T = \mathop{\mathbb{E}}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}} \left[\mathrm{Tr} \left(\underbrace{\nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x})^\top \left[\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I}\right]^{-1} \nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x})}_{\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}} \cdot \underbrace{\nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big) \nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)^\top}_{\mathbf{B}_{(\boldsymbol{x},\boldsymbol{y})}} \right) \right] \tag{17}$$

where, we have just rearranged the terms using the cyclic property of the trace. Now, we have two symmetric matrices inside the trace, $\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})} \in \mathbb{R}^{K\times K}$, $\mathbf{B}_{(\boldsymbol{x},\boldsymbol{y})} \in \mathbb{R}^{K\times K}$, where $\mathbf{B}_{(\boldsymbol{x},\boldsymbol{y})}$ is trivially positive semi-definite. Also, the matrix $\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}$ is positive semi-definite since by assumptions A1, A2, the Hessian at the optimum $\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) = \mathbf{H}_o^S(\boldsymbol{\theta}^\star)$ and $\mathbf{H}_o^S(\boldsymbol{\theta}^\star)$ is positive semi-definite and $\lambda > 0$.

Next, using the Lemma 11, we get the following lower-bound for T:

$$\begin{aligned}
T &\geq \mathop{\mathbb{E}}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}} \left[\mathrm{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right) \cdot \lambda_{\min} \left(\nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big) \nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)^\top \right) \right] \\
&= \mathop{\mathbb{E}}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}} \left[\mathrm{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right) \cdot \mathrm{Tr} \left(\nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big) \nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)^\top \right) \right] \\
&= \mathop{\mathbb{E}}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}} \left[\mathrm{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right) \cdot \|\nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)\|^2 \right] \\
&\geq \sigma_{\min}^2 \mathop{\mathbb{E}}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}} \left[\mathrm{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right) \right] \\
&= \sigma_{\min}^2 \mathrm{Tr} \left(\left[\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I}\right]^{-1} \mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^\star) \right)
\end{aligned} \tag{18}$$

In the second line, we have used the fact that the trace of the rank-1 matrix $\mathbf{B}_{(\boldsymbol{x},\boldsymbol{y})} = \nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big) \nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)^\top$ will be the same as its non-zero eigenvalue. In the fourth line, we take the lower bound by considering the minimum over all non-zero $\sigma_{(\boldsymbol{x},\boldsymbol{y})}^2 = \|\nabla_{\boldsymbol{f}} \ell\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)\|^2$, the existence of which is guaranteed by assumption A3, i.e.,

$$\sigma_{\min}^2 = \min_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}:\, \sigma_{(\boldsymbol{x},\boldsymbol{y})}^2 > 0} \sigma_{(\boldsymbol{x},\boldsymbol{y})}^2$$

Finally, in the last line, we use the cyclic property of the trace once again and move the expectation inside the trace, obtaining the covariance of function gradients $\mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^\star)$.

Proceeding further, we again make use of Lemma 11 in the eq. 18, since these matrices are also symmetric positive semi-definite. This yields,

$$T \geq \sigma_{\min}^2 \; \lambda_{\min} \left(\mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^\star) \right) \; \mathrm{Tr} \left(\left[\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I}\right]^{-1} \right) \tag{19}$$

Next, we use the assumption A5 to lower bound the minimum non-zero eigenvalue of the covariance of function gradients at the optimum by the corresponding one at the initialization.

$$T \geq \sigma_{\min}^2 \; A_{\boldsymbol{\theta}^0} \, \lambda_{\min} \left(\mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^0) \right) \; \mathrm{Tr} \left(\left[\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I}\right]^{-1} \right) \tag{20}$$

However, by assumption A4, the columns of $\mathbf{Z}_S := [\nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^0}(\boldsymbol{x}_1) \cdots \nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^0}(\boldsymbol{x}_{|S|})]^\top \in \mathbb{R}^{p\times K|S|}$ are isotropic vectors in $\mathbb{R}^p$. This means that $\mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^0) = \mathbb{E}_{(\boldsymbol{x},\boldsymbol{y})\sim\mathcal{D}} \left[\nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}}^0(\boldsymbol{x}) \nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}}^0(\boldsymbol{x})^\top \right] = \mathbf{I}$ (identity matrix) and thus $\lambda_{\min} \left(\mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^0) \right) = 1$, which results in the following lower bound,

$$
\begin{aligned}
T &\geq \sigma_{\min}^2 \; A_{\boldsymbol{\theta}^0} \operatorname{Tr}\left(\left[\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I}\right]^{-1}\right) \\
&= \sigma_{\min}^2 \; A_{\boldsymbol{\theta}^0} \sum_{i=1}^{p} \frac{1}{\lambda_i\left(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star)\right) + \lambda}
\end{aligned}
$$

where, the notation $\lambda_i(\cdot)$ denotes the i-th eigenvalue of the corresponding matrix and we will use the convention that for some matrix in $\mathbb{R}^{m \times m}$,

$$
\lambda_1 \geq \cdots \geq \lambda_m \, .
$$

Let us suppose r denotes the rank of the Hessian at the optimum, i.e., $r := \operatorname{rank}(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^0))$. Since the Hessian at the optimum is positive semi-definite, all the eigenvalues are non-negative and we can further lower bound the previous expression to T as follows:

$$
\begin{aligned}
T &\geq \sigma_{\min}^2 \; A_{\boldsymbol{\theta}^0} \sum_{i=1}^{r} \frac{1}{\lambda_i\left(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star)\right) + \lambda} \\
&\geq \sigma_{\min}^2 \; A_{\boldsymbol{\theta}^0} \frac{1}{\lambda_r\left(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star)\right) + \lambda}
\end{aligned}
$$

where in the second line, we have used the fact that a sum of non-negative numbers can be lower bounded by the maximum summand. The maximum here will correspond to using the inverse of the minimum non-zero eigenvalue λ_r.

Now, substituting the following lower bound together with expression of population risk and taking the limit of $\lambda \to 0$, finishes the proof.

$$
\mathcal{L}(\boldsymbol{\theta}^\star) \; \geq \; \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) \; + \; \frac{1}{n+1} \cdot \frac{\sigma_{\min}^2 \; A_{\boldsymbol{\theta}^0}}{\lambda_r\left(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star)\right)} \; . \tag{21}
$$

□

Note. For the purposes of this lower bound, we ignored the $\mathcal{O}\left(\frac{1}{n^2}\right)$ part in Theorem 3. This is because they take the form, e.g. for second-order influences as shown in Section A.5, and by the assumptions A1 and A2, they are non-negative since the Hessian at the optimum is positive semi-definite.

A.2.1 Empirical approximations of TIC like expressions

Consider the case when both the Hessian and covariance in the complexity term that shows up in our lower-bound expression,

$$
\operatorname{Tr}\left(\left[\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I}\right]^{-1} \mathbf{C}_{\boldsymbol{f}}^{\mathcal{D}}(\boldsymbol{\theta}^\star)\right)
$$

are based/approximated on the training set.

Let us further assume the case of MSE loss and that we are at the optimum, where by A2, $\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) = \mathbf{C}_{\boldsymbol{f}}^S(\theta^\star)$. Under these set of assumptions, the term above reduces to,

$$
\operatorname{Tr}\left((\mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}^\star) + \lambda \mathbf{I})^{-1} \mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}^\star)\right)
$$

Since we can always express a positive semi-definite matrix as some $\mathbf{Z}\mathbf{Z}^\top$. Let us substitute this in the above expression and consider the limit of regularization $\lambda \to 0$.

$$\lim_{\lambda \to 0} \text{Tr}\left((\mathbf{Z}\mathbf{Z}^\top + \lambda \mathbf{I})^{-1}\mathbf{Z}\mathbf{Z}^\top\right) = \text{Tr}\left(\mathbf{Z}^\dagger \mathbf{Z}\right) \tag{22}$$

where in the last line we have used the result from Tikhonov regularization, and $\mathbf{Z}^\dagger$ denotes the pseudo-inverse of $\mathbf{Z}$. But this expression is nothing but the $\text{rank}(\mathbf{Z})$ and thus shows the ineffectiveness of such an analysis where both the Hessian and the covariance of function gradients are based on the training set.

A.3 Double descent behaviour

Corollary 13. *For MSE loss, under the setting of Theorem 4 and the additional assumption A6, the population risk takes the following form and diverges almost surely to ∞ at $p = \frac{1}{\sqrt{c}}Kn$,*

$$\mathcal{L}(\boldsymbol{\theta}^\star) \geq \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) + \frac{\sigma_{min}^2 \; A_{\boldsymbol{\theta}^0}}{B_{\boldsymbol{\theta}^0} \left(\sqrt{n} - \frac{c}{\sqrt{K}}\sqrt{p}\right)^2}, \quad a.s.$$

in the asymptotic regime of $p, n \to \infty$ but their ratio is a fixed constant, and where $c > 0$ is a constant that depends only on the sub-Gaussian norm of columns of $\mathbf{Z}_S(\boldsymbol{\theta}^0)$.

Proof. Let us start from the lower bound shown in Theorem 4.

$$\mathcal{L}(\boldsymbol{\theta}^\star) \geq \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) + \frac{1}{n+1} \cdot \frac{\sigma_{\text{min}}^2 \; A_{\boldsymbol{\theta}^0}}{\lambda_r\left(\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star)\right)}. \tag{23}$$

For MSE loss, we have the Hessian $\mathbf{H}_{\mathcal{L}}^S(\boldsymbol{\theta}^\star) = \mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}^\star)$. We then use the assumption A5 to bound the minimum non-zero eigenvalue (here, λ_r) of $\mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}^\star)$ to the corresponding minimum non-zero eigenvalue at initialization. This results in the following lower bound,

$$\mathcal{L}(\boldsymbol{\theta}^\star) \geq \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) + \frac{1}{n+1} \cdot \frac{\sigma_{\text{min}}^2 \; A_{\boldsymbol{\theta}^0}}{B_{\boldsymbol{\theta}^0} \; \lambda_r\left(\mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}^0)\right)}. \tag{24}$$

Now notice that, via assumptions A4 and A6, $\mathbf{C}_{\boldsymbol{f}}^S(\boldsymbol{\theta}^0) = \frac{1}{|S|}\mathbf{Z}_S\mathbf{Z}_S^\top$ is a covariance matrix with the columns of $\mathbf{Z}_S$ containing independent, sub-gaussian random vectors in $\mathbb{R}^p$.

Thus, we can leverage the results of Vershynin (2010) on the extremal eigenvalues of covariance matrices. Specifically, given a random matrix, $\mathbf{Z} \in \mathbb{R}^{m \times n}$ whose columns are independent, sub-gaussian random vectors in $\mathbb{R}^m$, Vershynin (2010) states that the extremal eigenvalues of $\frac{1}{n}\mathbf{Z}\mathbf{Z}^\top$, in the asymptotic regime where $m, n \to \infty$ but their ratio $\frac{m}{n} \to \gamma \in (0, 1]$, we have:

$$\lambda_{\text{min}}\left(\frac{1}{n}\mathbf{Z}\mathbf{Z}^\top\right) \to \left(1 - c\sqrt{\frac{m}{n}}\right)^2 \quad \text{and} \quad \lambda_{\text{max}}\left(\frac{1}{n}\mathbf{Z}\mathbf{Z}^\top\right) \to \left(1 + c\sqrt{\frac{m}{n}}\right)^2 \quad \text{a.s.}, \tag{25}$$

where c is a constant that depends on the subgaussian norm.

Applying this result to our case, for $m = p$ and $n = Kn$. In the asymptotic regime where $n \to \infty$, we can safely consider the $n + 1$ in the denominator of the left fraction to be simply n.

$$\mathcal{L}(\boldsymbol{\theta}^\star) \geq \widetilde{\mathcal{L}}_S(\boldsymbol{\theta}^\star) + \frac{1}{n} \cdot \frac{\sigma_{\text{min}}^2 \; A_{\boldsymbol{\theta}^0}}{B_{\boldsymbol{\theta}^0} \; \left(1 - c\sqrt{\frac{p}{Kn}}\right)^2}. \tag{26}$$

Finally, by a simple rearrangement, we recover our desired lower bound, thus finishing the proof.

□

A.4 LEAVE-ONE-OUT DERIVATION

Influence calculation Let us recall the equation which the estimator $\widehat{\boldsymbol{\theta}}_{\tilde{\mathcal{D}}}$ should satisfy:

$$-\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})] + (1-\epsilon)\mathop{\mathbb{E}}_{\boldsymbol{z}\sim\mathcal{D}}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\boldsymbol{\theta})]\frac{d\boldsymbol{\theta}}{d\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_0,\boldsymbol{\theta}) - \epsilon\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_0,\boldsymbol{\theta})\frac{d\boldsymbol{\theta}}{d\epsilon} \tag{27}$$

Here, $G := (1-\epsilon)F + \epsilon\delta_{\boldsymbol{z}_0}$. For leave-one-out, we have that $G = D_{n-1}^{\setminus i}$ and $F = D_n$. Without loss of generality, assume that the removed sampled index $i = n$ and so $\boldsymbol{z}_i = \boldsymbol{z}_n$, and consider the shorthand $\tilde{D} = D_{n-1}^{\setminus n}$. Overall, then we are considering the contamination: $\tilde{D} := (1-\epsilon)D_n + \epsilon\delta_{\boldsymbol{z}_n}$, with $\epsilon = \frac{-1}{n-1}$. Let us substitute all of this back into the equation above:

$$-\mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})] + (1-\epsilon)\mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})]\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) - \epsilon\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}})\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} \tag{28}$$

$$-\mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})] + \mathop{\mathbb{E}}_{\boldsymbol{z}\sim \tilde{D}}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})]\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) \tag{29}$$

$$\begin{aligned}
&-\mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})] + \mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})]\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) - \epsilon\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}})\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} + \epsilon\mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})]\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} \\
\implies &-\mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})] + \mathop{\mathbb{E}}_{\boldsymbol{z}\sim D_n}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})]\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) + \epsilon\mathop{\mathbb{E}}_{\boldsymbol{z}\sim \tilde{D}}[\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z},\widehat{\boldsymbol{\theta}}_{\tilde{D}})]\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon}
\end{aligned}$$

The last equation holds because notice that in the second term $\boldsymbol{z} \sim \widehat{\boldsymbol{\theta}}_{\tilde{D}}$ (where $\tilde{D} = (1-\epsilon)D_n + \epsilon\delta_{\boldsymbol{z}_n}$) and not $\boldsymbol{z} \sim D_n$. Since, both D_n and $\tilde{D}$ are empirical distributions, we can compute the expectation as a finite sum,

$$-\sum_{i=1}^{n}\frac{1}{n}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_i,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) + \left(\frac{1}{n-1}\sum_{i=1}^{n-1}\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_i,\widehat{\boldsymbol{\theta}}_{\tilde{D}})\right)\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) \tag{30}$$

We can split the term involving gradient of the loss into two parts: one based on the $n-1$ samples (of which $\widehat{\boldsymbol{\theta}}_{\tilde{D}}$ is also the parameter estimator) and the other based on the left-out sample with index n.

$$\underbrace{-\frac{n-1}{n}\sum_{i=1}^{n-1}\frac{1}{n-1}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_i,\widehat{\boldsymbol{\theta}}_{\tilde{D}})}_{=0} - \frac{1}{n}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) + \left(\frac{1}{n-1}\sum_{i=1}^{n-1}\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_i,\widehat{\boldsymbol{\theta}}_{\tilde{D}})\right)\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}}) \tag{31}$$

The first term is zero since $\widehat{\boldsymbol{\theta}}_{\tilde{D}}$ is the parameter estimator of the first $n-1$ samples, and will satisfy the first-order stationarity conditions. Now, rearranging terms results in:

$$\left(\frac{1}{n-1}\sum_{i=1}^{n-1}\nabla^2_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_i,\widehat{\boldsymbol{\theta}}_{\tilde{D}})\right)\frac{\Delta\boldsymbol{\theta}}{\Delta\epsilon} = -\frac{n-1}{n}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n,\widehat{\boldsymbol{\theta}}_{\tilde{D}})$$

The term on the left is nothing but the Hessian computed over $\tilde{D}$ or in other words, the first $n-1$ samples. Let us denote it by $\mathbf{H}_{\mathcal{L}}^{\setminus n}(\widehat{\boldsymbol{\theta}}_{\tilde{D}})$, where the $\setminus n$ in the superscript emphasizes that the Hessian is computed over samples excluding the n-th sample. Further, $\Delta\epsilon = \epsilon - 0 = \epsilon = -\frac{1}{n-1}$. Also,

$\Delta\boldsymbol{\theta} = \widehat{\boldsymbol{\theta}}_{\tilde{D}} - \widehat{\boldsymbol{\theta}}_{D}$. Thus we obtain that the change in parameter estimator (i.e., influence on the parameter estimator) is,

$$\Delta\boldsymbol{\theta} = \frac{1}{n}\mathbf{H}_{\mathcal{L}}^{\setminus n}(\widehat{\boldsymbol{\theta}}_{\tilde{D}})^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}}) \tag{32}$$

Linearizing the loss at the left-out sample. In order to look at the change in loss evaluated on the n-th sample (left-out sample), we can consider linearizing the loss at $\widehat{\boldsymbol{\theta}}_{\tilde{D}}$.

$$\Delta\boldsymbol{\ell}_n = \frac{1}{n}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}})^{\top}\mathbf{H}_{\mathcal{L}}^{\setminus n}(\widehat{\boldsymbol{\theta}}_{\tilde{D}})^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}})$$

But, $\Delta\boldsymbol{\ell}_n = \boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}}) - \boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{D})$. And, so we have:

$$\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}}) = \boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{D}) + \frac{1}{n}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}})^{\top}\mathbf{H}_{\mathcal{L}}^{\setminus n}(\widehat{\boldsymbol{\theta}}_{\tilde{D}})^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}})$$

Now, we would like to average over the choice of the left-out sample n. However, we have to be careful and remember that $\tilde{D} = D_{n-1}^{\setminus i}$ depends on the particular left-out sample i. Thus, the full expression we get is the following:

$$\mathrm{LOO}_S = \mathcal{L}_S(\widehat{\boldsymbol{\theta}}_D) \; + \; \frac{1}{n^2}\sum_{i=1}^{n}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\left(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_{D_{n-1}^{\setminus i}}\right)^{\top}\mathbf{H}_{\mathcal{L}}^{\setminus i}\left(\widehat{\boldsymbol{\theta}}_{D_{n-1}^{\setminus i}}\right)^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}\left(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_{D_{n-1}^{\setminus i}}\right)$$

Let us recall the change in parameters with the leave-one-out contamination. From eq. 32, this was

$$\Delta\boldsymbol{\theta} = \frac{1}{n}\mathbf{H}_{\mathcal{L}}^{\setminus n}(\widehat{\boldsymbol{\theta}}_{\tilde{D}})^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{\tilde{D}})$$

where, $\tilde{D} := D_{n-1}^{\setminus n} = (1-\epsilon)D_n + \epsilon\delta_{z_n}$, with $\epsilon = \frac{-1}{n-1}$ and assuming n is large enough.

Under this asymptotic n scenario, it is reasonable to estimate the Hessian $\mathbf{H}_{\mathcal{L}}^{\setminus n}$ and $\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}$ at the distribution D_n instead of $\tilde{D} := D_{n-1}^{\setminus n}$. We thereby get the following (where we also make the dependence of $\Delta\boldsymbol{\theta}$ explicit on the particular left-out sample index),

$$\begin{aligned}
\Delta\boldsymbol{\theta}^{(n)} = \widehat{\boldsymbol{\theta}}_{D_{n-1}^{\setminus n}} - \widehat{\boldsymbol{\theta}}_S &= \frac{1}{n}\mathbf{H}_{\mathcal{L}}^{\setminus n}(\widehat{\boldsymbol{\theta}}_S)^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{D_n}) \\
&= \frac{1}{n}\left(\frac{1}{n-1}\sum_{i=1}^{n-1}\nabla_{\boldsymbol{\theta}}^2\boldsymbol{\ell}\left(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S\right)\right)^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{D_n}) \\
&= \left(\sum_{i=1}^{n-1}\nabla_{\boldsymbol{\theta}}^2\boldsymbol{\ell}\left(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S\right)\right)^{\dagger}\nabla_{\boldsymbol{\theta}}\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{D_n})
\end{aligned} \tag{33}$$

Now, we have a nice expression — in the sense that everything on the right-hand side is in terms of the parameters $\widehat{\boldsymbol{\theta}}_S$ obtained by training with the usual (empirical) distribution D_n. Since, we are interested in the leave-one-out loss, let us analyze how the above change in the estimated parameters (i.e., those obtained at convergence) brings about a change in the loss incurred on the n-th sample itself $\Delta\boldsymbol{\ell}^{(n)}$. To this end, we consider a **quadratic approximation** of the n-th sample loss at $\widehat{\boldsymbol{\theta}}_S$, as shown below (this can alternatively be thought of :

$$\boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_{D_{n-1}^{\setminus n}}) = \boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_S) + \nabla_{\boldsymbol{\theta}} \boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_S)^\top \Delta\boldsymbol{\theta}^{(n)} + \frac{1}{2} \Delta\boldsymbol{\theta}^{(n)\top} \nabla^2_{\boldsymbol{\theta}} \boldsymbol{\ell}(\boldsymbol{z}_n, \widehat{\boldsymbol{\theta}}_S) \Delta\boldsymbol{\theta}^{(n)} + \mathcal{O}(\|\Delta\boldsymbol{\theta}^{(n)}\|^3) \,. \tag{34}$$

Next, we consider this procedure for every choice of the left-out-sample, and then average out to get the expression of overall leave-one-out loss. We will also assume $\|\Delta\boldsymbol{\theta}^{(n)}\|^3$ and thus ignore the third-order term.

$$\begin{aligned}
\text{LOO} &= \frac{1}{n} \sum_{i=1}^{n} \boldsymbol{\ell}\left(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_{D_{n-1}^{\setminus i}}\right) \\
&= \frac{1}{n} \sum_{i=1}^{n} \boldsymbol{\ell}(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S) + \nabla_{\boldsymbol{\theta}} \boldsymbol{\ell}(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S)^\top \Delta\boldsymbol{\theta}^{(i)} + \frac{1}{2} \Delta\boldsymbol{\theta}^{(i)\top} \nabla^2_{\boldsymbol{\theta}} \boldsymbol{\ell}(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S) \Delta\boldsymbol{\theta}^{(i)} \\
&= \mathcal{L}(\widehat{\boldsymbol{\theta}}_S) + \frac{1}{n} \sum_{i=1}^{n} \underbrace{\nabla_{\boldsymbol{\theta}} \boldsymbol{\ell}(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S)^\top \Delta\boldsymbol{\theta}^{(i)}}_{=:B_i} + \frac{1}{2n} \sum_{i=1}^{n} \underbrace{\Delta\boldsymbol{\theta}^{(i)\top} \nabla^2_{\boldsymbol{\theta}} \boldsymbol{\ell}(\boldsymbol{z}_i, \widehat{\boldsymbol{\theta}}_S) \Delta\boldsymbol{\theta}^{(i)}}_{=:C_i}
\end{aligned} \tag{35}$$

To go further, we need to plug in the change in parameters computed in eq. 33 above. However, we first establish the validity of this general expression via the following result.

A.5 Leave-one-out formula proofs

Theorem 14. *Consider the particular case of the ordinary least-squares, where* $\boldsymbol{\ell}(\boldsymbol{z}, \boldsymbol{\theta}) := \boldsymbol{\ell}((\boldsymbol{x}, y), \boldsymbol{\theta}) = (y - \boldsymbol{\theta}^\top \boldsymbol{x})^2$. *We assume that we are given a training set* $S = \{(\boldsymbol{x}_i, y_i)\}_{i=1}^n$ *of points sampled i.i.d from the uniform empirical distribution* D_n. *Let the inputs be gathered into the data matrix* $\mathbf{X} \in \mathbb{R}^{n \times d}$ *where* d *is the input dimension and the targets are collected in the vector* $\boldsymbol{y} \in \mathbb{R}^n$. *Under the assumption that the number of samples* n *is large enough such that* $n \approx n - 1$, *we have that the leave-one-out expression from eq. 35 (with parameter change given in eq. 33) is equal to the widely-known closed-form formula of leave-one-out for least-squares* LOO^{LS}. *Mathematically, we have that,*

$$\text{LOO} = \text{LOO}^{LS} = \frac{1}{n} \sum_{i=1}^{n} \left(\frac{y_i - \boldsymbol{\theta}^\top \boldsymbol{x}_i}{1 - \mathbf{A}_{ii}} \right)^2 .$$

where, $\mathbf{A}_{ii} = \boldsymbol{x}_i^\top (\mathbf{X}^\top \mathbf{X})^{-1} \boldsymbol{x}_i$ *denotes the* i*-th diagonal entry of the matrix* $\mathbf{A} = \mathbf{X}(\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top$ *(i.e., the so-called 'hat-matrix') and* $\boldsymbol{\theta}$ *denotes the usual solution of* $\boldsymbol{\theta} = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \boldsymbol{y}$ *obtained via ordinary least-squares.*

Proof. We will first separately analyze the terms in the summation corresponding to the first-order and second-order parts in the eq. 35, which have been accorded the shorthand B_i and C_i respectively. Also, to make expressions simpler, we will define the residual as $r_i := \boldsymbol{\theta}^\top \boldsymbol{x}_i - y_i$ and denote the matrix $\mathbf{X}^\top \mathbf{X}$ by $\boldsymbol{\Sigma}$.

Let us write down the individual loss gradients and Hessian in this case. We have

$$\nabla \boldsymbol{\ell}_i := \nabla_{\boldsymbol{\theta}} \boldsymbol{\ell}((\boldsymbol{x}_i, y_i), \boldsymbol{\theta}) = 2 r_i \, \boldsymbol{x}_i \,,$$

$$\text{and, } \nabla^2 \boldsymbol{\ell}_i := \nabla^2_{\boldsymbol{\theta}} \boldsymbol{\ell}((\boldsymbol{x}_i, y_i), \boldsymbol{\theta}) = 2 \, \boldsymbol{x}_i \boldsymbol{x}_i^\top \,.$$

The term B_i in summation corresponding to the first-order part can be computed as follows.

$$
\begin{aligned}
B_i := \nabla \boldsymbol{\ell}_i{}^\top \Delta \boldsymbol{\theta}^{(i)} = \nabla \boldsymbol{\ell}_i{}^\top \left(\sum_{j \neq i} \nabla^2 \boldsymbol{\ell}_j \right)^{-1} \nabla \boldsymbol{\ell}_i &= 2 r_i \, \boldsymbol{x}_i^\top \left(2 \sum_{j \neq i} \boldsymbol{x}_j \boldsymbol{x}_j^\top \right)^{-1} 2 r_i \, \boldsymbol{x}_i \\
&= 2 r_i^2 \, \boldsymbol{x}_i^\top \left(\mathbf{X}^\top \mathbf{X} - \boldsymbol{x}_i \boldsymbol{x}_i^\top \right)^{-1} \boldsymbol{x}_i \\
&= 2 r_i^2 \, \boldsymbol{x}_i^\top \left(\boldsymbol{\Sigma} - \boldsymbol{x}_i \boldsymbol{x}_i^\top \right)^{-1} \boldsymbol{x}_i \\
&= 2 r_i^2 \, \boldsymbol{x}_i^\top \left(\boldsymbol{\Sigma}^{-1} + \frac{\boldsymbol{\Sigma}^{-1} \boldsymbol{x}_i \boldsymbol{x}_i^\top \boldsymbol{\Sigma}^{-1}}{1 - \boldsymbol{x}_i^\top \boldsymbol{\Sigma}^{-1} \boldsymbol{x}_i} \right) \boldsymbol{x}_i \\
&= 2 r_i^2 \left(\mathbf{A}_{ii} + \frac{(\mathbf{A}_{ii})^2}{1 - \mathbf{A}_{ii}} \right) \\
&= 2 r_i^2 \, \frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \, .
\end{aligned}
$$

Essentially, we have used the Sherman-Morrison-Woodbury formula in the fourth line above and rest is mere manipulation. Similarly, we compute the term C_i from the second-order part:

$$
\begin{aligned}
C_i := \Delta \boldsymbol{\theta}^{(i)}{}^\top \, \nabla^2 \boldsymbol{\ell}_i \, \Delta \boldsymbol{\theta}^{(i)} &= \nabla \boldsymbol{\ell}_i{}^\top \left(\sum_{j \neq i} \nabla^2 \boldsymbol{\ell}_j \right)^{-1} \nabla^2 \boldsymbol{\ell}_i \left(\sum_{j \neq i} \nabla^2 \boldsymbol{\ell}_j \right)^{-1} \nabla \boldsymbol{\ell}_i \\
&= 2 r_i^2 \, \boldsymbol{x}_i^\top \left(\boldsymbol{\Sigma} - \boldsymbol{x}_i \boldsymbol{x}_i^\top \right)^{-1} \boldsymbol{x}_i \boldsymbol{x}_i^\top \left(\boldsymbol{\Sigma} - \boldsymbol{x}_i \boldsymbol{x}_i^\top \right)^{-1} \boldsymbol{x}_i \\
&= 2 r_i^2 \left(\boldsymbol{x}_i^\top \left(\boldsymbol{\Sigma} - \boldsymbol{x}_i \boldsymbol{x}_i^\top \right)^{-1} \boldsymbol{x}_i \right)^2 \\
&= 2 r_i^2 \left(\frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \right)^2
\end{aligned}
$$

In the last line, we just reuse the computation that we already did in the previous part for $\boldsymbol{x}_i^\top \left(\boldsymbol{\Sigma} - \boldsymbol{x}_i \boldsymbol{x}_i^\top \right)^{-1} \boldsymbol{x}_i$. Finally, let us combine everything we have got with eq. 35

$$
\begin{aligned}
\text{LOO} &= \mathcal{L}(\boldsymbol{\theta}) \, + \, \frac{1}{n} \sum_{i=1}^n B_i \, + \, \frac{1}{2n} \sum_{i=1}^n C_i \\
&= \frac{1}{n} \sum_{i=1}^n r_i^2 \, + \, \frac{1}{n} \sum_{i=1}^n 2 r_i^2 \, \frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \, + \, \frac{1}{n} \sum_{i=1}^n r_i^2 \left(\frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \right)^2 \\
&= \frac{1}{n} \sum_{i=1}^n r_i^2 \left(1 \, + \, 2 \, \frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \, + \, \left(\frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \right)^2 \right) \\
&= \frac{1}{n} \sum_{i=1}^n r_i^2 \left(1 \, + \, \frac{\mathbf{A}_{ii}}{1 - \mathbf{A}_{ii}} \right)^2 \\
&= \frac{1}{n} \sum_{i=1}^n \left(\frac{r_i}{1 - \mathbf{A}_{ii}} \right)^2 .
\end{aligned}
$$

Thus, remembering our shorthand $r_i := \boldsymbol{\theta}^\top \boldsymbol{x}_i - y_i$, we conclude our proof.

□

Corollary 15. *The first-order influence function underestimates leave-one-out error for least-squares regression by a factor of* $\frac{1}{4}$.

Proof.

$$
\begin{aligned}
\mathrm{LOO}^{(1)} &= \mathcal{L}(\boldsymbol{\theta}) \,+\, \frac{1}{n}\sum_{i=1}^{n} B_i \\
&= \frac{1}{n}\sum_{i=1}^{n} r_i^2 \left(1 \,+\, 2\,\frac{\mathbf{A}_{ii}}{1-\mathbf{A}_{ii}}\right) \\
&= \frac{1}{n}\sum_{i=1}^{n} r_i^2 \,\frac{\mathbf{A}_{ii}}{1-\mathbf{A}_{ii}} \\
&= \frac{1}{n}\sum_{i=1}^{n} \left(\frac{r_i}{1-\mathbf{A}_{ii}}\right)^2 \mathbf{A}_{ii}(1-\mathbf{A}_{ii}) \\
&\leq \frac{1}{4}\,\mathrm{LOO}
\end{aligned}
$$

Here, the last line follows from using the AM (arithmetic mean) - GM (geometric mean) inequality. □

Corollary 16. *For any finite-width neural network, the first and second-order influence function give similar formulas for LOO like that in Theorem 7, but with* $\mathbf{A} = \mathbf{Z}_S(\boldsymbol{\theta}^\star)^\top \left(\mathbf{Z}_S(\boldsymbol{\theta}^\star)\mathbf{Z}_S(\boldsymbol{\theta}^\star)^\top\right)^{-1}\mathbf{Z}_S(\boldsymbol{\theta}^\star)$*, where* $\mathbf{Z}_S(\boldsymbol{\theta}^\star) := [\nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}_1), \cdots, \nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}_n)]$ *and* $\theta^\star$ *are the parameters at convergence for MSE loss.*

Proof. Simply replace $\boldsymbol{x}_i$ to $\nabla_{\boldsymbol{\theta}} \boldsymbol{f}_{\boldsymbol{\theta}}(\boldsymbol{x}_i)$ in the proof of Theorem 14 and repeat the procedure, since all the steps hold under the assumption of MSE loss and considering the Hessian $\mathbf{H}^S_{\mathcal{L}}(\boldsymbol{\theta}^\star) = \mathbf{H}^S_o(\boldsymbol{\theta}^\star)$ provided for by the assumption A2.

□

B Influence Functions: a primer

In this primer on influence functions, we closely follow the textbook of Hampel et al. (1986). The key objective of influence function is to investigate the infinitesimal behaviour of functionals, such as $T(\mathcal{D}_n)$. In particular, this is defined when the change in underlying distribution can be expressed in the form of a Dirac distribution. Formally, this is defined as follows:

Definition 17. *The influence function* IF *of the estimator* T*, at some distribution* F*, evaluated at a point* $\boldsymbol{z}$ *(where such a limit exists) is given by,*

$$
\mathrm{IF}(\boldsymbol{z}; T, F) = \lim_{\epsilon \to 0} \frac{T\left((1-\epsilon)F + \epsilon\delta_{\boldsymbol{z}}\right) - T(F)}{\epsilon}
$$

Essentially, the influence function (IF) involves a directional derivative of T at F along the direction of $\delta_{\boldsymbol{z}}$. Further, in order to interpret the above definition, substitute F by $\mathcal{D}_{n-1}$ and put $\epsilon = \frac{1}{n}$. This implies that (IF) measures n times the change of statistic T due to an additional observation $\boldsymbol{z}$. In other words, it describes the (standardized) effect of an infinitesimal contamination on the estimate T. E.g., in the case of parameter estimator, the change in estimated parameters due to presence of an additional datapoint. By now, the astute reader can already see its natural application for leave-one-out error, but we ask for some patience so as to discuss some other important aspects of influence functions.

More generally, one can view influence functions from the perspective of a Taylor series expansion (to be precise, the first-order von Mises expansion) of T at F, evaluated on some distribution G "close" to F,

$$
T(G) = T(F) \,+\, \int \mathrm{IF}(\boldsymbol{z}; T, F) d(G-F)(\boldsymbol{z}) \,+\, \text{remainder terms}\,. \tag{36}
$$

So, as evident from this, it is also possible to consider higher-order influence functions and use them in a combined manner, as considered in Debruyne et al. (2008). However, since the first-order term is often the dominating term — as well as to ensure tractability when we later consider neural networks — we will restrict our attention to only first-order influence functions hereafter.

B.1 PROPERTIES OF INFLUENCE FUNCTIONS

(i) Zero expectation. The expectation of influence function over the same distribution is zero, i.e., $\int \mathrm{IF}(\boldsymbol{z}; T, F) dF(\boldsymbol{z}) = 0$. This should actually be quite intuitive as we are basically averaging out all possible deviations of the estimator. But, the formal reasoning is that influence function is essentially akin to Gâteaux derivative at the distribution F,

$$\lim_{\epsilon \to 0} \frac{T\left((1-\epsilon)F + \epsilon G\right) - T(F)}{\epsilon} = \int \mathrm{IF}(\boldsymbol{z}) dG(\boldsymbol{z}) \,.$$

Now, replace G by F in the above equation and the stated property follows.

(ii) Variance of IF provides asymptotic variance of corresponding estimator. When the observations $\boldsymbol{z}_i$ are sampled i.i.d. according to F, then the empirical distribution $\mathcal{D}_n$ converges to F, for n sufficiently large, by Glivenko-Cantelli theorem. So, replacing G by $\mathcal{D}_n$ in eq. 36 and using the first property, we get:

$$T_n(\mathcal{D}_n) \approx T(F) \,+\, \int \mathrm{IF}(\boldsymbol{z}; T, F) d\mathcal{D}_n(\boldsymbol{z}) \,+\, \text{remainder terms} \,.$$

Integrating over the empirical distribution $\mathcal{D}_n$, we obtain:

$$\sqrt{n}(T_n(\mathcal{D}_n) - T(F)) \,\approx\, \frac{1}{\sqrt{n}} \sum_{i=1}^{n} \mathrm{IF}(\boldsymbol{z}_i; T, F) \,+\, \text{remainder terms} \,.$$

The term on the right-hand side involving IF is asymptotically normal by (multi-variate) Central Limit Theorem. Further, the remainder terms can often be neglected for $n \to \infty$, and thereby the estimator T_n is also asymptotically normal. Thus, $\sqrt{n}(T_n(\mathcal{D}_n) - T(F)) \xrightarrow{\mathrm{d}} \mathcal{N}_p(0, V(T, F))$, with the asymptotic variance (more accurately, covariance matrix) $V(T, F)$ as follows:

$$V(T, F) \,=\, \int \mathrm{IF}(\boldsymbol{z}; T, F)\, \mathrm{IF}(\boldsymbol{z}; T, F)^{\top}\, dF(\boldsymbol{z}) \,.$$

For the one-dimensional case, i.e., $T(F) \in \mathbb{R}$, we simply get $V(T, F) \,=\, \int \mathrm{IF}(\boldsymbol{z}; T, F)^2\, dF(\boldsymbol{z})$.

(iii) Chain rule for influence functions. Suppose our estimator $\widehat{\boldsymbol{\theta}}(F)$ depends on some other estimators, i.e., $\widehat{\boldsymbol{\theta}}(F) := T\left(\widehat{\boldsymbol{\theta}}_1(F), \cdots, \widehat{\boldsymbol{\theta}}_k(F)\right)$, then we can expression the influence function for $\widehat{\boldsymbol{\theta}}(F)$ as,

$$\mathrm{IF}(\boldsymbol{z}; \widehat{\boldsymbol{\theta}}, F) \,=\, \sum_{j=1}^{k} \frac{\partial T}{\partial \widehat{\boldsymbol{\theta}}_j}\, \mathrm{IF}(\boldsymbol{z}; \widehat{\boldsymbol{\theta}}_j, F) \,. \tag{37}$$

Remark. We refer the mathematically-oriented reader to Huber (2004) for details on the regularity conditions needed to ensure the existence of IF and the like.

C EMPIRICAL RESULTS AND DETAILS

C.1 EMPIRICAL DETAILS

We train all the networks via SGD with learning rate 0.5 and learning rate decay by a factor of 0.75 after each quarter of the target number of epochs.

For the experiments based on MNIST1D (Greydanus, 2020), the input dimension size is $d = 40$ while number of classes is $K = 10$.

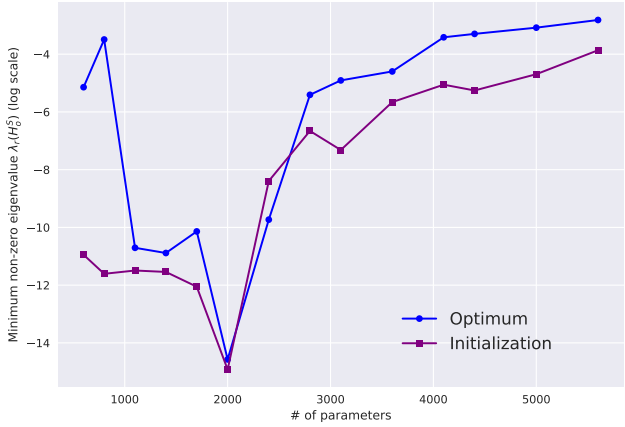
Double-descent model sizes (a) For the 3-layer double descent plot, we consider hidden widths $m_1,\ m_2 \in \{10, 20, 30, 40, 50\}$ and choose all those pairs as experiments where $|m_1 - m_2| \leq 20$.

(b) For the 2-layer plot in MSE, we take the hidden layer widths from $[1, 151]$ at an interval of 10.

(c) For the 2-layer plot corresponding to CE, we take hidden layer sizes from $[2, 32]$ at intervals of 2 and then for reducing the computation load when the network sizes increase, we take coarser intervals with a gap 5 and for even bigger sizes, at a gap of later 10. But this is purely to reduce computational load that comes with Hessian computation.

In all the double descent curves, we ensure that near the interpolation threshold all models are at interpolation, in the sense that the training accuracy is 100%.

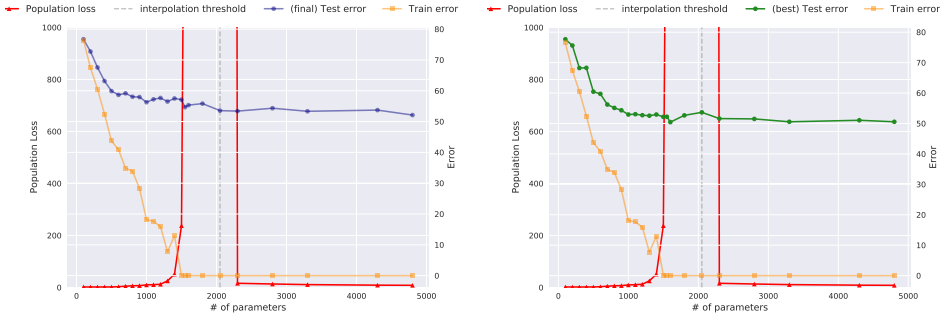
C.2 VERIFICATION OF ASSUMPTION A5



(a) 2-hidden layer case

Figure 4: Additional results on verifying the assumption A5

C.3 CROSS ENTROPY DOUBLE DESCENT

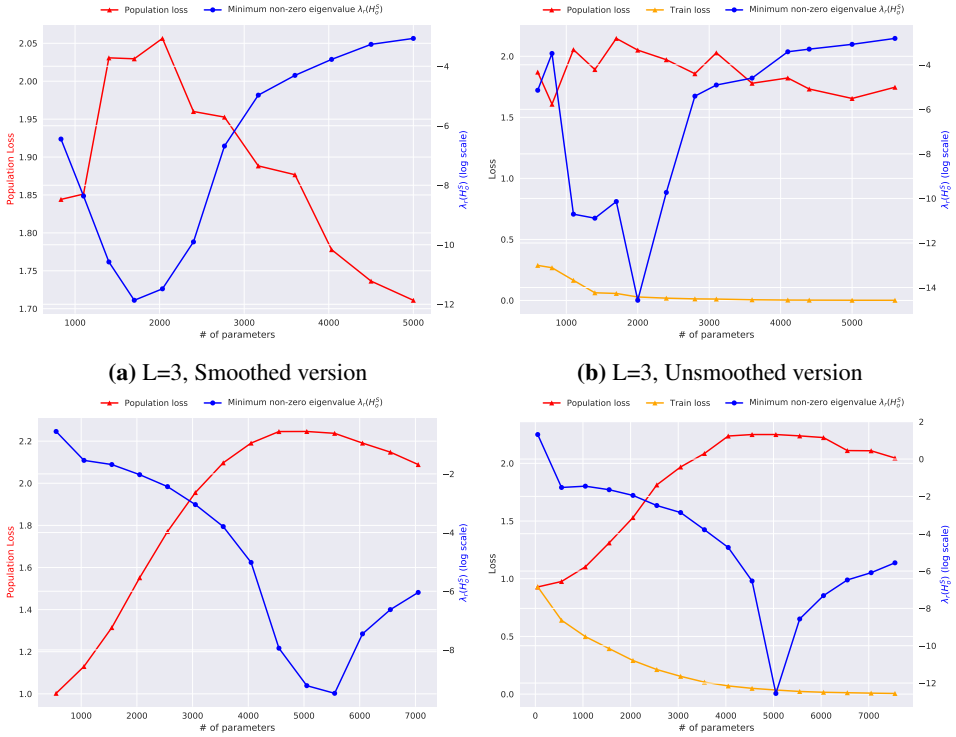


(a) Final Test Error **(b)** Best Test Error

Figure 5: Both test error at the final epoch as well as the 'best' test error in terms of the one selected based on a small validation set.

C.4 MSE DOUBLE DESCENT

For the sake of visualization in the 3-layer case, we smooth the quantities involved by considering a moving average of them over three successive model sizes.



(a) L=3, Smoothed version **(b)** L=3, Unsmoothed version

(c) L=2, Smoothed version **(d)** L=2, Unsmoothed version

Figure 6: We see that the same trend holds in both the cases — In fact, the unsmoothed version perhaps even more starkly shows the minimum non zero eigenvalue vanishes at the maximum population loss.

Also, in the unsmoothed version of the above plots, we also show the corresponding train loss. It must be emphasized that near interpolation the training accuracy is 100%, even though there might be some non-zero training loss.

C.5 LOWER BOUND COMPUTATION

At an initial glance, it might seem that the computed lower bounds could be very small in magnitude to be of use — as the bound is scaled by σ^2_{min}.

However, while practically computing one can consider a tighter lower bound, by using a particular tolerance τ.

Let us illustrate by considering evaluating this bound on a test set S' (just like we actually do). This is because in the lower bound we have the expression,

$$LHS = \frac{1}{|S|} \sum_{(\boldsymbol{x},\boldsymbol{y}) \in S'}^{|S|} \left[\text{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right) \cdot \|\nabla_{\boldsymbol{f}} \boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)\|^2 \right]$$

Now, since $\text{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right)$ is non-negative, if we are given a tolerance τ, we filter out those samples with $\|\nabla_{\boldsymbol{f}} \boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)\|^2 < \tau$

$$LHS \geq \frac{1}{|S|} \sum_{(\boldsymbol{x},\boldsymbol{y}) \in S'}^{|S|} \left[\text{Tr}\left(\mathbf{A}_{(\boldsymbol{x},\boldsymbol{y})}\right) \cdot \tau \cdot 1\{\|\nabla_{\boldsymbol{f}} \boldsymbol{\ell}\big(\boldsymbol{f}_{\boldsymbol{\theta}^\star}(\boldsymbol{x}), \boldsymbol{y}\big)\|^2\} \right]$$

where, 1 is the indicator function. This helps us compute lower bounds which can be evaluated in practice.

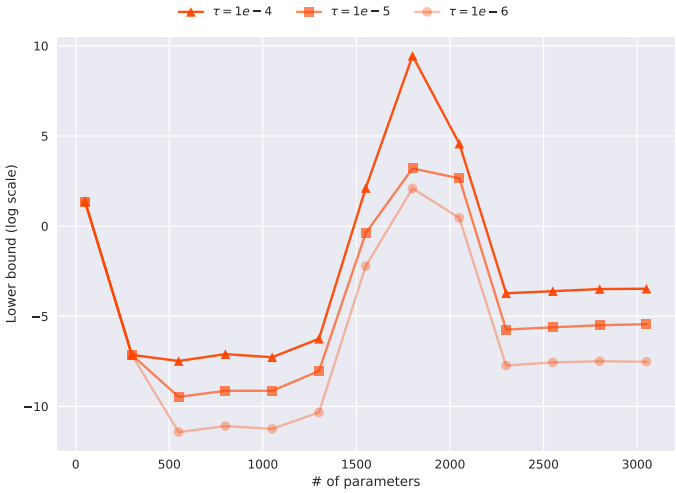
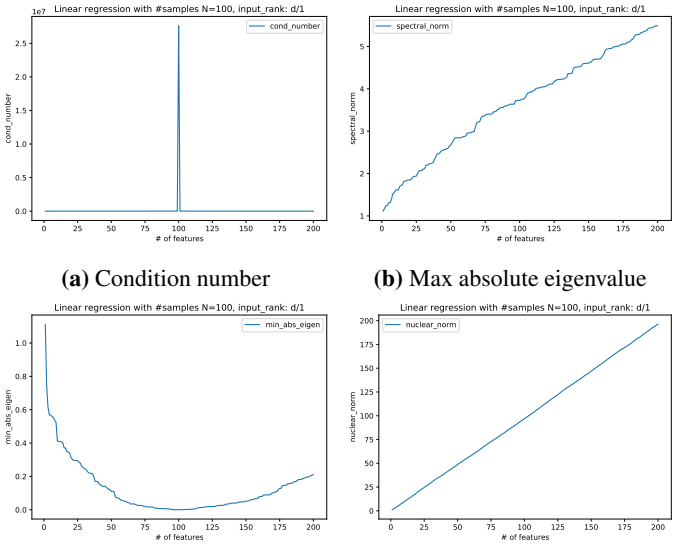


Figure 7: Comparison of various tolerance τ values used for lower bound computation. The lower bound values are displayed in log scale for the sake of presentation. In the figure, shown in the main text, we utilized tolerance $\tau = 1e - 6$. Overall, this illustrates that our bounds are not sensitive to choice of the tolerance parameter used.

C.6 EMPIRICAL OBSERVATIONS FROM LINEAR REGRESSION

C.6.1 HESSIAN STATISTICS ON TRAINING SET

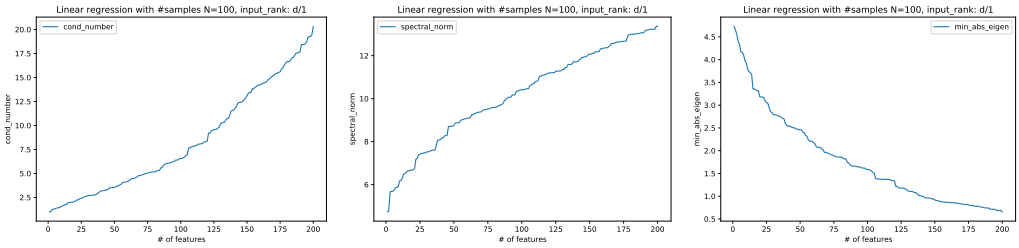


(a) Condition number **(b)** Max absolute eigenvalue

(c) Min absolute eigenvalue **(d)** Nuclear norm (or Trace)

Figure 8: Hessian statistics computed over the training set. We observe that condition number diverges at the double descent peak, owing to minimum absolute eigenvalue becoming close to zero at this threshold.

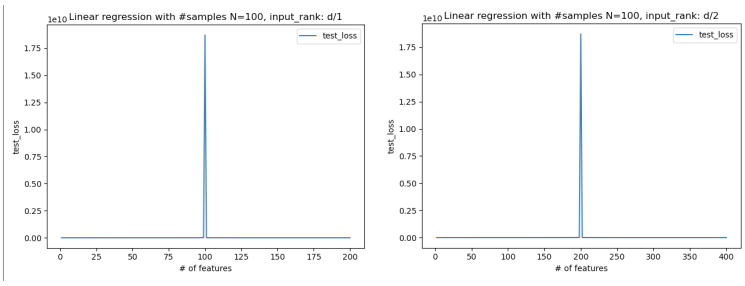
C.6.2 HESSIAN STATISTICS ON TEST SET



(a) Condition number **(b)** Max absolute eigenvalue **(c)** Min absolute eigenvalue

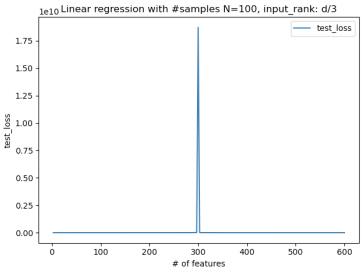
Figure 9: Hessian statistics computed over the test set (size 500).

C.6.3 DRAWING DOUBLE-DESCENT PLOTS AT ARBITRARY THRESHOLDS

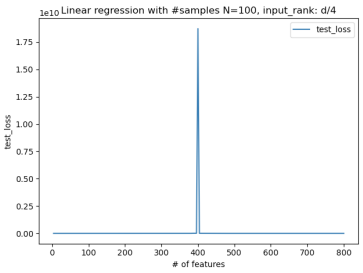


(a) $n = 100$, $d^\star = 100$, $\alpha = 0$

(b) $n = 100$, $d^\star = 200$, $\alpha = 1$



(c) $n = 100$, $d^\star = 300$, $\alpha = 2$



(d) $n = 100$, $d^\star = 400$, $\alpha = 3$

Figure 10: Changing the position of the interpolation threshold by adjusting the rank of the underlying input-covariance matrix.

The input design matrix is constructed in such a way so that given some $d^\star$ many input features, the number of base features is d. But we additionally have $\alpha\, d$ many redundant features (composed of linear combinations of the base features). Also, as the number of features is increased in the above test loss curve, these redundant features follow in succession to the base features. Thus, overall, we have $d^\star = (\alpha + 1)d$.