From Mean to Extreme: Formal Differential Privacy Bounds on the Success of Real-World Data Reconstruction Attacks

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Abstract

Data reconstruction attacks against machine learning (ML) models pose a strong risk of leaking sensitive data. When training ML models with differential privacy (DP), the success of such reconstruction attacks can be upper bounded. However, so far, these bounds have been formulated under worst-case assumptions, which in practice are overly pessimistic and, by that, result in unnecessary privacy-utility trade-offs. In this work, we provide formal bounds on the protection of DP against reconstruction attacks that do not necessitate prior knowledge of the data. We measure the reconstruction success with different metrics and determine its probabilistic behaviour from expectation (mean) to tail behaviour (extreme). By providing bounds for these metrics in a real-world scenario, we equip practitioners with a larger wealth of information to decide upon a certain set of privacy parameters.

1 Introduction and Related Work

Machine learning (ML) techniques and innovations have revolutionised a multitude of research and application areas, such as computer vision and natural language processing. However, the performance of ML methods is contingent on the availability of large real-world datasets. In certain domains, such as finance or medicine, data sets contain highly sensitive information about individuals and, therefore, are subject to (stronger) privacy legislation that can restrict their potentially already limited accessibility. This particularly concerns ML models since, for instance, they and/or their gradient updates store information about the training data, allowing for an (almost) perfect reconstruction of data samples (Fowl et al., 2022; Boenisch et al., 2023; Feng & Tramèr, 2024). For the acceptance of stakeholders and, by that, the availability of their data for ML training, it is thus imperative to implement protective measures. This has sparked the research area of privacy-preserving ML.

The gold standard technique for providing data owners with theoretical privacy protection is differential privacy (DP) (Dwork et al., 2014). DP is a formal framework that provides mathematically provable privacy guarantees for individuals whose data is being used, e.g., to train ML models. The strength of the guarantee depends on the chosen set of parameters. Utilising DP inherently enforces theoretical upper bounds on the success of attacks that aim to reconstruct the input data (Guo et al., 2022; Balle et al., 2022; Hayes et al., 2023; Kaissis et al., 2023a). Providing such bounds by applying DP allows data owners to control the risks of sharing their data. However, to do so, practitioners must be able to make an informed decision about privacy parameters, which effectively protects against data reconstruction while at the same time not being overly pessimistic, which would result in imprecise outputs and, in the context of machine learning, performance losses of the resulting models. Therefore, in addition to these abstract privacy budgets, assessing privacy guarantees in terms of the risk associated with specific attacks in concrete scenarios allows for a more nuanced understanding of the provided protection for all parties involved. In the following, we refer to such guarantees concerning interpretable risks as semantic quarantees.

Previous works have investigated these semantic guarantees employing DP (e.g. Nasr et al. (2021)). However, they have almost exclusively analysed scenarios under "worst-case" assumptions, in which the most powerful conceivable adversary performs the attack. Such a worst-case scenario includes knowledge about the reconstruction target. Defining this *prior knowledge* of an adversary is not trivial. However, at worst, the

reconstruction attack in such a scenario is equivalent to matching an element of a set of candidates of the training data to a specific model or gradient. While this follows a Bayesian notion of updating a prior belief with a posterior distribution, in many practical applications, an adversary is rather interested in recovering the data sample that is not already in their possession. Evaluating this worst-case scenario is appealing since an upper bound on the success of such a threat is also a bound on the success of any other reconstruction attack under less strict settings. Moreover, the constructed bound cannot be deteriorated by post-processing as the worst-case has control of all parameters (including the input and function) except for the introduced randomness by the DP mechanism. Therefore, no side knowledge could systematically increase the adversary's success. However, only considering the worst-case scenario comes at a cost: if the adversary is weaker than presumed, the constructed bounds become loose, overrating the adversary's success. Overestimating privacy risks is problematic because it drives practitioners to implement stricter privacy conditions, which, in turn, reduces the quality of the result. The resulting compromise between privacy and performance is called the privacy-utility trade-off (De et al., 2022). It imposes a dilemma on practitioners, particularly in fields where highly accurate results are critical, such as the medical domain. For example, in oncological tasks, they must choose between correctly detecting tumours or accepting the risk of leaking patients' private information. Hence, in practice, it is crucial to determine and analyse the case-dependent most powerful but realistic adversary to enforce the "least amount of privacy" that sufficiently protects against such threat without overly compromising the model's performance.

Interestingly, state-of-the-art (SOTA) real-world reconstruction attacks do not even require all the capabilities of a worst-case adversary. For instance, the currently best performing real-world attacks can reconstruct the data perfectly without any assumptions about the input data or distribution beyond its dimensionality (Fowl et al., 2022; Boenisch et al., 2023; Feng & Tramèr, 2024). Hence, while the analysis of the worst case is crucial for understanding the risks of reconstruction at the worst, we believe it is necessary to augment it with risk models tailored to real-world attacks in order to achieve an optimal privacy-utility trade-off. These real-world attacks operate under a relaxed threat model, i.e., the considered adversaries have fewer capabilities and/or knowledge and are thus weaker. We note that bounds on such relaxed threat models are not immune to post-processing. However, we believe that only practitioners who can accurately estimate the risks for their specific threat model can make an informed decision about an appropriate choice of privacy parameters, which leads to an optimal trade-off between privacy and utility. We envision that such analyses can lead to full system cards (analogously to model cards (Mitchell et al., 2019)), where the privacy risks in different contexts are described. In this work, we formulate probabilistic bounds on the reconstruction success for this branch of attacks (Fowl et al., 2022; Boenisch et al., 2023; Feng & Tramèr, 2024). With this, we hope to provide tools for a more educated choice of privacy parameters and, by that, better privacy-utility trade-offs that allow for both private and highly performing AI models.

1.1 Contributions

In this work, we investigate semantic privacy guarantees of Gaussian mechanisms under a relaxed threat model, which aligns with real-world SOTA reconstruction attacks. Our contributions can be summarised as follows:

- In Section 3.1, we formulate the analytic gradient inversion attack that delivers the optimal minimum variance unbiased estimator for any target input.
- In Section 3.2.1, we measure the success of reconstruction attacks using MSE, PSNR, and we formulate probabilistic bounds for these metrics in terms of $(\eta, \gamma(\eta))$ -reconstruction robustness (ReRo).
- In Section 3.2.2, we measure the success of reconstruction attacks using NCC, and derive a theoretical upper bound for this quantity.
- In Section 4, we visualize the implications of our results and their correspondence to empirical measurements.

1.2 Assumptions

Our assumptions are based on prior work, which we discuss in-depth in the Appendix A.1. Risks in DP settings are typically analysed under a worst-case adversary, which is colloquially often also referred to as "DP threat model". In this threat model, an adversary knows the query function q and all parameters of the privacy mechanism \mathcal{M} . The query function is the function of which we are interested in the output; in the context of AI models, it is a forward/backward pass over the model and the output is a gradient. The architecture and parameters of the model are essential to this function output. The adversary can manipulate these settings in their favour, including the network architecture, weights and loss. In a worst-case threat model, the adversary aims to distinguish whether a privatised output is based on the input of two almost synonymous datasets: \mathcal{D} or \mathcal{D}' . These two datasets are known and adjacent datasets, meaning \mathcal{D}' is the same as \mathcal{D} , but one entry is either removed or added, and the adversary knows all entries in \mathcal{D} and also which entry is removed or added. This scenario corresponds to a worst-case analysis of membership inference risk analysis. Hayes et al. (2023) and Kaissis et al. (2023a) have empirically and theoretically analysed the success of reconstruction attacks under multiple relaxations of the DP threat model. Even though both analyses provide a "step closer" towards choosing DP guarantees tailored to more practicable contexts, their assumptions for constructing theoretical bounds can still be considered "too strict" and fail to describe the capabilities of adversaries of some of the current SOTA real-world reconstruction attacks.

Given these considerations, we choose our assumptions to be equivalent to those of several SOTA real-world reconstruction attacks, e.g., Fowl et al. (2022); Boenisch et al. (2023); Feng & Tramèr (2024). These are relaxed compared to Hayes et al. (2023): In particular, they differ by assuming an adversary without knowledge about the input data beyond the dimensionality. In all other regards, our threat model is congruent with the worst-case reconstruction threat by Hayes et al. (2023). However, we note that the goal of our work differs from Hayes et al. (2023) as they describe a posterior probability which is influenced by a prior. In contrast, we describe the probabilities of reconstructing the input data in a certain quality. In the taxonomy by Cummings et al. (2024), our model differs in the assumptions of population-level and dataset-level auxiliary information compared to the DP worst-case threat model.

1.3 Background

In this preliminary subsection, we recall the established concept of (η, γ) -reconstruction robustness (ReRo) (Balle et al., 2022) that allows us to express DP guarantees in terms of the reconstruction success of an adversary. We note that further basic concepts, specifically regarding the studied reconstruction metrics, can be found in Appendix A.2.

Defining the "robustness" of ML models against reconstruction attacks is central to our findings, particularly in translating the "reconstruction success" given by reconstruction error functions into properties of DP mechanisms. Arguably, the most general definition of reconstruction robustness was given by Balle et al. (2022):

Definition 1. A randomized mechanism $\mathcal{M}: \mathcal{Z}^n \to \Theta$ is (η, γ) -reconstruction robust with respect to a prior π over \mathcal{Z} and a reconstruction error function $l: \mathcal{Z} \times \mathcal{Z} \to \mathbb{R}_{\geq 0}$ if for any dataset $D_- \in \mathcal{Z}^{n-1}$ and any reconstruction attack $R: \Theta \to \mathcal{Z}$:

$$\mathbb{P}_{Z \sim \pi, \Theta \sim \mathcal{M}(D_- \cup \{Z\})}[l(Z, R(\Theta)) \leq \eta] \leq \gamma.$$

In Definition 1, the reconstruction error function l measures the similarity between the true data Z and its reconstruction $R(\Theta)$, where $R(\Theta)$ is obtained using the randomised mechanism's output Θ and the adversary's prior knowledge π . The closer Z and $R(\Theta)$ are, the lower the values of $l(Z, R(\Theta))$. Therefore, η can be chosen in such a way that a reconstruction error above this threshold η indicates a distorted, non-usable reconstruction, and an error below η characterises informative reconstructions. In turn, quantifying the probability of having an error lower than this threshold η offers an insight into the semantic meaning of the protection provided by the DP mechanism \mathcal{M} . Concretely, if said probability is upper bounded by γ for all reconstruction attacks R, then the probability of "successfully" reconstructing the input Z is upper bounded by γ , and we call \mathcal{M} (η, γ)-reconstruction robust with respect to π and l.

Methods

Analytic Gradient Inversion Attacks

Throughout this work, we consider an adversary who aims to recover the input data used to train a ML model. We assume this adversary has control over the model's architecture, hyperparameters and loss function, and knows the dimension N of the input data as stated in Section 1.2. The works by Boenisch et al. (2023); Fowl et al. (2022); Feng & Tramèr (2024) have shown that, under such assumptions, said adversary can conduct a successful reconstruction attack against a ML model. All of these attacks require performing an analytic gradient inversion attack based on inverting linear layers with an added bias term present in (or added into) the model.

For instance, consider a network whose first layer is a fully connected linear layer $f: \mathbb{R}^N \to \mathbb{R}^M$ with weight matrix $W \in \mathbb{R}^{M \times N}$ and bias scalar term $b \in \mathbb{R}^M$. Formally, the operation of a linear layer on an input sample $X \in \mathbb{R}^N$ can be written as

$$f(X) = WX + b. (1)$$

Typically, such a linear layer is succeeded by other neural network operations and a loss function, which we summarise in the term $g: \mathbb{R}^M \to \mathbb{R}$. Hence, g describes the part of the network starting from the linear layer and q(f(X)) denotes the network's loss function output. As customary, all network parameters are updated according to the loss function during each training step. For that purpose, the gradient of $g \circ f$ with respect to all network parameters is computed by a backward pass. We call this gradient the global, concatenated gradient and denote it by G_X . Naturally, the global gradient G_X is dependent on the training step and on the input sample point X of that specific training step, which is evident since model updates change from one training step to another. However, to ease notation, we do not additionally index G_X with the iteration step. We recall that the adversary knows G_X under the capabilities assumed by the threat model.

Therefore, for a fixed training step, besides other model updates concerning the parameters of the part of the network given by g, for all $j \in \{1, ..., M\}$, the adversary observes

$$\nabla_{W_j} g(f(X)) \in \mathbb{R}^N \quad \text{and} \quad \frac{\partial g(f(X))}{\partial b_j} \in \mathbb{R},$$
 (2)

namely, the gradient of $g \circ f$ with respect to jth row of the matrix W and the derivative of $g \circ f$ with respect to the j-th entry of the bias term b, respectively. Moreover, the adversary is aware that the gradient and the derivative in 2 are constructed by a backward pass in the following way:

$$\nabla_{W_j} g(f(X)) = \frac{\partial g(f(X))}{\partial f(X)_j} \nabla_{W_j} f(X)_j = \frac{\partial g(f(X))}{\partial f(X)_j} X,$$

$$\frac{\partial g(f(X))}{\partial b_j} = \frac{\partial g(f(X))}{\partial f(X)_j} \frac{\partial f(X)_j}{\partial b_j} = \frac{\partial g(f(X))}{\partial f(X)_j},$$
(4)

$$\frac{\partial g(f(X))}{\partial b_j} = \frac{\partial g(f(X))}{\partial f(X)_j} \frac{\partial f(X)_j}{\partial b_j} = \frac{\partial g(f(X))}{\partial f(X)_j},\tag{4}$$

for all $j \in \{1, ..., M\}$. Note that the gradient in Equation 3 is a scaled version of the input X and that the multiplicative factor $\frac{\partial g(f(X))}{\partial f(X)_j}$ on the right-hand side of Equation 3 equals the observed update with respect to the j-th entry of the bias given in Equation 4. Thus, if there exists $j' \in \{1, ..., M\}$, such that the update $\frac{\partial g(f(X))}{\partial b_{x'}} \neq 0$, the adversary can reconstruct the input sample X analytically by performing

$$\nabla_{W_{j'}} g(f(X)) \oslash \frac{\partial g(f(X))}{\partial b_{j'}} = X, \tag{5}$$

where \oslash denotes the entry-wise division. Computing 5 is possible for all $j \in \{1, ..., M\}$ such that $\frac{\partial g(f(X))}{\partial b_i} \neq 0$.

In the assumed setting, the adversary can modify the model to facilitate their analytic reconstruction attack. Specifically, if necessary, they can insert the fully connected linear layer f, as given in Equation 1, as a first layer into the network or even replace it entirely. Therefore, from now on, we consider neural networks of the above form $g \circ f$. Similarly, we assume the adversary chooses the batch size B to equal one to ensure that only one input sample X is used per iteration step and avoid contaminating the model updates with the information from different samples. For a more detailed discussion concerning the batch size B, we refer to Section A.1, specifically the paragraph referring to prior works on analytic gradient inversion attacks.

2.2 Analytic Gradient Inversion Attacks under DP

Fulfilling the requirements of DP can be achieved by randomising the output of a deterministic function $q: \mathcal{X} \to \mathbb{R}^N$, where \mathcal{X} denotes a universe of sensitive records. This randomisation can be conducted employing a DP additive noise mechanism \mathcal{M} , that adds noise drawn from an appropriate distribution, calibrated to the (global) sensitivity Δ_p of q, to the deterministic output $q(\mathcal{D})$ of the function q over a data set \mathcal{D} . The sensitivity Δ_p of q is calculated using the ℓ_p -norm: $\Delta_p(q) = \sup_{D, D' \in \mathcal{X}, D \simeq D'} \|q(D) - q(D')\|_p$, for $p \in [1, \infty)$, where $D \simeq D'$ denotes that D and D' are adjacent data sets, i.e., data sets which differ in exactly one entry. In this work, we consider the Gaussian additive noise mechanism or Gaussian mechanism, the predominant approach in applications with high-dimensional data. Let \mathcal{M} be the Gaussian mechanism over a query function $q: \mathcal{X} \to \mathbb{R}^N$ with sensitivity $\Delta_2 := \Delta_2(q)$, then the output of the DP mechanism \mathcal{M} is given by

$$\mathcal{M}(q(D)) = q(D) + \xi, \quad \xi \sim \mathcal{N}\left(0, \sigma^2 \Delta_2^2 I_N\right),$$

for I_N being the N-dimensional identity matrix and σ the noise scale. In particular, the choice of σ modulates the privacy guarantee provided by (a single execution of) the Gaussian mechanism \mathcal{M} .

From now on, we consider training networks of the form $g \circ f$, as stated in Section 2.1, with DP to protect the training data against analytic gradient inversion attacks and evaluate the provided protection. In the context of neural networks, applying DP is usually achieved by training with DP-SGD (Song et al., 2013). Since in DP-SGD the privatised quantity is the global gradient, the noise must be calibrated to the sensitivity of the global gradient, which is not (necessarily) bounded and can be hard to compute. Hence, the DP-SGD algorithm is based on two main steps: (1) Clipping the ℓ_2 -norm of the global, concatenated gradient to a predefined bound C in order to have an artificial bound on the sensitivity and (2) adding calibrated, zero-centered Gaussian noise to the gradient. The hyperparameter C is usually called the maximum gradient norm.

For a network $g \circ f$, for a fixed iteration step, the global gradient G_X has the following form:

$$G_X = \left[\nabla_{W_1} g(f(X))^T, \frac{\partial g(f(X))}{\partial b_1}, ..., \nabla_{W_M} g(f(X))^T, \frac{\partial g(f(X))}{\partial b_M}, G_{X,P}^T \right]^T, \tag{6}$$

where $G_{X,P}$ denotes the concatenated gradient of $g \circ f$ with respect to the rest of the parameters of the network, where all vectors are unrolled to scalars. To induce a bound on the norm of the global gradient G_X , G_X is multiplied by the clipping term:

$$\beta_C(X) := \frac{1}{\max\left(1, \frac{\|G_X\|_2}{C}\right)}.\tag{7}$$

We note that $\beta_C(X)$ is dependent on the iteration step by definition and it decreases with increasing norm of the global gradient G_X . Next, independent and identically distributed (i.i.d.) noise samples ξ are drawn from a multivariate Gaussian distribution $\mathcal{N}(\mathbf{0}, C^2\sigma^2I)$, where I is the identity matrix of the dimension of the global gradient G_X and $\mathbf{0}$ denotes zero vector also matching the dimension of G_X . Hence, when a network is trained with DP-SGD, the adversary observes *noisy* versions of the model updates in Equations 3 and 4 that could be used for the gradient inversion attack:

$$\widetilde{\nabla}_{W_j} := \beta_C(X) \nabla_{W_j} g(f(X)) + \xi_j = \beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j} X + \xi_j, \qquad \xi_j \sim \mathcal{N}(\mathbf{0}_N, C^2 \sigma^2 I_N), \tag{8}$$

$$\widetilde{\nabla}_{b_j} := \beta_C(X) \frac{\partial g(f(X))}{\partial b_j} + \xi_j' = \beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j} + \xi_j', \qquad \xi_j' \sim \mathcal{N}(0, C^2 \sigma^2), \tag{9}$$

for all $j \in \{1, ..., M\}$. Under non-DP training, the gradient in Equation 3 stored a scaled version of the target input point X (see Section 2.1). In contrast, when employing DP, the model updates have the form presented in 8 where the stored target X has not only been rescaled but also perturbed by adding Gaussian noise. Therefore, for each $j \in \{1, ..., M\}$, Expression 8 denotes a *privatised* version of the input X. It is easy to see that the introduced randomness impedes performing a simple division to recover the input sample X

analogously to the operation in 5. Using the distribution of the sampled noise, the noisy model updates in Equations 8 and 9 can be expressed as samples from random variables in the following way

$$\widetilde{\nabla}_{W_j} \stackrel{d}{=} Y_j, \quad \text{for} \quad Y_j \sim \mathcal{N}\left(\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j} X, C^2 \sigma^2 I_N\right),$$
(10)

and
$$\widetilde{\nabla}_{b_j} \stackrel{d}{=} z_j$$
, for $z_j \sim \mathcal{N}\left(\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j}, C^2 \sigma^2\right)$, (11)

for all $j \in \{1, ..., M\}$, where $\stackrel{d}{=}$ means equal in distribution. Even though the adversary cannot perform the division in 5 to reconstruct X, they can use the privatised global gradient, in particular, the observations $\widetilde{\nabla}_{W_1}, ..., \widetilde{\nabla}_{W_M}, \widetilde{\nabla}_{b_1}, ..., \widetilde{\nabla}_{b_M}$ and the knowledge about their distributions, as given in 10 and 11, to design an estimator for the target point X. Ultimately, this estimator serves as the reconstruction Y_X of X. We highlight that $\widetilde{\nabla}_{W_1}, ..., \widetilde{\nabla}_{W_M}$ are M observed, privatised versions of X, whereby M can be modulated by the adversary since it denotes the number of rows of the matrix W that specifies the linear layer f (see 1). In Section 3.1, we address parameter M's importance for the estimator and, thus, for the attack's success.

Without making assumptions on the part of the network given by g, it is difficult to determine whether, for all or for some iteration steps, the part of the gradient denoted by $G_{X,P}$ (see 6) and, thereby, its privatised version contain usable information to estimate the target X. Therefore, we first assume that the privatised model updates given by $\widetilde{\nabla}_{W_1}, ..., \widetilde{\nabla}_{W_M}, \widetilde{\nabla}_{b_1}, ..., \widetilde{\nabla}_{b_M}$ are sufficient to estimate X and ignore the remaining part of the privatised global gradient. Later, we address the influence $G_{X,P}$ has on formulating an estimator for X and, in Corollary 1, we specify the choice of the part of the network given by g that renders the analytic gradient inversion attack with the highest reconstruction success in terms of the MSE.

Reconstructing the target X using only the observations $\widetilde{\nabla}_{W_1},...,\widetilde{\nabla}_{W_M}$ is "not far from" solving a classical mean estimation problem. However, the means of the distributions of the samples $\widetilde{\nabla}_{W_1},...,\widetilde{\nabla}_{W_M}$ are not X but rescaled versions of X, namely $\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_1}X,...,\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_M}X$, respectively (see 10). Removing the dependency on the scaling factors $\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_1},...,\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_M}$, is necessary for creating an unbiased estimator for X (see Proposition B.1 in the appendix) such as the sample mean. Since the adversary also observes noisy versions of these scaling factors, namely $\widetilde{\nabla}_{b_1},...,\widetilde{\nabla}_{b_M}$ (see 11), we can differentiate between the cases where the adversary needs and does not need $\widetilde{\nabla}_{b_1},...,\widetilde{\nabla}_{b_M}$ to construct an unbiased estimator for X.

In the former case, the adversary can try to integrate the observed privatised gradients $\widetilde{\nabla}_{W_1},...,\widetilde{\nabla}_{W_M},\widetilde{\nabla}_{b_1},...,\widetilde{\nabla}_{b_M}$ by dividing entry-wise $\widetilde{\nabla}_{W_j}$ by $\widetilde{\nabla}_{b_j}$ for all $j \in \{1,...,M\}$ such that $\widetilde{\nabla}_{b_j} \neq 0$. Consequently, to determine the behaviour of $\widetilde{\nabla}_{W_j} \oslash \widetilde{\nabla}_{b_j}$, $j \in \{1,...,M\}$, they combine 10 and 11 and consider

$$\widetilde{\nabla}_{W_j} \oslash \widetilde{\nabla}_{b_j} \stackrel{d}{=} V_j, \quad \text{for} \quad V_{j,i} \sim \frac{\mathcal{N}\left(\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j} x_i, C^2 \sigma^2\right)}{\mathcal{N}\left(\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j}, C^2 \sigma^2\right)},$$
(12)

where $V_{j,i}$ denotes the *i*-th entry of V_j for all $i \in \{1,...,N\}$, and $V_{j,i}$ are pairwise independently distributed. The ratio of two Gaussian distributions, as the one in 12, follows the Cauchy distribution. Generally, this distribution has no defined statistical moments (Marsaglia, 2006), such as an expectation. Although, under certain assumptions these moments can be approximated (Marsaglia, 2006; Díaz-Francés & Rubio, 2013), no general statements about the behaviour of the samples $\widetilde{\nabla}_{W_j} \oslash \widetilde{\nabla}_{b_j}$, $j \in \{1,...,M\}$, or the asymptotic behaviour of estimators constructed using said samples, can be made if the statistical moments do not exist. This is particularly problematic considering that the distribution in 12 varies for all $j \in \{1,...,M\}$, $X \in \mathcal{D}$ and each iteration step. Therefore, we assume that the adversary refrains from employing 12 to estimate X. However, if necessary, the adversary uses $\widetilde{\nabla}_{b_1},...,\widetilde{\nabla}_{b_M}$ to estimate the scaling factors $\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_1},...,\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_M}$ first and then, separately, they use these estimators to rescale the observed privatised gradients 8, reformulate their distributions 10 and solve the mean estimation problem. Due to the inherent randomness of estimators, it is easy to see that a reconstruction Y_X of X is more accurate if the scaling factors $\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_1},...,\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_M}$ are known. In particular, this implies that upper bounding the success of a reconstruction Y_X when the scaling factors are known is sufficient to bound the reconstruction

success of Y_X when the scaling factors are estimated. As we are interested in the most precise reconstruction an adversary could obtain, in the following, we assume that these scaling factors are available to the attacker.

3 Theoretical results

3.1 Construction of the optimal analytic gradient inversion reconstruction attack under DP

Let $X = (x_1, ..., x_N)^T \in \mathcal{D} \subseteq \mathbb{R}^N$ be a fixed reconstruction target point, and also fix the training iteration step. Again, we emphasise that from now on, we assume the adversary can estimate the scales $\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_1}, ..., \beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_M}$ without requiring their noisy versions 9 and after observing the privatised gradients $\widetilde{\nabla}_{W_1}, ..., \widetilde{\nabla}_{W_M}$ (see 8), as explained in the last part of Section 2.2. However, we also consider this case as such scaling factors can be approximated by various strategies independently of 9, in the simplest case, by constraints on the data, e.g., for images by setting their pixels to range from 0 to 255.

Having access to $\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_1},...,\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_M}$ implies that the adversary can rescale the observed privatised gradients $\widetilde{\nabla}_{W_1},...,\widetilde{\nabla}_{W_M}$ (see 8), construct their sample average \hat{X}_M , and use \hat{X}_M as an unbiased estimator for X:

$$\hat{X}_M := \frac{1}{M} \sum_{j=1}^M \frac{1}{\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j}} \widetilde{\nabla}_{W_j}. \tag{13}$$

Consequently, using the distribution of the observations $\widetilde{\nabla}_{W_j}$, $j \in \{1, ..., M\}$, given in 10, the distribution of the sample mean \hat{X}_M is given by:

$$\hat{X}_M \sim \mathcal{N}\left(X, \frac{1}{M^2} \sum_{j=1}^M \frac{C^2 \sigma^2}{\beta_C(X)^2 \frac{\partial g(f(X))}{\partial f(X)_j}^2} I_N\right). \tag{14}$$

In this work, we have assumed the adversary is able to modify the model's architecture, hyperparameters and loss function to facilitate their attack. Therefore, they are capable of adjusting the neural network $g \circ f$ to minimise the coordinate-wise variance of the sample mean \hat{X}_M in 14, aiming to increase the probability that \hat{X}_M takes a value "close enough" to its expectation, namely to the target X. One potential strategy, could be to increase the number of rows M of the matrix W of the first (linear) layer f (see 1) of the network. However, for increasing M, the norm of the global gradient G_X (see 6) increases and the clipping term $\beta_C(X)$ (see 7) decreases requiring a careful consideration of the interaction between M and $\beta_C(X)$. In the appendix (see Proposition B.2), we show that the coordinate-wise variance of the sample mean \hat{X}_M is lower bounded by $\sigma^2 \|X\|_2^2$ for $M \to \infty$. Intuitively, this implies that, from the adversary's perspective, the variability of the estimator \hat{X}_M can only get "so good", but after a certain value for M it will not further decrease and, in particular, it will never be zero such that \hat{X}_M results in the undistorted target point.

The next proposition specifies the concrete modifications to the network that render the sample mean \hat{X}_M with the lowest variance:

Proposition 1. Let the part of neural network given by g be replaced by the loss function $\mathcal{L}: \mathbb{R}^M \to \mathbb{R}$ with $\mathcal{L}(f(X)) = \mathbf{1}_M^T f(X)$, where $\mathbf{1}_M$ is the M-dimensional 1-vector, and

$$M \ge \max\left(1, \left\lceil \frac{C}{\min_{X \in \mathcal{D} \setminus \{\theta_N\}} \|X\|_2} \right\rceil\right),\tag{15}$$

where $\lceil \cdot \rceil$ denotes the function that rounds up its argument to the nearest integer, then $\frac{1}{M^2} \sum_{j=1}^{M} \frac{C^2 \sigma^2}{\beta_C(X)^2 \frac{\partial g(f(X))}{\partial f(X)_j}^2}$ is minimal and takes the value $\sigma^2 \|X\|_2^2$.

Assume the adversary replaces the entire network by the linear function $f: \mathbb{R}^N \to \mathbb{R}^M$ given by f(X) = WX, $W \in \mathbb{R}^{M \times N}$ and the loss function $\mathcal{L}: \mathbb{R}^M \to \mathbb{R}$ given by $\mathcal{L}(f(X)) := \mathbf{1}_M^T f(X)$, then $\frac{\partial g(f(X))}{\partial f(X)_j} = \frac{\partial \mathcal{L}(f(X))}{\partial f(X)_j} = 1$ for all $j \in \{1, ..., M\}$. In particular, this implies that

$$\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_1} = \dots = \beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_M} = \beta_C(X)$$
(16)

reducing the approximation of M scales, presented in the first part of this section, to the estimation of only one scale, namely $\beta_C(X)$.

Moreover, by Proposition 1, for all M' satisfying 15, the sample mean $\hat{X}_{M'}$ is distributed in the following way:

$$\hat{X}_{M'} \stackrel{d}{=} \hat{X}, \quad \text{for} \quad \hat{X} \sim \mathcal{N}\left(X, \sigma^2 \|X\|_2^2 I_N\right).$$
 (17)

If $C/\min_{X\in\mathcal{D}\setminus\{\mathbf{0}_N\}}\|X\|_2>1$, Proposition 1 implies that setting M below $\left\lceil\frac{C}{\min_{X\in\mathcal{D}\setminus\{\mathbf{0}_N\}}\|X\|_2}\right\rceil$, results in a sample mean with a higher variability in terms of its variance. Therefore, from the adversary's perspective, engineering M appropriately can be viewed informally as counteracting the addition of "overproportional" noise to the observed privatised model updates, since no clipping is triggered and thus the noise can be calibrated to a smaller value. However, increasing M beyond the bound in 15 does not further minimise the variance of the estimator in 17. In particular, the choices of the network presented in Proposition 1 suggest that no privatised version of the gradient $G_{X,P}$ (see 6) can improve the estimator's variance in 17 or improve the performance of the analytic reconstruction attack.

Since the adversary cannot directly manipulate the randomisation introduced by the DP mechanism to their benefit, their strategy can only rely in increasing the number of observed privatised versions of the input X and aggregating them, as presented in Proposition 1. As mentioned before, this approach makes the analytic reconstruction attack a classic mean estimation problem whose success depends on the variance of the estimator in 17. Due to the nature of the clipping term $\beta_C(X)$ (see 6), any non-usable observed information contained in $G_{X,P}$, increases the variance of this estimator (see Proof of Proposition 1 in Appendix B). Moreover, usable observed information contained in $G_{X,P}$ can be optimally incorporated into the mean estimation problem if it is a privatised version of the input X of the form in 8, making $G_{X,P}$ the gradient of a linear layer. Thus, the (insertion of the) linear layer and choosing M as in Proposition 1 makes $G_{X,P}$ redundant at best. In other words, in such a case, the entire, necessary information to estimate the target sample X is already contained in the gradient of the linear layer and $G_{X,P} \neq \mathbf{0}$ can only decrease the reconstruction quality by adding to the gradient norm and triggering the addition of "overproportional" noise.

If the adversary chooses M according to Proposition 1, then the behaviour of the sample mean \hat{X}_M coincides with \hat{X} 's one (see 17) independently of the specific value of M. Moreover, it is worth noting that \hat{X} is independent of the clipping norm C and its variance is exactly calibrated to the norm of the target X. Next, we examine some properties of \hat{X} .

Proposition 2. \hat{X} is the minimum variance unbiased estimator (MVUE) for X. Moreover, the expected mean squared error between the target X and \hat{X} is given by:

$$\mathbb{E}_X[\mathrm{MSE}_X(X,\hat{X})] = \sigma^2 ||X||_2^2. \tag{18}$$

A minimum variance unbiased estimator (MVUE) for the target X is desirable when dealing with the statistical estimation problem because such an estimator achieves the lowest expected mean squared error (MSE) and has the lowest variability in terms of its variance compared to all other estimators for X constructed with the same observations $\widetilde{\nabla}_{W_1}, ..., \widetilde{\nabla}_{W_M}$. The following result further emphasises the relevance of the estimator \widehat{X} regarding reconstructing the target X:

Theorem 1. Using the MSE as an optimality criterion, \hat{X} is the best achievable estimator and, thus, reconstruction for the target point X.

In this work, we utilise the MSE as an optimality criterion, because the mean squared error is equal to zero, i.e., $\text{MSE}_X(X,\hat{X}) = 0$, if and only if $X = \hat{X}$, namely whenever the adversary perfectly reconstructs the target X. A perfect reconstruction denotes the best case scenario for the adversary or conversely the worst case scenario concerning privacy preservation. Therefore, by Theorem 1, analysing the error between X and \hat{X} , specifically lower bounding this error, is sufficient to upper bound the reconstruction success of the adversary:

Corollary 1. Assume the adversary replaces the original model by a single, fully connected linear layer f(X) = WX of a single input $X \in \mathbb{R}^N$ by a learnable matrix $W \in \mathbb{R}^{M \times N}$, $M = \max\left(1, \left\lceil \frac{C}{\min_{X \in \mathcal{D} \setminus \{a_N\}} \|X\|_2} \right\rceil\right)$, and sets the loss function $\mathcal{L} : \mathbb{R}^M \to \mathbb{R}$ to be $\mathcal{L}(W, X) := \mathbf{1}_M^T W X$. Then, \hat{X} , as defined in Equation 17, is the

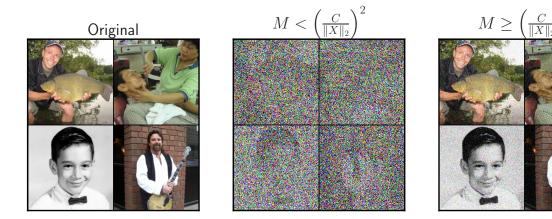


Figure 1: Demonstration of how an adversary can enforce to obtain the least amount of noise on the reconstruction for constant privacy parameters by adjusting the parameter M. If they increase M such that $M \geq \left(\frac{C}{\|X\|_2}\right)$, they will get the least amount of additive noise (right column). If they do not exceed this threshold, the additive noise is stronger ("overproportional") for the same privacy parameters C and σ (middle column). Middle column: M = 1, right column: M = 1000. All other parameters remain constant for both reconstructions: $\sigma = 5 \cdot 10^{-4}$, $C = 5.0 \cdot 10^3$, N = 150528. All images are from the ImageNet dataset (Deng et al., 2009).

MVUE with the lowest expected MSE and variability in terms of the variance compared to all other possible MVUE obtained under other choices regarding the model's architecture, hyperparameters and loss function.

Upper bounding the reconstruction success of an adversary that performs the attack described in Corollary 1 automatically bounds the success of an adversary that performs any other analytical gradient inversion attacks under DP. Therefore, we focus on the attack specified in Corollary 1 as well as on the optimal estimator \hat{X} as in 17. From now, we set \hat{X} to be the reconstruction of the adversary.

3.2 Determining the reconstruction success

In this section, we aim to determine the privacy effect DP has on the success of analytic reconstruction attacks performed against models privatised by the Gaussian Mechanism. To do so, we evaluate how well the adversary's reconstruction \hat{X} , as defined in 17, resembles its respective target input point X via the mean squared error (MSE), the peak signal-to-noise ratio (PSNR) and the normalised crossed-correlation (NCC). These metrics compare the input point X to its reconstructed version \hat{X} in different ways. However, we interpret it as high reconstruction success whenever they indicate high similarity between X and \hat{X} . We associate the reconstruction success of the adversary with the privacy guarantees provided by the DP framework. Additionally, we express the MSE and PSNR as random variables, examine their tail behaviour and provide guarantees in terms of the well-established notion of reconstruction robustness.

3.2.1 Reconstruction success measured by the MSE, PSNR and ReRo

First, we recover the value of the expected $\mathrm{MSE}_X(X,\hat{X})$ for all $X \neq \mathbf{0}_N$ stated in Equation 18. To interpret the result in 18, we consider the cases $C \leq \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$ and $C > \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$ separately.

Assume $C \leq \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$ holds. Then, the expected reconstruction error measured by the MSE between X and and the reconstruction \hat{X} can be lower bounded by the product of the squared maximum gradient norm C^2 and the squared noise scale σ^2 for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$, implying that the MSE between X and any of its unbiased estimators can be lower bounded by $C^2\sigma^2$. Hence, the higher the variance of the additive noise mechanism \mathcal{M} , the higher the expected error in terms of the MSE. It is, thus, easy to see that the privacy of the DP-mechanism and the expected MSE are positively correlated.

However, if we assume $C > \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$ holds, then the expected reconstruction error measured by the MSE between X and the reconstruction \hat{X} equals $\sigma^2 \|X\|_2^2$ for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$ and cannot be lower bounded solely using the DP parameters C and σ . In such a case, the result 18 illustrates the effect that increasing M, i.e., the number of reconstructions that are retrieved, has on the reconstruction success of the adversary, namely it cancels the addition of "overproportional" noise to all X such that $C > \|X\|_2$. Specifically, in this case, the error measured by the MSE is exactly calibrated to the norm of each target X. Due to the optimality of the estimator \hat{X} , we conclude as before, that the expected MSE between X and any of its unbiased estimators can be lower bounded by the product of the squared norm $\|X\|_2^2$ and σ^2 if $C > \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$.

It is difficult to make general claims regarding the occurrence of the cases $C \leq \min_{X \in \mathcal{X} \setminus \{\mathbf{0}_N\}} \|X\|_2$ and $C > \min_{X \in \mathcal{X} \setminus \{\mathbf{0}_N\}} \|X\|_2$ since, in practice, the maximum gradient norm C is usually an arbitrary hyperparameter, chosen on experience and not necessarily dependent on $\min_{X \in \mathcal{D} \setminus \mathbf{0}_N} \|X\|_2$. For instance, C = 1 is a very typical choice for classification tasks or whenever practitioners want to simplify all of their calculations.

To illustrate the behaviour of the $\mathrm{MSE}_X(X,\hat{X})$, we turn to the calculation of probabilistic bounds for this quantity. Concretely, fixing the target X, we can formulate the $\mathrm{MSE}_X(X,\hat{X})$ as a random variable determined by the randomness of the reconstruction \hat{X} and explore its tail behaviour. We are particularly interested in computing and bounding the probability of the error measured by the MSE falling below a certain threshold η . The choice of the threshold η indicates the interpretation of such a probabilistic bound since, for instance, η can be chosen such that an error measured by the $\mathrm{MSE}_X(X,\hat{X})$ below η characterises informative reconstructions.

Theorem 2.

$$MSE_X(X, \hat{X}) \stackrel{d}{=} \frac{\sigma^2 ||X||_2^2}{N} \cdot Y \quad with \quad Y \sim \chi_N^2, \tag{19}$$

where χ^2_N denotes the central chi-squared distribution with N degrees of freedom. In particular, for η given,

$$\mathbb{P}_{\hat{X}}(\mathrm{MSE}_X(X,\hat{X}) \le \eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 \|X\|_2^2}\right),\tag{20}$$

where Γ_R is the regularised gamma function.

By Theorem 2, the $\mathrm{MSE}_X(X,\hat{X})$ between any target point and its reconstruction is a chi-squared distributed random variable with N degrees of freedom multiplied by the product between 1/N and $\sigma^2 \|X\|_2^2$. Thus, it is no coincidence that the latter ratio equals the expected MSE given in Equation 18 because the expectation of a chi-squared distributed random variable with N degrees of freedom equals N. Therefore, we can recover the interpretation of $\sigma^2 \|X\|_2^2$ stated in the first part of this section, specifically the behaviour of this quantity for the cases $C \leq \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$ and $C > \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$. Moreover, Theorem 2 demonstrates that $\sigma^2 \|X\|_2^2$ determines not only the expectation of the MSE but also its distribution, highlighting the influence that the DP parameters C and σ and the parameter M, chosen by the adversary, have on the error measured by the MSE and consequently on the reconstruction success of the adversary measured by the MSE.

Informally, the chi-squared distribution with N degrees of freedom converges to the normal distribution for $N \to \infty$. Therefore, for cases where the dimension of the training data is "very high", it is pertinent to additionally analyse the asymptotic behaviour of the $\mathrm{MSE}_X(X,\hat{X})$ for $N \to \infty$. In the appendix, in Proposition B.3, we formally present the convergence in distribution of the $\mathrm{MSE}_X(X,\hat{X})$ for $N \to \infty$. However, we leave out an in-depth discussion of the result in Proposition B.3.

Expressing the MSE as a random variable enables the formulation of its tail behaviour as given in 20. We note that the result in Equation 20 resembles the notion of reconstruction robustness as given in Definition 1. In the following, we generalise this result to lose the dependency on the target point X and translate the probabilistic bound in 20 into reconstruction robustness:

Proposition 3. Let η given. Then, for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$,

$$\mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_{X}\left(X,\hat{X}\right) \leq \eta\right) \leq \Gamma_{R}\left(\frac{N}{2}, \frac{N\eta}{2\sigma^{2} \min_{X \in \mathcal{D}} \|X\|_{2}^{2}}\right),\tag{21}$$

where Γ_R is the regularised gamma function. Moreover, the DP-mechanism \mathcal{M} is $(\eta, \gamma(\eta))$ -reconstruction robust with respect to the MSE for any reconstruction and $\gamma(\eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 \min_{X \in \mathcal{D}} \|X\|_2^2}\right)$. If $C \leq \min_{X \in \mathcal{X}} \|X\|_2$ holds, then \mathcal{M} is $(\eta, \gamma'(\eta))$ -reconstruction robust with respect to the MSE for any reconstruction and $\gamma'(\eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2C^2}\right)$.

Depending on the training data, other reconstruction error functions might be more informative when assessing the similarity between the target X and the reconstruction \hat{X} . Concretely, we consider the peak signal-to-noise-ratio (PSNR) (see Definition 3) as a reconstruction quality function next. We note that the PSNR is defined via the MSE. In particular, as the MSE decreases, the PSNR increases, implying that if there exists one reconstruction that minimises the MSE and this minimum is non-zero, this reconstruction maximises the PSNR. Thus, by Theorem 1 and Corollary 1, the chosen reconstruction \hat{X} is also optimal with respect to the PSNR.

Next, we provide probabilistic bounds for the PSNR between X and \hat{X} to determine its tail behaviour. To do so, we assume $\max_{X \in \mathcal{D}} \max(X)$ and $\min_{X \in \mathcal{D}} \min(X)$ are known quantities. This information is known whenever the adversary has access to the range of the data, such as when the training data are images. Unlike the MSE, the PSNR increases as the similarity between X and \hat{X} grows. Therefore, we formulate a bound for the probability of the event when the PSNR exceeds a certain threshold η . In contrast to the MSE, η can be chosen such that a similarity measured by the PSNR above η characterises informative reconstructions.

Proposition 4. Assume $\max_{X \in \mathcal{D}} \max(X)$ and $\min_{X \in \mathcal{D}} \min(X)$ are known quantities. Then, for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$,

$$\mathbb{P}_{\hat{X}}\left(\mathrm{PSNR}_{X}\left(X,\hat{X}\right) \geq \eta\right) \leq \Gamma_{R}\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^{2} \min_{X \in \mathcal{D} \setminus \{\theta_{N}\}} \|X\|_{2}^{2}}\right),\tag{22}$$

for $\tilde{\eta}(\eta) := 10^{-\frac{\eta}{10}} \left(\max_{X \in \mathcal{D}} \max(X) - \min_{X \in \mathcal{D}} \min(X) \right)^2$, and for Γ_R being the regularised gamma function. If $C \leq \min_{X \in \mathcal{D}} \|X\|_2$ holds, then

$$\mathbb{P}_{\hat{X}}\left(\text{PSNR}_X\left(X,\hat{X}\right) \ge \eta\right) \le \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 C^2}\right),\tag{23}$$

for all $X \in \mathcal{D} \setminus \{\boldsymbol{o}_N\}$.

To reformulate the result in Proposition 4 into reconstruction robustness terms, we utilise the negative PSNR (-PSNR) as a reconstruction error function:

Proposition 5. Assume $\max_{X \in \mathcal{D}} \max(X)$ and $\min_{X \in \mathcal{D}} \min(X)$ are known quantities, and let Y_X be any possible reconstruction. Then, for all $X \in \mathcal{D} \setminus \{\boldsymbol{0}_N\}$,

$$\mathbb{P}_{Y_X}\left(\mathrm{PSNR}_X\left(X, Y_X\right) \ge \eta\right) \le \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 \min_{X \in \mathcal{D} \setminus \{\boldsymbol{\theta}_N\}} \|X\|_2^2}\right),\tag{24}$$

for Γ_R being the regularised gamma function and $\tilde{\eta}(\eta)$ as defined in Proposition 4. In particular, the DP-mechanism \mathcal{M} is $(-\eta, \tilde{\gamma}(\tilde{\eta}(\eta)))$ -reconstruction robust with respect to the negative PSNR (-PSNR) for any analytic reconstruction and $\tilde{\gamma}(\tilde{\eta}(\eta)) = \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 \min_{X \in \mathcal{D} \setminus \{\theta_N\}} \|X\|_2^2}\right)$. Moreover, if $C \leq \min_{X \in \mathcal{D}} \|X\|_2$ holds, then the DP-mechanism \mathcal{M} is $(-\eta, \tilde{\gamma}'(\tilde{\eta}(\eta)))$ -reconstruction robust with respect to -PSNR for any reconstruction, $\tilde{\gamma}'(\tilde{\eta}(\eta)) = \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2C^2}\right)$.

The CDF of the $MSE_X(X, \hat{X})$ and the survival function (SF) of the $PSNR_X(X, \hat{X})$ for different values of σ , N and $||X||_2$ are depicted in Figure 3. We remark that bounding the probability of a specific reconstruction loss allows for interpreting our results in terms of $(\eta, \gamma(\eta))$ -ReRo (compare Figure 3 to Propositions 3 and 5).

3.2.2 Reconstruction success measured by the NCC

Let us consider the linear correlation given by the NCC between a target X and its reconstruction \hat{X} . The NCC can be measured empirically using their entries $\{x_1,...,x_N\}$ and $\{\hat{x}_1,...,\hat{x}_N\}$ as sample sets, thereby

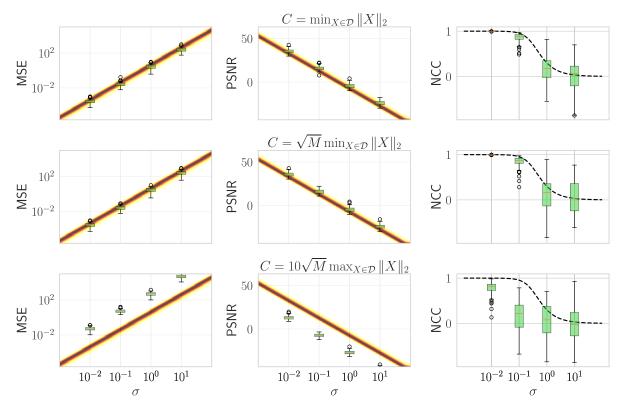


Figure 2: Empirical reconstruction results overlaid on our theoretical results. Each column shows one reconstruction metric: Left: MSE, Middle: PSNR, Right: NCC. For MSE and PSNR, the probability density function (PDF) of the theoretical distribution is encoded in colour, where white corresponds to low values and black to high values along each σ . For the NCC, the dashed line shows the theoretical bound. For all plots, we have performed empirical reconstructions of four-dimensional data samples (N=100) with M=100, which are sampled from a uniform distribution $\mathcal{U}(0,1)$. The resulting reconstruction metrics are displayed as boxplots, where the orange line shows the mean reconstruction result, the green box displays the interquartile range (IQR) from the first quartile to the third quartile, and the whiskers extend to the data point, which is the last within 1.5 of the IQR. Further outliers are displayed as circles. First row shows the case where $C=\min_{X\in\mathcal{D}}\|X\|_2$. Hence, the sensitivity is exhausted, and the resulting reconstruction cannot be further improved for this scenario even if M=1. In the second row, we have $C=\sqrt{M}\min_{X\in\mathcal{D}}\|X\|_2$. Here, several data samples only exceed the sensitivity threshold because multiple observations are combined. In the last row, no data sample is clipped as C is set to be larger than any data norm and M has not been chosen according to the optimal attack. Therefore, the sensitivity limit is not reached, and "overproportional" noise is applied. This decreases the reconstruction success.

computing the *sample* NCC. The calculation of the sample NCC and a brief comment regarding its usefulness can be found in the appendix in Proposition B.4 and Remark B.1, respectively. However, in this section, we concentrate on the (theoretical) correlation between X and \hat{X} .

To compute the NCC between X and \hat{X} , we apply the following strategy: First, let the target X be fixed. We assume there exists a continuous, one-dimensional random variable x distributed in such a way that $\{x_1, ..., x_N\}$ are probable samples from this distribution. In particular, we let $[\min_{i \in \{1,...,N\}} x_i, \max_{i \in \{1,...,N\}} x_i]$ be the support of x and set $\text{Var}(x) \leq \|X\|_2^2/N$. The latter assumption is motivated by the following fact: X is an N-dimensional vector, hence, drawing a random element from its entries $\{x_1, ..., x_N\}$ can be represented by a discrete, uniformly distributed random variable with support $\{x_1, ..., x_N\}$ and variance bounded by $\|X\|_2^2/N$. However, for our analysis, x needs to be a continuous random variable. Hence, we define x to be continuous, but maintain the range of the support and the variance of its discrete counterpart. Analogously, we construct the continuous, one-dimensional random variable \hat{x} such that $\hat{x} := x + \zeta$, for $\zeta \sim \mathcal{N}(0, \sigma^2 \|X\|_2^2)$ independent

of x. By definition, $\hat{x}_1, ..., \hat{x}_N$ are probable samples of the random variable \hat{x} . It is easy to see, that measuring the correlation between x and \hat{x} is equivalent to measuring the correlation between X and \hat{X} .

Proposition 6. Let x and \hat{x} be the two random variables as defined above. Then,

$$NCC(x,\hat{x}) = \sqrt{\frac{1}{1 + \sigma^2 ||X||_2^2 / Var(x)}} \le \sqrt{\frac{1}{1 + \sigma^2 N}}.$$
 (25)

By Proposition 6, the NCC (x, \hat{x}) is determined by the ratio between the entry-wise variance of the reconstruction \hat{X} given by $\sigma^2 \|X\|_2^2$ and variability within the target X given by $\operatorname{Var}(x)$. Therefore, if this ratio is not "high enough", we conclude from Equation 25 that the DP mechanism's noise is insufficient to disrupt the linear association between x and \hat{x} , or, conversely, between X and its reconstruction \hat{X} . Moreover, due to the nature of the reconstruction \hat{X} (see 17) a perfect (or "good enough") correlation between x and \hat{x} implies a perfect (or "good enough") reconstruction of the target X. In particular, Equation 25 equals one if and only if the noise scale σ equals zero, i.e., when no noise is added to model updates before release and the observed gradient G_X is not privatised during training. Since any non-affine transformation of the observed privatised gradient would distort the linear dependency between X and the non-privatised gradient, it is easy to see that \hat{X} is an optimal reconstruction with respect to the NCC as well.

If $\sigma \neq 0$, we observe that the NCC (x,\hat{x}) decreases with increasing noise scale σ . However, Proposition 6 demonstrates that the NCC between X and \hat{X} highly depends on the dimension N. In particular, the right-hand side of Expression $25 \in \mathcal{O}(\sqrt{1/N})$. Thus, for increasing dimension N, the correlation between X and \hat{X} measured by the NCC decreases independently of the noise of the DP mechanisms as long as $\sigma \neq 0$. Therefore, in the context of our work, for high values of N, the NCC is not necessarily a good metric to assess the reconstruction success of the adversary.

4 Visualisation and Interpretation

In this section, we show the implications of our bounds and their correspondence to empirical results. Specifically, we illustrate the effect of the parameters σ , C, M, N and $\|X\|_2$ on our results to allow for a more intuitive understanding of their behaviour. The code to reproduce our figures can be found on https://anonymous.4open.science/r/FromMeanToExtreme-46D1.

In Figure 1, we demonstrate the influence of M on the reconstruction result. Only by selecting an M large enough to exceed $\left(\frac{C}{\|X\|_2}\right)^2$ no overproportional noise is added and the reconstruction cannot be further improved. The same effect is present in Figure 2. Here, we display the behaviour of MSE, PSNR, and NCC under variation of the privacy parameters C and σ . For the MSE and PSNR, we plot the PDF for each value of σ as colour coding, where white areas have low probability and dark show high probability areas. For the NCC, we plot the bound on the expected reconstruction success. We create an artificial dataset consisting of 100 four-dimensional data points sampled from a uniform distribution $\mathcal{U}(0,1)$. An adversary attempts to reconstruct these data points using the attack specified in Section 2.2. The empirical reconstruction success measured by the respective metrics is displayed as boxplots. We observe, analogously to Figure 1, only if the clipping threshold is exceeded, either because the data norm is larger than C (first row) or because the adversary set M to be large (second row), the empirical reconstruction results overlap with the worst case distributions we have derived earlier. However, if this threshold is not exceeded then the additive noise is larger, which in turn increases the MSE and respectively decreases the PSNR and NCC.

In Figure 3, we visualize the influence of σ , N, and $||X||_2$ on the distribution of the MSE and PSNR. If a parameter is not varied in a specific row this parameter is also set to be 1. Unsurprisingly, we observe in the first row that increasing the noise scale σ leads to smaller probabilities for high-fidelity reconstructions. In the second row, we observe that the distributions become "steeper" for a changing data dimensionality, i.e., they are not spread out as much for high-dimensional data. This means that for high dimensional data it becomes increasingly unlikely to observe errors, which largely differ from the expected error. Of note, the expected error for the MSE remains constant, but changes for the PSNR. This is due to the conversion to an exponential scale, a transformation which the expected value is not robust against. In the last row, we see

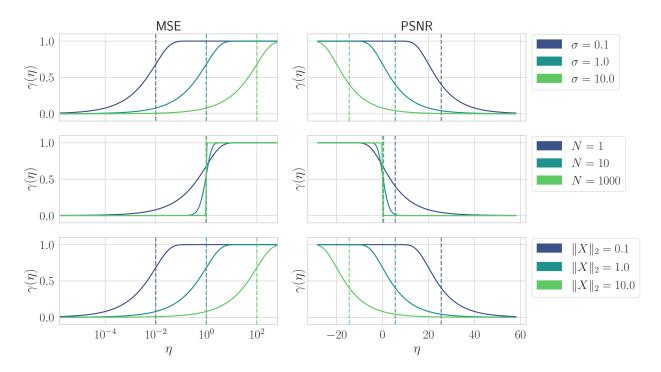


Figure 3: Cumulative Distribution Function of the MSE and Survival Function of the PSNR for varying parameters σ , N and $\min_{X \in \mathcal{X}} \|X\|_2$. If not varied then parameters are set to be $\sigma = 1$, N = 1, $\max_{X \in \mathcal{D}} \|X\|_2 = 1$. Vertical dashed lines show the corresponding bounds on the expected values. For a given error threshold η , we have a risk probability lower than or equal to $\gamma(\eta)$. Note that lower values of the PSNR denote better reconstruction results. Hence, for $(\eta, \gamma(\eta))$ -ReRo the negative PSNR needs to be considered (compare Proposition 5).

the impact of the ℓ_2 -norm of the data sample on the reconstruction error distribution. Here, we observe that if all other parameters remain unchanged, data samples with larger norms have higher errors. This is because the MSE is scaled linearly along with its input. More concretely, the same data sample and reconstruction have a different MSE if both are multiplied by the same constant, namely the multiplication of the constant and the previous MSE, although semantically the reconstruction contains the same information (compare Supplementary Figure 5). The PSNR is principally robust against this behaviour as it incorporates the data range. However, for the above experiment, the data range was kept constant.

Our work attempts to address a central question: How can practitioners choose the privacy parameters σ and C for their specific model to protect the training data? As explained in Section 3.1, the adversary can counteract the privatising effect of the maximum gradient norm C under the given threat model. Therefore, C does not (necessarily) influence the reconstruction success and can be set depending on the model application to retrieve the best-performing outcome. However, reconstruction success is determined by the selected noise scale σ and depends on the norm of the data points and their dimension N. Given a fixed data set with samples of the same dimension N, practitioners can choose some data samples and observe the impact various choices of the noise scale σ can have on their reconstruction success (see results in Sections 3.2.1 and 3.2.2). This is particularly important for assessing thresholds η (see Definition 1) that describe informative reconstructions in terms of the MSE or PSNR. Understanding what are appropriate choices for such thresholds, practitioners can calculate the probability γ of occurrence of such reconstructions and obtain the levels of protection determined by the reconstruction robustness. We highlight that no one general threshold for the MSE, PSNR and NCC determines a successful reconstruction for all cases as one can see, for instance, in Supplementary Figure 6.

5 Discussion and Conclusion

In this work, we formalise bounds over three reconstruction metrics for any data protected by the Gaussian mechanism. This is motivated by several state-of-the-art attacks, which found ways to extract such privatised data in settings and are designed to be real-world applicable. Providing bounds for the best current reconstruction attacks allows practitioners to make informed decisions to defend against this family of attacks based on mathematical guarantees. Considering the problem of the so-called privacy-utility trade-off, it is important to choose privacy parameters adjusted to the specific threat model in order to achieve the lowest utility penalty while maintaining an acceptable level of data security.

The MSE, PSNR, and NCC show different but complementary notions of reconstruction success. The MSE is a standard, wide-spread error metric. However, a notable drawback in practice is that it is only comparable to other MSE-values if it is zero and thus, the data is perfectly reconstructed. Namely, general assessments regarding the quality of a reconstruction based on the MSE can only by made on a comparative basis with the exception of an MSE equal to zero which signifies that the data is perfectly reconstructed. For many applications such as optimisation tasks, this is sufficient. However, for a metric measuring reconstruction success, it is desirable to be comparable to any other value. Moreover, the MSE is not robust to scaling, implying that if the data and its reconstruction are multiplied by the same scalar value, the resulting MSE changes (see Figure 5) even though no real development in the difference between the data and its reconstruction occurred. Therefore, if the privacy parameters are decided with respect to a certain bound on the MSE, practitioners must be aware of the exact scale of their dataset. PSNR can correct for this effect by setting the value range appropriately. The NCC is, by design, robust against any (positive) multiplication of the original data sample or its reconstructed version. This is because it only measures the linear correlation between these entities. Most importantly, for the MSE and PSNR, we can, for any set of privacy parameters, calculate the risk of being above a certain value, which corresponds to the notion of (η, γ) -ReRo. This gives any practitioner the maximum flexibility of choosing a reconstruction error and its risk, which is acceptable for their workflow. We argue that for a holistic decision for a certain set of privacy parameters, practitioners should consider all three metrics for their specific dataset in order to decide on bounds that sufficiently protect against reconstruction.

Even though an adversary cannot manipulate the DP mechanism, we showed they can improve their reconstruction result by setting M to be "large enough" so that even for large sensitivity thresholds C, the data is clipped and no overproportional noise is added. However, we demonstrate that the randomisation introduced by the DP mechanism still bounds the reconstruction success independently of the choice of M. Moreover, we note that increasing M also increases the computational complexity in $\mathcal{O}(M)$, whereas the enforced increase in the gradient norm is only in $\mathcal{O}(\sqrt{M})$. Therefore, in real-world applications, an adversary needs to trade off the likelihood of an optimal reconstruction against the computational requirements. At the same time, the AI practitioner has an interest in setting the clipping threshold C not too large, as this would again lead to overproportional additive noise and thus lead to stronger utility losses of a trained network architecture. Hence, for any defence evaluation, the optimal reconstruction under a specific set of privacy parameters should be considered as a realistic outcome.

We note that while our bounds are directly transferable to some of the currently best real-world reconstruction attacks, none of these attacks exploits data priors. Depending on the specific data, an adversary might have related data or prior knowledge that they can exploit to further improve their reconstruction success. This was, for example, empirically demonstrated by Schwethelm et al. (2024). In that case, our bounds also do not hold anymore, and other bounds, which consider prior knowledge, need to be applied, e.g., Hayes et al. (2023). However, years of research in generating artificial data are proof that modelling the distribution of a specific kind of data is not straightforward, especially not in providing a mathematical description of it. Hence, it is likely that only optimistic (our work) and pessimistic (Hayes et al., 2023) mathematical bounds can be provided.

One limitation of our work is that the NCC cannot be interpreted in a (η, γ) -ReRo notation. This is due to the fact that the NCC is an descriptive property of the interaction between two random variables or two data samples such as an image and its reconstructed version. The NCC is determined by the covariance between these quantities and their standard deviations, attributes that are deterministic in our

work. Therefore, although we have established a bound indicating that the NCC does not exceed a certain value, individual samples drawn from the distribution are likely to deviate from this bound due to the probabilistic nature of the random variable (see Remark B.1). Furthermore, we note that the NCC varies drastically for different dimensions and impedes comparison of the reconstruction success between scenarios with unequal dimensionality. Especially for $N \to \infty$, only perfect reconstructions will lead to NCC > 0, rendering its practical use in very high-dimensional settings to be limited.

Other well-established metrics, such as the normalized mutual information (NMI), structural similarity (SSIM) or even perceptual losses, also empirically seem to directly correspond to the influences we observed (see Figure 4). Deriving mathematical bounds on these metrics is left for future work. Analogously, in this work, we have considered the case for a single query of one data sample. However, a central field of research in the context of differentially private neural network training is accounting the privacy loss over repeated queries with subsampling amplification. Repeated queries would obviously allow for a better reconstruction as the result of these queries could be averaged, which does not affect the underlying signal but, in expectation, cancels out the additive noise. Subsampling, on the other hand, would impede this strategy, as the adversary needs to correctly match queries which are based on the same sample. It remains for future work to analyse how these counteracting effects affect the results of our work.

The overarching question of providing bounds for adversarial attacks remains: how can we optimally choose the *least* amount of privacy in order to not introduce utility losses but still provide reasonable protection? We argue that solely investigating worst-case bounds introduces stronger privacy-utility trade-offs than necessary. Our work provides the theoretical bounds on the risks for a branch of real-world reconstruction attacks and allows for a more precise evaluation of the potential privacy leakage in these settings. Specifically, we allow a practitioner to tune their privacy parameters from *mean* attack success to *extreme* attack success given by the expected reconstruction success and the tail behaviour of the reconstruction, respectively. However, we note that our bounds should be seen in context and as an augmentation to the worst-case bounds, as adversaries exploiting data priors outperform the reconstruction success our bounds suggest. We see our work as a first step towards a broad suite of threat model analyses as part of a full system model card, which provide a tailored risk report for practitioners for their specific settings in addition to a contextualisation of how changes in the capabilities of the adversary can influence the attack success probabilities.

Author Contributions

Acknowledgments

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A Background

A.1 Prior Work

Reconstruction Bounds DP is a standard privacy-preserving technique in ML. We assume the reader to be familiar with basic DP terminology and refer to Dwork et al. (2014) for a comprehensive background explanation. Additionally, we also refer to Mironov (2017) for a detailed introduction to Rényi-DP. Guo et al. (2022) laid the foundation for bounding the success of data reconstruction attacks by deriving a bound based on Rényi-DP. However, their work is limited to $\alpha = 2$, which leads to loose bounds, as shown by Hayes et al. (2023). Balle et al. (2022) further formalised and tightened the success upper-bound of reconstruction attacks by introducing the notion of reconstruction robustness (ReRo). Specifically, they defined (η, γ) -ReRo as the probability of an adversary having a reconstruction error lower or equal to η to be lower or equal than γ (see Definition 1). Intuitively, a practitioner here can define an error metric on the reconstruction and a threshold η where they think that if the error metric is above this threshold, the reconstruction is not usable anymore. Then an algorithm which is (η, γ) -ReRo gives the practitioner a guarantee that the probability of having a reconstruction below this threshold is at most γ . Based on this work, Hayes et al. (2023) empirically analysed reconstruction robustness of a worst-case adversary, with access to all intermediate outputs (gradients), all training data except the target sample as well as an additional collection of data samples, including the target sample (prior set). Kaissis et al. (2023a) extended this work by providing closed-form ReRo-bounds for the Gaussian and the Laplacian mechanism. Under the above-mentioned worst-case assumptions, the reconstruction attack is equivalent to matching the privatised output of a query function (in the case of neural network this corresponds to the intermediate gradients) to the correct sample in the prior set. Hence, the adversary's success can be represented as a binary variable as they either succeed, i.e., achieve perfect reconstruction, or fail. The authors show that using DP formally bounds the success rate of such an adversary and fulfils $(0, \gamma)$ -ReRo (Hayes et al., 2023; Kaissis et al., 2023a). Kaissis et al. (2023b) challenged the idea that the worst-case scenario is realistic and instead examined privacy guarantees (including the bounds for the success of a reconstruction attack) under a slightly relaxed threat model. However, the proposed assumptions still cannot be considered to describe a typical real-world workflow either, as, for example, the adversary requires a perfect reconstruction algorithm and only checks whether it succeeded. Cummings et al. (2024) recently proposed a taxonomy of threat models and alongside formulated bounds for a specific relaxed threat model. This analysed threat model does not impose the strict requirements to have knowledge about the concrete dataset. Furthermore, they assume that the distribution of the data is known and samples are drawn i.i.d. from this distribution. However, this implies that bounds are not over "most distinguishable" data samples. They can show that their bounds fulfil a strictly weaker notion of (η, γ) -ReRo, which they term distributional reconstruction robustness (DistReRo). More realistic threat models have also been investigated for Membership Inference Attacks (MIA), another type of attack where the adversary tries to infer whether a specific data point has been part of the training set. Nasr et al. (2021) construct empirical lower bounds on the probability that an MIA is successful depending on the adversary's capabilities and compare them to the success' upper bound given by the most powerful adversary. MIA is a much simpler attack than reconstruction attacks since only one bit of information is reconstructed (whether a subject was part of the training set or not). However, Nasr et al. (2021) conclude that relaxing the threat model considerably reduces the practical risk of attack success even for this simpler attack type.

Analytic Gradient Inversion Attacks Our work is inspired by a branch of reconstruction attacks, which have recently received substantial attention, as they (a) are conceivable and hard to detect in real-world scenarios, (b) allow for an (almost) perfect reconstruction success, and (c) are based on computationally low-cost analytical procedures, in contrast to optimisation based attacks (Geiping et al., 2020). Fowl et al. (2022) and Boenisch et al. (2023) concurrently proposed attacks on a decentralised, federated learning setup. They showed that potentially unnoticed modifications to a deep learning model by a malicious central server can lead to the input data being stored in the gradients of the model. In particular, they exploit that (a) linear layers are a common type of neural network component and (b) they have the property to encode their input in the gradients calculated by the chain rule (backward pass). The former makes them unsuspicious for any check, which is unaware of the latter. To recover the entire input such a linear layer is prepended to the original network architecture. Here, the input is encoded once for each projection, i.e., a layer projecting an

input $x \in \mathbb{R}^N$ to $b \in \mathbb{R}^M$, recovers M versions of x (bins). Specifically, this is achieved by performing an element-wise division between the gradient of the bias and the gradient of the weight of the linear layer, the exact input can be recovered. For batch sizes larger than one or multiple consecutive update steps, each bin encodes an arbitrary amount of inputs to the network. This is problematic, as any reconstruction more than one (meaningful) input, is an average of its inputs, which implies that the signals of these inputs overlap and may make them meaningless. While the authors propose several sophisticated strategies to avoid this effect, it cannot be impeded entirely. Hence, the worst-case is a single sample update step (i.e. batch size is one) where no data can overlap.

Based upon the same attack principle, Feng & Tramèr (2024) proposed an extension that no longer relies on the intermediate gradients but only requires access to the model before and after training. This strategy makes such attacks much more critical for real-world scenarios, as practitioners would only need to use a manipulated pretrained model in order for the training data to be stored in the weights. From there, the adversary can read it out after training. Feng & Tramèr (2024) carry out this attack by designing and inserting a so-called data trap into the model, i.e., a part of the network where the data is encoded and, with high probability, is not contaminated by any other gradients. By doing so, they circumvent the problem of overlapping data after multiple gradient update steps (yet not larger batch sizes), such that the attack is independent of the access to intermediate gradients. Furthermore, they show that these attacks can be performed without leading to substantial performance losses and with only minimal changes to the architecture, making it challenging to detect such an attack.

A.2 Metrics

Concretely, we introduce three metrics used in our analysis: the mean squared error (MSE), the peak signal-to-noise ratio (PSNR) and the normalised cross-correlation (NCC). These metrics allow us to assess the similarity and relation between an input point and its reconstruction and, consequently, determine the quality of said reconstruction.

In the following, let $X \in \mathbb{R}^N$, $N \in \mathbb{N}$, represent a fixed target input point and $Y_X \in \mathbb{R}^N$, be its estimator that serves as its reconstructed version. Moreover, we let x_i and $y_{i,X}$, for $i \in \{1, ..., N\}$, denote the *i*-th element of X and Y_X , respectively. First, we turn to the mean squared error (MSE), defined as average of the squared distances between the entries of the input vector and the entries of its reconstruction:

Definition 2. The mean squared error (MSE) between a fixed input X and its estimator Y_X is given by

$$MSE_X(X, Y_X) = \frac{1}{N}||X - Y_X||_2^2,$$

where $||\cdot||_2$ denotes the euclidean norm.

The MSE is one of the most common measures to quantify the reconstruction error. A decreasing MSE denotes a high similarity between X and Y_X or, conversely, a low reconstruction error. In particular, we note that $X = Y_X$ if and only if the $\text{MSE}_X(X, Y_X)$ equals zero. Nevertheless, while the MSE is widely used as an error function, it has unfavourable properties when assessing the reconstruction quality. Most importantly, it is not robust to scaling, hence, two reconstructions with the same $\text{MSE} \neq 0$ could have a very different reconstruction quality. This drawback is well-known and to overcome this, the PSNR, which is based on the MSE, has been established as a standard metric in signal processing applications:

Definition 3. The peak signal-to-noise-ratio (PSNR) between a fixed input point X and its estimator Y_X is given as

$$\mathrm{PSNR}_X(X,Y_X) = 10 \cdot \log_{10} \left(\frac{(\max(X) - \min(X))^2}{\mathrm{MSE}(X,Y_X)} \right),$$

with

$$\max(X) = \max_{i \in \{1,...,N\}} x_i$$
, and $\min(X) = \min_{i \in \{1,...,N\}} x_i$.

The PSNR contrasts the maximal range of the values of the entries of X to the MSE between X and its estimator Y_X . Thus, the PSNR puts the error measured in terms of the MSE into the context of the input

value range of X. By that, the PSNR is robust to a linear scaling of the data and and can be better compared across different scenarios. In this regard, the PSNR is a superior indicator of the reconstruction quality, and, although it is based on the same measurement, has advantages over the MSE (Wang & Bovik, 2009). Notably, unlike MSE, the PSNR is no error but a reconstruction quality function where larger values correspond to higher fidelity (and lower error) between X and Y_X .

Furthermore, we note that fixing the target input point X enables formulating the MSE and the PSNR in terms of the randomness of the estimator Y_X . In particular, if we view Y_X as the sample of a random variable, the MSE and the PSNR can be formulated as random variables and we can examine their tail behaviour (see Section 3.2).

Next, we introduce the NCC:

Definition 4. The normalised cross-correlation (NCC) (Rodgers & Nicewander, 1988) between a fixed input point X and its estimator Y_X is defined as

$$NCC(X, Y_X) = \frac{Cov(X, Y_X)}{\sigma_X \sigma_{Y_X}},$$

where

$$Cov(X, Y_X) = \frac{1}{N} \sum_{i=1}^{N} (x_i - \mathbb{E}_x[x])(y_{i,X} - \mathbb{E}_{y_X}[y_X]),$$

for $\mathbb{E}_x[x]$, $\mathbb{E}_{y_X}[y_X]$ denoting the numerically obtained expected values within a sample. Moreover, $\sigma_X = \sqrt{\operatorname{Var}(X)}$ and $\sigma_{Y_X} = \sqrt{\operatorname{Var}(Y_X)}$, for $\operatorname{Var}(X)$ and $\operatorname{Var}(Y_X)$ being the sample variances of X and Y_X , respectively.

We note that the NCC is equivalent to the Pearson's Correlation Coefficient, which is common in statistical testing. As both names suggest, this metric measures the linear correlation between a target point X and its estimated counterpart Y_X instead of measuring a difference. Hence, as opposed to metrics such as the MSE or PSNR, it is robust to linearly transformed inputs. High values of the NCC denote a high linear correlation between X and Y_X , which, depending on the context, can imply a high similarity between X and Y_X . For more information on interpreting the NCC results in the context of this work, we refer to Section 3.2.

B Proofs

In the following, we give the proofs for our theoretical results.

Proposition B.1. If the scaling factors $\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_1}, ..., \beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_M}$ are unknown, then there is no "realisable" unbiased estimator of the target X that can be constructed solely using the observed privatised gradients $\widetilde{\nabla}_{W_1}, ..., \widetilde{\nabla}_{W_M}$.

Proof. First, we show there is no deterministic transformation $T_X : \mathbb{R}^N \to \mathbb{R}^N$, such that $T_X \left(\widetilde{\nabla}_{W_j} \right)$, $j \in \{1, ..., M\}$, is a "realisable" unbiased estimator of X.

Let X be a fixed reconstruction target point and w.l.o.g. assume $X \neq \mathbf{0}_N$, for $\mathbf{0}_N$ the N-dimensional zero vector. Moreover, for a fixed $j \in \{1, ..., M\}$, let $T_X\left(\widetilde{\nabla}_{W_j}\right)$ be a realisable, unbiased estimator of X. Since

$$\widetilde{\nabla}_{W_j} = \beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j} X + \xi_j, \qquad \xi_j \sim \mathcal{N}(\mathbf{0}_N, C^2 \sigma^2 I_N),$$

 $T_X\left(\widetilde{\nabla}_{W_j}\right)$ must be an affine function, because that is the only transformation that can invert a multiplication and a sum. Hence, $T_X\left(\widetilde{\nabla}_{W_j}\right)$ has the following form:

$$T_X\left(\widetilde{\nabla}_{W_j}\right) = A\widetilde{\nabla}_{W_j} + b,\tag{26}$$

for $A \in \mathbb{R}^{N \times N}$ invertible and constant, and $b \in \mathbb{R}^{N \times N}$ constant. In particular, A and b are not functions of X, since $T_X\left(\widetilde{\nabla}_{W_j}\right)$ is a realisable estimator. Next, we compute the expectation of $T_X\left(\widetilde{\nabla}_{W_j}\right)$ for X fixed:

$$\mathbb{E}_{X}\left[T_{X}\left(\widetilde{\nabla}_{W_{j}}\right)\right] = \mathbb{E}_{X}\left[T_{X}\left(\beta_{C}(X)\frac{\partial g(f(X))}{\partial f(X)_{j}}X + \xi_{j}\right)\right]$$

$$= \mathbb{E}_{X}\left[A\beta_{C}(X)\frac{\partial g(f(X))}{\partial f(X)_{j}}X + A\xi_{j} + b\right]$$

$$= A\beta_{C}(X)\frac{\partial g(f(X))}{\partial f(X)_{j}}X + b.$$

 $T_X\left(\widetilde{\nabla}_{W_j}\right)$ is an unbiased estimator of X if and only if $\mathbb{E}_X\left[T_X\left(\widetilde{\nabla}_{W_j}\right)\right] = X$, which is equivalent to the following:

$$A\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_j}X + b = X \iff \left(\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_j}A - I_N\right)X + b = \mathbf{0}_N.$$
 (27)

Since b cannot depend on X and $\left(\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_j}A - I_N\right)X$ is a linear transformation of X, we cannot set $\left(\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_j}A - I_N\right)X = -b$ and solve the equation. Therefore, the equations in 27 are satisfied if and only if

$$\left(\beta_C(X)\frac{\partial g(f(X))}{\partial f(X)_j}A - I_N\right)X = \mathbf{0}_N \quad \text{and} \quad b = \mathbf{0}_N.$$
(28)

Since $X \neq \mathbf{0}$, it is easy to see that $A\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j} X + b = X$ holds if and only if

$$A = \frac{1}{\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_i}} I_N \quad \text{and} \quad b = \mathbf{0}_N,$$

implying that A depends on $\beta_C(X) \frac{\partial g(f(X))}{\partial f(X)_j}$, a quantity that can only be computed knowing X and contradicting the assumption that $T_X(\tilde{\nabla}_{W_j})$ is a realisable estimator. Therefore, for any $j \in \{1, ..., M\}$, there is no realisable unbiased estimator of X that can be constructed using the observed privatised gradient $\tilde{\nabla}_{W_j}$.

It is easy to see that, for this setting, if there was an unbiased estimator of X that can be constructed with $\widetilde{\nabla}_{W_1},...,\widetilde{\nabla}_{W_M}$, then there would exists at least one $j\in\{1,...,M\}$ such that an unbiased estimator of X can be constructed with $\widetilde{\nabla}_{W_j}$. However, since there is no $j\in\{1,...,M\}$ such that an unbiased estimator of X can be constructed with $\widetilde{\nabla}_{W_j}$, it follows by contraposition, that there is no unbiased estimator of X than can be constructed with $\widetilde{\nabla}_{W_1},...,\widetilde{\nabla}_{W_M}$.

Proposition B.2. The coordinte-wise variance of sample average \hat{X}_M stated in 13 is lower bounded by $\sigma^2 \|X\|_2^2$ for $M \to \infty$ and for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$.

Proof. Let the iteration step be fixed and observe the sample mean \hat{X}_M given in 13 with distribution described in 14. Then, consider the *i*th entry of \hat{X}_M , $i \in \{1, ..., N\}$, particularly, its variance given by

$$Var(\hat{X}_{M,i}) = \frac{1}{M^2} \sum_{j=1}^{M} \frac{C^2 \sigma^2}{\beta_C(X)^2 \frac{\partial g(f(X))}{\partial f(X)_i}^2}.$$
 (29)

Without loss of generality, we assume that $\frac{\partial g(f(X))}{\partial f(X)_j} > 0$ for all $j \in \{1, ..., M\}$.

It is easy to see that 29 decreases for decreasing $\sum_{j=1}^{M} \frac{1}{\beta_{C}(X)^{2} \frac{\partial g(f(X))^{2}}{\partial f(X)_{s}}}$. Therefore, minimising

$$\sum_{j=1}^{M} \frac{1}{\beta_C(X)^2 \frac{\partial g(f(X))}{\partial f(X)_j}^2} \tag{30}$$

with respect to $\frac{\partial g(f(X))}{\partial f(X)_1}$, ..., $\frac{\partial g(f(X))}{\partial f(X)_M}$ minimises the variance in 29. Note that doing so does not affect the multiplicative term $C^2\sigma^2/M$. By definition of the global concatenated gradient G_X (see 6), the squared norm of G_X is given by

$$||G_X||_2^2 = \sum_{j=1}^M \frac{\partial g(f(X))}{\partial f(X)_j}^2 ||X||_2^2 + \sum_{j=1}^M \frac{\partial g(f(X))}{\partial b_j}^2 + ||G_{X,P}||_2^2.$$
(31)

Set $\|\text{Rest}\|_2^2 := \sum_{j=1}^M \frac{\partial g(f(X))}{\partial b_j}^2 + \|G_{X,P}\|_2^2$. Without loss of generality, we assume the target $X \neq \mathbf{0}_N$, for $\mathbf{0}_N$ being the N dimensional vector with zeros everywhere. Then, we can reformulate 31 and obtain the following constraint regarding $\frac{\partial g(f(X))}{\partial f(X)_1}, ..., \frac{\partial g(f(X))}{\partial f(X)_M}$:

$$\sum_{j=1}^{M} \frac{\partial g(f(X))}{\partial f(X)_{j}}^{2} = \frac{\|G_{X}\|_{2}^{2} - \|\text{Rest}\|_{2}^{2}}{\|X\|_{2}^{2}}.$$
(32)

Minimising the variance in 29 with respect to $\frac{\partial g(f(X))}{\partial f(X)_1},...,\frac{\partial g(f(X))}{\partial f(X)_M}$ under the constraint given in 32, does not affect the norm of the global gradient G_X and, thus, it does not affect the value of $\beta_C(X)$. Therefore, minimising 30 with respect to $\frac{\partial g(f(X))}{\partial f(X)_1},...,\frac{\partial g(f(X))}{\partial f(X)_M}$ under the constraint given in 32 is equivalent to minimising $\sum_{j=1}^{M} \frac{1}{\frac{\partial g(f(X))}{\partial f(X)_j}}$ with respect to $\frac{\partial g(f(X))}{\partial f(X)_1},...,\frac{\partial g(f(X))}{\partial f(X)_M}$ under 32. Hence, setting $y_j = \frac{\partial g(f(X))}{\partial f(X)_j}^2$, for $j \in \{1,...,M\}$, we have an optimisation problem of the following form:

Minimise
$$\sum_{j=1}^{M} \frac{1}{y_j}$$
 for $\sum_{j=1}^{M} y_j = \frac{\|G_X\|_2^2 - \|\text{Rest}\|_2^2}{\|X\|_2^2}$ and $y_1, ..., y_M > 0$, (33)

33 is a well-known minimisation problem with solution given by $y_j = \frac{\|G_X\|_2^2 - \|\text{Rest}\|_2^2}{M\|X\|_2^2}$ for all $j \in \{1, ..., M\}$. However, if needed, a proof of the statement can be obtained using the gradient of the function in 33 to

construct the direction of the steepest descent and combining this with the given constraints in 33. Hence, setting

$$\frac{\partial g(f(X))}{\partial f(X)_j}^2 = \frac{\|G_X\|_2^2 - \|\text{Rest}\|_2^2}{M\|X\|_2^2}$$
(34)

for all $j \in \{1, ..., M\}$, minimises the variance given in 29 with respect to $\frac{\partial g(f(X))}{\partial f(X)_1}, ..., \frac{\partial g(f(X))}{\partial f(X)_M}$.

We insert the choice 34 into the variance 35 and obtain

$$\operatorname{Var}(\hat{X}_{M,i}) \ge \frac{1}{M^2} \sum_{i=1}^{M} \frac{C^2 \sigma^2 M \|X\|_2^2}{\beta_C(X)^2 (\|G_X\|_2^2 - \|\operatorname{Rest}\|_2^2))} = \frac{C^2 \sigma^2 \|X\|_2^2}{\beta_C(X)^2 (\|G_X\|_2^2 - \|\operatorname{Rest}\|_2^2))}, \tag{35}$$

for $i \in \{1, ..., N\}$.

Recall the definition of the clipping term $\beta_C(X)$ given in 7. Using 34, we can see that the norm of the global gradient $\|G_X\|_2$ is linearly increasing in M. Thus, there exist \hat{M} , such that for all $M \ge \hat{M}$, $\|G_X\|_2 \ge C$ and $\beta_C(X) = \frac{C}{\|G_X\|_2}$. Hence, by 35

$$\lim_{M \to \infty} \operatorname{Var}(\hat{X}_{M,i}) \ge C^2 \sigma^2 \|X\|_2^2 \cdot \lim_{M \to \infty} \frac{1}{\beta_C(X)^2 (\|G_X\|_2^2 - \|\operatorname{Rest}\|_2^2))}$$

$$= C^2 \sigma^2 \|X\|_2^2 \cdot \lim_{M \to \infty} \frac{1}{C^2 - \frac{\|\operatorname{Rest}\|_2^2}{\|G_X\|_2^2}}$$

$$= \sigma^2 \|X\|_2^2. \tag{36}$$

Equality 36 holds because $\|\text{Rest}\|_2$ is independent of M. Lastly, if $X \neq \mathbf{0}_N$, then $\sigma^2 \|X\|_2^2 > 0$.

Proposition 1. Let the part of neural network given by g be replaced by the loss function $\mathcal{L}: \mathbb{R}^M \to \mathbb{R}$ with $\mathcal{L}(f(X)) = \mathbf{1}_M^T f(X)$, where $\mathbf{1}_M$ is the M-dimensional 1-vector, and

$$M \ge \max\left(1, \left\lceil \frac{C}{\min_{X \in \mathcal{D} \setminus \{\boldsymbol{\varrho}_N\}} \|X\|_2} \right\rceil\right),\tag{15}$$

where $\lceil \cdot \rceil$ denotes the function that rounds up its argument to the nearest integer, then $\frac{1}{M^2} \sum_{j=1}^{M} \frac{C^2 \sigma^2}{\beta_C(X)^2 \frac{\partial g(f(X))}{\partial f(X)_j}^2}$ is minimal and takes the value $\sigma^2 \|X\|_2^2$.

Proof. Let the iteration step be fixed and observe the sample mean \hat{X}_M given in 13 with distribution described in 14. Then, consider the *i*th entry of \hat{X}_M , particularly, its variance given by

$$Var(\hat{X}_{M,i}) = \frac{1}{M^2} \sum_{j=1}^{M} \frac{C^2 \sigma^2}{\beta_C(X)^2 \frac{\partial g(f(X))}{\partial f(X)_i}^2},$$
(37)

for $i \in \{1, ..., N\}$. Without loss of generality, let $X \neq \mathbf{0}_N$. We have shown in the proof of Proposition B.1 that the choice

$$\frac{\partial g(f(X))^{2}}{\partial f(X)_{j}} = \frac{\|G_{X}\|_{2}^{2} - \|\text{Rest}\|_{2}^{2}}{M\|X\|_{2}^{2}}$$
(38)

for all $j \in \{1,...,M\}$, minimises the variance given in 37 with respect to $\frac{\partial g(f(X))}{\partial f(X)_1},...,\frac{\partial g(f(X))}{\partial f(X)_M}$. Let us set

$$\frac{\partial g(f(X))}{\partial f(X)_1}^2 = \frac{\|G_X\|_2^2 - \|\text{Rest}\|_2^2}{M\|X\|_2^2},\tag{39}$$

and insert this choice into the variance 37:

$$\operatorname{Var}(\hat{X}_{M,i}) \ge \frac{C^2 \sigma^2}{M} \frac{1}{\beta_C(X)^2 \frac{\partial g(f(X))^2}{\partial f(X)_1}},\tag{40}$$

for $i \in \{1, ..., N\}$. If $\|G_X\|_2$ is fixed, it follows from 38 that $\frac{\partial g(f(X))}{\partial f(X)_1}^2$ increases with decreasing norm $\|\text{Rest}\|_2$. However, $\|\text{Rest}\|_2$ cannot be bounded or quantified for any iteration step without specific knowledge of the neural network. Thus, the adversary cannot minimise $\|\text{Rest}\|_2$ without manipulating some layers of the network. If they manipulate these layers, we see 40 is minimal whenever $\frac{\partial g(f(X))}{\partial f(X)_1}$ is maximal, i.e., whenever $\|\text{Rest}\|_2 = 0$. $\|\text{Rest}\|_2 = 0$ occurs for all $X \in \mathcal{D}$ and all iteration steps when the adversary replaces the entire network by the linear layer f (see Section 2.1, specifically Equation 1) and sets the bias term b to be equal to $\mathbf{0}_M$. In such a case the neural network is given by the linear layer f(X) = WX and a loss function which we denote by $\mathcal{L}: \mathbb{R}^M \to \mathbb{R}$. As a consequence, $\frac{\partial g(f(X))}{\partial f(X)_1} = \frac{\partial \mathcal{L}(f(X))}{\partial f(X)_1}$. In particular, 38 implies

$$\frac{\partial g(f(X))}{\partial f(X)_1}^2 = \frac{\partial \mathcal{L}(f(X))}{\partial f(X)_1}^2 = \frac{\|G_X\|_2^2}{M\|X\|_2^2}.$$
 (41)

Inserting 41 into the right hand side of 40 further bounds the variance $Var(\hat{X}_{M,i})$:

$$\operatorname{Var}(\hat{X}_{M,i}) \ge \frac{C^2 \sigma^2 \|X\|_2^2}{\beta_C(X)^2 \|G_X\|_2^2},\tag{42}$$

for all $i \in \{1, ..., N\}$.

Now, we observe the lower bound in 42. Naturally, the right hand side of 42 is lowest when the denominator in 42 is highest. By definition of the clipping term $\beta_C(X)$ (see 7), the product $\beta_C(X)^2 ||G_X||_2^2$ is upper bounded by C^2 , delivering

$$\operatorname{Var}(\hat{X}_{M,i}) \ge \sigma^2 \|X\|_2^2,\tag{43}$$

for $i \in \{1, ..., N\}$. In particular, no change in the parameters or architecture of the network can increase the product $\beta_C(X)^2 \|G_X\|_2^2$ beyond C^2 to further decrease the lower bound given in 43. Therefore, we assume, the adversary chooses M and $\frac{\partial \mathcal{L}(f(X))}{\partial f(X)_1}$ such that that $\|G_X\|_2^2 \geq C^2$ for as many data points X as possible. Using 41, $\|G_X\|_2^2 \geq C^2$ implies

$$M\frac{\partial \mathcal{L}(f(X))}{\partial f(X)_1}^2 \ge \frac{C^2}{\|X\|_2^2}.$$
(44)

Next, we consider two cases, when $\min_{X\in\mathcal{D}} \|X\|_2 > 0$ and $\min_{X\in\mathcal{D}} \|X\|_2 = 0$. If $\min_{X\in\mathcal{D}} \|X\|_2 > 0$, then

$$\|G_X\|_2^2 \ge C^2 \quad \forall X \in \mathcal{D} \iff M \frac{\partial \mathcal{L}(f(X))}{\partial f(X)_1}^2 \ge \frac{C^2}{\min_{X \in \mathcal{D}} \|X\|_2^2} \quad \forall X \in \mathcal{D}. \tag{45}$$

M, $\min_{X\in\mathcal{D}}\|X\|_2$ and C fixed during training and do not changed from iteration to iteration. Thus, 45 holds for all $X\in\mathcal{D}$ if $\frac{\partial \mathcal{L}(f(X))}{\partial f(X)_1}$ is constant for all X and all iteration steps, implying \mathcal{L} is an affine function of f(X). If $C>\min_{X\in\mathcal{D}}\|X\|_2$, then choosing $M\geq \left\lceil \frac{C}{\min_{X\in\mathcal{D}}\|X\|_2}\right\rceil$ and the loss function $\mathcal{L}:\mathbb{R}^M\to\mathbb{R}$ to be $\mathcal{L}(f(X))=\mathbf{1}_M^Tf(X)$, where $\mathbf{1}_M$ is the M-dimensional 1-vector, delivers the sample average \hat{X}_M with the lowest variance per entry given by

$$Var(\hat{X}_{M,i}) = \sigma^2 ||X||_2^2, \tag{46}$$

where $\lceil \cdot \rceil$ denotes the function that rounds up its argument to the nearest integer. If $C \leq \min_{X \in \mathcal{D}} \|X\|_2$, then choosing $M \geq 1$ and the loss function to be $\mathcal{L}(f(X)) = f(X)$ delivers the sample average \hat{X}_M with the lowest variance per entry given by

$$Var(\hat{X}_{M,i}) = \sigma^2 ||X||_2^2 \ge \sigma^2 C^2.$$
(47)

All in all, we conclude that if $\min_{X \in \mathcal{D}} \|X\|_2 > 0$, replacing the subpart of the neural network given by g by the loss function $\mathcal{L}(f(X)) = \mathbf{1}_M^T f(X)$ and setting $M \ge \max\left(1, \left\lceil \frac{C}{\min_{X \in \mathcal{D}} \|X\|_2} \right\rceil\right)$ minimises the variance $\operatorname{Var}(\hat{X}_{M,i})$ for all $i \in \{1, ..., N\}$.

If $\min_{X\in\mathcal{D}}\|X\|_2=0$, then there is no choice for $\frac{\partial\mathcal{L}(f((X))}{\partial f(X)_1}$ or M such that 44 holds for all $X\in\mathcal{D}$. However, in such a case, without loss of generality, we assume that the adversary sets the loss function to be $\mathcal{L}(f(X))=f(X)$ and chooses M to ensure that 44 holds for all $X\in\mathcal{D}$ with $X\neq\mathbf{0}_N$. In such a case, the adversary sets $M=\left\lceil\frac{C}{\min_{X\in\mathcal{D}\setminus\{\mathbf{0}_N\}}\|X\|_2}\right\rceil$, analogously as the argumentation above.

The following lemma serves as an auxiliary result to obtain Proposition 2:

Lemma B.1. For all $j \in \{1,...,N\}$, the jth entry \hat{x}_j of the estimator \hat{X} is a (fully) efficient estimator for the jth entry x_j of the target X.

Proof. The estimator \hat{X} given in 17 is normally distributed with mean X and covariance matrix given by $\sigma^2 \|X\|_2^2 I_N$. Let \hat{x}_i , $i \in \{1, ..., N\}$, denote the *i*th entry of \hat{X} . Then, by distribution of \hat{X} , \hat{x}_1 , ..., \hat{x}_N are independent, normally distributed with mean $x_1, ..., x_N$, respectively, and same variance given by $\sigma^2 \|X\|_2^2$. Thus, for all $i \in \{1, ..., N\}$, \hat{x}_i is an unbiased estimator of the *i*th entry of the target X.

Moreover, applying the Cramér-Rao bound for scalar unbiased estimators, we compute a lower bound for the variance of any the estimator of \hat{x}_i , $i \in \{1, ..., N\}$:

$$\operatorname{Var}_{X}(\hat{x}_{i}) \ge I(x_{i})^{-1} = \sigma^{2} \|X\|_{2}^{2},\tag{48}$$

where $I(x_i)$ denotes the Fisher information matrix that measures the amount of information the rescaled, observable normally distributed random variables \hat{x}_i carries about its unknown mean x_i . Since this matrix is well-known in literature, we do not provide a proof for the right hand side of the equality in 48.

Since for all $i \in \{1, ..., N\}$, \hat{x}_i is an unbiased estimator of x_i that achieves the Cramér-Rao bound, it is a (fully) efficient estimator of x_i achieving the smallest variability in terms of the variance.

Proposition 2. \hat{X} is the minimum variance unbiased estimator (MVUE) for X. Moreover, the expected mean squared error between the target X and \hat{X} is given by:

$$\mathbb{E}_X[\mathrm{MSE}_X(X,\hat{X})] = \sigma^2 \|X\|_2^2. \tag{18}$$

Proof. Let $\hat{Y} = (\hat{y}_1, ..., \hat{y}_N)^T$ denote any estimator of X. Then,

$$\mathbb{E}_{X}[\text{MSE}_{X}(X,\hat{Y})] = \mathbb{E}_{X} \left[\frac{||X - \hat{Y}||_{2}^{2}}{N} \right]$$

$$= \frac{1}{N} \sum_{i=1}^{N} \mathbb{E}_{X} \left[(x_{i} - \hat{y}_{i})^{2} \right]$$

$$= \frac{1}{N} \sum_{i=1}^{N} \mathbb{E}_{x_{i}} \left[(x_{i} - \hat{y}_{i})^{2} \right]$$

$$= \frac{1}{N} \sum_{i=1}^{N} \left(\mathbb{E}_{x_{i}} \left[x_{i} - \hat{y}_{i} \right]^{2} + \text{Var}_{x_{i}} (x_{i} - \hat{y}_{i}) \right)$$

$$= \frac{1}{N} \sum_{i=1}^{N} \left(\text{Bias}_{x_{i}} (x_{i}, \hat{y}_{i})^{2} + \text{Var}_{x_{i}} (\hat{y}_{i}) \right). \tag{49}$$

For all unbiased estimators, the expected MSE, as given in 49, is solely determined by the sum of the variances of each entry \hat{y}_i . Therefore, by Lemma B.1, \hat{X} is the unbiased estimator that minimises 49. In other words, \hat{X} is the unbiased estimator that achieves the lowest expected MSE. Such estimators are called *minimum* variance unbiased estimators in the literature and achieve the smallest variability in terms of the variance.

Lastly, we compute the expected MSE between X and \hat{X} :

$$\mathbb{E}_{X}[\text{MSE}_{X}(X, \hat{X})] = \frac{1}{N} \sum_{i=1}^{N} \text{Var}_{x_{i}}(\hat{x}_{i}) = \sigma^{2} ||X||_{2}^{2}.$$

Theorem 1. Using the MSE as an optimality criterion, \hat{X} is the best achievable estimator and, thus, reconstruction for the target point X.

Proof. Since the adversary only observes one privatised version of the gradient $\tilde{\nabla}_W$ they can use to construct an estimator for X, it is easy to see that \hat{X} is a sufficient statistic for estimating X. Moreover, by Proposition 2, \hat{X} is the unbiased estimator, which uses the sufficient statistic $\hat{\nabla}_W$ as input, that achieves the lowest expected $\text{MSE}_X(X,\hat{X})$. Such unbiased estimators achieve the lowest possible MSE and have the smallest variability in terms of their variance. Thus, using the MSE as an optimality criterion, \hat{X} is the optimal estimator for X. Since lower values of $\text{MSE}_X(X,\hat{X})$ denote high similarity between X and \hat{X} , we conclude that \hat{X} is the best achievable reconstruction for X.

Theorem 2.

$$MSE_X(X, \hat{X}) \stackrel{d}{=} \frac{\sigma^2 ||X||_2^2}{N} \cdot Y \quad with \quad Y \sim \chi_N^2, \tag{19}$$

where χ^2_N denotes the central chi-squared distribution with N degrees of freedom. In particular, for η given,

$$\mathbb{P}_{\hat{X}}(\mathrm{MSE}_X(X,\hat{X}) \le \eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 \|X\|_2^2}\right),\tag{20}$$

where Γ_R is the regularised gamma function.

Proof. We can compute the MSE between X and its reconstruction \hat{X} as the mean error over their components:

$$MSE_{X}(X, \hat{X}) = \frac{1}{N} \sum_{i=1}^{N} (x_{i} - \hat{x}_{i})^{2}$$

$$= \frac{1}{N} \sum_{i=1}^{N} \tilde{\xi}_{i}^{2}, \qquad \tilde{\xi}_{i} \sim \mathcal{N}(0, \sigma^{2} || X ||_{2}^{2})$$

$$= \frac{1}{N} \sum_{i=1}^{N} \sigma^{2} || X ||_{2}^{2} \left(\frac{1}{\sigma || X ||_{2}} \tilde{\xi}_{i} \right)^{2}$$

$$= \frac{\sigma^{2} || X ||_{2}^{2}}{N} \sum_{i=1}^{N} \rho_{i}^{2}, \qquad (50)$$

where $\rho_i := \frac{1}{\sigma ||X||_2} \xi_i$, for $i \in \{1, ..., N\}$. If X is fixed, $\rho_1, ..., \rho_N$ are pairwise independent random variables with $\rho_i \sim \mathcal{N}(0, 1)$ for all $i \in \{1, ..., N\}$. Hence,

$$\sum_{i=1}^{N} \rho_i^2 \sim \chi_N^2,\tag{51}$$

where χ_N^2 denotes the central chi-squared distribution with N degrees of freedom. Thus,

$$MSE_X(X, \hat{X}) \stackrel{d}{=} \frac{\sigma^2 ||X||_2^2}{N} \cdot Y$$
 with $Y \sim \chi_N^2$.

Then, for η given

$$\mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_X(X,\hat{X}) \leq \eta\right) = \gamma \iff \mathbb{P}_Y\left(\frac{\sigma^2 \|X\|_2^2}{N} \cdot Y \leq \eta\right) = \gamma.$$

Since Y is a centered chi-squared distributed random variable with N degrees of freedom, its cumulative distribution function can be computed via the regularized gamma function Γ_R . Hence,

$$\mathbb{P}_{Y}\left(\frac{\sigma^{2}\|X\|_{2}^{2}}{N} \cdot Y \leq \eta\right) = P_{Y}\left(Y \leq \frac{N\eta}{\sigma^{2}\|X\|_{2}^{2}}\right) = \Gamma_{R}\left(\frac{N}{2}, \frac{N\eta}{2\sigma^{2}\|X\|_{2}^{2}}\right)$$

implies

$$\mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_X(X,\hat{X}) \le \eta\right) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 \|X\|_2^2}\right). \tag{52}$$

Proposition B.3. For $N \to \infty$, it holds

$$\sqrt{N}\left(\mathrm{MSE}_X(X,\hat{X}) - \sigma^2 \|X\|_2^2\right) \stackrel{d}{\to} \mathcal{N}\left(0, 2\sigma^4 \|X\|_2^4\right). \tag{53}$$

Proof. By Theorem 2, it holds

$$MSE_X(X, \hat{X}) = \sigma^2 ||X||_2^2 \frac{1}{N} \sum_{i=1}^N \rho_i^2 = \frac{1}{N} \sum_{i=1}^N \sigma^2 ||X||_2^2 \rho_i^2,$$
 (54)

where ρ_i are i.i.d. standard normally distributed random variables, for all $i \in \{1, ..., N\}$. Consider the behaviour of $\sigma^2 ||X||_2^2 \rho_i^2$ for $i \in \{1, ..., N\}$:

$$\mathbb{E}\left[\sigma^{2} \|X\|_{2}^{2} \rho_{i}^{2}\right] = \sigma^{2} \|X\|_{2}^{2} \mathbb{E}\left[\rho_{i}^{2}\right] = \sigma^{2} \|X\|_{2}^{2} \operatorname{Var}(\rho_{i}) = \sigma^{2} \|X\|_{2}^{2}$$

and

$$\mathrm{Var}\left(\sigma^{2}\|X\|_{2}^{2}\rho_{i}^{2}\right) = \sigma^{4}\|X\|_{2}^{4}\mathrm{Var}\left(\rho_{i}^{2}\right) = \sigma^{4}\|X\|_{2}^{4}\left(\mathbb{E}\left[\rho_{i}^{4}\right] - \mathbb{E}\left[\rho_{i}^{2}\right]^{2}\right) = \sigma^{4}\|X\|_{2}^{4}(3-1) = 2\sigma^{4}\|X\|_{2}^{4}$$

since $\mathbb{E}\left[\rho_i^4\right]$ corresponds to the 4-th moment of the standard normal distribution and can be, thus, computed easily. Therefore, using the central limit theorem, for $N \to \infty$, it follows

$$\sqrt{N} \left(\mathrm{MSE}_X(X, \hat{X}) - \sigma^2 \|X\|_2^2 \right) = \sqrt{N} \left(\frac{1}{N} \sum_{i=1}^N \sigma^2 \|X\|_2^2 \rho_i^2 - \sigma^2 \|X\|_2^2 \right) \overset{d}{\to} \mathcal{N} \left(0, 2\sigma^4 \|X\|_2^4 \right).$$

Proposition 3. Let η given. Then, for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$,

$$\mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_{X}\left(X,\hat{X}\right) \leq \eta\right) \leq \Gamma_{R}\left(\frac{N}{2}, \frac{N\eta}{2\sigma^{2} \min_{X \in \mathcal{D}} \|X\|_{2}^{2}}\right),\tag{21}$$

where Γ_R is the regularised gamma function. Moreover, the DP-mechanism \mathcal{M} is $(\eta, \gamma(\eta))$ -reconstruction robust with respect to the MSE for any reconstruction and $\gamma(\eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 \min_{X \in \mathcal{D}} \|X\|_2^2}\right)$. If $C \leq \min_{X \in \mathcal{X}} \|X\|_2$ holds, then \mathcal{M} is $(\eta, \gamma'(\eta))$ -reconstruction robust with respect to the MSE for any reconstruction and $\gamma'(\eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2C^2}\right)$.

Proof. Consider the CDF of the $\mathrm{MSE}_X(X,\hat{X})$ for $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$ given Equation 20 in Theorem 2. Since the regularised gamma function is increasing in its second argument, using Theorem 2, it follows that

$$\mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_{X}(X,\hat{X}) \leq \eta\right) = \Gamma_{R}\left(\frac{N}{2}, \frac{N\eta}{2\sigma^{2}\|X\|_{2}^{2}}\right) \leq \Gamma_{R}\left(\frac{N}{2}, \frac{N\eta}{2\sigma^{2}\min_{X \in \mathcal{X}} \|X\|_{2}^{2}}\right),\tag{55}$$

for all reconstruction target points $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$. Moreover, due to the optimality of the estimator \hat{X} , it holds

$$\mathbb{P}_{Y_X'}\left(\mathrm{MSE}_X(X, Y_X') \le \eta\right) \le \mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_X(X, \hat{X}) \le \eta\right) \tag{56}$$

for all possible reconstructions Y_X' . Therefore, combining 55 and 56, we conclude that \mathcal{M} is $(\eta, \gamma(\eta))$ -reconstruction robust with respect to the MSE for any reconstruction and $\gamma(\eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 \min_{X \in \mathcal{X}} \|X\|_2^2}\right)$.

Moreover, if $C \leq \min_{X \in \mathcal{X}} ||X||_2$, then

$$\mathbb{P}_{\hat{X}}\left(\mathrm{MSE}_X(X,\hat{X}) \le \eta\right) \le \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2 C^2}\right),\tag{57}$$

for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$. In particular, in such a case, \mathcal{M} is $(\eta, \gamma'(\eta))$ -reconstruction robust with respect to the MSE for any reconstruction and $\gamma'(\eta) = \Gamma_R\left(\frac{N}{2}, \frac{N\eta}{2\sigma^2C^2}\right)$.

Proposition 4. Assume $\max_{X \in \mathcal{D}} \max(X)$ and $\min_{X \in \mathcal{D}} \min(X)$ are known quantities. Then, for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$,

$$\mathbb{P}_{\hat{X}}\left(\mathrm{PSNR}_{X}\left(X,\hat{X}\right) \geq \eta\right) \leq \Gamma_{R}\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^{2} \min_{X \in \mathcal{D} \setminus \{\boldsymbol{\theta}_{N}\}} \|X\|_{2}^{2}}\right),\tag{22}$$

for $\tilde{\eta}(\eta) := 10^{-\frac{\eta}{10}} \left(\max_{X \in \mathcal{D}} \max(X) - \min_{X \in \mathcal{D}} \min(X) \right)^2$, and for Γ_R being the regularised gamma function. If $C \leq \min_{X \in \mathcal{D}} \|X\|_2$ holds, then

$$\mathbb{P}_{\hat{X}}\left(\text{PSNR}_X\left(X,\hat{X}\right) \ge \eta\right) \le \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 C^2}\right),\tag{23}$$

for all $X \in \mathcal{D} \setminus \{\boldsymbol{o}_N\}$.

Proof. The cumulative distribution function (CDF) of the PSNR can be calculated using the CDF of the MSE. In particular, this implies that we can also compute probabilistic bounds for the PSNR using Theorem 2. Let η be given. Then,

$$\begin{aligned} \operatorname{PSNR}_X(X,\hat{X}) &\geq \eta \\ \iff & 10 \log_{10} \left(\frac{(\max(X) - \min(X))^2}{\operatorname{MSE}(X,\hat{X})} \right) \geq \eta \\ \iff & \log_{10} ((\max(X) - \min(X))^2) - \log_{10} (\operatorname{MSE}(X,\hat{X}) \geq \frac{\eta}{10} \\ \iff & 2 \log_{10} [(\max(X) - \min(X))] - \frac{\eta}{10} \geq \log_{10} \left(\operatorname{MSE}(X,\hat{X}) \right) \\ \iff & (\max(X) - \min(X))^2 10^{-\frac{\eta}{10}} \geq \operatorname{MSE}(X,\hat{X}). \end{aligned}$$

Thus, setting

$$\hat{\eta}(\eta) = 10^{-\frac{\eta}{10}} (\max(X) - \min(X))^2, \tag{58}$$

it follows from Theorem 2 that

$$\mathbb{P}_{\hat{X}}(\mathrm{PSNR}_{X}(X,\hat{X}) \geq \eta)$$

$$= \mathbb{P}_{\hat{X}}(\mathrm{MSE}_{X}(X,\hat{X}) \leq \hat{\eta}(\eta)) \leq \Gamma_{R}\left(\frac{N}{2}, \frac{N\hat{\eta}(\eta)}{2} \frac{1}{\sigma^{2} \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_{N}\}} \|X\|_{2}^{2}}\right). \tag{59}$$

We note that the right hand side of the previous result is still dependent on the target value X due to $\hat{\eta}(\eta)$. To remove this dependency, we find an upper bound for $\hat{\eta}(\eta)$:

$$\hat{\eta}(\eta) = 10^{-\frac{\eta}{10}} (\max(X) - \min(X))^2 \le 10^{-\frac{\eta}{10}} \left(\max_{X \in \mathcal{D}} \max(X) - \min_{X \in \mathcal{D}} \min(X) \right)^2 := \tilde{\eta}(\eta).$$

Since the regularised gamma function Γ_R is increasing with respect to the second argument, it follows:

$$\mathbb{P}_{\hat{X}}(\mathrm{PSNR}_X(X,\hat{X}) \ge \eta) \le \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2^2}\right),\tag{60}$$

for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$. Using the same argument, if $C \leq \min_{X \in \mathcal{D} \setminus \{\mathbf{0}_N\}} \|X\|_2$, then

$$\mathbb{P}_{\hat{X}}(\mathrm{PSNR}_X(X,\hat{X}) \ge \eta) \le \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 C^2}\right). \tag{61}$$

Proposition 5. Assume $\max_{X \in \mathcal{D}} \max(X)$ and $\min_{X \in \mathcal{D}} \min(X)$ are known quantities, and let Y_X be any possible reconstruction. Then, for all $X \in \mathcal{D} \setminus \{\mathbf{0}_N\}$,

$$\mathbb{P}_{Y_X}\left(\text{PSNR}_X\left(X, Y_X\right) \ge \eta\right) \le \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 \min_{X \in \mathcal{D} \setminus \{\boldsymbol{\varrho}_N\}} \|X\|_2^2}\right),\tag{24}$$

for Γ_R being the regularised gamma function and $\tilde{\eta}(\eta)$ as defined in Proposition 4. In particular, the DP-mechanism \mathcal{M} is $(-\eta, \tilde{\gamma}(\tilde{\eta}(\eta)))$ -reconstruction robust with respect to the negative PSNR (-PSNR) for any analytic reconstruction and $\tilde{\gamma}(\tilde{\eta}(\eta)) = \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2\min_{X\in\mathcal{D}\setminus\{\theta_N\}}\|X\|_2^2}\right)$. Moreover, if $C \leq \min_{X\in\mathcal{D}}\|X\|_2$ holds, then the DP-mechanism \mathcal{M} is $(-\eta, \tilde{\gamma}'(\tilde{\eta}(\eta)))$ -reconstruction robust with respect to -PSNR for any reconstruction, $\tilde{\gamma}'(\tilde{\eta}(\eta)) = \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2C^2}\right)$.

Proof. On the one hand, by the optimality of the reconstruction \hat{X} with respect to the PSNR, it holds that

$$\mathbb{P}_{Y_X}(\mathrm{PSNR}_X(X, Y_X') \ge \eta) \le \mathbb{P}_{\hat{X}}(\mathrm{PSNR}_X(X, \hat{X}) \ge \eta),$$

for any analytic reconstruction Y'_X . On the other hand, by Proposition 4,

$$\mathbb{P}_{\hat{X}}(\mathrm{PSNR}_X(X,\hat{X}) \geq \eta) = \mathbb{P}_{\hat{X}}(-\mathrm{PSNR}_X(X,\hat{X}) \leq -\eta) \leq \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2 \min_{X \in \mathcal{D}\setminus \{\mathbf{0}_N\}} \|X\|_2^2}\right),$$

for $\tilde{\eta}(\eta)$ as defined in Proposition 4 and Γ_R being the regularised gamma function. Therefore, using Definition 1, we conclude that the DP mechanism \mathcal{M} is $(-\eta, \tilde{\gamma}(\tilde{\eta}(\eta)))$ -reconstruction robust for $\tilde{\gamma}(\tilde{\eta}(\eta)) = \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2\min_{X \in \mathcal{D}\setminus \{\mathbf{0}_N\}} \|X\|_2^2}\right)$ with respect to the negative PSNR, i.e., -PSNR.

Lastly, since the regularised gamma function Γ_R is increasing with respect to the second argument, if $C \leq \min_{X \in \mathcal{D}} \|X\|_2$, then the DP mechanism \mathcal{M} is $(-\eta, \tilde{\gamma}'(\tilde{\eta}(\eta)))$ -reconstruction robust for $\tilde{\gamma}'(\tilde{\eta}(\eta)) = \Gamma_R\left(\frac{N}{2}, \frac{N\tilde{\eta}(\eta)}{2\sigma^2C^2}\right)$ with respect to the negative PSNR, i.e., -PSNR.

Proposition B.4. Let σ_X and $\sigma_{\hat{X}}$ denote the sample standard deviations of $\{x_1, ..., x_N\}$ and $\{\hat{x}_1, ..., \hat{x}_N\}$, respectively. Then, the sample NCC between X and \hat{X} is given by

$$NCC(X, \beta_C(X)^{-1}\hat{X}) = \frac{\sigma_X}{\sigma_{\hat{X}}} + \frac{1}{\sigma_X \sigma_{\hat{X}}} \left(\frac{1}{N} \sum_{i=1}^N x_i \zeta_i - \bar{x}\bar{\zeta} \right), \tag{62}$$

for $\zeta \sim \mathcal{N}(\boldsymbol{\theta}_N, \sigma^2 ||X||_2^2 I_N)$.

Proof. We recall that the reconstruction \hat{X} of X is obtained by the adversary using the observed privatised gradients given in 8. Due to the from X independent random noise added to these gradients, they can be viewed as samples from random variables (see 10). Consequently, \hat{X} can also be viewed as a sample from a random variable (see 17):

$$\hat{X} \stackrel{d}{=} \hat{Y}_X$$
, for $\hat{Y}_X \sim \mathcal{N}\left(X, \sigma^2 ||X||_2^2 I_N\right)$.

For the target point X being fixed and, thus, $||X||_2$ being also fixed, we let $\zeta \sim \mathcal{N}\left(\mathbf{0}_N, \sigma^2 ||X||_2^2 I_N\right)$ be a random variable drawn independently of X. It is easy to see that

$$\hat{X} \stackrel{d}{=} X + \zeta.$$

Let

$$\bar{x} := \frac{1}{N} \sum_{i=1}^{N} x_i, \quad \text{and} \quad \bar{\hat{x}} := \frac{1}{N} \sum_{i=1}^{N} \hat{x}_i = \frac{1}{N} \sum_{i=1}^{N} (x_i + \zeta_i) = \bar{x} + \bar{\zeta}, \quad \text{for} \quad \bar{\zeta} := \frac{1}{N} \sum_{i=1}^{N} \zeta_i, \quad (63)$$

denote the sample means of $\{x_1,...,x_N\}$ and $\{\hat{x}_1,...,\hat{x}_N\}$, respectively. Moreover, let

$$\sigma_X = \sqrt{\frac{1}{N} \sum_{i=1}^{N} (x_i - \bar{x})^2}, \quad \text{and} \quad \sigma_{\hat{X}} = \sqrt{\frac{1}{N} \sum_{i=1}^{N} (\hat{x}_i - \bar{\hat{x}})^2},$$
 (64)

denote the sample standard deviations of $\{x_1, ..., x_N\}$ and $\{\hat{x}_1, ..., \hat{x}_N\}$, respectively. Then, by definition of the sample NCC (see Definition 4), it follows

$$NCC(X, \hat{X}) = \frac{1}{N} \sum_{i=1}^{N} \frac{(x_i - \bar{x})}{\sigma_X} \frac{(\hat{x}_i - \bar{x})}{\sigma_{\hat{X}}}$$

$$= \frac{1}{N} \sum_{i=1}^{N} \frac{(x_i - \bar{x})}{\sigma_X} \frac{(x_i - \bar{x} + \zeta_i - \bar{\zeta})}{\sigma_{\hat{X}}}$$

$$= \frac{1}{N} \sum_{i=1}^{N} \frac{(x_i - \bar{x})^2 + (x_i - \bar{x})(\zeta_i - \bar{\zeta})}{\sigma_X \sigma_{\hat{X}}}$$

$$= \frac{\sigma_X}{\sigma_{\hat{X}}} + \frac{1}{\sigma_X \sigma_{\hat{X}}} \left(\frac{1}{N} \sum_{i=1}^{N} (x_i - \bar{x})(\zeta_i - \bar{\zeta}) \right). \tag{65}$$

The multiplicative term $\frac{1}{N}\sum_{i=1}^{N}(x_i-\bar{x})(\zeta_i-\bar{\zeta})$ of 65 corresponds to the sample covariance between the entries of target X and the random noise, namely between the sets $\{x_1,...,x_N\}$ and $\{\zeta_1,...,\zeta_N\}$. Next, we simplify the sample covariance:

$$\frac{1}{N} \sum_{i=1}^{N} (x_i - \bar{x})(\zeta_i - \bar{\zeta}) = \frac{1}{N} \sum_{i=1}^{N} (x_i \zeta_i - \bar{x} \zeta_i - x_i \bar{\zeta} + \bar{x} \bar{\zeta})$$

$$= \frac{1}{N} \sum_{i=1}^{N} x_i \zeta_i - \bar{x} \frac{1}{N} \sum_{i=1}^{N} \xi_i - \bar{\zeta} \frac{1}{N} \sum_{i=1}^{N} x_i + \bar{x} \bar{\zeta}$$

$$= \frac{1}{N} \sum_{i=1}^{N} x_i \zeta_i - \bar{x} \bar{\zeta}.$$
(66)

Using 66, we conclude:

$$NCC(X, \beta_C(X)^{-1}\hat{X}) = \frac{\sigma_X}{\sigma_{\hat{X}}} + \frac{1}{\sigma_X \sigma_{\hat{X}}} \left(\frac{1}{N} \sum_{i=1}^N x_i \zeta_i - \bar{x}\bar{\zeta} \right).$$

Remark B.1. Since the entries of the target $\{x_1,...,x_N\}$ and the noise $\{\zeta_1,...,\zeta_N\}$ are independent, the sample covariance, i.e., multiplicative term $\frac{1}{N}\sum_{i=1}^N x_i\zeta_i - \bar{x}\bar{\zeta}$ on the right-hand side of 62, converges to its theoretical value, i.e., zero, for $N\to\infty$. Otherwise, for $N<\infty$, the sample covariance can be positive or negative, increasing or decreasing the right-hand side of 62 arbitrarily, hindering the computation of meaningful upper and lower bounds for the sample NCC.

Proposition 6. Let x and \hat{x} be the two random variables as defined above. Then,

$$NCC(x,\hat{x}) = \sqrt{\frac{1}{1 + \sigma^2 ||X||_2^2 / Var(x)}} \le \sqrt{\frac{1}{1 + \sigma^2 N}}.$$
 (25)

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Proof. Using the definition of the NCC (see Definition 4) it follows:

$$NCC(x, \hat{x}) = \frac{Cov(x, \hat{x})}{\sqrt{Var(x)}\sqrt{Var(\hat{x})}}$$

$$= \frac{Cov(x, x) + Cov(x, \zeta)}{\sqrt{Var(x)}\sqrt{Var(x) + Var(\zeta)}}$$

$$= \frac{Var(x)}{\sqrt{Var(x)}\sqrt{Var(x) + Var(\xi)}}$$

$$= \sqrt{\frac{Var(x)}{Var(x) + Var(\zeta)}}$$

$$= \sqrt{\frac{1}{1 + Var(\zeta)/Var(x)}}.$$
(67)

We recall that $\text{Var}(\zeta) = \sigma^2 ||X||_2^2$ by definition. Thus, using the assumption that $\text{Var}(x) \leq ||X||_2^2/N$, it follows from 67:

$$NCC(x, \hat{x}) = \sqrt{\frac{1}{1 + \sigma^2 ||X||_2^2 / Var(x)}} \le \sqrt{\frac{1}{1 + \sigma^2 N}}.$$
 (68)

C Additional Figures

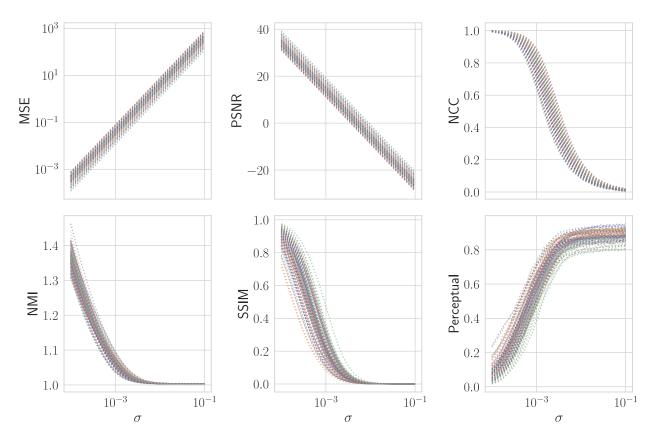


Figure 4: Reconstruction metrics for 100 reconstructed data samples from the ImageNet dataset under varying σ . Additionally to the MSE, PSNR, and NCC we also include the Normalized Mutual Information (NMI), Structural Similarity (SSIM) and a perceptual loss. We see that the NMI, SSIM and the perceptual loss show a very similar result as our bounded metrics, especially the NCC. Of note, the perceptual loss inherently incorporates an image prior due to the network being trained on image data. Although this violates our assumption for the adversary, the metric still aligns with all other reconstruction metrics.

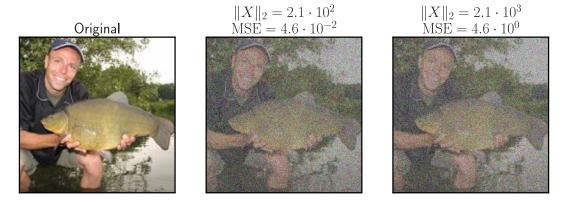


Figure 5: Demonstration how the ℓ_2 -norm of the data sample influences the MSE. The same image, which is rescaled by a constant factor a=10, is being reconstructed. We keep $\sigma=1.0\cdot 10^{-3}$ constant for both and set $C=\|X\|_2$. The resulting MSE is scaled by the factor a^2 .

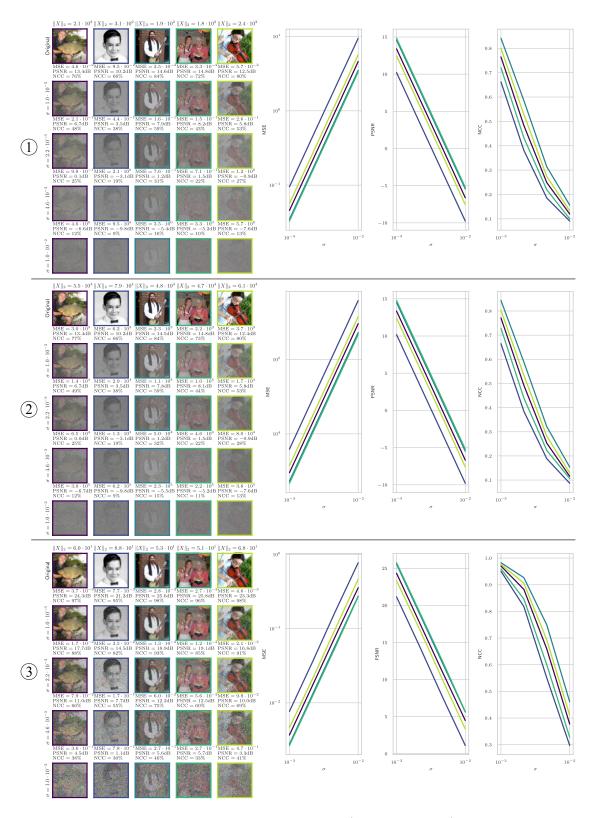


Figure 6: Exemplary reconstructions on the ImageNet dataset (Deng et al., 2009) under varying σ . Scenario (1): $N=224\times224\times3$, data range in (0,1). Scenario (2): $N=224\times224\times3$, data range in (0,255). Scenario (3): $N=64\times64\times3$, data range in (0,1). All scenarios: C=1, M=1. The colours of the image frames match the corresponding metric curves.