# DP-SGD Without Clipping: The Lipschitz Neural Network Way

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#### Abstract

State-of-the-art approaches for training Differentially Private (DP) Deep Neural 1 2 Networks (DNN) faces difficulties to estimate tight bounds on the sensitivity of 3 the network's layers, and instead rely on a process of per-sample gradient clipping. This clipping process not only biases the direction of gradients but also proves 4 5 costly both in memory consumption and in computation. To provide sensitivity bounds and bypass the drawbacks of the clipping process, our theoretical analysis 6 of Lipschitz constrained networks reveals an unexplored link between the Lipschitz 7 8 constant with respect to their input and the one with respect to their parameters. 9 By bounding the Lipschitz constant of each layer with respect to its parameters we guarantee DP training of these networks. This analysis not only allows the 10 computation of the aforementioned sensitivities at scale but also provides leads 11 on to how maximize the gradient-to-noise ratio for fixed privacy guarantees. To 12 facilitate the application of Lipschitz networks and foster robust and certifiable 13 learning under privacy guarantees, we provide a Python package that implements 14 building blocks allowing the construction and private training of such networks. 15

## 16 **1** Introduction

Machine learning relies more than ever on foundational models, and such practices raise questions 17 about privacy. Differential privacy allows to develop methods for training models that preserve 18 the privacy of individual data points in the training set. The field seeks to enable deep learning on 19 sensitive data, while ensuring that models do not inadvertently memorize or reveal specific details 20 about individual samples in their weights. This involves incorporating privacy-preserving mechanisms 21 into the design of deep learning architectures and training algorithms, whose most popular example 22 is Differentially Private Stochastic Gradient Descent (DP-SGD) [1]. One main drawback of classical 23 DP-SGD methods is that they require costly per-sample backward processing and gradient clipping. 24 In this paper, we offer a new method that unlocks fast differentially private training through the use 25 of Lipschitz constrained neural networks. Additionally, this method offers new opportunities for 26 practitioners that wish to easily "DP-fy" [2] the training procedure of a deep neural network. 27

**Differential privacy fundamentals.** Informally, differential privacy is a *definition* that quantifies how 28 much the change of a single sample in a dataset affects the range of a stochastic function (here the DP 29 training), called mechanism in this context. This quantity can be bounded in an inequality involving 30 two parameters  $\epsilon$  and  $\delta$ . A mechanism fulfilling such inequality is said  $(\epsilon, \delta)$ -DP (see Definition 1). 31 This definition is universally accepted as a strong guarantee against privacy leakages under various 32 scenarii, including data aggregation or post-processing [3]. A popular rule of thumb suggests using 33  $\epsilon \leq 10$  and  $\delta < \frac{1}{N}$  with N the number of records [2] for mild guarantees. In practice, most classic 34 algorithmic procedures (called queries in this context) do not readily fulfill the definition for useful 35 36 values of  $(\epsilon, \delta)$ , in particular the deterministic ones: randomization is mandatory. This randomization

```
model = DP_Sequential ( # step 1: use DP_Sequential to build a model
    Γ
        # step 2: add Lipschitz layers of known sensitivity
        DP_BoundedInput(input_shape=(28, 28, 1), upper_bound=20.),
        DP_SpectralConv2D(filters=16, kernel_size=3, use_bias=False),
        DP_GroupSort(2),
        DP_Flatten(),
        DP_SpectralDense(10),
    ],
    noise_multiplier = 1.2, # step 3: choose DP parameters
    sampling_probability = batch_size / dataset_size,
) # step 4: compile the model, and choose any first order optimizer
model.compile(loss=DP_Crossentropy(), optimizer=Adam(1e-3))
model.fit( # step 5: train the model and measure the DP guarantees
    train_dataset, validation_data=val_dataset,
    epochs=num_epochs, callbacks=[DP_Accountant()]
)
```

Figure 1: An example of usage of our framework, illustrating how to create a small Lipschitz VGG and how to train it under  $(\epsilon, \delta)$ -DP guarantees while reporting  $(\epsilon, \delta)$  values.

comes at the expense of "utility", i.e the usefulness of the output for downstream tasks [4]. The goal 37 is then to strike a balance between privacy and utility, ensuring that the released information remains 38 useful and informative for the intended purpose while minimizing the risk of privacy breaches. The 39 privacy/utility trade-off yields a Pareto front, materialized by plotting  $\epsilon$  against a measurement of 40 utility, such as validation accuracy for a classification task. 41 Private gradient descent. The SGD algorithm consists of a sequence of queries that (i) take the 42 dataset in input, sample a minibatch from it, and return the gradient of the loss evaluated on the 43 minibatch, before (ii) performing a descent step following the gradient direction. The sensitivity (see 44 Definition 2) of SGD queries is proportional to the norm of the per-sample gradients. DP-SGD turns 45 each query into a Gaussian mechanism by perturbing the gradients with a noise  $\zeta$ . The upper bound 46 on gradient norms is generally unknown in advance, which leads practitioners to clip it to C > 0, in 47 order to bound the sensitivity manually. This is problematic for several reasons: 1. Hyper-parameter 48 search on the broad-range clipping value C is required to train models with good privacy/utility trade-49 50 offs [5], 2. The computation of per-sample gradients is expensive: DP-SGD is usually slower and 51 consumes more memory than vanilla SGD, in particular for the large batch sizes often used in private training [6], **3.** Clipping the per-sample gradients biases their average [7]. This is problematic as the 52 average direction is mainly driven by misclassified examples, that carry the most useful information 53 for future progress. 54

An unexplored approach: Lipschitz constrained networks. We propose to train neural networks for which the parameter-wise gradients are provably and analytically bounded during the whole training procedure, in order to get rid of the clipping process. This allows for rapid training of models without a need for tedious hyper-parameter optimization.

The main reason why this approach has not been experimented much in the past is that upper bounding
the gradient of neural networks is often intractable. However, by leveraging the literature of Lipschitz
constrained networks [8], we show that these networks allows to estimate their gradient bound.
This yields tight bounds on the sensitivity of SGD steps, making their transformation into Gaussian
mechanisms inexpensive - hence the name Clipless DP-SGD.
Informally, the Lipschitz constant quantifies the rate at which the function's output varies with respect

to changes in its input. A Lipschitz constrained network is one in which its weights and activations
are constrained such that it can only represent *l*-Lipschitz functions. In this work, we will focus our
attention on feed-forward networks (refer to Definition 3). Note that the most common architectures,
such as Convolutional Neural Networks (CNNs), Fully Connected Networks (FCNs), Residual
Networks (ResNets), or patch-based classifiers (like MLP-Mixers), all fall under the category of
feed-forward networks. We will also tackle the particular case of Gradient Norm Preserving (GNP)
networks, a subset of Lipschitz networks that enjoy tighter bounds (see appendix).

#### 72 Contributions

73 While the properties of Lipschitz constrained networks regarding their inputs are well explored, the

<sup>74</sup> properties with respect to its parameters remain non-trivial. This work provides a first step to fill this

- <sup>75</sup> gap: our analysis shows that under appropriate architectural constraints, a *l*-Lipschitz network has a
- <sup>76</sup> tractable, finite Lipschitz constant with respect to its parameters. We prove that this Lipschitz constant
- allows for easy estimation of the sensitivity of the gradient computation queries. The prerequisite and
   details of the method to compute the sensitivities are explained in Section 2.
- 79 Our contributions are the following:
- We extend the field of applications of Lipschitz constrained neural networks. So far the
   literature focused on Lipschitzness with respect to the *inputs*: we extend the framework to
   compute the Lipschitzness with respect to the parameters. This is exposed in Section 2.
- We propose a general framework to handle layer gradient steps as Gaussian mechanisms
   that depends on the loss and the model structure. Our framework covers widely used
   architectures, including VGG and ResNets.
- 3. We show that SGD training of deep neural networks can be achieved without gradient
   clipping using Lipschitz layers. This allows the use of larger networks and larger batch
   sizes, as illustrated by our experiments in Section 4.
- 4. We establish connections between Gradient Norm Preserving (GNP) networks and improved privacy/utility trade-offs (Section 3.1).
- 5. Finally, a **Python package**<sup>1</sup> companions the project, with pre-computed Lipschitz constant and noise for each layer type, ready to be forked on any problem of interest (Section 3.2).

#### 93 1.1 Differential Privacy and Lipschitz Networks

The definition of DP relies on the notion of neighboring datasets, i.e datasets that vary by at most one example. We highlight below the central tools related to the field, inspired from [9].

**Definition 1**  $((\epsilon, \delta)$ -Differential Privacy). A labeled dataset  $\mathcal{D}$  is a finite collection of input/label pairs  $\mathcal{D} = \{(x_1, y_1), (x_2, y_2), \dots, (x_N, y_N)\}$ . Two datasets  $\mathcal{D}$  and  $\mathcal{D}'$  are said to be neighboring for the "replace-one" relation if they differ by at most one sample:  $\mathcal{D}' = \mathcal{D} \cup \{(x'_i, y'_i)\} \setminus \{(x_i, y_i)\}$ . Let  $\epsilon$  and  $\delta$  be two non-negative scalars. A mechanism  $\mathcal{A}$  is  $(\epsilon, \delta)$ -DP if for any two neighboring datasets  $\mathcal{D}$  and  $\mathcal{D}'$ , and for any  $S \subseteq range(\mathcal{A})$ :

$$\mathbb{P}[\mathcal{A}(\mathcal{D}) \in S] \le e^{\epsilon} \times \mathbb{P}[\mathcal{A}(\mathcal{D}') \in S] + \delta.$$
(1)

A cookbook to create a  $(\epsilon, \delta)$ -DP mechanism from a query is to compute its *sensitivity*  $\Delta$  (see Definition 2), and to perturb its output by adding a Gaussian noise of predefined variance  $\zeta^2 = \Delta^2 \sigma^2$ , where the  $(\epsilon, \delta)$ -DP guarantees depends on  $\sigma$ . This yields what is called a *Gaussian mechanism* [3].

**Definition 2** ( $l_2$ -sensitivity). Let  $\mathcal{M}$  be a query mapping from the space of the datasets to  $\mathbb{R}^p$ . Let  $\mathcal{N}$ 

be the set of all possible pairs of neighboring datasets  $\mathcal{D}, \mathcal{D}'$ . The  $l_2$  sensitivity of  $\mathcal{M}$  is defined by:

$$\Delta(\mathcal{M}) = \max_{\mathcal{D}, \mathcal{D}' \in \mathcal{N}} \|\mathcal{M}(D) - \mathcal{M}(D')\|_2.$$
(2)

**Differentially Private SGD.** The classical algorithm keeps track of  $(\epsilon, \delta)$ -DP values with a *moments accountant* [1] which allows to keep track of privacy guarantees at each epoch, by composing different sub-mechanisms. For a dataset with N records and a batch size b, it relies on two parameters: the sampling ratio  $p = \frac{b}{N}$  and the "noise multiplier"  $\sigma$  defined as the ratio between effective noise strength  $\zeta$  and sensitivity  $\Delta$ . Bounds on gradient norm can be turned into bounds on sensitivity of SGD queries. In "replace-one" policy for  $(\epsilon, \delta)$ -DP accounting, if the gradients are bounded by K > 0, the sensitivity of the gradients averaged on a minibatch of size b is  $\Delta = 2K/b$ .

Crucially, the algorithm requires a bound on  $\|\nabla_{\theta} \mathcal{L}(\hat{y}, y)\|_2 \leq K$ . The whole difficulty lies in bounding tightly this value in advance for neural networks. Currently, gradient clipping serves as a patch to circumvent the issue [1]. Unfortunately, clipping individual gradients in the batch is costly

and will bias the direction of their average, which may induce underfitting [7].

<sup>&</sup>lt;sup>1</sup>Code and documentation are given as supplementary material during review process.

**Lipschitz constrained networks.** Our proposed solution comes from the observation that the norm 117 of the gradient and the Lipschitz constant are two sides of the same coin. The function  $f: \mathbb{R}^m \to \mathbb{R}^n$ 118 is said *l*-Lipschitz for  $l_2$  norm if for every  $x, y \in \mathbb{R}^m$  we have  $||f(x) - f(y)||_2 \le l ||x - y||_2$ . Per 119 Rademacher's theorem [10], its gradient is bounded:  $\|\nabla_x f\| \le l$ . Reciprocally, continuous functions 120 gradient bounded by *l* are *l*-Lipschitz. 121

In Lipschitz networks, the literature has predominantly concentrated on investigating the control 122 of Lipschitzness with respect to the inputs (i.e bounding  $\nabla_x f$ ), primarily motivated by concerns 123 of robustness [11]. However, in this work, we will demonstrate that it is also possible to control 124 Lipschitzness with respect to parameters (i.e bounding  $\nabla_{\theta} f$ ), which is essential for ensuring privacy. 125 Our first contribution will point out the tight link that exists between those two quantities. 126 Definition 3 (Lipschitz feed-forward neural network). A feedforward neural network of depth D, 127

with input space  $\mathcal{X} \subset \mathbb{R}^n$ , output space  $\mathcal{Y} \subset \mathbb{R}^K$  (e.g logits), and parameter space  $\Theta \subset \mathbb{R}^p$ , is a 128

parameterized function  $f: \Theta \times \mathcal{X} \to \mathcal{Y}$  defined by the sequential composition of layers  $f_d$ : 129

$$f(\theta, x) := \left(f_D(\theta_d) \circ \dots \circ f_2(\theta_2) \circ f_1(\theta_1)\right)(x).$$
(3)

The parameters of the layers are denoted by  $\theta = (\theta_d)_{1 \le d \le D} \in \Theta$ . For affine layers, it corresponds to bias and weight matrix  $\theta_d = (W_d, b_d)$ . For activation functions, there is no parameters:  $\theta_d = \emptyset$ . 130 131

Lipschitz networks are feed-forward networks, with the additionnal constraint that each 132 layer  $x_d \mapsto f_d(\theta_d, x_d) := y_d$  is  $l_d$ -Lipschitz for all  $\theta_d$ . Consequently, the function  $x \mapsto f(\theta, x)$ 133

is *l*-Lipschitz with  $l = l_1 \times \ldots \times l_d$  for all  $\theta \in \Theta$ . 134

In practice, this is enforced by using activations with Lipschitz constant  $l_d$ , and by applying a con-135

straint  $\Pi: \mathbb{R}^p \to \Theta$  on the weights of affine layers. This corresponds to spectrally normalized matri-136 137

ces [12, 13], since for affine layers we have  $l_d = ||W_d||_2 := \max_{\|x\| \le 1} ||W_d x||_2$  hence  $\Theta = \{||W_d|| \le l_q\}$ .

The seminal work of [8] proved that universal approximation in the set of *l*-Lipschitz functions was 138 achievable by this family of architectures. Concurrent approaches are based on regularization (like

139 in [14, 15, 16]) but they fail to produce formal guarantees. While they have primarily been studied in

140 the context of adversarial robustness [11, 17], recent works have revealed additional properties of 141

these networks, such as improved generalization [13, 18]. However, the properties of their parameter 142

gradient  $\nabla_{\theta} f(\theta, x)$  remain largely unexplored. 143

#### Clipless DP-SGD with *l*-Lipschitz networks 2 144

Our framework consists of 1. a method that computes the maximum gradient norm of a network with 145 respect to its parameters to obtain a *per-layer* sensitivity  $\Delta_d$ , 2. a moments accountant that relies on 146 the per-layer sensitivities to compute  $(\epsilon, \delta)$ -DP guarantees. The method 1. is based on the recursive 147 formulation of the chain rule involved in backpropagation, while 2. keeps track of  $(\epsilon, \delta)$ -DP values 148 with RDP accounting. It requires some natural assumptions that we highlight below. 149

**Requirement 1** (Lipschitz loss.). The loss function  $\hat{y} \mapsto \mathcal{L}(\hat{y}, y)$  must be L-Lipschitz with respect to 150

the logits  $\hat{y}$  for all ground truths  $y \in \mathcal{Y}$ . This is notably the case of Categorical Softmax-Crossentropy. 151

The Lipschitz constants of common classification losses can be found in the appendix. 152

**Requirement 2** (Bounded input). *There exists*  $X_0 > 0$  *such that for all*  $x \in \mathcal{X}$  *we have*  $||x|| \leq X_0$ . 153

While there exist numerous approaches for the parametrization of Lipschitz networks (e.g differen-154 tiable re-parametrization [19, 8], optimization over matrix manifolds [20] or projections [21]), our 155 framework only provides sensitivity bounds for projection-based algorithms (see appendix). 156

**Requirement 3** (Lipschitz projection). The Lipschitz constraints must be enforced with a projection 157 operator  $\Pi: \mathbb{R}^p \to \Theta$ . This corresponds to Tensorflow [22] constraints and Pytorch [23] hooks. 158

Projection is a post-processing of private gradients: it induces no privacy leakage [3]. 159

To compute the per-layer sensitivities, our framework mimics the backpropagation algorithm, where 160

Vector-Jacobian products (VJP) are replaced by Scalar-Scalar products of element-wise bounds. For 161 an arbitrary layer  $x_d \mapsto f_d(\theta_d, x_d) := y_d$  the operation is sketched below: 162

$$\nabla_{x_d} \mathcal{L} := (\nabla_{y_d} \mathcal{L}) \frac{\partial f_d}{\partial x_d} \implies \|\nabla_{x_d} \mathcal{L}\|_2$$

$$\Rightarrow \|\nabla_{x_d} \mathcal{L}\|_2 \le \|\nabla_{y_d} \mathcal{L}\|_2 \times \left\|\frac{\partial f_d}{\partial x_d}\right\|_2.$$
(4)

Vector-Jacobian product: backpropagate gradients

Scalar-Scalar product: backpropagate bounds



Figure 2: **Backpropagation for bounds**, Algorithm 1. Compute the per-layer sensitivity  $\Delta_d$ .

The notation  $\|\cdot\|_2$  must be understood as the spectral norm for Jacobian matrices, and the Euclidean norm for gradient vectors. The scalar-scalar product is inexpensive. For Lipschitz layers the spectral norm of the Jacobian  $\|\frac{\partial f}{\partial x}\|$  is kept constant during training with projection operator  $\Pi$ . The bound of the gradient with respect to the parameters then takes a simple form:

$$\|\nabla_{\theta_d} \mathcal{L}\|_2 = \|\nabla_{y_d} \mathcal{L}\|_2 \times \left\|\frac{\partial f_d}{\partial \theta_d}\right\|_2.$$
<sup>(5)</sup>

Once again the operation is inexpensive. The upper bound  $\left\|\frac{\partial f}{\partial \theta}\right\|_2$  typically depends on the supremum of  $\|x_d\|_2$ , that can also be analytically bounded, as exposed in the following section.

#### 169 2.1 Backpropagation for bounds

The pseudo-code of **Clipless DP-SGD** is sketched in Algorithm 2. The algorithm avoids clipping by computing a *per-layer* bound on the element-wise gradient norm. The computation of this *per-layer* bound is described by Algorithm 1 (graphically explained in Figure 2). Crucially, it requires to compute the spectral norm of the Jacobian of each layer with respect to input and parameters.

**Input bound propagation (line 2).** We compute  $X_d = \max_{\|x\| \le X_{d-1}} \|f_d(x)\|_2$ . For activation functions it depends on their range. For linear layers, it depends on the spectral norm of the operator itself. This quantity can be computed with SVD or Power Iteration [24, 19], and constrained during training using projection operator  $\Pi$ . In particular, it covers the case of convolutions, for which tight bounds are known [25]. For affine layers, it additionally depends on the amplitude of the bias  $\|b_d\|$ .

**Remark 1** (Tighter bounds in literature.). Although libraries such as Decomon [26] or auto-LiRPA [27] provide tighter bounds  $X_d$  via linear relaxations [28, 29], our approach is capable of delivering practically tighter bounds than worst-case scenarios thanks to the projection operator  $\Pi$ , while also being significantly less computationally expensive. Moreover, hybridizing our method with scalable certification methods can be a path for future extensions.

**Computing maximum gradient norm (line 6).** We bound the Jacobian  $\frac{\partial f_d(\theta_d, x)}{\partial \theta_d}$ . In neural networks, the parameterized layers  $f(\theta, x)$  (fully connected, convolutions) are bilinear operators. Hence we typically obtain bounds of the form:

$$\left\| \frac{\partial f_d(\theta_d, x)}{\partial \theta_d} \right\|_2 \le K(f_d, \theta_d) \|x\|_2 \le K(f_d, \theta_d) X_{d-1}, \tag{6}$$

where  $K(f_d, \Theta_d)$  is a constant that depends on the nature of the operator.  $X_{d-1}$  is obtained in line 2 with input bound propagation. Values of  $K(f_d, \theta_d)$  for popular layers are pre-computed in the library.

**Backpropagate cotangeant vector bounds (line 7).** We bound the Jacobian  $\frac{\partial f_d(\theta_d, x)}{\partial x}$ . For activation functions this value can be hard-coded, while for affine layers it is the spectral norm of the linear operator. Like before, this value is constrained with projection operator II.

#### 192 2.2 Privacy accounting for Clipless DP-SGD

<sup>193</sup> Two strategies are available to keep track of  $(\epsilon, \delta)$  values as the training progresses, based on <sup>194</sup> accounting either a per-layer "local" sensitivity, either by aggregating them into a "global" sensitivity.

#### Algorithm 1 Backpropagation for Bounds(f, X)

**Input**: Feed-forward architecture  $f(\theta, \cdot) = f_D(\theta_D, \cdot) \circ \ldots \circ f_1(\theta_1, \cdot)$ **Input**: Weights  $\theta = (\theta_1, \theta_2, \dots, \theta_D)$ , input bound  $X_0$ 1: for all layers  $1 \le d \le D$  do  $X_d \leftarrow \max \|f_d(\theta_d, x)\|_2.$ 2: ▷ Input bounds propagation  $\|x\| \leq X_{d-1}$ 3: end for 4:  $G \leftarrow L/b$ . Lipschitz constant of the loss for batchsize b 5: for all layers  $D \ge d \ge 1$  do  $\Delta_d \leftarrow G \max_{\|x\| \leq X_{d-1}} \| \frac{\partial f_d(\theta_d, x)}{\partial \theta_d} \|_2.$ Compute sensitivity from gradient norm 6:  $G \leftarrow G \max_{\|x\| \leq X_{d-1}} \| \frac{\partial f_d(\theta_d, x)}{\partial x} \|_2 = G l_d.$ 7: Backpropagate cotangeant vector bounds 8: end for 9: **return** sensitivities  $\Delta_1, \Delta_2, \ldots, \Delta_D$ 

#### Algorithm 2 Clipless DP-SGD with local sensitivity accounting

**Input**: Feed-forward architecture  $f(\theta, \cdot) = f_D(\theta_D, \cdot) \circ \ldots \circ f_1(\theta_1, \cdot)$ **Input**: Initial weights  $\theta = (\theta_1, \theta_1, \dots, \theta_D)$ , learning rate  $\eta$ , noise multiplier  $\sigma$ . 1: repeat  $\Delta_1, \Delta_2 \dots \Delta_D \leftarrow \text{Backpropagation for Bounds}(f, X).$ 2: Update Moment Accountant state with **local** sensitivities  $\Delta_1, \Delta_2, \dots \Delta_d$ . 3: Sample a batch  $\mathcal{B} = \{(x_1, y_1), (x_2, y_2), \dots, (x_b, y_b)\}.$ Compute per-layer averaged gradient:  $g_d := \frac{1}{b} \sum_{i=1}^b \nabla_{\theta_d} \mathcal{L}(f(\theta, x_i), y_i)).$ 4: 5: Sample local noise:  $\zeta_d \sim \mathcal{N}(0, \sigma \Delta_d)$ . 6: 7: Perform noisified gradient step:  $\theta_d \leftarrow \theta_d - \eta(g_d + \zeta_d)$ . Enforce Lipschitz constraint with projection:  $\theta_d \leftarrow \Pi(\theta_d)$ . 8: 9: **until** privacy budget  $(\epsilon, \delta)$ -DP budget has been reached.

**The "global" strategy.** Illustrated in the appendix,this strategy simply aggregates the individual sensitivities  $\Delta_d$  of each layer to obtain the global sensitivity of the whole gradient vector  $\Delta = \sqrt{\sum_d \Delta_d^2}$ . The origin of the clipping-based version of this strategy can be traced back to [30]. With noise variance  $\sigma^2 \Delta^2$  we recover the accountant that comes with DP-SGD. It tends to overestimate the true sensitivity (in particular for deep networks), but its implementation is straightforward with existing tools.

**The "local" strategy.** Recall that we are able to characterize the sensitivity  $\Delta_d$  of every layer of the network. Hence, we can apply a different noise to each of the gradients. We dissect the whole training procedure in Figure 3. At same noise multiplier  $\sigma$ , it tends to produce a higher value of  $\epsilon$  per epoch than "global" strategy, but has the advantage over the latter to add smaller effective noise  $\zeta$  to each weight.

We rely on the autodp<sup>2</sup> library [32, 33, 34] as it uses the Renyi Differential Privacy (RDP) adaptive composition theorem [35, 36], that ensures tighter bounds than naive DP composition.

# **3** From theory to practice

Beyond the application of Algorithms 1 and 2, our framework provides numerous opportunities to enhance our understanding of prevalent techniques identified in the literature. An in-depth exploration of these is beyond the scope of this work, so we focus on giving insights on promising tracks based on our theoretical analysis. In particular, we discuss how the tightness of the bound provided by Algorithm 1 can be influenced by working on the architecture, the input pre-processing and the loss post-processing.

<sup>&</sup>lt;sup>2</sup>https://github.com/yuxiangw/autodp distributed under Apache License 2.0 licence.



Figure 3: Accountant for locally enforced differential privacy. (i) The gradient query for each layer is turned into a Gaussian mechanism [9], (ii) their composition at the scale of the whole network is a non isotropic Gaussian mechanism, (iii) that benefits from amplification via sub-sampling [31], (iv) the train steps are composed over the course of training.

#### 215 3.1 Gradient Norm Preserving networks

- We can manually derive the bounds obtained from Algorithm 2 across diverse configurations. Below, we conduct a sensitivity analysis on *l*-Lipschitz networks.
- **Theorem (informal) 1. Gradient Norm of Lipschitz Networks.** Assume that every layer  $f_d$  is *K-Lipschitz, i.e*  $l_1 = \cdots = l_D = K$ . Assume that every bias is bounded by *B*. We further assume that
- each activation is centered in zero (e.g ReLU, tanh, GroupSort). We recall that  $\theta = [\theta_1, \theta_2, \dots, \theta_D]$ .
- 221 Then the global upper bound of Algorithm 2 can be expanded analytically.
- 222 I. If K < 1 we have:  $\|\nabla_{\theta} \mathcal{L}(f(\theta, x), y)\|_2 = \mathcal{O}\left(L\left(K^D(X_0 + B) + 1\right)\right)$ .
- 223 Due to the  $K^D \ll 1$  term this corresponds to a vanishing gradient phenomenon [37]. The output of 224 the network is essentially independent of its input, and the training is nearly impossible.
- 225 **2.** If K > 1 we have:  $\|\nabla_{\theta} \mathcal{L}(f(\theta, x), y)\|_2 = \mathcal{O}(LK^D(X_0 + B))$ .
- 226 Due to the  $K^D \gg 1$  term this corresponds to an exploding gradient phenomenon [38]. The upper 227 bound becomes vacuous for deep networks: the added noise  $\zeta$  is at risk of being too high.

228 3. If 
$$K = 1$$
 we have:  $\|\nabla_{\theta} \mathcal{L}(f(\theta, x), y)\|_2 = \mathcal{O}\left(L\left(X_0 + \sqrt{D} + \sqrt{BX_0}D + BD^{3/2}\right)\right)$ 

which for linear layers without biases further simplify to  $\mathcal{O}(L(X_0 + \sqrt{D}))$ .

The formal statement can be found in appendix. From Theorem 1 we see that most favorable bounds 230 are achieved by 1-Lipschitz neural networks with 1-Lipschitz layers. In classification tasks, they are 231 not less expressive than conventional networks [18]. Hence, this choice of architecture is not at the 232 expense of utility. Moreover an accuracy/robustness trade-off exists, determined by the choice of 233 loss function [18]. However, setting K = 1 merely ensures that  $\|\nabla_x f\| \le 1$ , and in the worst-case 234 scenario we have  $\|\nabla_x f\| < 1$  almost everywhere. This could result in a situation where the bound of 235 case 3 in Theorem 1 is not tight, leading to an underfitting regime as in case K < 1. With Gradient 236 Norm Preserving (GNP) networks [17], we expect to mitigate this issue. 237

**Controlling** *K* with Gradient Norm Preserving (GNP) networks. GNP networks are 1-Lipschitz neural networks with the additional constraint that the Jacobian of layers consists of orthogonal matrices. They fulfill the Eikonal equation  $\left\|\frac{\partial f_d(\theta_d, x_d)}{\partial x_d}\right\|_2 = 1$  for any intermediate activation  $f_d(\theta_d, x_d)$ . Without biases these networks are also norm preserving:  $\|f(\theta, x)\| = \|x\|$ .

As a consequence, the gradient of the loss with respect to the parameters is easily bounded by

$$\|\nabla_{\theta_d} \mathcal{L}\| = \|\nabla_{y_D} \mathcal{L}\| \times \left\| \frac{\partial f_d(\theta_d, x_d)}{\partial \theta_d} \right\|,\tag{7}$$

which for weight matrices  $W_d$  further simplifies to  $\|\nabla_{W_d} \mathcal{L}\| \le \|\nabla_{y_D} \mathcal{L}\| \times \|f_{d-1}(\theta_{d-1}, x_{d-1})\|$ . We see that this upper bound crucially depends on two terms than can be analyzed separately. On one hand,  $\|f_{d-1}(\theta_{d-1}, x_{d-1})\|$  depends on the scale of the input. On the other,  $\|\nabla_{y_D} \mathcal{L}\|$  depends on the loss, the predictions and the training stage. We show below how to intervene on these two quantities. **Remark 2** (Implementation of GNP Networks). *In practice, GNP are parametrized with GroupSort* activation [8, 39], Householder activation [40], and orthogonal weight matrices [17, 41]. Strict orthogonality is challenging to enforce, especially for convolutions for which it is still an active research area (see [42, 43, 44, 45, 46] and references therein). Our line of work traces an additional motivation for the development of GNP and the bounds will strengthen as the field progresses.

**Controlling**  $X_0$  with input pre-processing. The weight gradient norm  $\|\nabla_{\theta_d} \mathcal{L}\|$  indirectly depends on the norm of the inputs. This observation implies that the pre-processing of input data significantly influences the bounding of sensitivity. Multiple strategies are available to keep the input's norm under control: projection onto the ball ("norm clipping"), or projection onto the sphere ("normalization"). In the domain of natural images for instance, this result sheds light on the importance of color space such as RGB, HSV, YIQ, YUV or Grayscale. These strategies are natively handled by our library.

Controlling L with the hybrid approach, loss gradient clipping. As training progresses, the 258 magnitude of  $\|\nabla_f \mathcal{L}\|$  tends to diminish when approaching a local minima, quickly falling below the 259 upper bound and diminishing the gradient norm to noise ratio. To circumvent the issue, the gradient 260 clipping strategy is still available in our framework. Crucially, instead of clipping the parameter 261 gradient  $\nabla_{\theta} \mathcal{L}$ , any intermediate gradient  $\nabla_{f_d} \mathcal{L}$  can be clipped during backpropagation. This can be achieved with a special "*clipping layer*" that behaves like the identity function at the forward 262 263 pass, and clips the gradient during the backward pass. The resulting cotangeant vector is not a true 264 gradient anymore, but rather a descent direction [47]. In vanilla DP-SGD the clipping is applied on 265 the batched gradient  $\nabla_{W_d} \mathcal{L}$  of size  $b \times h^2$  for matrix weight  $W_d \in \mathbb{R}^{h \times h}$  and clipping this vector can cause memory issues or slowdowns [6]. In our case,  $\nabla_{y_D} \mathcal{L}$  is of size  $b \times h$  which reduces overhead. 266 267

#### 268 3.2 Lip-dp library

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To foster and spread accessibility, we provide an opensource tensorflow library for Clipless DP-SGD training, named lip-dp. It provides an exposed Keras API for seamless usability. It is implemented as a wrapper over the Lipschitz layers of deel-lip<sup>3</sup> library [48]. Its usage is illustrated in Figure 1.

#### 272 **4 Experimental results**

We validate our implementation with a speed benchmark against competing approaches, and we present the privacy/utility Pareto front that can be obtained with GNP networks.

Speed and memory consumption. We bench-275 marked the median runtime per epoch of vanilla 276 DP-SGD against the one of Clipless DP-SGD, 277 on a CNN architecture and its Lipschitz equiv-278 alent respectively. The experiment was run on 279 a GPU with 48GB video memory. We compare 280 against the implementation of tf\_privacy, 281 opacus and optax. In order to allow a fair com-282 parison, when evaluating Opacus, we reported 283 the runtime with respect to the logical batch size, 284 while capping the physical batch size to avoid 285 Out Of Memory error (OOM). Although our li-286 brary does not implement logical batching yet, 287 it is fully compatible with this feature. 288

An advantage of projection  $\Pi$  over per-sample

gradient clipping is that the projection cost is



Figure 4: Our approach outperforms concurrent frameworks in terms of runtime and memory: we trained CNNs (ranging from 130K to 2M parameters) on CIFAR-10, and report the median batch processing time (including noise, and constraints application  $\Pi$  or gradient clipping).

independent of the batch size. Fig 4 validates that our method scales much better than vanilla DP-SGD, and is compatible with large batch sizes. It offers several advantages: firstly, a larger batch size contributes to a decrease of the sensitivity  $\Delta \propto 1/b$ , which diminishes the ratio between noise and gradient norm. Secondly, as the batch size *b* increases, the variance decreases at the parametric rate  $O(\sqrt{b})$  (as demonstrated in appendix), aligning with expectations. This observation does not apply to DP-SGD: gradient clipping biases the direction of the average gradient, as noticed by [7].

<sup>3</sup>https://github.com/deel-ai/deel-lip distributed under MIT License (MIT).



Figure 5: **Our framework paints a clearer picture of the privacy/utility trade-off.** We trained models in an "out of the box setting" (no pre-training, no data augmentation and no handcrafted features) on multiple tasks. While our results align with the baselines presented in other frameworks, we recognize the importance of domain-specific engineering. In this regard, we find the innovations introduced in [49, 50, 51] and references therein highly relevant. These advancements demonstrate compatibility with our framework and hold potential for future integration.

**Pareto front of privacy/utility trade-off.** We performed a search over a broad range of hyper-297 parameters values to cover the Pareto front between utility and privacy. Results are reported in 298 Figure 5. We emphasize that our experiments did not use the elements behind the success of most 299 recent papers (pre-training, data preparation, or handcrafted feature are examples). Hence our 300 results are more representative of the typical performance that can be obtained in an "out of the 301 box" setting. Future endeavors or domain-specific engineering can enhance the performance even 302 further, but such improvements currently lie beyond the scope of our work. We also benchmarked 303 architectures inspired from VGG [52], Resnet [53] and MLP\_Mixers [54] see appendix for more 304 details. Following standard practices of the community [2], we used sampling without replacement at 305 each epoch (by shuffling examples), but we reported  $\epsilon$  assuming *Poisson sampling* to benefit from 306 privacy amplification [31]. We also ignore the privacy loss that may be induced by hyper-parameter 307 308 search, which is a limitation per recent studies [5], but is common practice.

# **309 5 Limitations and future work**

Although this framework offers a novel approach to address differentially private training, it introduces 310 new challenges. We primary rely on GNP networks, where high performing architectures are 311 quite different from the usual CNN architectures. As emphasized in Remark 2, we anticipate that 312 progress in these areas would greatly enhance the effectiveness of our approach. Additionally, to 313 meet requirement 3, we rely on projections, necessitating additional efforts to incorporate recent 314 advancements associated with differentiable reparametrizations [42, 43]. It is worth noting that 315 our methodology is applicable to most layers. Another limitation of our approach is the accurate 316 computation of sensitivity  $\Delta$ , which is challenging due to the non-associativity of floating-point 317 arithmetic and its impact on numerical stability [55]. This challenge is exacerbated on GPUs, where 318 operations are inherently non-deterministic [56]. Finally, as mentioned in Remark 1, our propagation 319 bound method can be refined. 320

### 321 6 Concluding remarks and broader impact

Besides its main focus on differential privacy, our work provides (1) a motivation to further develop Gradient Norm Preserving architectures. Furthermore, the development of networks with known Lipschitz constant with respect to parameters is a question of independent interest, (2) a useful tool for the study of the optimization dynamics in neural networks. Finally, Lipschitz networks are known to enjoy certificates against adversarial attacks [17, 57], and from generalization guarantees [13], without cost in accuracy [18]. We advocate for the spreading of their use in the context of robust and certifiable learning.

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