SHARPNESS AWARE MINIMIZATION: GENERAL ANALYSIS AND IMPROVED RATES

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ABSTRACT

Sharpness-Aware Minimization (SAM) has emerged as a powerful method for improving generalization in machine learning models by minimizing the sharpness of the loss landscape. However, despite its success, several important questions regarding the convergence properties of SAM in non-convex settings are still open, including the benefits of using normalization in the update rule, the dependence of the analysis on the restrictive bounded variance assumption, and the convergence guarantees under different sampling strategies. To address these questions, in this paper, we provide a unified analysis of SAM and its unnormalized variant (USAM) under one single flexible update rule (Unified SAM), and we present convergence results of the new algorithm under a relaxed and more natural assumption on the stochastic noise. Our analysis provides convergence guarantees for SAM under different step size selections for non-convex problems and functions that satisfy the Polyak-Lojasiewicz (PL) condition (a non-convex generalization of strongly convex functions). The proposed theory holds under the arbitrary sampling paradigm, which includes importance sampling as special case, allowing us to analyze variants of SAM that were never explicitly considered in the literature. Experiments validate the theoretical findings and further demonstrate the practical effectiveness of Unified SAM in training deep neural networks for image classification tasks.

1 Introduction

Consider the classical finite-sum optimization problem

$$\min_{x \in \mathbb{R}^d} \left[f(x) = \frac{1}{n} \sum_{i=1}^n f_i(x) \right]. \tag{1}$$

where each f_i is differentiable, L_i -smooth and lower bounded. Let X^* be the set of minimizers of f, which we assume is non-empty. In practical scenarios, the variable x represents the model parameters, n is the total number of training instances and the functions f_i are loss functions that measure how close our model is to the i-th training data point. The goal is to minimize the average loss of all training instances.

Understanding the generalization capabilities of overparameterized deep neural networks is a central concern in machine learning research, (Zhang et al., 2021). The training objective function f has numerous global optima that perfectly fit the training data, (Liu et al., 2020); however, these different global optima can lead to dramatically varying generalization performances. Empirical observations indicate that stochastic gradient descent (SGD) tends to converge to solutions with good generalization properties, even in the absence of regularization methods (Zhang et al., 2021). This phenomenon suggests that minimizing the training objective using a specific algorithm and initialization strategy can lead to convergence toward a solution with better generalization.

Recent studies have observed that the sharpness of the training loss, that is how rapidly it changes in a neighborhood around the model's parameters, correlates strongly with the generalization error (Keskar et al., 2016; Jiang et al., 2019). Additionally, generalization bounds related to sharpness have been derived (Dziugaite & Roy, 2018). This observation has motivated recent works (Foret et al., 2021; Zheng et al., 2021; Wu et al., 2020) aiming to minimize sharpness to improve generalization. More specifically, building on these ideas Foret et al. (2021), proposed reformulating the

optimization problem in (1) into a min-max problem of the following form:

$$\min_{x \in \mathbb{R}^d} \max_{\|\varepsilon\| \le \rho} f(x + \varepsilon)$$

where ε represents the radius of the desired neighborhood. The merits of such a formulation reside in the fact that essentially we minimize the empirical sharpness measure $\max_{\|\varepsilon\| \le \rho} f(x+\varepsilon) - f(x)$, which inevitably will lead to flatter minima. The objective now is to find x that minimizes f not just at a specific point but across the entire ε -neighborhood. By taking the first-order Taylor expansion of f around x and solving for the optimal ε^* , the (Normalized) Sharpness-Aware Minimization (SAM) update rule is obtained:

$$x^{t+1} = x^t - \gamma_t \nabla f_{S_t} \left(x^t + \rho_t \frac{\nabla f_{S_t}(x^t)}{\|\nabla f_{S_t}(x^t)\|} \right), \tag{SAM}$$

where $S_t \subseteq [n]$ is a random subset of data points (mini-batch) with cardinality τ sampled independently at each iteration t. The normalization of the inner gradient ensures that the point $\tilde{x}^t = x^t + \rho_t \frac{\nabla f_{S_t}(x^t)}{\|\nabla f_{S_t}(x^t)\|}$ is a good approximation of x^t , since $\|\tilde{x}^t - x^t\| = \rho_t$. This leads to a more stable optimization process, as explained in Dai et al. (2023).

In an orthogonal direction and building upon SAM, Unnormalized Sharpness-Aware Minimization (USAM) was introduced by Andriushchenko & Flammarion (2022) and further investigated in Shin et al. (2024); Dai et al. (2023). The update rule for USAM is defined as follows:

$$x^{t+1} = x^t - \gamma_t \nabla f_{S_t} \left(x^t + \rho_t \nabla f_{S_t} (x^t) \right). \tag{USAM}$$

In contrast to SAM, USAM omits the normalization thus the point \tilde{x}^t can be potentially far from x^t making the \tilde{x}^t 's of USAM updates much larger. This means that the removal of normalization can lead to much more aggressive steps making the USAM potentially more unstable.

Although the two variants appear closely related, the proof techniques and upper bounds used in the convergence analysis of SAM are substantially different from those in USAM. Furthermore, the convergence guarantees of the two variants vary significantly. For example, in the deterministic setting (full-batch), SAM guarantees convergence only to a neighborhood of the solution, whereas USAM does not. Additionally, the step sizes γ_t and ρ_t used in the two update rules to guarantee convergence are very different. All of these differences motivate the importance and necessity of a novel general analysis of SAM-type algorithms, unifying the two main variants (SAM and USAM) and providing the ability to design and analyze new SAM-like methods filling existing gaps in the theoretical understanding of the update rules.

In this work we develop such unified framework that allows the combination of the two approaches and, at the same time, obtains the best-known convergence guarantees under relaxed assumptions.

Main Contributions. Our main contributions are summarized below.

♦ **Unified Framework.** We propose the Unified SAM, an update rule that is a convex combination of SAM and USAM, given by:

$$x^{t+1} = x^t - \gamma_t \nabla f_{S_t} \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|\nabla f_{S_t}(x^t)\|} \right) \nabla f_{S_t}(x^t) \right)$$
 (Unified SAM)

where $\lambda \in [0,1]$. The new formulation captures both USAM and SAM as special cases ($\lambda = 0$ and $\lambda = 1$, respectively), but more importantly, it opens up a wide range of possible update rules beyond these traditional settings. The unified framework offers the flexibility to adjust the degree of normalization (using different values for λ) based on specific model needs, offering a more versatile approach to SAM.

 \diamond Technical Assumptions on the Stochastic Noise. Existing convergence analyses of stochastic SAM rely heavily on the bounded variance assumption, that is, there exists a $\sigma \geq 0$ such that $\mathbb{E} \|\nabla f_{S_t}(x) - \nabla f(x)\|^2 \leq \sigma^2$, (Andriushchenko & Flammarion, 2022; Si & Yun, 2023; Li & Giannakis, 2023; Harada & Iiduka, 2024; Mi et al., 2022; Zhuang et al., 2022) or sometimes to the much stronger bounded gradient condition $\mathbb{E} \|\nabla f_{S_t}(x)\|^2 \leq q^2$, where $q \geq 0$ (Mi et al., 2022; Zhuang et al., 2022). While these assumptions have been crucial in previous analyses, they can be overly restrictive. In the literature of convergence analysis for stochastic gradient descent (SGD),

| Work | Assumptions | Arbitrary Sampling? | SAM Variant |
|--|--|---------------------|---|
| PL functions | | | |
| (Andriushchenko & Flammarion, 2022) (Shin et al., 2024) (Dai et al., 2023) Theorems 3.2 and 3.5 | BV Interpolation Deterministic ER | X X X | USAM USAM USAM Unified SAM |
| General Non-convex functions | | | |
| (Mi et al., 2022) (Zhuang et al., 2022) (Andriushchenko & Flammarion, 2022) (Li & Giannakis, 2023) (Si & Yun, 2023) Theorem 3.7 | BV, BG BV, BG BV BV BV | x x x x | SAM/SSAM GSAM USAM SAM SAM Unified SAM |

Table 1: Summary of the convergence results in the SAM literature. In all works, smoothness is assumed. The top part of the table is for PL functions and the lower part is for general non-convex functions. Here BV = Bounded Variance, BG = Bounded Gradients.

there have been a lot of efforts recently on relaxing such assumptions (Gower et al., 2019; Khaled & Richtárik, 2020; Gower et al., 2021), but, to date, no work has successfully used similar ideas for the analysis of SAM. In our analysis, we relax the bounded gradients/variance assumptions by utilizing the recently proposed Expected Residual (ER) condition Gower et al. (2021); Khaled & Richtárik (2020). As we explain later, in several scenarios, including smooth non-convex problems, ER holds for free and allows us to provide step sizes for SAM related to the sampling strategies.

- ♦ Convergence guarantees for Unified SAM. We provide tight convergence guarantees for Unified SAM, for smooth functions satisfying the Polyak-Lojasiewicz (PL) condition (Polyak, 1987; Lojasiewicz, 1963; Karimi et al., 2016) and for general non-convex functions. See also Table 1 for a summary of our results and comparison with closely related works.
- *PL functions:* For constant step-sizes γ and ρ we prove linear convergence for Unified SAM to a neighborhood of the solution. Our theorem holds without requiring the much stronger assumptions of interpolation condition or the bounded variance assumption of previous works (Andriushchenko & Flammarion, 2022; Shin et al., 2024). Additionally, we prove that for decreasing step sizes γ_t and ρ_t , the Unified SAM converges to the exact solution with a sublinear O(1/t) rate (under the ER condition). Our theoretical results hold under the arbitrary sampling paradigm and, as such, can capture tight convergence guarantees in the deterministic setting. For PL functions in the deterministic setting (full-batch SAM), we show that USAM converges to the exact solution, while SAM does not. This observation was first noted by Si & Yun (2023) for deterministic algorithms. To the best of our knowledge, our work is the first that provides tight convergence guarantees, showing this behaviour as a special case of stochastic algorithms.
- Non-convex functions: Under the ER condition, we show that for general non-convex functions Unified SAM with step sizes that depend on T (the total number of iterations) achieves $\mathbb{E} \|\nabla f(x^T)\| < \varepsilon$ for a given ε at a sublinear rate. This is the first result that drops the bounded variance assumption for both USAM and SAM, (Andriushchenko & Flammarion, 2022; Li & Giannakis, 2023), and substitutes it with the Expected Residual condition.

Finally, as corollaries of the main theorems for the above two classes of problems, we obtain the state-of-the-art convergence guarantees for SGD (a special case of SAM with $\rho=0$), showing the tightness of our analysis.

 \diamond **Arbitrary Sampling.** Via a stochastic reformulation of the finite sum problem (1), firstly introduced in Gower et al. (2019), we explain how our convergence guarantees of Unified SAM hold under the arbitrary sampling paradigm. This allows us to cover a wide range of samplings for USAM and SAM (and their convex combination via $\lambda \in [0,1]$) that were never considered in the literature

before, including uniform sampling and importance sampling as special cases. In this sense, our analysis of Unified SAM is unified for different sampling strategies.

♦ Numerical Evaluation. In Section 4, we present extensive experiments validating different aspects of our theoretical results (behavior of methods in the deterministic setting, importance sampling, and different step-size selections). We also assess the performance of Unified SAM in training deep neural networks for multi-class image classification problems. The code for reproducing all experiments is included as supplementary material with our submission.

2 Unified SAM with Arbitrary Sampling

In this work, we provide a theoretical analysis of Unified SAM that allows us to obtain convergence guarantees of any minibatch and reasonable sampling selection.

We are able to do that by leveraging the recently proposed "stochastic reformulation" of the minimization problem (1) from Gower et al. (2019; 2021). Following an identical setting to Gower et al. (2021), we assume that we have access to unbiased gradient estimates $g(x) \in \mathbb{R}^d$ such that $\mathbb{E}[g(x)] = \nabla f(x)$. For example, we can have $g(x) = \frac{1}{\tau} \sum_{i \in S} \nabla f_i(x)$ to be a mini-batch, where $S \subseteq [n]$ is chosen uniformly at random with $|S| = \tau$. To accommodate any form of mini-batching, we utilize the arbitrary sampling notation $g(x) = \nabla f_v(x) := \frac{1}{n} \sum_{i=1}^n v_i \nabla f_i(x)$, where $v \in \mathbb{R}^n_+$ is a random sampling vector drawn from a distribution \mathcal{D} such that $\mathbb{E}_{\mathcal{D}}[v_i] = 1$, for $i = 1, \ldots, n$. Then the original problem (1) can be reformulated as $\min_{x \in \mathbb{R}^d} \mathbb{E}_{\mathcal{D}}\left[f_v(x) := \frac{1}{n} \sum_{i=1}^n v_i f_i(x)\right]$. Note that it follows immediately from the definition of sampling vector that $\mathbb{E}[g(x)] = \frac{1}{n} \sum_{i=1}^n \mathbb{E}[v_i] \nabla f_i(x) = \nabla f(x)$. Using this reformulation of the original problem, the update rule of Unified SAM can be rewritten as follows:

$$x^{t+1} = x^t - \gamma_t g\left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|}\right) g(x^t)\right), \tag{Unified SAM}$$

where $g(x^t) \sim \mathcal{D}$ is sampled i.i.d at each iteration and $\rho_t \geq 0$, $\gamma_t > 0$ and $\lambda_t \in [0,1]$. The name unified stems from the fact that the update rule indeed unifies both USAM and SAM, however, we acknowledge that there exist other SAM-like variants that our approach does not cover.

Arbirtary Sampling. Using the stochastic reformulation, the update rule of Unified SAM includes several variants of the algorithm, each related to different sampling, by simply varying the distribution \mathcal{D} (that satisfies $\mathbb{E}_{\mathcal{D}}[v_i] = 1, \forall i \in [n]$). This flexibility implies that different choices of \mathcal{D} lead to distinct SAM-type methods (never proposed in the literature before) for solving the original problem (1). In this work we focus on two representative sampling distributions, without aiming to be exhaustive:

- 1. Single element sampling: We choose only singleton sets $S = \{i\}$ for $i \in [n]$, i.e. $\mathbb{P}[|S| = 1] = 1$. Each number i is sampled with probability $p_i \in [0,1]$ or more formally the vector $v \in \mathbb{R}^n$ is defined via $\mathbb{P}[v = e_i/p_i] = p_i$, where $\sum_{i=1}^n p_i = 1$. It is clear that $\mathbb{E}[v_i] = 1$. For example, when $p_i = 1/n$ for all i then this reduces to the well-known uniform sampling.
- 2. au-nice Sampling: Let $au \in [n]$. We generate a random subset $S \subseteq [n]$ by choosing uniformly from all subsets of size au. More formally, the vector $v \in \mathbb{R}^n$ is defined by $\mathbb{P}[v = \frac{n}{\tau} \sum_{i \in S} e_i] := 1/\binom{n}{\tau} = \frac{\tau!(n-\tau)!}{n!}$ for any subset $S \subseteq [n]$ with $|S| = \tau$. Using a double counting argument, one can show that $\mathbb{E}[v_i] = 1$, see Gower et al. (2019).

Importantly, our analysis applies to all forms of mini-batching and supports various choices of sampling vectors v. Later in Section 3.4, we provide additional details on non-uniform single element sampling strategies. In addition, it is clear that if $\tau=n$ in the τ -nice sampling then we recover the full batch or deterministic regime. Later in Section 3.2, we further demonstrate how our analysis encompasses deterministic SAM and USAM as special cases.

3 Convergence Analysis

In this section, we present our main convergence results. Firstly, we introduce the main assumption for our results, namely the Expected Residual (ER) condition. Then we focus on PL functions,

where we demonstrate a linear convergence rate using constant step sizes, and also provide a variant with decreasing step sizes for convergence to the exact solution. Moreover, we extend the analysis to general non-convex functions. Lastly, we discuss the use of importance sampling.

3.1 MAIN ASSUMPTION

In all our theoretical results, we rely on the Expected Residual condition.

Assumption 3.1 (Expected Residual Condition). Let $f^{\inf} = \inf_{x \in \mathbb{R}^n} f(x)$. We say the Expected Residual condition holds if there exist parameters $A, B, C \geq 0$ such that for an unbiased estimator g(x) of $\nabla f(x)$, we have that for all $x \in \mathbb{R}^d$

$$\mathbb{E}_{\mathcal{D}} \|g(x)\|^{2} \le 2A[f(x) - f^{\inf}] + B\|\nabla f(x)\|^{2} + C.$$
 (ER)

Most prior works in the SAM literature assume either bounded gradients (e.g., Mi et al. (2022); Zhuang et al. (2022)) or bounded variance (e.g., Andriushchenko & Flammarion (2022); Harada & Iiduka (2024)). Both conditions are stronger assumptions than ER. Note that the bounded gradients assumption is captured by ER for A = B = 0, C > 0, while the bounded variance is obtained for A = 0, B = 1, and C > 0. For a detailed analysis of other conditions that automatically satisfy ER, see Gower et al. (2021) and Khaled & Richtárik (2020). Finally, when each f_i is L_i -smooth under mild assumptions on the distribution \mathcal{D} , one can show that ER holds immediately (not an assumption but property of the problem) and has closed-from expressions for the constants A, B, and C. For more details on the expressions A, B, and C in this scenario, please check Appendix B.

3.2 PL FUNCTIONS

One of the popular generalizations of strong convexity in the literature is the Polyak-Lojasiewicz (PL) condition, (Karimi et al., 2016; Lei et al., 2019). We formally define this condition in Definition A.2 and here we establish linear convergence of Unified SAM for functions that satisfy it.

Theorem 3.2. Assume that each f_i is L_i -smooth, f is μ -PL and the ER is satisfied. Set $L_{\max} = \max_{i \in [n]} L_i$ and let

$$\rho_t = \rho \le \frac{\mu}{L_{\max}(\mu + 2[B\mu + A](1 - \lambda)^2)}, \gamma_t = \gamma \le \frac{\mu - L_{\max}\rho\left(\mu + 2[B\mu + A](1 - \lambda)^2\right)}{2L_{\max}(B\mu + A)\left[2L_{\max}^2\rho^2(1 - \lambda)^2 + 1\right]}.$$

Then for Unified SAM it holds

$$\mathbb{E}[f(x^t) - f(x^*)] \le (1 - \gamma \mu)^t \left[f(x^0) - f(x^*) \right] + N,$$

where
$$N = \frac{L_{\max}}{\mu} \left(C \gamma + \rho (1 + 2 \gamma L_{\max}^2 \rho) \left[\lambda^2 + C (1 - \lambda)^2 \right] \right)$$
.

As an immediate corollary of Theorem 3.2 we get the following guarantees for USAM and SAM.

Corollary 3.3. *Make the same assumptions as Theorem 3.2.*

• USAM: Let
$$\rho \leq \frac{\mu}{L_{\max}(\mu + 2[B\mu + A])}$$
 and $\gamma \leq \frac{\mu - L_{\max}\rho(\mu + 2[B\mu + A])}{2L_{\max}(B\mu + A)[2L_{\max}^2\rho^2 + 1]}$. Then for USAM it holds

$$\mathbb{E}[f(x^t) - f(x^*)] \le (1 - \gamma \mu)^t \left[f(x^0) - f(x^*) \right] + \frac{L_{\max}C}{\mu} \left(\gamma + \rho(1 + 2\gamma L_{\max}^2 \rho) \right).$$

• SAM: Let $\rho \leq \frac{1}{L_{\max}}$ and $\gamma \leq \frac{\mu(1-L_{\max}\rho)}{2L_{\max}(B\mu+A)}$. Then for SAM it holds

$$\mathbb{E}[f(x^t) - f(x^*)] \le (1 - \gamma\mu)^t \left[f(x^0) - f(x^*) \right] + \frac{L_{\max}}{\mu} \left(C\gamma + \rho(1 + 2\gamma L_{\max}^2 \rho) \right).$$

To the best of our knowledge, all prior convergence results for (stochastic) SAM have relied on the strong assumption of bounded variance, as seen in works like Andriushchenko & Flammarion

(2022), Si & Yun (2023), Li & Giannakis (2023) and Harada & Iiduka (2024). In contrast, the above theorem is the first to establish convergence for SAM without relying on this assumption. The closest related works on the convergence of constant step size SAM for PL functions are Shin et al. (2024) and Dai et al. (2023). The former provides a linear convergence rate for USAM in the *interpolated* regime, while the latter establishes a linear convergence rate for USAM in the *deterministic* regime. Our result is the first to demonstrate linear convergence in the fully stochastic regime. Additionally, when $\rho = 0$, Unified SAM reduces to SGD, and Theorem 3.2 recovers the step sizes and rates (up to constants) of Theorem 4.6 from Gower et al. (2021), demonstrating the tightness of our results.

Notice that the constant C from the ER and the parameter λ control the neighborhood of the convergence. In particular, in the deterministic regime, the ER holds trivially with A=0, B=1 and C=0. Under these conditions, Theorem 3.2 simplifies as follows.

Corollary 3.4 (Deterministic SAM). Let f be L-smooth and μ -PL function. Let $\rho \leq \frac{1}{L(1+2(1-\lambda)^2)}$ and $\gamma \leq \frac{1-L\rho\left(1+2(1-\lambda)^2\right)}{2L[2L_{\max}^2\rho^2(1-\lambda)^2+1]}$. Then for Unified SAM it holds

$$\mathbb{E}[f(x^t) - f(x^*)] \le (1 - \gamma\mu)^t \left[f(x^0) - f(x^*) \right] + \frac{L\rho(1 + 2\gamma L^2\rho)\lambda^2}{\mu}.$$

First, observe that the PL parameter μ no longer appears in the step sizes ρ and γ . Additionally, when $\lambda=0$, i.e. USAM, the method converges to the exact solution at a linear rate. However, for $\lambda>0$, and in particular when $\lambda=1$ (SAM), convergence is only up to a neighborhood. This suggests that even in the deterministic setting, SAM does not fully converge to the minimum. This was first investigated in Si & Yun (2023) and a similar result appear in Dai et al. (2023). We illustrate this phenomenon experimentally in Section 4.1.

Finally, as an extension of Theorem 3.2, we also show how to obtain convergence to exact solution with an O(1/t) rate for Unified SAM using decreasing step sizes.

Theorem 3.5. Assume that each f_i is L_i -smooth, f is μ -PL and the ER is satisfied. Let $\rho_t = \min\left\{\frac{1}{2t+1}, \rho^*\right\}$ and $\gamma_t = \min\left\{\frac{2t+1}{(t+1)^2\mu}, \gamma^*\right\}$, where ρ^* and γ^* are the upper bounds of ρ and γ , respectively, in Theorem 3.2. Then for Unified SAM it holds $E[f(x^t) - f(x^*)] \leq O\left(\frac{1}{t}\right)$.

The detailed expression hidden under the big O notation can be found in Appendix C. A similar result appears in Andriushchenko & Flammarion (2022) where they provide decreasing step size selection for USAM for PL functions and prove a convergence rate of O(1/t). However, their result relies on the additional assumption of bounded variance. In contrast, our theorem does not require this assumption and is valid for any $\lambda \in [0,1]$. Notably, to the best of our knowledge, this is the first decreasing step size result for SAM.

3.3 General non-convex functions

In this section, we remove the PL assumption and work with general non-convex functions for Unified SAM. First, we start with a general proposition that upper bounds the quantity $\mathbb{E} \|\nabla f(x^t)\|$. Our approach follows a similar derivation to the analysis of SGD in the same setting by Khaled & Richtárik (2020).

Proposition 3.6. Assume that each
$$f_i$$
 is L_i -smooth and the ER is satisfied. Let $\rho \leq \min\left\{\frac{1}{8L_{\max}}, \frac{1}{BL_{\max}(1-\lambda)^2}\right\}$ and $\gamma \leq \frac{1}{8BL_{\max}}$. Then for Unified SAM it holds
$$\min_{t=0,\dots,T-1} \mathbb{E} \|\nabla f(x^t)\|^2 \leq \frac{2\left(1+2A\gamma L_{\max}\left[\rho(1-\lambda)^2(1+2\gamma\rho L_{\max}^2)+\gamma\right]\right)^T}{T\gamma} [f(x^0)-f^{\inf}] + 2L_{\max}\left[C\gamma + \rho(1+2\gamma\rho L_{\max}^2)(\lambda^2 + C(1-\lambda)^2)\right].$$

In order to control the exponential convergence term of the previous theorem we need to carefully select the step sizes ρ and γ . This is what the following theorem achieves.

Theorem 3.7. Let $\varepsilon > 0$ and set $\delta_0 = f(x_0) - f^{\inf} \ge 0$. For any $\rho \le \overline{\rho}$ and $\gamma \le \overline{\gamma}$. Provided that $T \ge \frac{\delta_0 L_{\max}}{\varepsilon^2} \max \left\{ 96B, 24(1-\lambda)\sqrt{3L_{\max}A}, \frac{5184L_{\max}A^2(1-\lambda)^4\delta_0}{\varepsilon^2}, \frac{864\delta_0A}{\varepsilon^2}, \frac{144C}{\varepsilon^2}, \frac{288L_{\max}^2(1-\lambda)^2}{\varepsilon^2} \right\}$

the iterates of Unified SAM satisfy $\min_{t=0,...,T-1} \mathbb{E} \|\nabla f(x^t)\| \leq \varepsilon$.

For the precise expressions of $\overline{\rho}$ and $\overline{\gamma}$ we refer to Theorem C.4. The results in Proposition 3.6 and Theorem 3.7 are tight, as setting $\rho=0$ Unified SAM reduces in SGD and these simplify to the step sizes and rates (up to constants) of Theorem 2 and Corollary 1 from Khaled & Richtárik (2020). Other results for general non-convex functions can be found in Mi et al. (2022), Zhuang et al. (2022) and Li & Giannakis (2023) for SAM and in Andriushchenko & Flammarion (2022) for USAM. However, as mentioned earlier, all these analyses rely on the strong assumption of bounded variance and/or bounded gradients. In contrast, our result uses weaker assumptions and offers guarantees for any $\lambda \in [0,1]$. Furthermore, Khanh et al. (2024) have results for general non-convex functions in the deterministic setting though all their results are asymptotic. Another closely related work is Nam et al. (2023), where they also assume the expected residual condition, however, they additionally assume bounded gradient of f and their results are only asymptotic and hold almost surely.

3.4 BEYOND UNIFORM SAMPLING: IMPORTANCE SAMPLING

In our contributions, we highlighted that this work is the first to offer a theoretical justification for applying importance sampling in SAM. Importance sampling refers to selecting probabilities that optimize convergence rates, with our focus being on single-element sampling. To derive the importance sampling probabilities, we substitute the bounds for A, B, and C from Proposition B.2 into the sample complexity obtained in Theorem 3.7, resulting in the following:

$$T \ge \frac{\delta_0 L_{\max}}{\varepsilon^2} \max \left\{ 24(1-\lambda) \sqrt{\frac{3L_{\max}}{n\tau} \max_i \frac{L_i}{p_i}}, \frac{\frac{5184L_{\max}}{n^2\tau^2} (\max_i \frac{L_i}{p_i})^2 (1-\lambda)^4 \delta_0}{\varepsilon^2}, \frac{\frac{864\delta_0}{n\tau} \max_i \frac{L_i}{p_i}}{\varepsilon^2}, \frac{\frac{288}{n\tau} \max_i \frac{L_i}{p_i} \sigma^*}{\varepsilon^2}, \frac{288L_{\max}^2 (1-\lambda)^2}{\varepsilon^2} \right\}$$

Now we need to optimize the quantity $\max_i \frac{L_i}{np_i}$ over the probabilities (p_i) . Thus we get

$$p_i = \frac{L_i}{\sum_{j=1}^n L_j},\tag{2}$$

which is precisely the *Importance Sampling*, see also Gower et al. (2019).

4 NUMERICAL EXPERIMENTS

In this section, we evaluate our proposed step sizes for Unified SAM on both deterministic and stochastic PL problems, with experiments designed to illustrate our theoretical findings. Additionally, we explore different values of λ when training deep neural networks to improve accuracy.

4.1 VALIDATION OF THE THEORY

In this part, we empirically validate our theoretical results and illustrate the main properties of Unified SAM that our theory suggests in Section 3. In these experiments, we focus on ℓ_2 -regularized regression problems (problems with strongly convex objective f and components f_i and thus PL) and we evaluate the performance of Unified SAM on synthetic data. The loss function of the ℓ_2 -regularized *ridge regression* is given by

$$f(x) = \frac{1}{2} \|\mathbf{A}x - b\|^2 + \lambda_r \|x\|^2 = \frac{1}{2n} \sum_{i=1}^n (\mathbf{A}[i,:]x - b_i)^2 + \frac{\lambda_r}{2} \|x\|^2$$

 and the loss function of the ℓ_2 -regularized *logistic regression* is given by

$$f(x) = \frac{1}{2n} \sum_{i=1}^{n} \log (1 + \exp(-b_i \mathbf{A}[i,:]x)) + \frac{\lambda_r}{2} ||x||^2.$$

In both problems $A \in \mathbb{R}^{n \times d}$, $b \in \mathbb{R}^n$ are the given data and $\lambda_r \geq 0$ is the regularization parameter.

Normalized SAM converges only in neighborhood. In Corollary 3.4 we highlight that our theoretical results indicate that in the deterministic case, USAM achieves linear convergence to the exact solution. In contrast, when $\lambda > 0$, and specifically for SAM ($\lambda = 1$), convergence is only to a neighborhood of the solution, an unusual outcome for deterministic optimization methods. We validate this observation experimentally in Figure 1. We run a ridge regression problem with $n = 100, d = 100, \lambda_r = 0$. The matrix A has been generated according to Lenard & Minkoff (1984) such that the condition number of A is 10 and the vector b has been sampled from the standard Gaussian distribution. We have used the deterministic Unified SAM for $\lambda = 0.0, \dots, 1.0$ and we run each algorithm for 50 epochs. Indeed we can see that USAM $(\lambda = 0)$ converges all the way to the exact solution while the other choices of λ converge to a neighborhood. It is also noteworthy that the neighborhood increases as λ approaches 1.

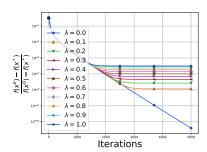


Figure 1: Deterministic Unified SAM for various values of λ applied to the ridge regression problem. USAM ($\lambda = 0$) converges to the exact solution while the other variants $\lambda > 0$ converge to a neighborhood of the solution.

Constant vs Decreasing Step size. In this part, we compare the performance of Unified SAM under both constant and decreasing step size regimes, as described in Theorem 3.2 and Theorem 3.5, respectively. For this experiment, we consider a logistic regression task with n=100, d=100, and $\lambda_r=3/n$. As before the matrix A has been generated such that its condition number is 10 and the vector b has been sampled from the standard Gaussian distribution. We run Unified SAM with $\lambda=0.0,0.5,1.0$, using uniform single-element sampling for 10,000 epochs across 5 trials. The average results of these trials, along with one standard deviation, are shown in Figure 2. Initially, the trajectory of the decreasing step size follows that of the constant step size. However, as the constant step size version of Theorem 3.2 approaches a neighborhood near optimality, it stagnates. In contrast, the decreasing step size from Theorem 3.5 allows for continued improvement, leading to better overall accuracy.

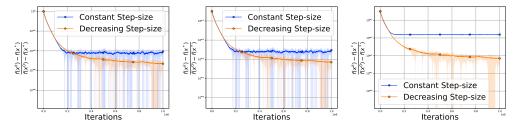


Figure 2: Comparison between constant and decreasing step size regimes of Unified SAM. From left to right we have $\lambda = 0.0, 0.5, 1.0$

Uniform vs Importance Sampling. In this experiment, we demonstrate the benefit of using importance sampling compared to uniform sampling. We consider a ridge regression problem with n=100, d=100, and $\lambda_r=3/n$. The eigenvalues of matrix \boldsymbol{A} are sampled uniformly from the interval [1.0,10.0], and the vector b is drawn from a standard Gaussian distribution. We run Unified SAM with $\lambda=0.0,0.5,1.0$, employing single-element uniform sampling for 3,000 epochs across five trials. For importance sampling, the probabilities are set as $p_i=L_i/\sum_{j=1}^n L_j$. The averaged results with one standard deviation are presented in Figure 3. The results clearly show that importance sampling enhances the convergence of Unified SAM for all values of λ , which is consistent with the discussion in Section 3.4.

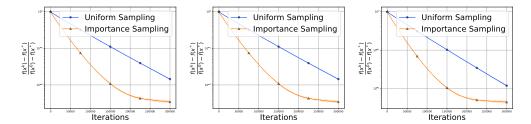


Figure 3: Comparison between uniform and importance sampling for Unified SAM. From left to right we have $\lambda=0.0,0.5,1.0$

4.2 IMAGE CLASSIFICATION

In this section, we evaluate the generalization performance of Unified SAM using the classic image classification task. The models are trained on the CIFAR-10 and CIFAR-100 datasets, (Krizhevsky et al., 2009). For data augmentation, we apply standard techniques such as random crop, random horizontal flip, and normalization (DeVries, 2017). All experiments are conducted on NVIDIA RTX 6000 Ada GPUs.

Models. We use several models in our experiments, including ResNet-18 (RN-18) (He et al., 2016a), and PreActResNet-18 (PRN-18) (He et al., 2016b). To demonstrate the scalability of Unified SAM, we also include Wider ResNet (WRN-28-10) (Zagoruyko & Komodakis, 2016).

Different values of λ . This subsection explores how varying λ in the Unified SAM update rule affects generalization performance. We experiment with the PRN-18 and WRN-28-10 models trained on CIFAR-10 and CIFAR-100. The Unified SAM method is trained using $\rho=0.1,0.2,0.3,0.4$ and $\lambda_t=0.0,0.5,1.0,1/t,1-1/t$. The last two choices of λ_t can be intuitively explained as follows: for $\lambda_t=1-1/t$, the algorithm starts as USAM ($\lambda_1=0$) and gradually transitions toward SAM as $\lambda_t\to 1$ when $t\to\infty$, meaning it begins with USAM and approaches SAM. Conversely, for $\lambda_t=1/t$, the algorithm behaves the other way around, starting closer from SAM and converging to USAM over time. Following Pang et al. (2021) and Zhang et al. (2024), we set the weight decay to 5×10^{-4} , the momentum to 0.9, and train for 100 epochs. The step size γ is initialized at 0.1 and reduced by a factor of 10 at the 75-th and 90-th epochs.

Each experiment is repeated three times, and we report the averages and standard errors in Tables 2 and 3 and Tables 8 and 9 in appendix. The results indicate that, for a fixed ρ , the optimal λ value is not always $\lambda=0$ or $\lambda=1$. Overall, $\lambda_t=1-1/t$ appears to be a reliable choice. In particular, we observe that USAM ($\lambda=0$) does not have the best performance in any of these experiments. On the contrary in the CIFAR10 dataset the value $\lambda_t=1-1/t$ is a good choice while in the CIFAR100 dataset both $\lambda_t=1$ and $\lambda_t=1-1/t$ offer strong performance.

Table 2: Test accuracy (%) of Unified SAM for WRN-28-10 on CIFAR10, evaluated across different values of ρ and λ . With bold we highlight the best performance for fixed ρ .

| Unified SAM | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|--------------|-------------------|--------------------|--------------------|--------------------|---------------------|
| $\rho = 0.1$ | $95.7_{\pm 0.01}$ | $95.68_{\pm0.11}$ | $95.9_{\pm 0.08}$ | $95.84_{\pm0.07}$ | $95.81_{\pm 0.03}$ |
| $\rho = 0.2$ | $95.8_{\pm 0.05}$ | $95.77_{\pm 0.09}$ | $95.93_{\pm 0.07}$ | $95.71_{\pm 0.13}$ | $95.98_{\pm0.1}$ |
| $\rho = 0.3$ | $95.35_{\pm0.3}$ | $95.88_{\pm0.1}$ | $95.95_{\pm 0.09}$ | $95.68_{\pm0.02}$ | $95.99_{\pm 0.06}$ |
| $\rho = 0.4$ | $95.46_{\pm0.02}$ | $95.76_{\pm0.1}$ | $95.62_{\pm 0.05}$ | $95.46_{\pm0.27}$ | $95.79_{\pm 0.07}$ |
| SGD | | | $95.35_{\pm 0.06}$ | | |

Unified VaSSO. The Variance-Suppressed Sharpness-Aware Optimization (VaSSO) method, introduced in Li & Giannakis (2023), is an extension of SAM designed to reduce variance in gradient estimates. VaSSO adjusts the direction of gradient updates using a combination of past gradients and the current gradient, controlled by a parameter θ , aiming to suppress noise during training and enhance generalization, particularly in overparameterized models. In their work, they prove that VaSSO converges at the same asymptotic rate as SAM, under the bounded variance assumption. The update rule

Table 3: Test accuracy (%) of Unified SAM for WRN-28-10 on CIFAR100, evaluated across different values of ρ and λ . With bold we highlight the best performance for fixed ρ .

| Unified SAM | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|--------------------|-------------------|--------------------|----------------------|--------------------|----------------------|
| $\rho = 0.1$ | $80.84_{\pm0.08}$ | $81.01_{\pm 0.11}$ | $80.69_{\pm0.11}$ | $80.81_{\pm0.24}$ | $80.88_{\pm0.31}$ |
| $\rho = 0.2$ | $81.12_{\pm0.18}$ | $81.45_{\pm0.23}$ | $f 81.53_{\pm 0.09}$ | $81.31_{\pm 0.21}$ | $81.22_{\pm 0.19}$ |
| $\rho = 0.3$ | $80.94_{\pm0.13}$ | $81.64_{\pm0.16}$ | $81.62_{\pm0.16}$ | $81.03_{\pm0.14}$ | $f 81.71_{\pm 0.17}$ |
| $\rho = 0.4$ | $80.1_{\pm 0.22}$ | $81.31_{\pm 0.16}$ | $f 81.70_{\pm 0.06}$ | $80.39_{\pm 0.07}$ | $81.59_{\pm 0.05}$ |
| SGD | | | $79.79_{\pm 0.18}$ | | |

of VaSSO is defined as follows: $d_t = (1-\theta)d_{t-1} + \theta \nabla f_i(x^t), x^{t+1} = x^t - \gamma_t \nabla f_i \left(x^t + \rho_t \frac{d_t}{\|d_t\|}\right)$. We can incorporate our unification approach, initially designed for SAM, into VaSSO. This leads to the following modified update rule:

$$d_t = (1 - \theta)d_{t-1} + \theta \nabla f_i(x^t)$$

$$x^{t+1} = x^t - \gamma_t \nabla f_i \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|d_t\|} \right) d_t \right)$$
(Unified VaSSO)

Note that when $\lambda_t = 1$ then we recover VaSSO as introduced in Li & Giannakis (2023).

In this section, we conduct experiments to evaluate the performance of Unified VaSSO. All models are trained for 200 epochs with a batch size of 128. A cosine scheduler is employed in all cases, with an initial step size of 0.05. The weight decay is set to 0.001. For VaSSO, we use $\theta=0.4$, as this value provides the best accuracy according to Li & Giannakis (2023). For the CIFAR-10 dataset, we set $\rho=0.1$, while for CIFAR-100, we use $\rho=0.2$. Each experiment is repeated three times, and we report the average of the maximum test accuracy along with the standard error. The numerical results are presented in Tables 4 and 5.

The results show that, for the WideResNet-28-10 model the value $\lambda_t=1-1/t$ produces the best accuracy. For the ResNet-18 model the best values for λ appear to be $\lambda=0.5$ and $\lambda=1-1/t$. Thus, the choice $\lambda_t=1-1/t$ appears to be a strong overall choice. In any cases, a λ value different from 1 achieves the best performance.

Table 4: Test accuracy (%) of Unified VaSSO on various neural networks trained on CIFAR10.

| Model | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|------------------|--------------------|--------------------|--------------------|--------------------|---------------------|
| ResNet-18 | $96.12_{\pm 0.02}$ | $96.10_{\pm 0.05}$ | $96.22_{\pm 0.11}$ | $96.03_{\pm 0.04}$ | $96.34_{\pm 0.01}$ |
| WideResNet-28-10 | $96.77_{\pm 0.07}$ | $96.93_{\pm0.03}$ | $97.03_{\pm 0.06}$ | $96.71_{\pm 0.06}$ | $97.06_{\pm 0.09}$ |

Table 5: Test accuracy (%) of Unified VaSSO on various neural networks trained on CIFAR100.

| Model | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | / | $\lambda = 1 - 1/t$ |
|------------------|--------------------|--------------------|--------------------|-------------------|-----------------------|
| ResNet-18 | $79.82_{\pm 0.15}$ | $80.01_{\pm 0.07}$ | $79.76_{\pm0.03}$ | $79.93_{\pm0.13}$ | $79.86_{\pm0.04}$ |
| WideResNet-28-10 | $83.07_{\pm0.25}$ | $83.29_{\pm0.33}$ | $83.51_{\pm 0.09}$ | $82.83_{\pm0.18}$ | ${f 83.66}_{\pm0.19}$ |

5 CONCLUSION

In this paper, we introduced the Unified SAM framework, which generalizes both SAM and USAM by using a convex combination parameter λ . We also relaxed the common assumption of bounded variance used in previous analyses by using the ER condition, allowing us to prove convergence under weaker assumptions. In particular, we established convergence guarantees for Unified SAM for both PL and general non-convex functions under this condition. Our analysis also covered importance sampling for SAM. Future work could explore extending this unified approach to adaptive step size choices for SAM or analyzing Unified SAM in the distributed and/or federated learning settings. Another interesting direction is to explore how our results help minimize the sharpness-aware loss (using for example the hessian-regularized version proposed in Wen et al. (2023) and Bartlett et al. (2023).

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Supplementary Material

The Supplementary Material is organized as follows: In Appendix A, we give the basic definitions and lemmas that we need for the proofs. Appendix C presents the proofs of the theoretical guarantees from the main paper. In Appendix D, we provide additional experiments.

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A TECHNICAL PRELIMINARIES

A.1 BASIC DEFINITIONS

In this section, we present some basic definitions we use throughout the paper.

Definition A.1 (L-smooth). A differentiable function $f: \mathbb{R}^d \to \mathbb{R}$ is L-smooth if there exists a constant L > 0 such that

$$\|\nabla f(x) - \nabla f(y)\| \le L\|x - y\|,$$

or equivalently

$$f(x) - f(y) \le \langle \nabla f(y), x - y \rangle + \frac{L}{2} ||x - y||^2$$

for all $x, y \in \mathbb{R}^d$.

Definition A.2 (μ -PL). We say that a differentiable function $f: \mathbb{R}^d \to \mathbb{R}$ satisfies the Polyak-Lojasiewicz condition if

$$\|\nabla f(x)\|^2 \ge 2\mu(f(x) - f(x^*)),$$

for all $x \in \mathbb{R}^d$.

It can be proven that PL functions admit a unique minimizer x^* and moreover, it holds that $x^* \in \arg\min f \Leftrightarrow \nabla f(x^*) = 0$, see Lemma 2.18, 2.13 and 2.22 in Garrigos & Gower (2023). Thus $f^{\inf} = f(x^*)$.

Definition A.3 (Interpolation). We say that the interpolation condition holds if there exists $x^* \in \mathcal{X}^*$ such that

$$\min_{x \in \mathbb{R}^n} f_i(x) = f_i(x^*)$$

for all $i = 1, \ldots, n$.

A.2 BASIC LEMMAS

Lemma A.4. Let f be an L-smooth function. Then for all $x \in \mathbb{R}^n$ we have

$$\|\nabla f(x)\|^2 \le 2L[f(x) - f^{\inf}].$$

Lemma A.5 (Young's Inequality). For any $a, b \in \mathbb{R}^n$ and $\beta \neq 0$ we have

$$\langle a, b \rangle \le \frac{1}{2\beta} ||a||^2 + \frac{\beta}{2} ||b||^2.$$

Completing the square we have

$$\|a+b\|^2 \leq (1+\beta^{-1})\|a\|^2 + (1+\beta)\|b\|^2.$$

For $\beta = 1$ we get

$$||a+b||^2 \le 2||a||^2 + 2||b||^2$$
.

Lemma A.6. Assume that each f_i is L_i -smooth and consider the iterates of Unified SAM. Then it holds

$$\mathbb{E}_{\mathcal{D}} \left\| g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right) \right\|^2$$

$$\leq 4L_{\max}^2 \rho_t^2 \lambda_t^2 + 2 \left[2L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + 1 \right] \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2.$$

Proof. We have

$$\begin{split} & \mathbb{E}_{\mathcal{D}} \left\| g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right) \right\|^2 \\ & = \mathbb{E}_{\mathcal{D}} \left\| g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right) - g(x^t) + g(x^t) \right\|^2 \\ & \stackrel{Lemma A.5}{\leq} 2 \mathbb{E}_{\mathcal{D}} \left\| g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right) - g(x^t) \right\|^2 + 2 \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 \\ & \stackrel{smoothness}{\leq} 2 L_{\max}^2 \rho_t^2 \mathbb{E}_{\mathcal{D}} \left\| (1 - \lambda_t) g(x^t) + \frac{\lambda_t}{\|g(x^t)\|} g(x^t) \right\|^2 + 2 \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 \\ & \stackrel{\leq}{\leq} 4 L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 + 4 L_{\max}^2 \rho_t^2 \lambda_t^2 + 2 \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 \\ & = 4 L_{\max}^2 \rho_t^2 \lambda_t^2 + 2 \left[2 L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + 1 \right] \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 \end{split}$$

Lemma A.7. Assume that each f_i is L_i -smooth and consider the iterates of Unified SAM. Then for any $\beta > 0$ it holds

$$\begin{split} & \mathbb{E}_{\mathcal{D}} \left\langle g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right), \nabla f(x^t) \right\rangle \\ & \geq \left(1 - \frac{\beta}{2} \right) \|\nabla f(x^t)\|^2 - \frac{L_{\max}^2 \rho_t^2 \lambda_t^2}{\beta} - \frac{L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2}{\beta} \, \mathbb{E}_{\mathcal{D}} \, \|g(x^t)\|^2. \end{split}$$

In particular, for $\beta = L_{\max} \rho_t$ we get

$$\begin{split} & \mathbb{E}_{\mathcal{D}} \left\langle g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right), \nabla f(x^t) \right\rangle \\ & \geq \left(1 - \frac{L_{\max} \rho_t}{2} \right) \|\nabla f(x^t)\|^2 - L_{\max} \rho_t \lambda_t^2 - L_{\max} \rho_t (1 - \lambda_t)^2 \, \mathbb{E}_{\mathcal{D}} \, \|g(x^t)\|^2. \end{split}$$

Proof. We have

$$\begin{split} & \mathbb{E}_{\mathcal{D}} \left\langle g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right), \nabla f(x^t) \right\rangle \\ &= \mathbb{E}_{\mathcal{D}} \left\langle g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right) - g(x^t), \nabla f(x^t) \right\rangle \\ &+ \mathbb{E}_{\mathcal{D}} \left\langle g(x^t), \nabla f(x^t) \right\rangle \\ &= \mathbb{E}_{\mathcal{D}} \left\langle g \left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|} \right) g(x^t) \right) - g(x^t), \nabla f(x^t) \right\rangle + \|\nabla f(x^t)\|^2. \end{split}$$

Now

$$-\mathbb{E}_{\mathcal{D}}\left\langle g\left(x^{t} + \rho_{t}\left(1 - \lambda_{t} + \frac{\lambda_{t}}{\|g(x^{t})\|}\right)g(x^{t})\right) - g(x^{t}), \nabla f(x^{t})\right\rangle$$

$$\stackrel{Lemma}{\leq} \frac{A.5}{2\beta} \mathbb{E}_{\mathcal{D}} \left\| g\left(x^{t} + \rho_{t}\left(1 - \lambda_{t} + \frac{\lambda_{t}}{\|g(x^{t})\|}\right)g(x^{t})\right) - g(x^{t})\right\|^{2}$$

$$+ \frac{\beta}{2} \mathbb{E}_{\mathcal{D}} \left\|\nabla f(x^{t})\right\|^{2}$$

$$\stackrel{smoothness}{\leq} \frac{L_{\max}^{2}\rho_{t}^{2}}{2\beta} \mathbb{E}_{\mathcal{D}} \left\| (1 - \lambda_{t})g(x^{t}) + \frac{\lambda_{t}}{\|g(x^{t})\|}g(x^{t})\right\|^{2} + \frac{\beta}{2} \|\nabla f(x^{t})\|^{2}$$

$$\stackrel{Lemma}{\leq} \frac{L_{\max}^2 \rho_t^2}{2\beta} 2(1 - \lambda_t)^2 \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 + \frac{L_{\max}^2 \rho_t^2}{2\beta} 2\lambda_t^2 + \frac{\beta}{2} \|\nabla f(x^t)\|^2 \\
= \frac{L_{\max}^2 \rho_t^2 \lambda_t^2}{\beta} + \frac{L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2}{\beta} \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 + \frac{\beta}{2} \|\nabla f(x^t)\|^2,$$

thus

$$\begin{split} & \mathbb{E}_{\mathcal{D}} \langle g(x^{t} + \rho_{t}g(x^{t})), \nabla f(x^{t}) \rangle \\ & \geq -\frac{L_{\max}^{2} \rho_{t}^{2} \lambda_{t}^{2}}{\beta} - \frac{L_{\max}^{2} \rho_{t}^{2} (1 - \lambda_{t})^{2}}{\beta} \, \mathbb{E}_{\mathcal{D}} \, \|g(x^{t})\|^{2} - \frac{\beta}{2} \|\nabla f(x^{t})\|^{2} + \|\nabla f(x^{t})\|^{2} \\ & = \left(1 - \frac{\beta}{2}\right) \|\nabla f(x^{t})\|^{2} - \frac{L_{\max}^{2} \rho_{t}^{2} \lambda_{t}^{2}}{\beta} - \frac{L_{\max}^{2} \rho_{t}^{2} (1 - \lambda_{t})^{2}}{\beta} \, \mathbb{E}_{\mathcal{D}} \, \|g(x^{t})\|^{2}. \end{split}$$

Lemma A.8. Let $(r_t)_{t\geq 0}$ and $(\delta_t)_{t\geq 0}$ be sequences of non-negative real numbers and let g>1 and $N\geq 0$. Assume that the following recursive relationship holds:

$$r_t \le g\delta_t - \delta_{t+1} + N \tag{3}$$

Then it holds

$$\min_{0 \le t \le T-1} r_t \le \frac{g^T}{T} \delta_0 + N.$$

Proof. Set $\gamma_t = g^{-t}$ for any $t \in \mathbb{Z}$. Then multiply both sides of (3) with g_t . This yields

$$g_t r_t \le g_{t-1} \delta_t - g_t \delta_{t+1} + N g_t.$$

Summing for t = 0, ..., T and telescoping we get

$$\sum_{t=0}^{T-1} g_t r_t \le g_{-1} \delta_0 - g_{T-1} \delta_T + N \sum_{t=0}^{T-1} g_t$$

$$\le g_{-1} \delta_0 + N \sum_{t=0}^{T-1} g_t. \tag{4}$$

Now let $G = \sum_{t=0}^{T-1} g_t$. Note that the sequence (g_t) is decreasing for $t \ge 0$, thus $G \ge Tg_{T-1}$. Now dividing (4) by G we get

$$\min_{0 \le t \le T-1} r_t \le \frac{1}{G} \sum_{t=0}^{T-1} g_t r_t$$

$$\le \frac{g_{-1} \delta_0}{G} + N$$

$$\le \frac{g_{-1} \delta_0}{T g_{T-1}} + N$$

$$= \frac{g^T \delta_0}{T} + N,$$

as wanted.

B More on the Expected Residual Condition

In this section, we introduce several assumptions under which the ER is automatically satisfied. Specifically, we consider the following cases:

• Assuming bounded gradients, i.e.

$$\mathbb{E}_{\mathcal{D}} \|g(x)\|^2 \le \sigma^2, \quad \forall x \in \mathbb{R}^n,$$

then the ER holds with A = 0, B = 0, $C = \sigma^2$.

• Assuming bounded variance, i.e.

$$\mathbb{E}_{\mathcal{D}} \| g(x) - \nabla f(x) \| \le \sigma^2, \quad \forall x \in \mathbb{R}^n,$$

then the ER holds with A = 0, B = 1, $C = \sigma^2$.

• Assuming expected smoothness, i.e.

$$\mathbb{E}_{\mathcal{D}} \|g(x) - \nabla f(x)\| \le 2\mathcal{L}[f(x) - f^{\inf}], \quad \forall x \in \mathbb{R}^n,$$

then the ER holds with $A = 2\mathcal{L}$, B = 0, C = 0.

• Assuming the relaxed strong growth condition, i.e.

$$\mathbb{E}_{\mathcal{D}} \|g(x)\| \le \rho \|\nabla f(x)\| + \sigma^2, \quad \forall x \in \mathbb{R}^n,$$

then the ER holds with A = 0, $B = \rho$, $C = \sigma^2$.

• Assuming the relaxed strong growth condition, i.e.

$$\mathbb{E}_{\mathcal{D}} \|g(x)\| \le \alpha [f(x) - f^{\inf}] + \sigma^2, \quad \forall x \in \mathbb{R}^n,$$

then the ER holds with $A = \alpha$, B = 0, $C = \sigma^2$.

Additionally, if we have more information about the problem and the distribution \mathcal{D} , we can derive stronger bounds on A, B, and C. The more assumptions we make, the tighter these bounds become, as demonstrated in the following propositions.

Proposition B.1 (Prop. 2, (Khaled & Richtárik, 2020)). Assume that each f_i is L_i -smooth and that $\mathbb{E}_{\mathcal{D}}[v_i^2] < \infty$ for all $i \in [n]$. Let $\sigma^* = \frac{1}{n} \sum_{i=1}^n f_i(x^*) - f_i^* \geq 0$, where $f_i^* = \inf_{x \in \mathbb{R}^n} f_i(x)$. Then ER holds with $A = \max_i L_i \mathbb{E}_{\mathcal{D}}[v_i^2]$, B = 0 and $C = 2A\sigma^*$.

This proposition indicates that if each f_i is L_i -smooth and minimal assumptions hold for the distribution \mathcal{D} , then ER is satisfied. As an immediate corollary of Proposition B.1, if the problem further satisfies the interpolation assumption (see Definition A.3), then C = 0. The next proposition gives much tighter constants for A, B, and C in the context of the sampling strategies considered in this paper.

Proposition B.2 (Prop. 3, (Khaled & Richtárik, 2020)). Assume that each f_i is L_i -smooth.

- 1. For the single element sampling ER holds with $A = \frac{1}{n\tau} \max_i \frac{L_i}{p_i}$, B = 0 and $C = 2A\sigma^*$.
- 2. For the τ -nice sampling ER holds with $A = \frac{n-\tau}{\tau(n-1)}L_{\max}$, $B = \frac{n(\tau-1)}{\tau(n-1)}$ and $C = 2A\sigma^*$, where $L_{\max} = \max_i L_i$.

Lastly, if we assume x^* -convexity with τ -nice sampling, we obtain the following constants:

Proposition B.3 (Prop. 3.3, (Gower et al., 2021)). Assume that each f_i is L_i -smooth and that there exists $x^* \in \mathcal{X}^*$ such that f_i is x^* -convex. In addition, assume that \mathcal{D} is the τ -nice sampling. Then ER holds with $A = \frac{n-\tau}{\tau(n-1)}L_{\max}$, B = 1 and $C = \frac{2(n-\tau)}{\tau(n-1)}\sigma_1$, where $\sigma_1 = \sup_{x^* \in \mathcal{X}^*} \frac{1}{n} \sum_{i=1}^n \|\nabla f_i(x^*)\|^2$.

C PROOFS OF THE MAIN RESULTS

In this section we present the proofs of the main theoretical results presented in the main paper, i.e., the convergence guarantees of Unified SAM for PL and smooth functions and general non-convex and smooth functions. We restate the main theorems here for completeness.

In all cases, we use the convention $1/0 = \infty$.

C.1 PROOF OF THEOREM 3.2

 Theorem C.1. Assume that each f_i is L_i -smooth, f is μ -PL and the ER is satisfied. Set $L_{\max} = \max_{i \in [n]} L_i$ and let

$$\rho_t = \rho \le \frac{\mu}{L_{\max}(\mu + 2[B\mu + A](1 - \lambda)^2)}, \gamma_t = \gamma \le \frac{\mu - L_{\max}\rho\left(\mu + 2[B\mu + A](1 - \lambda)^2\right)}{2L_{\max}(B\mu + A)\left[2L_{\max}^2\rho^2(1 - \lambda)^2 + 1\right]}.$$

Then for Unified SAM it holds

$$\mathbb{E}[f(x^t) - f(x^*)] \le (1 - \gamma \mu)^t \left[f(x^0) - f(x^*) \right] + \frac{L_{\max}}{\mu} \left(C\gamma + \rho (1 + 2\gamma L_{\max}^2 \rho) \left[\lambda^2 + C(1 - \lambda)^2 \right] \right).$$

Proof of Theorem 3.2. By combining the smoothness of function f with the update rule of Unified SAM we obtain:

$$f(x^{t+1}) \leq f(x^t) + \langle \nabla f(x^t), x^{t+1} - x^t \rangle + \frac{L_{\max}}{2} \|x^{t+1} - x^t\|^2$$

$$= f(x^t) - \gamma_t \left\langle \nabla f(x^t), g\left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|}\right) g(x^t)\right) \right\rangle$$

$$+ \frac{L_{\max} \gamma_t^2}{2} \left\| g\left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|}\right) g(x^t)\right) \right\|^2.$$

By taking expectation conditioned on x^t we obtain:

$$\mathbb{E}[f(x^{t+1}) - f(x^*)|x^t] - [f(x^t) - f(x^*)]$$

$$\leq -\gamma_t \, \mathbb{E}_{\mathcal{D}} \left\langle g\left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|}\right) g(x^t)\right), \nabla f(x^t) \right\rangle$$

$$+ \frac{L_{\max} \gamma_t^2}{2} \, \mathbb{E}_{\mathcal{D}} \left\| g\left(x^t + \rho_t \left(1 - \lambda_t + \frac{\lambda_t}{\|g(x^t)\|}\right) g(x^t)\right) \right\|^2$$

$$L^{emmas} \, \frac{A.6and \, A.7}{\leq} - \gamma_t \left[\left(1 - \frac{L_{\max} \rho_t}{2}\right) \|\nabla f(x^t)\|^2 - L_{\max} \rho_t \lambda_t^2 - L_{\max} \rho_t (1 - \lambda_t)^2 \, \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 \right]$$

$$+ \frac{\gamma_t^2 L_{\max}}{2} \left[4L_{\max}^2 \rho_t^2 \lambda_t^2 + 2 \left[2L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + 1 \right] \, \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2 \right]$$

$$= -\gamma_t \left(1 - \frac{L_{\max} \rho_t}{2}\right) \|\nabla f(x^t)\|^2 + \gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t \right] \, \mathbb{E}_{\mathcal{D}} \|g(x^t)\|^2$$

$$+ \gamma_t L_{\max} \rho_t \lambda_t^2 (1 + 2\gamma_t L_{\max}^2 \rho_t)$$

$$\leq -\gamma_t \left(1 - \frac{L_{\max} \rho_t}{2}\right) \|\nabla f(x^t)\|^2 + 2A\gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t \right] \left[f(x^t) - f(x^*) \right]$$

$$+ B\gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t \right] \|\nabla f(x^t)\|^2$$

$$+ C\gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t \right] + \gamma_t L_{\max} \rho_t \lambda_t^2 (1 + 2\gamma_t L_{\max}^2 \rho_t)$$

$$= -\gamma_t \left(1 - \frac{L_{\max} \rho_t}{2} - BL_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t \right] \right) \|\nabla f(x^t)\|^2$$

$$+ 2A\gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t \right] \left[f(x^t) - f(x^*) \right]$$

$$+ \gamma_t L_{\max} \left[C\rho_t (1 - \lambda_t)^2 + 2C\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + C\gamma_t + \rho_t \lambda_t^2 + 2\gamma_t L_{\max}^2 \rho_t^2 \lambda_t^2 \right]. \tag{5}$$

Now by the condition on ρ_t and γ_t it is easy to show that the coefficient of $\|\nabla f(x^t)\|^2$ is non-negative so we can use the fact that f is μ -PL to get

$$\mathbb{E}[f(x^{t+1}) - f(x^*)|x^t] - [f(x^t) - f(x^*)]$$

$$\leq -2\gamma_t \mu \left(1 - \frac{L_{\max}\rho_t}{2} - BL_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t\right]\right) [f(x^t) - f(x^*)]$$

$$+ 2A\gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t\right] [f(x^t) - f(x^*)]$$

$$+ \gamma_t L_{\max} \left[C\rho_t (1 - \lambda_t)^2 + 2C\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + C\gamma_t + \rho_t \lambda_t^2 + 2\gamma_t L_{\max}^2 \rho_t^2 \lambda_t^2\right]$$

SO

$$\mathbb{E}[f(x^{t+1}) - f(x^*)|x^t] \leq \left(1 - 2\gamma_t \mu \left(1 - \frac{L_{\max}\rho_t}{2} - BL_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t\right]\right) + 2A\gamma_t L_{\max} \left[\rho_t (1 - \lambda_t)^2 + 2\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + \gamma_t\right]\right) \left[f(x^t) - f(x^*)\right] + \gamma_t L_{\max} \left[C\rho_t (1 - \lambda_t)^2 + 2C\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + C\gamma_t + \rho_t \lambda_t^2 + 2\gamma_t L_{\max}^2 \rho_t^2 \lambda_t^2\right] \leq (1 - \gamma_t \mu) [f(x^t) - f(x^*)] + \gamma_t L_{\max} \left[C\rho_t (1 - \lambda_t)^2 + 2C\gamma_t L_{\max}^2 \rho_t^2 (1 - \lambda_t)^2 + C\gamma_t + \rho_t \lambda_t^2 + 2\gamma_t L_{\max}^2 \rho_t^2 \lambda_t^2\right],$$

where the last inequality follows from the conditions posed on ρ_t and γ_t Taking expectation again and using the tower property we get

$$\mathbb{E}[f(x^{t+1}) - f(x^*)] \\
\leq (1 - \gamma \mu) \, \mathbb{E}[f(x^t) - f(x^*)] \\
+ \gamma L_{\max} \left[C\rho(1 - \lambda)^2 + 2C\gamma L_{\max}^2 \rho^2 (1 - \lambda)^2 + C\gamma + \rho \lambda^2 + 2\gamma L_{\max}^2 \rho^2 \lambda^2 \right] \\
= (1 - \gamma \mu) \, \mathbb{E}[f(x^t) - f(x^*)] + \gamma L_{\max} \left(C\gamma + \rho (1 + 2\gamma L_{\max}^2 \rho) \left[\lambda^2 + C(1 - \lambda)^2 \right] \right).$$
(6)

Recursively applying the above and summing up the resulting geometric series gives:

$$\mathbb{E}[f(x^{t+1}) - f(x^*)] \le (1 - \gamma \mu)^t [f(x^0) - f(x^*)]$$

$$+ \gamma L_{\max} \left(C\gamma + \rho (1 + 2\gamma L_{\max}^2 \rho) \left[\lambda^2 + C(1 - \lambda)^2 \right] \right) \sum_{i=0}^t (1 - \gamma \mu)^j.$$

Using the fact that $\sum_{j=0}^{t} (1 - \gamma \mu)^j \leq \frac{1}{\gamma \mu}$ we get what we wanted.

C.2 Proof of Theorem 3.5

Theorem C.2. Assume that each f_i is L_i -smooth, f is μ -PL and the ER is satisfied. Let $\rho_t = \min\left\{\frac{1}{2t+1}, \rho^*\right\}$ and $\gamma_t = \min\left\{\frac{2t+1}{(t+1)^2\mu}, \gamma^*\right\}$, where ρ^* and γ^* are the upper bounds of ρ and γ , respectively, in Theorem 3.2. Then for Unified SAM it holds

$$E[f(x^t) - f(x^*)] \le O\left(\frac{1}{t}\right).$$

Proof of Theorem 3.5. Since $\gamma_t = \min\left\{\frac{2t+1}{(t+1)^2\mu}, \gamma^*\right\} \leq \gamma^*$ and $\rho_t = \min\left\{\frac{1}{2t+1}, \rho^*\right\} \leq \rho^*$ we get that inequality (6) holds for any $t \geq 0$, hence we have:

$$\mathbb{E}[f(x^{t+1}) - f(x^*)] \leq (1 - \gamma_t \mu) \, \mathbb{E}[f(x^t) - f(x^*)] + \gamma_t L_{\max} \left(C \gamma_t + \rho_t (1 + 2 \gamma_t \rho_t L_{\max}^2) \left[\lambda^2 + C (1 - \lambda)^2 \right] \right) \\
= (1 - \gamma_t \mu) \, \mathbb{E}[f(x^t) - f(x^*)] + C L_{\max} \gamma_t^2 + \gamma_t \rho_t L_{\max}^2 (1 + 2 \gamma_t \rho_t L_{\max}^2) \left[\lambda^2 + C (1 - \lambda)^2 \right] \\
\leq (1 - \gamma_t \mu) \, \mathbb{E}[f(x^t) - f(x^*)] + C L_{\max} \gamma_t^2 + \gamma_t \rho_t L_{\max}^2 \left(1 + \frac{2L_{\max}^2}{\mu} \right) \left[\lambda^2 + C (1 - \lambda)^2 \right], \tag{7}$$

where the last inequality follows from $\gamma_t \leq \frac{2t+1}{(t+1)^2\mu}$ and $\rho_t \leq \frac{1}{2t+1}$, and thus $\gamma_t \rho_t \leq \frac{1}{(t+1)^2\mu} \leq \frac{1}{\mu}$.

Now set $\delta_t = \mathbb{E}[f(x^t) - f(x^*)]$, $R = CL_{\max}$ and $Q = L_{\max}^2 \left(1 + \frac{2L_{\max}^2}{\mu}\right) \left[\lambda^2 + C(1-\lambda)^2\right]$.

Then the inequality (7) takes the following form:

$$\delta_{t+1} \le (1 - \gamma_t \mu) \delta_t + Q \gamma_t \rho_t + R \gamma_t^2. \tag{8}$$

Now since the sequences $\frac{2t+1}{(t+1)^2\mu}$ and $\frac{1}{2t+1}$ are clearly decreasing there exists an integer $t^* \in \mathbb{N}$ such that for any $t \geq t^*$ we have $\gamma_t = \frac{2t+1}{(t+1)^2\mu}$ and $\rho_t \frac{1}{2t+1}$. Substituting in (8) we get that for any $t \geq t^*$ we have

$$\delta_{t+1} \le \frac{t^2}{(t+1)^2} \delta_t + \frac{Q}{\mu(t+1)^2} + \frac{R(2t+1)^2}{\mu^2(t+1)^4}$$
$$\le \frac{t^2}{(t+1)^2} \delta_t + \frac{Q\mu + 4R}{\mu^2(t+1)^2},$$

because $\frac{(2t+1)^2}{(t+1)^4} \leq \frac{4(t+1)^2}{(t+1)^4} = \frac{4}{(t+1)^2}$. Multiplying both sides with $(t+1)^2$ and rearranging we have

$$(t+1)^2 \delta_{t+1} - t^2 \delta_t \le \frac{Q\mu + 4R}{u^2}.$$

Summing for $t = t^*, \dots, T-1$ and telescoping we have

$$T^2 \delta_T \le (t^*)^2 \delta_{t^*} + \frac{Q\mu + 4R}{\mu^2} (T - t^* - 1).$$

Changing notation from T to t we get

$$\mathbb{E}[f(x^t) - f(x^*)] \le \frac{(t^*)^2 \delta_{t^*}}{t^2} + \frac{Q\mu + 4R}{\mu^2} \frac{t - t^* - 1}{t^2}$$
$$\le \frac{(t^*)^2 \delta_{t^*}}{t^2} + \frac{Q\mu + 4R}{\mu^2} \frac{1}{t}$$
$$= O\left(\frac{1}{t}\right),$$

as wanted.

C.3 Proof of Proposition 3.6

Proposition C.3. Assume that each f_i is L_i -smooth and the ER is satisfied. Let

$$\rho \leq \min\left\{\frac{1}{8L_{\max}}, \frac{1}{BL_{\max}(1-\lambda)^2}\right\} \text{ and } \gamma \leq \frac{1}{8BL_{\max}},$$

where $L_{\max} = \max_{i \in [n]} L_i$. Then for Unified SAM it holds

$$\min_{t=0,...,T-1} \mathbb{E} \|\nabla f(x^t)\|^2 \le \frac{2\left(1 + 2A\gamma L_{\max}\left[\rho(1-\lambda)^2(1 + 2\gamma\rho L_{\max}^2) + \gamma\right]\right)^T}{T\gamma} [f(x^0) - f^{\inf}] + 2L_{\max}\left[C\gamma + \rho(1 + 2\gamma\rho L_{\max}^2)(\lambda^2 + C(1-\lambda)^2)\right].$$

Proof of Proposition 3.6. From Equation (5) we have

$$\begin{split} &\mathbb{E}[f(x^{t+1}) - f^{\inf}|x^{t}] - [f(x^{t}) - f^{\inf}] \\ &\leq -\gamma_{t} \left(1 - \frac{L_{\max}\rho_{t}}{2} - BL_{\max} \left[\rho_{t}(1 - \lambda_{t})^{2} + 2\gamma_{t}L_{\max}^{2}\rho_{t}^{2}(1 - \lambda_{t})^{2} + \gamma_{t} \right] \right) \|\nabla f(x^{t})\|^{2} \\ &\quad + 2A\gamma_{t}L_{\max} \left[\rho_{t}(1 - \lambda_{t})^{2} + 2\gamma_{t}L_{\max}^{2}\rho_{t}^{2}(1 - \lambda_{t})^{2} + \gamma_{t} \right] \left[f(x^{t}) - f^{\inf} \right] \\ &\quad + \gamma_{t}L_{\max} \left[C\rho_{t}(1 - \lambda_{t})^{2} + 2C\gamma_{t}L_{\max}^{2}\rho_{t}^{2}(1 - \lambda_{t})^{2} + C\gamma_{t} + \rho_{t}\lambda_{t}^{2} + 2\gamma_{t}L_{\max}^{2}\rho_{t}^{2}\lambda_{t}^{2} \right] \\ &\leq -\frac{\gamma_{t}}{2} \|\nabla f(x^{t})\|^{2} + 2A\gamma_{t}L_{\max} \left[\rho_{t}(1 - \lambda_{t})^{2} + 2\gamma_{t}L_{\max}^{2}\rho_{t}^{2}(1 - \lambda_{t})^{2} + \gamma_{t}L_{\max}^{2}\rho_{t}^{2}(1 - \lambda_{t})^{2} + \gamma_{t}L_{\max}^{2}\rho_{t}^{2}\lambda_{t}^{2} \right] \\ &\quad + \gamma_{t}L_{\max} \left[C\rho_{t}(1 - \lambda_{t})^{2} + 2C\gamma_{t}L_{\max}^{2}\rho_{t}^{2}(1 - \lambda_{t})^{2} + C\gamma_{t} + \rho_{t}\lambda_{t}^{2} + 2\gamma_{t}L_{\max}^{2}\rho_{t}^{2}\lambda_{t}^{2} \right], \end{split}$$

where the last inequality follows by the conditions posed in ρ_t and γ_t . After taking expectation, using the tower property and some algebraic manipulations we get

$$\begin{split} \frac{\gamma}{2} \, \mathbb{E} \, \| \nabla f(x^t) \|^2 & \leq \left(1 + 2A\gamma L_{\max} \left[\rho (1-\lambda)^2 + 2\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + \gamma \right] \right) \mathbb{E}[f(x^t) - f^{\inf}] \\ & - \mathbb{E}[f(x^{t+1}) - f^{\inf}] + \gamma L_{\max} \left[C\rho (1-\lambda)^2 + 2C\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + C\gamma + \rho \lambda^2 + 2\gamma L_{\max}^2 \rho^2 \lambda^2 \right]. \end{split}$$

Now set

$$\begin{split} \delta_t &= \frac{2}{\gamma} \, \mathbb{E}[f(x^t) - f^{\inf}] \ge 0 \\ r_t &= \mathbb{E} \, \|\nabla f(x^t)\|^2 \ge 0 \\ g &= \left(1 + 2A\gamma L_{\max} \left[\rho(1-\lambda)^2 + 2\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + \gamma\right]\right) > 1 \\ N &= 2L_{\max} \left[C\rho(1-\lambda)^2 + 2C\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + C\gamma + \rho\lambda^2 + 2\gamma L_{\max}^2 \rho^2 \lambda^2\right] \ge 0 \end{split}$$

and the previous inequality takes the following form:

$$\frac{\gamma}{2}r_t \le g\delta_t - \delta_{t+1} + N.$$

Now apply Lemma A.8 and we get what we wanted.

C.4 Proof of Theorem 3.7

Theorem C.4. Let $\varepsilon > 0$ and set $\delta_0 = f(x_0) - f^{\inf} \ge 0$. For any

$$\rho \leq \min\left\{\frac{1}{8L_{\max}}, \frac{1}{BL_{\max}(1-\lambda)^2}, \frac{1}{\sqrt{T}}, \frac{\varepsilon^2}{12L_{\max}(C(1-\lambda)^2 + \lambda^2)}\right\}$$

and

$$\begin{split} \gamma \leq \min \left\{ \frac{1}{8BL_{\max}}, \frac{1}{2L(1-\lambda)\sqrt{3AL_{\max}}}, \frac{1}{6L_{\max}A(1-\lambda)^2\sqrt{T}}, \frac{1}{\sqrt{6ALT}}, \right. \\ \left. \frac{\varepsilon^2}{24L_{\max}^3\left(C(1-\lambda)^2 + \lambda^2\right)}, \frac{\varepsilon^2}{12L_{\max}C} \right\} \end{split}$$

Then provided that

$$T \ge \frac{\delta_0 L_{\max}}{\varepsilon^2} \max \left\{ 96B, 24(1-\lambda)\sqrt{3L_{\max}A}, \frac{5184L_{\max}A^2(1-\lambda)^4\delta_0}{\varepsilon^2}, \frac{864\delta_0 A}{\varepsilon^2}, \frac{144C}{\varepsilon^2}, \frac{288L_{\max}^2(1-\lambda)^2}{\varepsilon^2} \right\}$$

we have $\min_{t=0,...,T-1} \mathbb{E} \|\nabla f(x^t)\| \leq \varepsilon$.

Proof of Theorem 3.7. From Proposition C.3 under the condition that $\rho \leq \min\left\{\frac{1}{8L_{\max}}, \frac{1}{BL_{\max}(1-\lambda)^2}\right\}$ and $\gamma \leq \frac{1}{8BL_{\max}}$ we have

$$\min_{t=0,\dots,T-1} \mathbb{E} \|\nabla f(x^t)\|^2 \leq \frac{2\left(1 + 2A\gamma L_{\max}\left[\rho(1-\lambda)^2 + 2\gamma L_{\max}^2 \rho^2(1-\lambda)^2 + \gamma\right]\right)^T}{T\gamma} [f(x^0) - f^{\inf}] \\
+ 2L_{\max}\left[C\rho(1-\lambda)^2 + 2C\gamma L_{\max}^2 \rho^2(1-\lambda)^2 + C\gamma + \rho\lambda^2 + 2\gamma L_{\max}^2 \rho^2\lambda^2\right].$$

Using the fact that $1 + x \le \exp(x)$, we have that

$$(1 + 2A\gamma L_{\max} \left[\rho(1-\lambda)^2 + 2\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + \gamma\right])^T$$

$$\leq \exp\left(2L_{\max}AT\gamma \left[\rho(1-\lambda)^2 + 2\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + \gamma\right]\right) \leq \exp(1) < 3.$$

For the second inequality to hold it is enough to assume that

$$2TAL_{\max}\gamma \left[\rho(1-\lambda)^2 + 2\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + \gamma\right] \le 1$$

For this to hold it is enough to have

1134
1135 •
$$2TAL_{\max}\gamma\rho(1-\lambda)^2 \le 1/3$$

1136 •
$$4TAL_{\text{max}}^3 \gamma \rho^2 (1-\lambda)^2 < 1/3$$

•
$$2TAL_{\max}\gamma^2 \le 1/3$$

Posing the restriction $\rho \leq \frac{1}{\sqrt{T}}$ it is enough to have (solving for γ):

•
$$\gamma \le \frac{1}{6L_{\max}A(1-\lambda)^2\sqrt{T}}$$

•
$$\gamma \le \frac{1}{2L(1-\lambda)\sqrt{3AL_{\max}}}$$

•
$$\gamma \leq \frac{1}{\sqrt{6ALT}}$$

Then substituting in (9) we get

$$\min_{t=0,\dots,T-1} \mathbb{E} \|\nabla f(x^t)\|^2 \le \frac{6\delta_0}{T\gamma} + 2L_{\max} \left[C\rho(1-\lambda)^2 + 2C\gamma L_{\max}^2 \rho^2 (1-\lambda)^2 + C\gamma + \rho\lambda^2 + 2\gamma L_{\max}^2 \rho^2 \lambda^2 \right].$$
(11)

To make the right hand side of (11) smaller than ε^2 , we require that the second term satisfies

$$2L_{\max}\left[C\rho(1-\lambda)^2 + 2C\gamma L_{\max}^2\rho^2(1-\lambda)^2 + C\gamma + \rho\lambda^2 + 2\gamma L_{\max}^2\rho^2\lambda^2\right] \le \frac{\varepsilon^2}{2} \Leftrightarrow 2L_{\max}\left[\rho\left(C(1-\lambda)^2 + \lambda^2\right) + 2L_{\max}^2\gamma\rho^2\left(C(1-\lambda)^2 + \lambda^2\right) + C\gamma\right] \le \frac{\varepsilon^2}{2}$$

For this to hold it is enough to have

•
$$2L_{\max}\rho(C(1-\lambda)^2+\lambda^2) \leq \frac{\varepsilon^2}{6} \Leftarrow \rho \leq \frac{\varepsilon^2}{12L_{\max}(C(1-\lambda)^2+\lambda^2)}$$

•
$$4L_{\max}^3 \gamma \rho^2 \left(C(1-\lambda)^2 + \lambda^2 \right) \le \frac{\varepsilon^2}{6} \stackrel{\rho \le 1}{\Leftarrow} \gamma \le \frac{\varepsilon^2}{24L_{\max}^3 \left(C(1-\lambda)^2 + \lambda^2 \right)}$$

•
$$2L_{\max}C\gamma \leq \frac{\varepsilon^2}{6} \Leftarrow \gamma \leq \frac{\varepsilon^2}{12L_{\max}C}$$

Similarly, for the first term, we get that the number of iterations must satisfy:

$$\frac{6\delta_0}{T\gamma} \le \frac{\varepsilon^2}{2} \Leftrightarrow T \ge \frac{12\delta_0}{\gamma\varepsilon^2} \tag{12}$$

Hence so far we need the following restrictions on ρ and γ :

•
$$\rho \leq \min\left\{\frac{1}{8L_{\max}}, \frac{1}{BL_{\max}(1-\lambda)^2}\right\}$$
 and $\gamma \leq \frac{1}{8BL_{\max}}$ (by the restrictions of Proposition C.3)

•
$$\rho \leq \frac{1}{\sqrt{T}}$$
 and $\rho \leq \frac{\varepsilon^2}{12L_{\max}(C(1-\lambda)^2 + \lambda^2)}$

$$\begin{array}{lll} \bullet \ \gamma & \leq & \frac{1}{6L_{\max}A(1-\lambda)^2\sqrt{T}} \ \ \text{and} \ \ \gamma & \leq & \frac{1}{2L(1-\lambda)\sqrt{3}AL_{\max}} \ \ \text{and} \ \ \gamma & \leq & \frac{1}{\sqrt{6}ALT} \ \ \text{and} \ \ \gamma & \leq & \frac{\varepsilon^2}{12L_{\max}C} \end{array}$$

Plugging each of the previous bounds of γ into (12) we get

•
$$T \ge \frac{96BL_{\max}\delta_0}{\varepsilon^2}$$

•
$$T \ge \frac{24L_{\max}\delta_0(1-\lambda)\sqrt{3L_{\max}A}}{\varepsilon^2}$$

•
$$T \ge \frac{5184L_{\max}^2 A^2 (1-\lambda)^4 \delta_0^2}{\varepsilon^4}$$

1188 •
$$T \geq \frac{864\delta_0^2AL_{\max}}{\varepsilon^4}$$
1190 • $T \geq \frac{144L_{\max}C\delta_0}{\varepsilon^4}$

•
$$T \ge \frac{288L_{\max}^3 \delta_0 (1-\lambda)^2}{\varepsilon^4}$$

Finally, collecting all the terms into a single bound we have:

$$\begin{split} \rho &\leq \min \left\{ \frac{1}{8L_{\max}}, \frac{1}{BL_{\max}(1-\lambda)^2}, \frac{1}{\sqrt{T}}, \frac{\varepsilon^2}{12L_{\max}(C(1-\lambda)^2 + \lambda^2)} \right\} \\ \gamma &\leq \min \left\{ \frac{1}{8BL_{\max}}, \frac{1}{2L(1-\lambda)\sqrt{3AL_{\max}}}, \frac{1}{6L_{\max}A(1-\lambda)^2\sqrt{T}}, \frac{1}{\sqrt{6ALT}}, \frac{\varepsilon^2}{24L_{\max}^3\left(C(1-\lambda)^2 + \lambda^2\right)}, \frac{\varepsilon^2}{12L_{\max}C} \right\} \\ T &\geq \frac{\delta_0 L_{\max}}{\varepsilon^2} \max \left\{ 96B, 24(1-\lambda)\sqrt{3L_{\max}A}, \frac{5184L_{\max}A^2(1-\lambda)^4\delta_0}{\varepsilon^2}, \frac{864\delta_0A}{\varepsilon^2}, \frac{144C}{\varepsilon^2}, \frac{288L_{\max}^2(1-\lambda)^2}{\varepsilon^2} \right\}, \end{split}$$

as wanted.

D ADDITIONAL EXPERIMENTS

In this section, we present additional experimental evaluations of Unified SAM, following the same setup as in Li & Giannakis (2023). Specifically, we train ResNet-18 and WRN-28-10 on CIFAR10 and CIFAR100 datasets. Standard data augmentation techniques, including random cropping, random horizontal flipping, and normalization DeVries (2017), are employed. The models are trained for 200 epochs with a batch size of 128, using a cosine scheduler starting from 0.05. Weight decay is set to 0.001. For SAM, we use $\rho=0.1$ for CIFAR10 and $\rho=0.2$ for CIFAR100. Each experiment is repeated three times, and we report the average of the maximum test accuracy along with the standard error. The numerical results are presented in Tables 6 and 7, demonstrate that Unified SAM consistently improves the test accuracy over both USAM and SAM across all tested models. Lastly, we also observe that careful tuning of the parameter λ is essential for achieving optimal performance.

Table 6: Test accuracy (%) of Unified SAM on various neural networks trained on CIFAR10.

| Model | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|------------------|--------------------|-------------------|--------------------|--------------------|---------------------|
| ResNet-18 | $96.13_{\pm 0.05}$ | $96.16_{\pm0.03}$ | $96.33_{\pm0.03}$ | $96.20_{\pm 0.05}$ | $96.22_{\pm 0.09}$ |
| WideResNet-28-10 | $97.26_{\pm0.44}$ | $96.98_{\pm0.08}$ | $97.05_{\pm 0.05}$ | $96.73_{\pm 0.04}$ | $96.63_{\pm 0.35}$ |

Table 7: Test accuracy (%) of Unified SAM on various neural networks trained on CIFAR100.

| Model | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|------------------|--------------------|--------------------|--------------------|--------------------|---------------------|
| ResNet-18 | $80.21_{\pm 0.02}$ | $80.28_{\pm0.16}$ | $80.14_{\pm 0.07}$ | $80.27_{\pm 0.09}$ | $80.19_{\pm 0.06}$ |
| WideResNet-28-10 | $83.49_{\pm0.23}$ | $83.71_{\pm 0.03}$ | $83.55_{\pm0.19}$ | $83.55_{\pm0.10}$ | $83.62_{\pm0.16}$ |

Table 8: Test accuracy (%) of Unified SAM for PRN-18 on CIFAR10, evaluated across different values of ρ and λ . With bold we highlight the best performance for fixed ρ .

| Unified SAM | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|--------------------|--------------------|--------------------|--------------------|-------------------|---------------------|
| $\rho = 0.1$ | $95.29_{\pm 0.09}$ | $95.32_{\pm 0.02}$ | $95.59_{\pm 0.09}$ | $95.24_{\pm0.03}$ | $95.53_{\pm 0.11}$ |
| $\rho = 0.2$ | $95.25_{\pm0.14}$ | $95.48_{\pm 0.05}$ | $95.5_{\pm 0.02}$ | $95.38_{\pm0.11}$ | $95.58_{\pm0.07}$ |
| $\rho = 0.3$ | $95.25_{\pm0.11}$ | $95.24_{\pm0.02}$ | $95.18_{\pm0.04}$ | $95.12_{\pm0.19}$ | $95.26_{\pm0.1}$ |
| $\rho = 0.4$ | $94.76_{\pm 0.09}$ | $94.98_{\pm 0.07}$ | $94.7_{\pm 0.02}$ | $94.64_{\pm0.03}$ | $94.61_{\pm 0.09}$ |
| SGD | | | $94.82_{\pm 0.02}$ | | |

Table 9: Test accuracy (%) of Unified SAM for PRN-18 on CIFAR100, evaluated across different values of ρ and λ . With bold we highlight the best performance for fixed ρ .

| Unified SAM | $\lambda = 0.0$ | $\lambda = 0.5$ | $\lambda = 1.0$ | $\lambda = 1/t$ | $\lambda = 1 - 1/t$ |
|--------------------|-------------------|--------------------|--------------------|--------------------|------------------------|
| $\rho = 0.1$ | $78.28_{\pm0.15}$ | $78.28_{\pm 0.06}$ | $78.32_{\pm0.22}$ | $78.33_{\pm0.32}$ | $78.39_{\pm0.31}$ |
| $\rho = 0.2$ | $78.98_{\pm0.18}$ | $78.68_{\pm0.13}$ | $78.96_{\pm0.12}$ | $78.87_{\pm 0.02}$ | $78.79_{\pm 0.1}$ |
| $\rho = 0.3$ | $79.0_{\pm 0.05}$ | $78.95_{\pm 0.07}$ | $79.21_{\pm 0.08}$ | $78.73_{\pm 0.06}$ | ${f 79.27}_{\pm 0.08}$ |
| $\rho = 0.4$ | $78.57_{\pm0.26}$ | $78.76_{\pm0.3}$ | $79.05_{\pm 0.16}$ | $78.36_{\pm0.09}$ | $78.79_{\pm 0.13}$ |
| SGD | | | $76.9_{\pm 0.23}$ | | |