
GeoClip: Geometry-Aware Clipping for Differentially Private SGD

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Abstract

Differentially private stochastic gradient descent (DP-SGD) is the most widely used method for training machine learning models with provable privacy guarantees. A key challenge in DP-SGD is setting the per-sample gradient clipping threshold, which significantly affects the trade-off between privacy and utility. While recent adaptive methods improve performance by adjusting this threshold during training, they operate in the standard coordinate system and fail to account for correlations across the coordinates of the gradient. We propose GeoClip, a geometry-aware framework that clips and perturbs gradients in a transformed basis aligned with the geometry of the gradient distribution. GeoClip adaptively estimates this transformation using only previously released noisy gradients, incurring no additional privacy cost. We provide convergence guarantees for GeoClip and derive a closed-form solution for the optimal transformation that minimizes the amount of noise added while keeping the probability of gradient clipping under control. Experiments on both tabular and image datasets demonstrate that GeoClip consistently outperforms existing adaptive clipping methods under the same privacy budget.

1 Introduction

As machine learning models are increasingly trained on sensitive user data, ensuring strong privacy guarantees during training is essential to reduce the risk of misuse, discrimination, or unintended data exposure. Differential privacy (DP) [Dwork et al., 2006a,b] offers a principled framework for protecting individual data, and has become a cornerstone of privacy-preserving machine learning. In deep learning, the most widely used approach for DP is differentially private stochastic gradient descent (DP-SGD) [Abadi et al., 2016], which clips per-sample gradients and adds calibrated noise to their average.

Despite its widespread use, standard DP-SGD has a key limitation: it relies on a fixed clipping threshold to bound the sensitivity of individual gradients. Selecting this threshold poses a challenging privacy-utility tradeoff—setting it too low discards useful gradient information, while setting it too high increases sensitivity and necessitates injecting more noise, ultimately degrading model performance. This trade-off was observed empirically by McMahan et al. [2018] and later analyzed theoretically and shown to be a fundamental limitation of differentially private learning by Amin et al. [2019]. Moreover, the optimal threshold can vary over the course of training, across tasks, and between datasets, limiting the effectiveness of a fixed setting.

To address this, recent work has proposed adaptive strategies that dynamically adjust the clipping threshold during training. One class of methods uses decay schedules to reduce the threshold over time. Yu et al. [2018] and Du et al. [2021] propose linear and near-linear decay rules, respectively, where the schedule is predefined and does not depend on the dataset—hence, no privacy budget is required. Lin et al. [2022] introduce a nonlinear decay schedule, along with a transfer strategy that leverages public data to guide threshold selection. Recently, methods have been introduced which set the clipping level based on the data during the training process. These include AdaClip [Pichapati et al., 2019], which applies coordinate-wise clipping based on estimated gradient variances. Yet another is quantile-based clipping [Andrew et al., 2021], which sets the threshold using differentially private quantiles of per-sample gradient norms. Although [Andrew et al., 2021] was designed for federated learning, it can be adapted to centralized DP-SGD. These adaptive strategies have been shown to improve model utility while preserving privacy guarantees.

Despite these advances, existing adaptive clipping methods remain agnostic to the geometry of the gradient distribution. They operate in the standard basis—treating each coordinate independently. This overlooks dependencies between coordinates, especially when gradients exhibit strong correlations across dimensions. In such cases, independently clipping and perturbing each coordinate can introduce redundant noise without improving privacy, ultimately degrading model utility. To address this, we propose *GeoClip*, a method that transforms gradients into a decorrelated basis that better reflects their underlying geometry. By applying DP mechanisms in this transformed space, GeoClip allocates noise more effectively, achieving a better privacy-utility tradeoff.

A different approach to correlations in gradients from the literature considers introducing correlations—dependencies across iterations and between entries of the noise vector—into the noise, rather than injecting i.i.d. Gaussian noise. This has been shown to improve the utility of private training [Denisov et al., 2022, Kairouz et al., 2021]. Choquette-Choo et al. [2024] strengthen this direction by analytically characterizing near-optimal spatio-temporal correlation structures that lead to provably tighter privacy-utility tradeoffs. However, these approaches have been developed independently of adaptive clipping, and the effect of combining both methods remains unexplored. In addition, [Choquette-Choo et al., 2024] uses pre-determined correlations in the noise, rather than being tailored to the data, as our approach is.

GeoClip is data-driven but does not require any additional privacy budget to compute the clipping transformation. Instead, it reuses the noisy gradients already released during training to estimate the mean and the covariance of the gradient distribution. By reusing these privatized gradients, GeoClip adapts its basis over time without accessing raw data or incurring additional privacy cost.

We list our main contributions below:

1. We propose GeoClip, a novel framework that applies differential privacy in a transformed basis rather than the standard coordinate system. To guide the choice of transformation, we derive a convergence theorem (Theorem 1) showing how the basis impacts convergence under DP-SGD, providing theoretical guidance for selecting transformations that improve utility.
2. Building on this insight, we formulate a convex optimization problem to find the transformation, and derive a closed-form solution (Theorem 2).
3. We introduce two algorithms to estimate the transformation using only previously released noisy gradients. The first is based on a moving average to estimate the gradient covariance matrix.
4. For large-scale models, the full gradient covariance matrix is prohibitively large to store. Thus, our second algorithm uses a streaming low-rank approximation of the covariance matrix. This second algorithm is thus suitable for deep models with large parameter counts.
5. We validate GeoClip through experiments on synthetic, tabular, and image datasets, showing that it consistently outperforms existing adaptive clipping methods under the same privacy budget.

Notation. We denote the $d \times d$ identity matrix by I_d . The notation $N \sim \mathcal{N}(0, \sigma^2 I_d)$ denotes a d -dimensional Gaussian with zero mean and covariance $\sigma^2 I_d$. We use $\|x\|_2$ for the Euclidean norm of a vector x , A^{-1} for the inverse of matrix A , and A^\top for its transpose. The trace of a matrix A is denoted by $\text{Tr}(A)$, and $\text{Cov}(x \mid y)$ refers to the conditional covariance of x given y .

2 General Framework

Let $\mathcal{D} = \{x_k\}_{k=1}^N$ be a dataset of N examples, and let $f : \mathbb{R}^d \rightarrow \mathbb{R}$ denote the empirical loss function defined as the average of per-sample losses:

$$f(\theta) = \frac{1}{N} \sum_{k=1}^N f_k(\theta), \quad (1)$$

where $\theta \in \mathbb{R}^d$ is the model parameter vector and each f_k corresponds to the loss on the k -th data point x_k . In DP-SGD, the algorithm updates θ using a noisy clipped stochastic gradient to ensure privacy. Let $g_t \in \mathbb{R}^d$ denote the stochastic gradient at iteration t . The update rule is:

$$\theta_{t+1} = \theta_t - \eta \tilde{g}_t, \quad (2)$$

where \tilde{g}_t is the privatized version of g_t , obtained by clipping and adding noise.

GeoClip builds on the DP-SGD framework but changes how gradients are processed before clipping and adding noise. It begins by shifting and projecting the gradient into a new coordinate system:

$$\omega_t = M_t(g_t - a_t), \quad (3)$$

where $a_t \in \mathbb{R}^d$ is a reference point and $M_t \in \mathbb{R}^{d \times d}$ is a full-rank transformation matrix that defines the new basis. To enforce differential privacy, we clip the transformed gradient ω_t to unit norm and add Gaussian noise:

$$\tilde{\omega}_t = \frac{\omega_t}{\max(1, \|\omega_t\|_2)} + N_t, \quad N_t \sim \mathcal{N}(0, \sigma^2 I_d), \quad (4)$$

where σ is set based on the desired privacy guarantee. We then map the noisy, clipped gradient back to the original space:

$$\tilde{g}_t = M_t^{-1} \tilde{\omega}_t + a_t. \quad (5)$$

Remark 1. *GeoClip generalizes AdaClip of Pichapati et al. [2019], which itself extends standard DP-SGD. AdaClip essentially assumes that M_t is diagonal for per-coordinate scaling, whereas GeoClip allows M_t to be any full-rank matrix. This added flexibility allows GeoClip to account for correlations between gradient components and inject noise along more meaningful directions.*

A key feature of our framework is that the privacy guarantees remain unaffected by the choice of M_t and a_t , as long as no privacy budget is used to compute them. This allows M_t and a_t to be chosen entirely based on utility, without compromising privacy. The main challenge, then, is how to select these parameters effectively. To address this, we first define a performance metric for GeoClip. Inspired by Pichapati et al. [2019], we derive the following convergence bound for our framework and use it to guide the design of M_t and a_t , ultimately improving both convergence and training efficiency.

Theorem 1 (Convergence of GeoClip). *Assume f has an L -Lipschitz continuous gradient. Further, assume the stochastic gradients are bounded, i.e., $\|\nabla f_k(\theta)\| \leq G$, and have bounded variance, i.e., $\mathbb{E}_k \|\nabla f_k(\theta) - \nabla f(\theta)\|^2 \leq \sigma_g^2$. Let $\theta^* = \arg \min_{\theta \in \mathbb{R}^d} f(\theta)$ denote the optimal solution, and suppose the learning rate satisfies $\eta < \frac{2}{3L}$. Then, for the iterates $\{\theta_t\}_{t=0}^{T-1}$ produced by GeoClip with batch size 1 using the update rule $\theta_{t+1} = \theta_t - \eta \tilde{g}_t$, where \tilde{g}_t is defined in (5), the average squared gradient norm satisfies:*

$$\begin{aligned} \frac{1}{T} \sum_{t=0}^{T-1} \mathbb{E} \|\nabla f(\theta_t)\|^2 &\leq \underbrace{\frac{f(\theta_0) - f(\theta^*)}{T \left(\eta - \frac{3L\eta^2}{2} \right)}}_{\text{Optimization gap}} + \underbrace{\frac{3L\eta}{2-3L\eta} \sigma_g^2}_{\text{Gradient variance term}} + \underbrace{\frac{L\eta\sigma^2}{T(2-3L\eta)} \sum_{t=0}^{T-1} \mathbb{E} \text{Tr} \left[(M_t^\top M_t)^{-1} \right]}_{\text{Noise-injection term}} \\ &\quad + \underbrace{\frac{2}{T(2-3L\eta)} \sum_{t=0}^{T-1} \mathbb{E} [\beta(a_t) (\text{Tr} (M_t^\top M_t \Sigma_t) + \|M_t(\mathbb{E}[g_t | \theta^t] - a_t)\|^2)]}_{\text{Clipping error term}}, \end{aligned} \quad (6)$$

where $\theta^t = (\theta_0, \dots, \theta_t)$ represents the history of parameter values up to iteration t , $\Sigma_t = \text{Cov}(g_t | \theta^t)$, and

$$\beta(a_t) = (G + \|a_t\|) \left(G + \frac{3L\eta}{2} (G + \|a_t\|) \right). \quad (7)$$

The above result generalizes Theorem 2 in [Pichapati et al. \[2019\]](#); proof details are in the supplement.

[Theorem 1](#) provides insight into how we should choose the transformation parameters a_t and M_t . In particular, we want to choose the transformation parameters a_t and M_t , for all t , to minimize the right-hand side of (6). The reference point a_t directly affects the clipping error by setting the center around which gradients are clipped, thereby influencing how much of each gradient is truncated. In contrast, the noise injection term remains independent of a_t , as noise is added regardless of the gradient's position relative to a_t . From [Theorem 1](#), the clipping error at iteration t takes the form

$$\beta(a_t) (\text{Tr}(M_t^\top M_t \Sigma_t) + \|M_t(\mathbb{E}[g_t | \theta^t] - a_t)\|^2), \quad (8)$$

where $\beta(a_t)$ is a scale factor that grows with $\|a_t\|$. A natural choice is to set $a_t = \mathbb{E}[g_t | \theta^t]$ which eliminates the bias term $\|M_t(\mathbb{E}[g_t | \theta^t] - a_t)\|^2$. Given this choice, since gradients are norm-bounded by G , we can apply Jensen's inequality to also bound the norm of their mean, to obtain the bound

$$\beta(a_t) \leq 2G^2(1 + 3L\eta). \quad (9)$$

We now show that the remaining contribution to the clipping error term $\text{Tr}(M_t^\top M_t \Sigma_t)$ in fact serves as an upper bound on the probability that the gradient is clipped (i.e., $|\omega| > 1$). We do so using Markov's inequality as detailed below:

$$\text{Tr}(M_t^\top M_t \Sigma_t) = \mathbb{E}[\|M_t(g_t - \mathbb{E}[g_t | \theta^t])\|^2 | \theta^t] \quad (10)$$

$$= \mathbb{E}[\|\omega\|^2 | \theta^t] \quad (11)$$

$$\geq \Pr(\|\omega\| > 1 | \theta^t). \quad (12)$$

We can now interpret the clipping error term as the likelihood that clipping occurs—ideally, the lower the better as the gradient information will be better preserved. However, in setting a clipping level, we must also be aware of the amount of noise: if M_t is scaled down, there is effectively more noise, as captured by the noise-injection term in (6). We handle this tradeoff via the following optimization problem for the transformation matrix M_t :

$$\begin{aligned} & \underset{M_t}{\text{minimize}} \quad \text{Tr}(M_t^\top M_t)^{-1} \\ & \text{subject to} \quad \text{Tr}(M_t^\top M_t \Sigma_t) \leq \gamma. \end{aligned} \quad (13)$$

Theorem 2. Let $\Sigma_t = \text{Cov}(g_t | \theta^t)$ be a positive definite matrix. The optimal transformation matrix $M_t^* \in \mathbb{R}^{d \times d}$ at iteration t for the optimization problem in (13), along with its corresponding objective value, are given by:

$$M_t^* = \left(\frac{\gamma}{\sum_{i=1}^d \sqrt{\lambda_i}} \right)^{1/2} \Lambda_t^{-1/4} U_t^\top, \quad \text{Tr}(M_t^{*\top} M_t^*)^{-1} = \frac{\left(\sum_{i=1}^d \sqrt{\lambda_i} \right)^2}{\gamma}, \quad (14)$$

where $\Sigma_t = U_t \Lambda_t U_t^\top$ is the eigendecomposition of the covariance matrix, with $\Lambda_t = \text{diag}(\lambda_1, \dots, \lambda_d)$ containing its eigenvalues.

Remark 2. Applying the optimal transformation M_t^* to the gradient g_t at iteration t , the covariance of the transformed gradient $\tilde{\omega}_t = M_t^*(g_t - \mathbb{E}[g_t | \theta^t])$, conditioned on the history θ^t , is

$$\text{Cov}(\tilde{\omega}_t | \theta^t) = M_t^* \Sigma_t M_t^{*\top} = \frac{\gamma \Lambda_t^{1/2}}{\sum_{i=1}^d \sqrt{\lambda_i}} = \gamma \text{diag} \left(\frac{\sqrt{\lambda_1}}{\sum_{i=1}^d \sqrt{\lambda_i}}, \dots, \frac{\sqrt{\lambda_d}}{\sum_{i=1}^d \sqrt{\lambda_i}} \right).$$

Thus, the optimal transformation decorrelates and scales down the gradients while preserving the relative ordering of variance across directions, in contrast to traditional whitening, which eliminates all variance structure. Note that setting $a_t = \mathbb{E}[g_t | \theta^t]$ and choosing $M_t = \sqrt{\frac{\gamma}{d}} \Lambda_t^{-1/2} U_t^\top$

would amount to a whitening of the gradients and ensures that the constraint in (13) is active, i.e., $\text{Tr}(M_t^\top M_t \Sigma_t) = \gamma$. Under this choice, the objective becomes:

$$\text{Tr}(M_t^\top M_t)^{-1} = \frac{d}{\gamma} \text{Tr}(\Lambda_t) = \frac{d}{\gamma} \sum_{i=1}^d \lambda_i. \quad (15)$$

Comparing our objective in (14) with the objective resulting from the whitening transformation, we observe that applying the Cauchy–Schwarz inequality yields:

$$\frac{d}{\gamma} \sum_{i=1}^d \lambda_i \geq \frac{1}{\gamma} \left(\sum_{i=1}^d \sqrt{\lambda_i} \right)^2, \quad (16)$$

with equality if and only if $\lambda_1 = \dots = \lambda_d$. This shows that our solution achieves a strictly smaller objective than whitening in all non-isotropic cases, where the gradient distribution exhibits unequal variance along the eigenbasis directions.

The proof of Theorem 2 is in the supplement.

3 Algorithm Overview

Algorithm 1 outlines our proposed GeoClip method. We explain its key steps below.

Moving average for mean and covariance estimation. While the theoretical results from Section 2 assume access to the true gradient distribution, in practice this distribution is unknown, and so it must be estimated using only privatized gradients. We estimate the mean and covariance using exponential moving averages computed from those privatized gradients. This enables us to estimate the geometry of the gradients without consuming additional privacy budget. Specifically, we maintain estimates a_t of the mean, and Σ_t of the covariance matrix, which are updated according to:

$$a_{t+1} \leftarrow \beta_1 a_t + (1 - \beta_1) \tilde{g}_t \quad (17)$$

$$\Sigma_{t+1} \leftarrow \beta_2 \Sigma_t + (1 - \beta_2)(\tilde{g}_t - a_t)(\tilde{g}_t - a_t)^\top \quad (18)$$

where β_1 and β_2 are constants close to 1 (e.g., $\beta_1 = 0.99$, $\beta_2 = 0.999$). The eigenvalues and eigenvectors used for the transformation are then computed from the estimated covariance.

Clamping eigenvalues. The covariance matrix is positive semi-definite and may contain zero eigenvalues, which can cause numerical instability. To address this, we clamp eigenvalues from below at a small threshold h_1 (e.g., 10^{-15}). Since we only observe privatized gradients, which may be noisy and unstable, we also clamp from above at h_2 to prevent extreme scaling.

Covariance update with mini-batch. Let B denote a mini-batch of training examples sampled at each iteration, with $|B|$ indicating the batch size. When using mini-batch gradient descent with $|B| > 1$, we must estimate the mean and covariance from the privatized batch averages of the per-sample gradients. Let g_i be the random variable representing the gradient of the i -th sample in the batch. Let $\bar{g} = \frac{1}{|B|} \sum_{i \in B} g_i$ be the batch average gradient. Assuming that the g_i are i.i.d., with the same distribution as g , the covariance of the average gradient \bar{g} satisfies

$$\text{Cov}(\bar{g}) = \text{Cov}\left(\frac{1}{|B|} \sum_{i \in B} g_i\right) = \frac{1}{|B|} \text{Cov}(g). \quad (19)$$

The same principle applies when we observe only the privatized average gradient \tilde{g} . To account for this averaging effect, the covariance update is scaled by the batch size; i.e., line 11 of GeoClip in Algorithm 1 becomes

$$\Sigma_{t+1} \leftarrow \beta_2 \Sigma_t + |B|(1 - \beta_2)(\tilde{g}_t - a_t)(\tilde{g}_t - a_t)^\top. \quad (20)$$

Low-Rank PCA. When the dimensionality is high, computing and storing the full gradient covariance matrix becomes impractical. To address this, we propose a method to maintain a low-rank approximation using a simple and efficient procedure we refer to as *Streaming Rank- k PCA* (Algorithm 2). Specifically, we maintain an approximate eigendecomposition of the covariance in the form $U \Lambda U^\top$,

Algorithm 1 GeoClip

Require: Dataset \mathcal{D} , model f_θ , loss \mathcal{L} , learning rate η , noise scale σ , steps T , hyperparameters $h_1, h_2, \beta_1, \beta_2$

- 1: Initialize θ , mean vector $a_0 = 0$, covariance $\Sigma_0 = I_d$, transform $M_0 = M_0^{\text{inv}} = I_d$
- 2: **for** $t = 0$ to T **do**
- 3: Sample a data point (x_t, y_t)
- 4: Compute gradient $g_t \leftarrow \nabla_\theta \mathcal{L}(f_\theta(x_t), y_t)$
- 5: Center and transform: $\omega_t \leftarrow M_t(g_t - a_t)$
- 6: Clip: $\bar{\omega}_t \leftarrow \omega_t / \max(1, \|\omega_t\|_2)$
- 7: Add noise: $\tilde{\omega}_t \leftarrow \bar{\omega}_t + N$, where $N \sim \mathcal{N}(0, \sigma^2 I_d)$
- 8: Map back: $\tilde{g}_t \leftarrow M_t^{\text{inv}} \tilde{\omega}_t + a_t$
- 9: Update model: $\theta_{t+1} \leftarrow \theta_t - \eta \tilde{g}_t$
- 10: Update mean: $a_{t+1} \leftarrow \beta_1 a_t + (1 - \beta_1) \tilde{g}_t$
- 11: Update covariance: $\Sigma_{t+1} \leftarrow \beta_2 \Sigma_t + (1 - \beta_2)(\tilde{g}_t - a_t)(\tilde{g}_t - a_t)^\top$
- 12: Eigendecompose: $\Sigma_{t+1} = U_t \Lambda_t U_t^\top$
- 13: Clamp eigenvalues: $\lambda_i \leftarrow \text{Clamp}(\lambda_i, \min = h_1, \max = h_2)$
- 14: Set $M_{t+1} \leftarrow (\gamma / \sum_i \sqrt{\lambda_i})^{1/2} \Lambda_t^{-1/4} U_t^\top$
- 15: Set $M_{t+1}^{\text{inv}} \leftarrow (\gamma / \sum_i \sqrt{\lambda_i})^{-1/2} U_t \Lambda_t^{1/4}$
- 16: **end for**
- 17: **return** Final parameters θ

Algorithm 2 STREAMING RANK- k PCA

Require: Eigenvectors $V \in \mathbb{R}^{d \times k}$, eigenvalues $\Lambda \in \mathbb{R}^{k \times k}$, gradient \tilde{g} , mean a , factor β_3 , rank k

- 1: Center: $z \leftarrow \tilde{g} - a$
- 2: Form augmented matrix: $U_{\text{aug}} \leftarrow [V \ z]$
- 3: Compute: $Z \leftarrow U_{\text{aug}} \cdot \text{diag}(\sqrt{\beta_3 \lambda_1}, \dots, \sqrt{\beta_3 \lambda_k}, \sqrt{1 - \beta_3})$
- 4: Perform SVD: $Z = USR^\top$
- 5: Set $V_{\text{new}} \leftarrow$ first k columns of U
- 6: Set $\Lambda_{\text{new}} \leftarrow$ squares of the first k singular values in S
- 7: Return: $V_{\text{new}}, \Lambda_{\text{new}}$

where $U \in \mathbb{R}^{d \times k}$ contains the top- k eigenvectors and $\Lambda \in \mathbb{R}^{k \times k}$ the corresponding eigenvalues, where $k \ll d$. Upon receiving a new gradient \tilde{g} , we center it using the running mean a , yielding $z = \tilde{g} - a$, and perform a weighted update to the covariance:

$$\Sigma_{\text{new}} = \beta_3 U \Lambda U^\top + (1 - \beta_3) z z^\top = [U \ z] \begin{bmatrix} \beta_3 \Lambda & 0 \\ 0 & 1 - \beta_3 \end{bmatrix} [U \ z]^\top \quad (21)$$

Rather than forming this full matrix, we compute its square root:

$$Z = [U \ z] \cdot \text{diag}(\sqrt{\beta_3 \lambda_1}, \dots, \sqrt{\beta_3 \lambda_k}, \sqrt{1 - \beta_3}) \in \mathbb{R}^{d \times (k+1)} \quad (22)$$

We then perform an SVD on Z and retain the top k singular vectors and squared singular values as the updated eigenvectors and eigenvalues. We note that this computation takes $\mathcal{O}(dk^2 + k^3)$ time [Li et al. \[2019\]](#), highlighting that it is only linear in d . The rest of the procedure follows Algorithm 1 by replacing the moving average over the full covariance matrix in line 11 of Algorithm 1 with the low-rank approximation described above. Another necessary change to the algorithm is the following: since U_t is no longer a square matrix, the transformation M_t takes the gradient into the lower k -dimensional space to clip and add noise, such that the resulting M_t^{inv} returns to the full d -dimensional space. The complete version of this variant is provided as Algorithm 3 in the supplementary material.

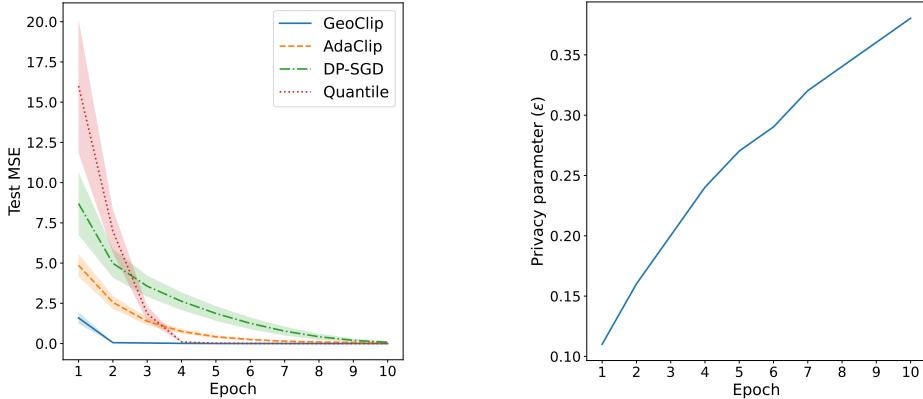


Figure 1: GeoClip results for the synthetic Gaussian dataset with 10 features. The left plot shows the average test MSE for each method over 10 epochs, with shaded regions representing the standard deviation across 20 random seeds. We observe that GeoClip achieves the fastest convergence and lowest average test MSE. The right plot shows the overall privacy budget ε expended for $\delta = 10^{-5}$. This plot applies to all four algorithms, as they are tuned to achieve the same privacy level for a given number of epochs.

4 Experimental Results

We present empirical results demonstrating the benefits of our proposed GeoClip framework compared to AdaClip, quantile-based clipping, and standard DP-SGD. For quantile-based clipping, we use the median of per-sample norms, which has been shown to perform well across various learning tasks [Andrew et al. \[2021\]](#). All (ε, δ) -DP guarantees are computed using the Connect-the-Dots accountant [Doroshenko et al. \[2022\]](#). All experiments were conducted on Google Colab using CPU resources.

We start with a synthetic dataset to demonstrate how GeoClip accelerates convergence, reflecting its design motivation from Theorem 1, and then evaluate its performance on real-world datasets. All datasets are split into 80-10-10 train-validation-test sets for consistent evaluation. The results in Sections 4.1, 4.2, and 4.3 are derived from Algorithm 1. We present the results using the low-rank PCA variant in Section 4.4.

Our results indicate that our framework in Algorithm 1 is robust to the choice of hyperparameters β_1 , β_2 , and h_1 . Standard values commonly used in optimization, such as $\beta_1 = 0.99$ and $\beta_2 = 0.999$, work well in our setting. The parameter h_1 only needs to be a small positive constant (e.g., 10^{-15}) to ensure numerical stability. Since the eigenvalues are clamped to the range $[h_1, h_2]$ (line 13 in Algorithm 1), the trace term $\sum_i \sqrt{\lambda_i}$ is bounded between $d\sqrt{h_1}$ and $d\sqrt{h_2}$, where d is the dimensionality. This allows us to set the parameter γ to 1, as its effect can be absorbed by tuning h_2 . For h_2 , we have observed that the values 1 and 10 perform consistently well across datasets. Throughout all experiments, we tune only h_2 to select between these two options. A similar setup is used in Algorithm 2 with β_2 replaced by β_3 (set to 0.99), which is also robust across experiments.

4.1 Synthetic Dataset

GeoClip is designed to improve convergence, particularly in the presence of feature correlation. To empirically demonstrate this, we evaluate it on a synthetic Gaussian dataset with 20,000 samples and 10 features—five of which are correlated, while the remaining five are independent. We train a linear regression model using various private training methods for 10 epochs with a batch size of 1024, tuning the learning rate for each method to ensure stable convergence. As shown in Figure 1 (left), GeoClip converges as early as epoch 2, while the next best method—quantile-based clipping—requires nearly twice as many epochs. Figure 1 (right) plots the privacy cost (ε) versus epoch, showing how faster convergence helps minimize overall privacy cost.

As illustrated by the standard deviation bands in Figure 1, GeoClip not only achieves faster convergence but also demonstrates more stable training with reduced variance—an important property when

privacy constraints limit training to a single run. In such scenarios, lower variability across runs enhances the reliability of the final model without requiring additional privacy budget.

4.2 Tabular Datasets

In addition to the synthetic dataset, we run experiments to compare the performance of different clipping strategies on three real-world datasets: Diabetes Efron et al. [2004], Breast Cancer Wolberg et al. [1993], and Android Malware Borah and Bhattacharyya [2020]. We briefly describe each dataset and its corresponding learning task below.

The Diabetes dataset contains 442 samples and 10 standardized features. The continuous valued target variable indicates disease progression, making this a regression task. The Breast Cancer dataset contains 569 samples and 30 numerical features, which are standardized to zero mean and unit variance before training. The target is to predict a binary label indicating the presence or absence of cancer. The Android Malware dataset contains 4465 samples and 241 integer attributes. The target is to classify whether a program is malware or not.

For each of the three datasets, we perform a grid search over the relevant hyperparameters to identify the best model under a given (ε, δ) privacy budget. We train for 5 epochs using 20 random seeds and report the average performance along with the standard deviation in Tables 1, 2, and 3. We observe that our proposed GeoClip framework consistently outperforms all baseline methods across both regression and classification tasks. GeoClip achieves better performance with noticeably smaller standard deviations, indicating greater stability across random seeds.

Table 1: Diabetes dataset test MSE comparison for $\delta = 10^{-5}$, batch size = 32, and model dimension $d = 11$.

Framework	$\varepsilon = 0.50$	$\varepsilon = 0.86$	$\varepsilon = 0.93$
GeoClip (ours)	0.073± 0.015	0.044± 0.003	0.039± 0.009
AdaClip	0.077±0.027	0.062±0.028	0.055±0.014
Quantile	0.090±0.027	0.083±0.044	0.072±0.014
DP-SGD	0.108±0.040	0.095±0.047	0.072±0.040

Table 2: Breast Cancer dataset test accuracy (%) comparison for $\delta = 10^{-5}$, batch size = 64, and model dimension $d = 62$.

Framework	$\varepsilon = 0.67$	$\varepsilon = 0.8$	$\varepsilon = 0.87$
GeoClip (ours)	87.87±3.32	88.57±3.37	93.63±1.63
AdaClip	84.90±5.91	85.42±5.34	87.71±5.87
Quantile	81.41±12.53	81.63±10.71	92.28±2.68
DP-SGD	77.32±6.17	79.42±9.71	85.95±4.45

Table 3: Malware dataset test accuracy (%) comparison for $\delta = 10^{-5}$, batch size = 512, and model dimension $d = 484$.

Framework	$\varepsilon = 0.26$	$\varepsilon = 0.49$	$\varepsilon = 0.67$
GeoClip (ours)	90.77±1.83	91.64±1.26	92.67±1.63
AdaClip	88.35±3.27	90.25±1.33	90.23±3.11
Quantile	77.84±1.29	78.84±1.27	81.86±1.31
DP-SGD	88.04±2.21	90.55±1.55	90.57±1.61

4.3 Final Layer Fine-Tuning

In many transfer learning scenarios, fine-tuning only the final layer is standard practice due to both its computational efficiency and minimal privacy cost. Last-layer fine-tuning is a well-suited application

of GeoClip, as it involves a small number of trainable parameters, making covariance estimation more tractable. This setup also benefits from GeoClip’s faster convergence, which is particularly valuable in privacy-constrained settings where only limited training iterations are feasible.

To demonstrate this, we design an experiment where a convolutional neural network (CNN) is first trained on MNIST LeCun et al. [1998] using the Adam optimizer and then transferred to Fashion-MNIST Xiao et al. [2017] by freezing all layers except the final fully connected layer. The CNN consists of two convolutional and pooling layers followed by a linear compression layer that reduces the feature size to 50, resulting in a total of only 510 trainable parameters. We fine-tune this layer using different methods under varying privacy budgets, and present the results in Table 4.

Table 4: Final layer DP fine-tuning on Fashion-MNIST for 4 epochs and $\delta = 10^{-6}$ over 5 seeds.

Framework	$\varepsilon = 0.6$	$\varepsilon = 1$
GeoClip (Ours)	73.09±0.72	73.09±0.63
AdaClip	68.35±0.41	69.24±0.28
Quantile-based	71.78±1.28	72.09±1.12
DP-SGD	69.40±0.82	69.83±0.83

4.4 Low-Rank PCA Results

To evaluate our method on low-rank PCA algorithm (Algorithm 2), we construct a synthetic binary classification dataset with 20,000 samples and 400 Gaussian features, where 50 are correlated and 350 are uncorrelated. Labels are generated by applying a linear function to the features, adding Gaussian noise, and thresholding the sigmoid output. We train a logistic regression model with 802 trainable parameters. We then apply our method using a low-rank PCA approximation with rank 50. As shown in the left panel of Figure 2, our approach converges faster than competing methods even with this low-rank approximation.

We also evaluate our method on the USPS dataset Hull [1994] using logistic regression with 2,570 trainable parameters (256 input features \times 10 classes + 10 biases). The USPS dataset contains 9,298 grayscale handwritten digit images (0–9), each of size 16×16 pixels. This compact benchmark is commonly used for evaluating digit classification models. For this dataset, we apply a low-rank PCA approximation with rank 100. Results are shown in the right panel of Figure 2. As with the synthetic dataset, GeoClip with low-rank PCA also achieves faster convergence on USPS compared to baseline methods. We provide additional ε -vs-iteration plots for both datasets, illustrating how faster convergence reduces overall privacy cost in the supplementary material.

5 Conclusions

We have introduced GeoClip, a geometry-aware framework for differentially private SGD that leverages the structure of the gradient distribution to improve both utility and convergence. By operating in a basis adapted to the estimated noisy gradients, GeoClip injects noise more strategically, thereby reducing distortion without incurring additional privacy cost. We have provided a formal analysis of convergence guarantees which characterizes the optimal transformation. Our empirical results on synthetic and real-world datasets show that GeoClip consistently converges faster and outperforms existing adaptive clipping methods, improving both the mean and standard deviation of the performance metrics over multiple runs. Via low-rank approximation method, we have shown that GeoClip scales to the high-dimensional data setting, thus making it suitable for practical deployment in large, privacy-sensitive models.

Limitations. The linear transformation involved in GeoClip requires an additional computation via an eigendecomposition. Our low-rank approximation addresses that to some extent. The algorithm introduces additional hyperparameters (most notably h_2) compared to standard DP-SGD, which must be tuned for optimal performance. Our experiments have been performed on a limited collection of datasets; additional testing is needed to see how our algorithm performs in more generality.

Broader impact. As discussed in the Introduction, our work is motivated by societal concerns, with a focus on improving the theoretical limits of differentially private optimization.

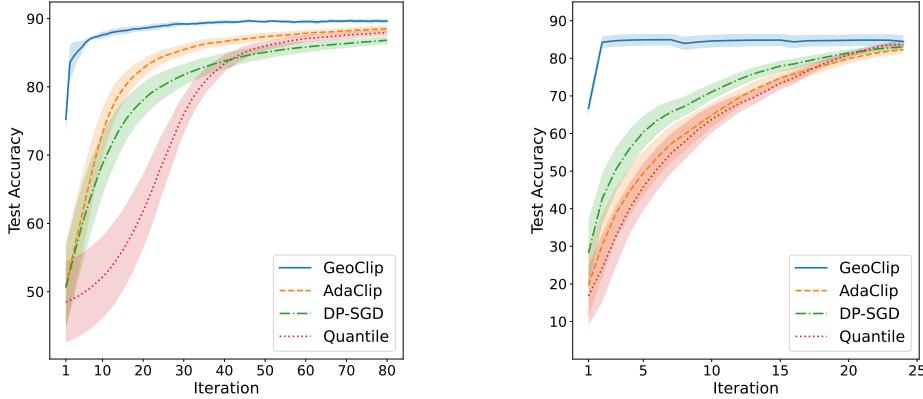


Figure 2: The left panel shows results on the synthetic Gaussian dataset with 400 features using a rank-50 PCA approximation for GeoClip. The plot displays average test accuracy (%) over 80 iterations with a batch size of 1024. GeoClip achieves the fastest convergence and highest average accuracy. The right panel shows results on the USPS dataset using a rank-100 approximation over 25 iterations with a batch size of 1024, where a similar convergence trend is observed. Shaded regions represent standard deviation across 20 random seeds.

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A Proof of Theorem 1

Theorem (Convergence of GeoClip). *Assume f has an L -Lipschitz continuous gradient. Further, assume the stochastic gradients are bounded, i.e., $\|\nabla f_k(\theta)\| \leq G$, and have bounded variance, i.e., $\mathbb{E}_k \|\nabla f_k(\theta) - \nabla f(\theta)\|^2 \leq \sigma_g^2$. Let $\theta^* = \arg \min_{\theta \in \mathbb{R}^d} f(\theta)$ denote the optimal solution, and suppose the learning rate satisfies $\eta < \frac{2}{3L}$. Then, for the iterates $\{\theta_t\}_{t=0}^{T-1}$ produced by GeoClip with batch size 1 using the update rule $\theta_{t+1} = \theta_t - \eta \tilde{g}_t$, where \tilde{g}_t is defined in (5), the average squared gradient norm satisfies:*

$$\begin{aligned} \frac{1}{T} \sum_{t=0}^{T-1} \mathbb{E} \|\nabla f(\theta_t)\|^2 &\leq \underbrace{\frac{f(\theta_0) - f(\theta^*)}{T \left(\eta - \frac{3L\eta^2}{2} \right)}}_{\text{Optimization gap}} + \underbrace{\frac{3L\eta}{2-3L\eta} \sigma_g^2}_{\text{Gradient variance term}} + \underbrace{\frac{L\eta\sigma^2}{T(2-3L\eta)} \sum_{t=0}^{T-1} \mathbb{E} \text{Tr} \left[(M_t^\top M_t)^{-1} \right]}_{\text{Noise-injection term}} \\ &\quad + \underbrace{\frac{2}{T(2-3L\eta)} \sum_{t=0}^{T-1} \mathbb{E} \left[\beta(a_t) \left(\text{Tr} (M_t^\top M_t \Sigma_t) + \|M_t(\mathbb{E}[g_t | \theta^t] - a_t)\|^2 \right) \right]}_{\text{Clipping error term}}, \end{aligned} \quad (23)$$

where $\theta^t = (\theta_0, \dots, \theta_t)$ represents the history of parameter values up to iteration t , $\Sigma_t = \text{Cov}(g_t | \theta^t)$, and

$$\beta(a_t) = (G + \|a_t\|) \left(G + \frac{3L\eta}{2} (G + \|a_t\|) \right). \quad (24)$$

Proof. We can express the noisy gradient as:

$$\tilde{g}_t = M_t^{-1} \tilde{\omega}_t + a_t \quad (25)$$

$$= \frac{M_t^{-1} \omega_t}{\max\{1, \|\omega_t\|\}} + M_t^{-1} N_t + a_t \quad (26)$$

$$= \frac{g_t - a_t}{\max\{\|M_t(g_t - a_t)\|, 1\}} + M_t^{-1} N_t + a_t. \quad (27)$$

Thus, the parameter update takes the form

$$\theta_{t+1} = \theta_t - \eta \tilde{g}_t \quad (28)$$

$$= \theta_t - \eta \left[\frac{g_t - a_t}{\max\{\|M_t(g_t - a_t)\|, 1\}} + M_t^{-1} N_t + a_t \right]. \quad (29)$$

We define the following quantities:

$$c_t = \frac{g_t - a_t}{\max\{\|M_t(g_t - a_t)\|, 1\}}, \quad (30)$$

$$\Delta_t = c_t - (g_t - a_t). \quad (31)$$

c_t is the clipped version of $g_t - a_t$. Δ_t quantifies the distortion due to clipping, being zero when no clipping occurs and negative when clipping is applied. By the L -Lipschitz continuity of the gradient of $f(\theta)$, we have:

$$f(\theta_{t+1}) \leq f(\theta_t) + \langle \nabla f(\theta_t), \theta_{t+1} - \theta_t \rangle + \frac{L}{2} \|\theta_{t+1} - \theta_t\|^2. \quad (32)$$

Let $\theta^t = (\theta_0, \dots, \theta_t)$ describe the entire history of states. Taking the expectation conditioned on θ^t , we obtain:

$$\begin{aligned} & \mathbb{E}[f(\theta_{t+1})|\theta^t] \\ & \leq f(\theta_t) + \mathbb{E}[\langle \nabla f(\theta_t), \theta_{t+1} - \theta_t \rangle | \theta^t] + \frac{L}{2} \mathbb{E}[\|\theta_{t+1} - \theta_t\|^2 | \theta^t] \end{aligned} \quad (33)$$

$$= f(\theta_t) - \eta \mathbb{E}[\langle \nabla f(\theta_t), c_t + M_t^{-1}N_t + a_t \rangle | \theta^t] + \frac{L\eta^2}{2} \mathbb{E}[\|c_t + M_t^{-1}N_t + a_t\|^2 | \theta^t] \quad (34)$$

$$= f(\theta_t) - \eta \mathbb{E}[\langle \nabla f(\theta_t), c_t + a_t \rangle | \theta^t] + \frac{L\eta^2}{2} \mathbb{E}[\|c_t + a_t\|^2 | \theta^t] + \frac{L\eta^2}{2} \mathbb{E}[\|M_t^{-1}N_t\|^2 | \theta^t] \quad (35)$$

$$= f(\theta_t) - \eta \mathbb{E}[\langle \nabla f(\theta_t), c_t + a_t \rangle | \theta^t] + \frac{L\eta^2}{2} \mathbb{E}[\|c_t + a_t\|^2 | \theta^t] + \frac{L\eta^2\sigma^2}{2} \|M_t^{-1}\|_F^2 \quad (36)$$

$$= f(\theta_t) - \eta \mathbb{E}[\langle \nabla f(\theta_t), g_t + \Delta_t \rangle | \theta^t] + \frac{L\eta^2}{2} \mathbb{E}[\|g_t + \Delta_t\|^2 | \theta^t] + \frac{L\eta^2\sigma^2}{2} \|M_t^{-1}\|_F^2 \quad (37)$$

$$\begin{aligned} & = f(\theta_t) - \eta \|\nabla f(\theta_t)\|^2 - \eta \mathbb{E}[\langle \nabla f(\theta_t), \Delta_t \rangle | \theta^t] \\ & \quad + \frac{L\eta^2}{2} \mathbb{E}[\|g_t - \nabla f(\theta_t) + \nabla f(\theta_t) + \Delta_t\|^2 | \theta^t] + \frac{L\eta^2\sigma^2}{2} \|M_t^{-1}\|_F^2 \end{aligned} \quad (38)$$

where (35) follows from $\mathbb{E}[N_t] = 0$ and the independence of N_t from g_t (and thus from $c_t + a_t$). The equality (36) follows because $\mathbb{E}[NN^\top] = \sigma^2 I_d$, and $\|M_t^{-1}\|_F$ represents the Frobenius norm of M_t^{-1} . The equality (37) follows from $c_t + a_t = \Delta_t + g_t$. The last equality follows because

$$\mathbb{E}[\langle \nabla f(\theta_t), g_t + \Delta_t \rangle | \theta^t] = \mathbb{E}[\langle \nabla f(\theta_t), \Delta_t \rangle | \theta^t] + \mathbb{E}[\langle \nabla f(\theta_t), g_t \rangle | \theta^t] \quad (39)$$

$$= \mathbb{E}[\langle \nabla f(\theta_t), \Delta_t \rangle | \theta^t] + \nabla f(\theta_t)^\top \mathbb{E}[g_t | \theta^t] \quad (40)$$

$$= \mathbb{E}[\langle \nabla f(\theta_t), \Delta_t \rangle | \theta^t] + \|\nabla f(\theta_t)\|^2, \quad (41)$$

where (41) follows because g_t is an unbiased estimator of the true gradient, i.e., $\mathbb{E}[g_t | \theta^t] = \nabla f(\theta_t)$. From Jensen's inequality, we have:

$$\begin{aligned} & \mathbb{E}[\|g_t - \nabla f(\theta_t) + \nabla f(\theta_t) + \Delta_t\|^2 | \theta^t] \\ & \leq 3 \left(\mathbb{E}[\|g_t - \nabla f(\theta_t)\|^2 | \theta^t] + \|\nabla f(\theta_t)\|^2 + \mathbb{E}[\|\Delta_t\|^2 | \theta^t] \right). \end{aligned} \quad (42)$$

From the Cauchy–Schwarz inequality, we have:

$$\mathbb{E}[\langle \nabla f(\theta_t), \Delta_t \rangle | \theta^t] \leq \|\nabla f(\theta_t)\| \mathbb{E}[\|\Delta_t\| | \theta^t]. \quad (43)$$

Plugging (42) and (43) into (38), we obtain:

$$\begin{aligned} \mathbb{E}[f(\theta_{t+1}) | \theta^t] & \leq f(\theta_t) - \eta \|\nabla f(\theta_t)\|^2 + \eta \|\nabla f(\theta_t)\| \mathbb{E}[\|\Delta_t\| | \theta^t] \\ & \quad + \frac{3L\eta^2}{2} \left[\mathbb{E}[\|g_t - \nabla f(\theta_t)\|^2 | \theta^t] + \|\nabla f(\theta_t)\|^2 + \mathbb{E}[\|\Delta_t\|^2 | \theta^t] \right] + \frac{L\eta^2\sigma^2}{2} \|M_t^{-1}\|_F^2. \end{aligned} \quad (44)$$

To bound $\mathbb{E}[\|\Delta_t\| | \theta^t]$ and $\mathbb{E}[\|\Delta_t\|^2 | \theta^t]$, we first bound $\Pr(\|\Delta_t\| > 0 | \theta^t)$ and $\|\Delta_t\|$ given θ^t and $\|\Delta_t\| > 0$. Using Markov's inequality, we obtain:

$$\Pr(\|\Delta_t\| > 0 | \theta^t) = \Pr(\|M_t(g_t - a_t)\|^2 > 1 | \theta^t) \quad (45)$$

$$\leq \Pr(\|M_t(g_t - a_t)\|^2 \geq 1 | \theta^t) \quad (46)$$

$$\leq \mathbb{E}[\|M_t(g_t - a_t)\|^2 | \theta^t] \quad (47)$$

$$= \mathbb{E}[\|M_t(g_t - \mathbb{E}[g_t | \theta^t] + \mathbb{E}[g_t | \theta^t] - a_t)\|^2 | \theta^t] \quad (48)$$

$$= \mathbb{E}[\|M_t(g_t - \mathbb{E}[g_t | \theta^t])\|^2 | \theta^t] + \|M_t(\mathbb{E}[g_t | \theta^t] - a_t)\|^2 \quad (49)$$

$$= \text{Tr}(\text{Cov}(M_t g_t | \theta^t)) + \|M_t(\mathbb{E}[g_t | \theta^t] - a_t)\|^2. \quad (50)$$

Using (30) and (31), we obtain:

$$\|\Delta_t\| = \left\| \left(\frac{1}{\max\{\|M_t(g_t - a_t)\|, 1\}} - 1 \right) (g_t - a_t) \right\| \quad (51)$$

$$\leq \|g_t - a_t\| \quad (52)$$

$$\leq \|g_t\| + \|a_t\| \quad (53)$$

$$\leq G + \|a_t\|, \quad (54)$$

where the final inequality follows from $\|\nabla f_k(\theta)\| \leq G$. Therefore,

$$\mathbb{E} [\|\Delta_t\| \mid \theta^t] = \Pr (\|\Delta_t\| > 0 \mid \theta^t) \mathbb{E} [\|\Delta_t\| \mid \|\Delta_t\| > 0, \theta^t] \quad (55)$$

$$\leq (G + \|a_t\|) \left(\text{Tr} (\text{Cov}(M_t g_t \mid \theta^t)) + \|M_t(\mathbb{E}[g_t \mid \theta^t] - a_t)\|^2 \right). \quad (56)$$

Similarly, we obtain the following bound:

$$\mathbb{E} [\|\Delta_t\|^2 \mid \theta^t] \leq (G + \|a_t\|)^2 \left(\text{Tr} (\text{Cov}(M_t g_t \mid \theta^t)) + \|M_t(\mathbb{E}[g_t \mid \theta^t] - a_t)\|^2 \right). \quad (57)$$

By plugging (56) and (57) into (44) and rearranging the terms, we obtain

$$\begin{aligned} & \mathbb{E} [f(\theta_{t+1}) \mid \theta^t] \\ & \leq f(\theta_t) + \left(\frac{3L\eta^2}{2} - \eta \right) \|\nabla f(\theta_t)\|^2 + \eta(G + \|a_t\|) \left(G + \frac{3L\eta}{2}(G + \|a_t\|) \right) \\ & \quad \left[\text{Tr} (\text{Cov}(M_t g_t \mid \theta^t)) + \|M_t(\mathbb{E}[g_t \mid \theta^t] - a_t)\|^2 \right] + \frac{3L\eta^2}{2} \mathbb{E} [\|g_t - \nabla f(\theta_t)\|^2 \mid \theta^t] + \frac{L\eta^2\sigma^2}{2} \|M_t^{-1}\|_F^2. \end{aligned} \quad (58)$$

Above, we have used the fact that $\|\nabla f(\theta_t)\| \leq G$, which itself follows from Jensen's inequality and the assumption that $\|\nabla f_k(\theta)\| \leq G$. Let

$$\beta(a_t) = (G + \|a_t\|) \left(G + \frac{3L\eta}{2}(G + \|a_t\|) \right). \quad (59)$$

Rearranging the terms, applying the law of total expectation (now taking expectation over the entire history of states θ^t), and using the bound $\mathbb{E}_k \|\nabla f_k(\theta) - \nabla f(\theta)\|^2 \leq \sigma_g^2$, we obtain:

$$\begin{aligned} \left(\eta - \frac{3L\eta^2}{2} \right) \mathbb{E} \|\nabla f(\theta_t)\|^2 & \leq \mathbb{E} f(\theta_t) - \mathbb{E} f(\theta_{t+1}) + \frac{3L\eta^2\sigma_g^2}{2} + \frac{L\eta^2\sigma^2}{2} \mathbb{E} \|M_t^{-1}\|_F^2 \\ & \quad + \eta \mathbb{E} [\beta(a_t) (\text{Tr} (\text{Cov}(M_t g_t \mid \theta^t)) + \|M_t(\mathbb{E}[g_t \mid \theta^t] - a_t)\|^2)]. \end{aligned} \quad (60)$$

Summing over $t = 0$ to $T - 1$ and applying the telescoping sum, we get:

$$\begin{aligned} \left(\eta - \frac{3L\eta^2}{2} \right) \sum_{t=0}^{T-1} \mathbb{E} \|\nabla f(\theta_t)\|^2 & \leq f(\theta_0) - \mathbb{E} f(\theta_T) + \frac{3L\eta^2\sigma_g^2}{2} T + \frac{L\eta^2\sigma^2}{2} \sum_{t=0}^{T-1} \mathbb{E} \|M_t^{-1}\|_F^2 \\ & \quad + \eta \sum_{t=0}^{T-1} \mathbb{E} [\beta(a_t) (\text{Tr} (\text{Cov}(M_t g_t \mid \theta^t)) + \|M_t(\mathbb{E}[g_t \mid \theta^t] - a_t)\|^2)] \end{aligned} \quad (61)$$

Dividing by T , assuming $\eta \leq \frac{2}{3L}$, applying the identity $\|M_t^{-1}\|_F^2 = \text{Tr} ((M_t^\top M_t)^{-1})$, and using the fact that $\mathbb{E} f(\theta_T) \geq f(\theta^*)$, we obtain:

$$\begin{aligned} & \frac{1}{T} \sum_{t=0}^{T-1} \mathbb{E} \|\nabla f(\theta_t)\|^2 \\ & \leq \frac{f(\theta_0) - f(\theta^*)}{T \left(\eta - \frac{3L\eta^2}{2} \right)} + \frac{3L\eta}{2 - 3L\eta} \sigma_g^2 + \frac{L\eta\sigma^2}{T(2 - 3L\eta)} \sum_{t=0}^{T-1} \mathbb{E} [\text{Tr} ((M_t^\top M_t)^{-1})] \\ & \quad + \frac{2}{T(2 - 3L\eta)} \sum_{t=0}^{T-1} \mathbb{E} [\beta(a_t) (\text{Tr} (M_t^\top M_t \text{Cov}(g_t \mid \theta^t)) + \|M_t(\mathbb{E}[g_t \mid \theta^t] - a_t)\|^2)]. \end{aligned} \quad (62)$$

□

B Proof of Theorem 2

We would like to solve the following optimization

$$\begin{aligned} & \underset{M_t}{\text{minimize}} \quad \text{Tr}(M_t^\top M_t)^{-1} \\ & \text{subject to} \quad \text{Tr}(M_t^\top M_t \Sigma_t) \leq \gamma. \end{aligned} \quad (63)$$

Defining A_t as the Gram matrix of M_t , i.e., $A_t = M_t^\top M_t$, we can reformulate the optimization problem as:

$$\begin{aligned} & \underset{A_t}{\text{minimize}} \quad \text{Tr}(A_t^{-1}) \\ & \text{subject to} \quad \text{Tr}(A_t \Sigma_t) \leq \gamma. \end{aligned} \quad (64)$$

Both the objective and the constraint are convex in A , so to solve the problem, we introduce the Lagrangian function:

$$\ell(A_t, \mu) = \text{Tr}(A_t^{-1}) + \mu(\text{Tr}(A_t \Sigma_t) - \gamma), \quad (65)$$

Taking the derivative of $\ell(A_t, \mu)$ with respect to A_t and setting the derivative to zero for optimality, we get

$$A_t^{-2} = \mu \Sigma_t. \quad (66)$$

and so

$$A_t = \frac{1}{\sqrt{\mu}} \Sigma_t^{-\frac{1}{2}}. \quad (67)$$

Substituting A_t in the constraint $\text{Tr}(A_t \Sigma_t) = \gamma$, we get

$$\frac{1}{\sqrt{\mu}} \text{Tr}\left(\Sigma_t^{\frac{1}{2}}\right) = \gamma. \quad (68)$$

Solving for μ :

$$\sqrt{\mu} = \frac{1}{\gamma} \text{Tr}\left(\Sigma_t^{\frac{1}{2}}\right). \quad (69)$$

Thus, the optimal A_t is:

$$A_t = \frac{\gamma}{\text{Tr}\left(\Sigma_t^{\frac{1}{2}}\right)} \Sigma_t^{-\frac{1}{2}}. \quad (70)$$

Using the eigen decomposition, we write Σ_t as:

$$\Sigma_t = U_t \Lambda_t U_t^\top, \quad (71)$$

where U_t is an orthogonal matrix whose columns are the eigenvectors of Σ_t , and

$$\Lambda_t = \text{diag}(\lambda_1, \dots, \lambda_d)$$

is a diagonal matrix containing the corresponding eigenvalues. We now have

$$\Sigma_t^{\frac{1}{2}} = U_t \Lambda_t^{\frac{1}{2}} U_t^\top, \quad \Sigma_t^{-\frac{1}{2}} = U_t \Lambda_t^{-\frac{1}{2}} U_t^\top, \quad (72)$$

Thus, the final expression for A_t is:

$$A_t = \frac{\gamma}{\text{Tr}(\Lambda_t^{\frac{1}{2}})} U_t \Lambda_t^{-\frac{1}{2}} U_t^\top. \quad (73)$$

Therefore, we have

$$M_t^\top M_t = \frac{\gamma}{\text{Tr}(\Lambda_t^{\frac{1}{2}})} U_t \Lambda_t^{-\frac{1}{2}} U_t^\top, \quad (74)$$

and

$$M_t = \left(\frac{\gamma}{\text{Tr}(\Lambda_t^{\frac{1}{2}})} \right)^{1/2} \Lambda_t^{-\frac{1}{4}} U_t^\top. \quad (75)$$

Since $\text{Tr}(\Lambda_t^{1/2}) = \sum_{i=1}^d \sqrt{\lambda_i}$, we can simplify M_t as follows:

$$M_t = \left(\frac{\gamma}{\sum_{i=1}^d \sqrt{\lambda_i}} \right)^{1/2} \Lambda_t^{-1/4} U_t^\top. \quad (76)$$

Algorithm 3 GEOCLIP WITH RANK- k PCA

Require: Dataset \mathcal{D} , model f_θ , loss \mathcal{L} , learning rate η , noise scale σ , steps T , rank k , hyperparameters $h_1, h_2, \beta_1, \beta_3$

- 1: Initialize θ , mean vector $a_0 = 0$, $U_0 = [e_1, \dots, e_k]$ where $e_i \in \mathbb{R}^d$ is the i -th standard basis vector, $\Lambda_0 = I_k$
- 2: Compute transform $M_0 \leftarrow (\gamma/k)^{1/2} \Lambda_0^{-1/4} U_0^\top$ and $M_0^{\text{inv}} \leftarrow (\gamma/k)^{-1/2} U_0 \Lambda_0^{1/4}$
- 3: **for** $t = 0$ to T **do**
- 4: Sample a data point (x_t, y_t)
- 5: Compute gradient $g_t \leftarrow \nabla_\theta \mathcal{L}(f_\theta(x_t), y_t)$
- 6: Center and transform: $\omega_t \leftarrow M_t(g_t - a_t)$
- 7: Clip: $\bar{\omega}_t \leftarrow \omega_t / \max(1, \|\omega_t\|_2)$
- 8: Add noise: $\tilde{\omega}_t \leftarrow \bar{\omega}_t + N$, where $N \sim \mathcal{N}(0, \sigma^2 I_k)$
- 9: Map back: $\tilde{g}_t \leftarrow M_t^{\text{inv}} \tilde{\omega}_t + a_t$
- 10: Update model: $\theta_{t+1} \leftarrow \theta_t - \eta \tilde{g}_t$
- 11: Update mean: $a_{t+1} \leftarrow \beta_1 a_t + (1 - \beta_1) \tilde{g}_t$
- 12: Update eigenspace: $(U_{t+1}, \Lambda_{t+1}) \leftarrow \text{STREAMING RANK-}k \text{ PCA}(U_t, \Lambda_t, \tilde{g}_t, a_{t+1}, \beta_3, k)$
- 13: Clamp eigenvalues: $\lambda_i \leftarrow \text{Clamp}(\lambda_i, \min = h_1, \max = h_2)$
- 14: Set $M_{t+1} \leftarrow (\gamma / \sum_i \sqrt{\lambda_i})^{1/2} \Lambda_{t+1}^{-1/4} U_{t+1}^\top$
- 15: Set $M_{t+1}^{\text{inv}} \leftarrow (\gamma / \sum_i \sqrt{\lambda_i})^{-1/2} U_{t+1} \Lambda_{t+1}^{1/4}$
- 16: **end for**
- 17: **return** Final parameters θ

C GeoClip with Low-Rank PCA

Computing and storing the full gradient covariance matrix becomes infeasible in high dimensions. To address this, we propose a low-rank approximation method in Algorithm 3, which incorporates a rank- k PCA step described in the **STREAMING RANK- k PCA** algorithm. This procedure efficiently maintains an approximate eigendecomposition $U_t \Lambda_t U_t^\top$, where $U_t \in \mathbb{R}^{d \times k}$ contains the top- k eigenvectors and $\Lambda_t = \text{diag}(\lambda_1, \dots, \lambda_k) \in \mathbb{R}^{k \times k}$ holds the corresponding eigenvalues at iteration t , with $k \ll d$.

Given a new gradient \tilde{g}_t , we center it using the running mean a_{t+1} to obtain $z = \tilde{g}_t - a_{t+1}$, and update the covariance as:

$$\Sigma_{\text{new}} = \beta_3 U_t \Lambda_t U_t^\top + (1 - \beta_3) z z^\top = [U_t \ z] \begin{bmatrix} \beta_3 \Lambda_t & 0 \\ 0 & 1 - \beta_3 \end{bmatrix} [U_t \ z]^\top. \quad (77)$$

Rather than forming this matrix directly, we compute its square root:

$$Z = [U_t \ z] \cdot \text{diag}(\sqrt{\beta_3 \lambda_1}, \dots, \sqrt{\beta_3 \lambda_k}, \sqrt{1 - \beta_3}), \quad (78)$$

and perform an SVD on Z , retaining the top k components to obtain U_{t+1} and Λ_{t+1} .

Since U_t is not square, $M_t \in \mathbb{R}^{k \times d}$ projects gradients into a k -dimensional subspace for clipping and noise addition, while $M_t^{\text{inv}} \in \mathbb{R}^{d \times k}$ maps them back to \mathbb{R}^d .

To compute M_0 and M_0^{inv} in line 2 of Algorithm 3, we use the simplification $\text{Tr}(\Lambda_0^{1/2}) = \sum_{i=1}^k 1 = k$. Also, in line 13, each λ_i depends on t , but we omit the subscript for notational simplicity.

D Privacy Cost vs. Iteration (Corresponding to Figure 2)

Figure 4 shows the privacy cost (ε) versus iteration curves associated with the accuracy-vs-iteration plots in Figure 2 (Section 4.4, Low-Rank PCA Results), which is repeated in this PDF as Figure 3 for easier reference. These plots highlight how faster convergence reduces overall privacy cost. For example, on the USPS dataset (Figure 3, right), GeoClip achieves high accuracy within the first few iterations, requiring only $\varepsilon \approx 0.15$, whereas quantile-based clipping takes about 24 iterations to reach similar accuracy, incurring a higher privacy cost of $\varepsilon \approx 0.5$.

STREAMING RANK- k PCA

Require: Eigenvectors $U_t \in \mathbb{R}^{d \times k}$, eigenvalues $\Lambda_t \in \mathbb{R}^{k \times k}$, gradient $\tilde{g}_t \in \mathbb{R}^d$, mean $a_{t+1} \in \mathbb{R}^d$, factor $\beta_3 \in \mathbb{R}$, rank k

- 1: Center: $z \leftarrow \tilde{g}_t - a_{t+1}$
- 2: Form augmented matrix: $U_{\text{aug}} \leftarrow [U_t \ z]$
- 3: Compute: $Z \leftarrow U_{\text{aug}} \text{ diag}(\sqrt{\beta_3 \lambda_1}, \dots, \sqrt{\beta_3 \lambda_k}, \sqrt{1 - \beta_3})$
- 4: Perform SVD: $Z = VSR^\top$
- 5: Set $U_{t+1} \leftarrow$ first k columns of V
- 6: Set $\Lambda_{t+1} \leftarrow$ squares of the first k singular values in S
- 7: Return: U_{t+1}, Λ_{t+1}

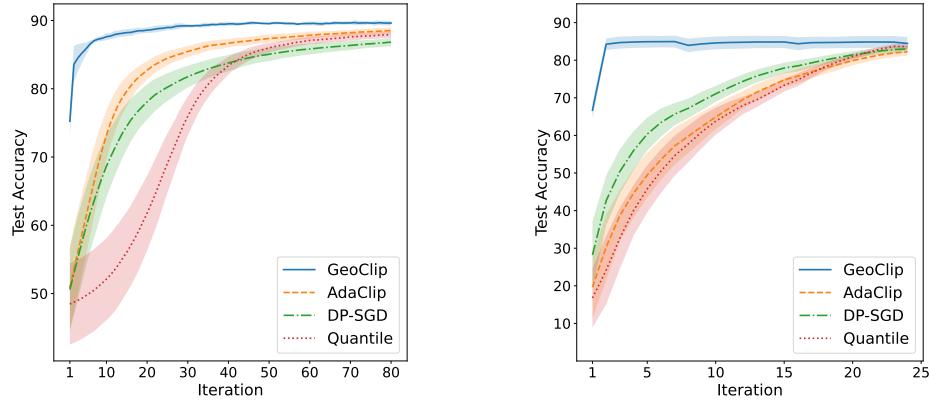


Figure 3: The left panel shows results on the synthetic Gaussian dataset with 400 features using a rank-50 PCA approximation for GeoClip. The plot displays average test accuracy (%) over 80 iterations with a batch size of 1024. GeoClip achieves the fastest convergence and highest average accuracy. The right panel shows results on the USPS dataset using a rank-100 approximation over 25 iterations with a batch size of 1024, where a similar convergence trend is observed. Shaded regions represent standard deviation across 20 random seeds.

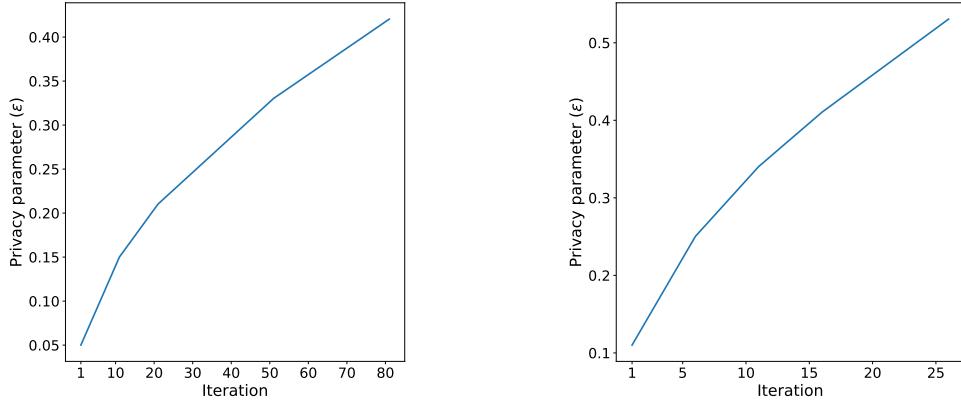


Figure 4: The left plot shows the overall privacy budget ε spent on training the synthetic Gaussian dataset with 400 features for $\delta = 10^{-5}$, while the right plot shows the same for the USPS dataset with $\delta = 10^{-5}$. These plots apply to all four algorithms, which are tuned to achieve the same privacy level for a given number of iterations.