ON THE GLOBAL CONVERGENCE OF TRAINING DEEP LINEAR RESNETS

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ABSTRACT

We study the convergence of gradient descent (GD) and stochastic gradient descent (SGD) for training $L$-hidden-layer linear residual networks (ResNets). We prove that for training deep residual networks with certain linear transformations at input and output layers, which are fixed throughout training, both GD and SGD with zero initialization on all hidden weights can converge to the global minimum of the training loss. Moreover, when specializing to appropriate Gaussian random linear transformations, GD and SGD provably optimize wide enough deep linear ResNets. Compared with the global convergence result of GD for training standard deep linear networks (Du & Hu, 2019), our condition on the neural network width is sharper by a factor of $O(\kappa Lq)$, where $\kappa$ denotes the condition number of the covariance matrix of the training data. We further propose a modified identity input and output transformations, and show that a $(d + k)$-wide neural network is sufficient to guarantee the global convergence of GD/SGD, where $d, k$ are the input and output dimensions respectively.

1 INTRODUCTION

Despite the remarkable power of deep neural networks (DNNs) trained using stochastic gradient descent (SGD) in many machine learning applications, theoretical understanding of the properties of this algorithm, or even plain gradient descent (GD), remains limited. Many key properties of the learning process for such systems are also present in the idealized case of deep linear networks. For example, (a) the objective function is not convex; (b) errors back-propagate; and (c) there is potential for exploding and vanishing gradients. In addition to enabling study of systems with these properties in a relatively simple setting, analysis of deep linear networks also facilitates the scientific understanding of deep learning because using linear networks can control for the effect of architecture choices on the expressiveness of networks (Arora et al., 2018; Du & Hu, 2019). For these reasons, deep linear networks have received extensive attention in recent years.

One important line of theoretical investigation of deep linear networks concerns optimization landscape analysis (Kawaguchi, 2016; Hardt & Ma, 2016; Freeman & Bruna, 2016; Lu & Kawaguchi, 2017; Yun et al., 2018; Zhou & Liang, 2018), where major findings include that any critical point of a deep linear network with square loss function is either a global minimum or a saddle point, and identifying conditions on the weight matrices that exclude saddle points. Beyond landscape analysis, another research direction aims to establish convergence guarantees for optimization algorithms (e.g. GD, SGD) for training deep linear networks. Arora et al. (2018) studied the trajectory of gradient flow and showed that depth can help accelerate the optimization of deep linear networks. Ji & Telgarsky (2019); Gunasekar et al. (2018) investigated the implicit bias of GD for training deep linear networks and deep linear convolutional networks respectively. More recently, Bartlett et al. (2019); Arora et al. (2019a); Shamir (2018); Du & Hu (2019) analyzed the optimization trajectory of
GD for training deep linear networks and proved global convergence rates under certain assumptions on the training data, initialization, and neural network structure.

Inspired by the great empirical success of residual networks (ResNets), Hardt & Ma (2016) considered identity parameterizations in deep linear networks, i.e., parameterizing each layer’s weight matrix as $I + W$, which leads to the so-called deep linear ResNets. In particular, Hardt & Ma (2016) established the existence of small norm solutions for deep residual networks with sufficiently large depth $L$, and proved that there are no critical points other than the global minimum when the maximum spectral norm among all weight matrices is smaller than $O(1/L)$. Motivated by this intriguing finding, Bartlett et al. (2019) studied the convergence rate of GD for training deep linear networks with identity initialization, which is equivalent to zero initialization in deep linear ResNets. They assumed whitened data and showed that GD can converge to the global minimum if (i) the training loss at the initialization is very close to optimal or (ii) the regression matrix $\Phi$ is symmetric and positive definite. (In fact, they proved that, when $\Phi$ is symmetric and has negative eigenvalues, GD for linear ResNets with zero-initialization does not converge.) Arora et al. (2019a) showed that GD converges under substantially weaker conditions, which can be satisfied by random initialization schemes. The convergence theory of stochastic gradient descent for training deep linear ResNets is largely missing; it remains unclear under which conditions SGD can be guaranteed to find the global minimum.

In this paper, we establish the global convergence of both GD and SGD for training deep linear ResNets without any condition on the training data. More specifically, we consider the training of $L$-hidden-layer deep linear ResNets with fixed linear transformations at input and output layers. We prove that under certain conditions on the input and output linear transformations, GD and SGD can converge to the global minimum of the training loss function. Moreover, when specializing to appropriate Gaussian random linear transformations, we show that, as long as the neural network is wide enough, both GD and SGD with zero initialization on all hidden weights can find the global minimum. There are two main ingredients of our proof: (i) establishing restricted gradient bounds and a smoothness property; and (ii) proving that these properties hold along the optimization trajectory and further lead to global convergence. We point out the second aspect is challenging especially for SGD due to the uncertainty of its optimization trajectory caused by stochastic gradients. We summarize our main contributions as follows:

- We prove the global convergence of GD and SGD for training deep linear ResNets. Specifically, we derive a generic condition on the input and output linear transformations, under which both GD and SGD with zero initialization on all hidden weights can find global minima. Based on this condition, one can design a variety of input and output transformations for training deep linear ResNets.

- When applying appropriate Gaussian random linear transformations, we show that as long as the neural network width satisfies $m = \Omega(kr\kappa^2)$, with high probability, GD can converge to the global minimum up to an $\epsilon$-error within $O(kr \log(1/\epsilon))$ iterations, where $k, r$ are the output dimension and the rank of training data matrix $X$ respectively, and $\kappa = \|X\|_2^2/\sigma^2(X)$ denotes the condition number of the covariance matrix of the training data. Compared with previous convergence results for training deep linear networks from Du & Hu (2019), our condition on the neural network width is independent of the neural network depth $L$, and is strictly better by a factor of $O(L\kappa)$.

- Using the same Gaussian random linear transformations, we also establish the convergence guarantee of SGD for training deep linear ResNets. We show that if the neural network width satisfies $m = \tilde{\Omega}(k r n^2)\log^2(1/\epsilon) \cdot n^2/B^2$, with constant probability, SGD can converge to the global minimum up to an $\epsilon$-error within $\tilde{O}(k^2\epsilon^{-1}\log(1/\epsilon) \cdot n/B)$ iterations, where $n$ is the training sample size and $B$ is the minibatch size of stochastic gradient. This is the first global convergence rate of SGD for training deep linear networks. Moreover, when the global minimum of the training loss is 0, we prove that SGD can further achieve linear rate of global convergence, and the condition on the neural network width does not depend on the target error $\epsilon$.

As alluded to above, we analyze networks with $d$ inputs, $k$ outputs, and $m \geq \max(d, k)$ nodes in each hidden layer. Linear transformations that are fixed throughout training map the inputs to the first hidden layer, and the last hidden layer to the outputs. We prove that our bounds hold with high probability when these input and output transformations are randomly generated by Gaussian distributions. If, instead, the input transformation simply copies the inputs onto the first $d$ compo-
nents of the first hidden layer, and the output transformation takes the first $k$ components of the last hidden layer, then our analysis does not provide a guarantee. There is a good reason for this: a slight modification of a lower bound argument from Bartlett et al. (2019) demonstrates that GD may fail to converge in this case. However, we describe a similarly simple, deterministic, choice of input and output transformations such that wide enough networks always converge. The resulting condition on the network width is weaker than that for Gaussian random transformations, and thus improves on the corresponding convergence guarantee for linear networks, which, in addition to requiring wider networks, only hold with high probability for random transformations.

1.1 ADDITIONAL RELATED WORK

In addition to what we discussed above, a large bunch of work focusing on the optimization of neural networks with nonlinear activation functions has emerged. We will briefly review them in this subsection.

It is widely believed that the training loss landscape of nonlinear neural networks is highly nonconvex and nonsmooth (e.g., neural networks with ReLU/LeakyReLU activation), thus it is fundamentally difficult to characterize the optimization trajectory and convergence performance of GD and SGD. Some early work (Andoni et al., 2014; Daniely, 2017) showed that wide enough (polynomial in sample size $n$) neural networks trained by GD/SGD can learn a class of continuous functions (e.g., polynomial functions) in polynomial time. However, those works only consider training some of the neural network weights rather than all of them (e.g., the input and output layers) \(^1\). In addition, a series of papers investigated the convergence of gradient descent for training shallow networks (typically 2-layer networks) under certain assumptions on the training data and initialization scheme (Tian, 2017; Du et al., 2018b; Brutzkus et al., 2018; Zhong et al., 2017; Li & Yuan, 2017; Zhang et al., 2018). However, the assumptions made in these works are rather strong and not consistent with practice. For example, Tian (2017); Du et al. (2018b); Zhong et al. (2017); Li & Yuan (2017); Zhang et al. (2018) assumed that the label of each training data is generated by a teacher network, which has the same architecture as the learned network. Brutzkus et al. (2018) assumed that the training data is linearly separable. Li & Liang (2018) addressed this drawback; they proved that for two-layer ReLU network with cross-entropy loss, as long as the neural network is sufficiently wide, under mild assumptions on the training data SGD with commonly-used Gaussian random initialization can achieve nearly zero expected error. Du et al. (2018c) proved the similar results of GD for training two-layer ReLU networks with square loss. Beyond shallow neural networks, Allen-Zhu et al. (2019); Du et al. (2019); Zou et al. (2019) generalized the global convergence results to multi-layer over-parameterized ReLU networks. Chizat et al. (2019) showed that training over-parameterized neural networks actually belongs to a so-called “lazy training” regime, in which the model behaves like its linearization around the initialization. Furthermore, the parameter scaling is more essential than over-paramterization to make the model learning within the “lazy training” regime. Along this line of research, several follow up works have been conducted. Oymak & Sołtanolkotabi (2019); Zou & Gu (2019); Su & Yang (2019); Kawaguchi & Huang (2019) improved the convergence rate and over-paramterization condition for both shallow and deep networks. Arora et al. (2019b) showed that training a sufficiently wide deep neural network is almost equivalent to kernel regression using neural tangent kernel (NTK), proposed in Jacot et al. (2018). Allen-Zhu et al. (2019); Du et al. (2019); Zhang et al. (2019) proved the global convergence for training deep ReLU ResNets. Frei et al. (2019) proved the convergence of GD for training deep ReLU ResNets under an over-parameterization condition that is only logarithmic in the depth of the network, which partially explains why deep residual networks are preferable to fully connected ones. However, all the results in Allen-Zhu et al. (2019); Du et al. (2019); Zhang et al. (2019); Frei et al. (2019) require a very stringent condition on the network width, which typically has a high-degree polynomial dependence on the training sample size $n$. Besides, the results in Allen-Zhu et al. (2019); Zhang et al. (2019) also require that all data points are separated by a positive distance and have unit norm. As shown in Du & Hu (2019) and will be proved in this paper, for deep linear (residual) networks, there is no assumption on the training data, and the condition on the network width is significantly milder, which is independent of the sample size $n$. While achieving a stronger result for linear networks than for nonlinear ones is not surprising, we believe that our analysis, conducted in the idealized deep linear case, can provide useful insights to understand optimization in the nonlinear case.

\(^{1}\)In Daniely (2017), the weight changes in all hidden layers make negligible contribution to the final output, thus can be approximately treated as only training the output layer.
Two concurrent works analyze gradient descent applied to deep linear (residual) networks (Hu et al., 2020; Wu et al., 2019). Hu et al. (2020) consider deep linear networks with orthogonal initialization, and Wu et al. (2019) consider zero initialization on the last layer and identity initialization for the rest of the layers, which are similar to our setting. However, there are several differences between their work and ours. One major difference is that Hu et al. (2020) and Wu et al. (2019) only prove global convergence for GD, but our results cover both GD and SGD. In addition, Hu et al. (2020) focuses on proving the global convergence of GD for sufficiently wide networks, while we provide a generic condition on the input and output linear transformations for ensuring global convergence. Wu et al. (2019) assumes whitened data and proves a $O(L^3 \log(1/\epsilon))$ bound on the number of iterations required for GD to converge, where we establish a $O(\log(1/\epsilon))^2$ bound.

1.2 Notation.

We use lower case, lower case bold face, and upper case bold face letters to denote scalars, vectors and matrices respectively. For a positive integer, we denote the set $\{1, \ldots, k\}$ by $[k]$. Given a vector $x$, we use $|x|_2$ to denote its $\ell_2$ norm. We use $N(\mu, \sigma^2)$ to denote the Gaussian distribution with mean $\mu$ and variance $\sigma^2$. Given a matrix $X$, we denote $\|X\|_{F, p}$, $\|X\|_2$, and $\|X\|_{2, \infty}$ as its Frobenious norm, spectral norm and $\ell_2, \infty$ norm (maximum $\ell_2$ norm over its columns), respectively. In addition, we denote by $\sigma_{\min}(X)$, $\sigma_{\max}(X)$ and $\sigma_r(X)$ the smallest, largest and $r$-th largest singular values of $X$ respectively. For a square matrix $A$, we denote by $\lambda_{\min}(A)$ and $\lambda_{\max}(A)$ the smallest and largest eigenvalues of $A$ respectively. For two sequences $\{a_k\}_{k=0}^\infty$ and $\{b_k\}_{k=0}^\infty$, we say $a_k = O(b_k)$ if $a_k \leq C_1 b_k$ for some absolute constant $C_1$, and use $a_k = \Omega(b_k)$ if $a_k \geq C_2 b_k$ for some absolute constant $C_2$. Except the target error $\epsilon$, we use $\tilde{O}(\cdot)$ and $\Omega(\cdot)$ to hide the logarithmic factors in $O(\cdot)$ and $\Omega(\cdot)$ respectively.

2 Problem Setup

Model. In this work, we consider deep linear ResNets defined as follows:

$$f_W(x) = B(I + W_L) \cdots (I + W_1)Ax,$$

where $x \in \mathbb{R}^d$ is the input, $f_W(x) \in \mathbb{R}^k$ is the corresponding output, $A \in \mathbb{R}^{m \times d}$, $B \in \mathbb{R}^{k \times m}$ denote the weight matrices of input and output layers respectively, and $W_1, \ldots, W_L \in \mathbb{R}^{m \times m}$ denote the weight matrices of all hidden layers. The formulation of ResNets in our paper is different from that in Hardt & Ma (2016); Bartlett et al. (2019), where the hidden layers have the same width as the input and output layers. In our formulation, we allow the hidden layers to be wider by choosing the dimensions of $A$ and $B$ appropriately.

Loss Function. Let $\{(x_i, y_i)\}_{i=1, \ldots, n}$ be the training dataset, $X = (x_1, \ldots, x_n) \in \mathbb{R}^{d \times n}$ be the input data matrix and $Y = (y_1, \ldots, y_n) \in \mathbb{R}^{k \times n}$ be the corresponding output label matrix. We assume the data matrix $X$ is of rank $r$, where $r$ can be smaller than $d$. Let $W = \{W_1, \ldots, W_L\}$ be the collection of weight matrices of all hidden layers. For an example $(x, y)$, we consider the square loss defined by

$$\ell(W; x, y) = \frac{1}{2} \| f_W(x) - y \|_2^2.$$

Then the training loss over the training dataset takes the following form

$$L(W) := \sum_{i=1}^n \ell(W; x_i, y_i) = \frac{1}{2} \| B(I + W_L) \cdots (I + W_1)AX - Y \|_F^2.$$

Algorithm. Similar to Allen-Zhu et al. (2019); Zhang et al. (2019), we consider algorithms that only train the weights $W$ for hidden layers while leaving the input and output weights $A$ and $B$ unchanged throughout training. For hidden weights, we follow the similar idea in Bartlett et al. (2019) and adopt zero initialization (which is equivalent to identity initialization for standard linear network). We would also like to point out that at the initialization, all the hidden layers automatically satisfy the so-called balancedness condition (Arora et al., 2018; 2019a; Du et al., 2018a). The optimization algorithms, including GD and SGD, are summarized in Algorithm 1.

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2 Considering whitened data immediately gives $\kappa = 1$. 

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may not be able to guarantee global convergence. Therefore, it is natural to ask in which setting \( L \) becomes \( I \) so that with appropriate input and output transformations, both GD and SGD can converge to the global minimum. It what follows, we will show that GD/SGD can converge to a point which implies that the global minima of deep linear ResNets cannot be smaller than that of linear model.

It is clear that the expressive power of deep linear ResNets is identical to that of simple linear model, which implies that the global minima of deep linear ResNets cannot be smaller than that of linear model. Therefore, our focus is to show that GD/SGD can converge to a point which is exactly the global minimum of the linear regression problem. It what follows, we will show that with appropriate input and output transformations, both GD and SGD can converge to the global minimum.

3 MAIN THEORY

The following theorem establishes the global convergence of GD for training deep linear ResNets.

**Theorem 3.1.** There are absolute constants \( C \) and \( C_1 \) such that, if the input and output weight matrices satisfy

\[
\frac{\sigma_{\min}^2(A) \sigma_{\min}^2(B)}{\|A\|_2 \|B\|_2} \geq C \frac{\|X\|_2 (L(W^{(0)}) - L(W^*))^{1/2}}{\sigma_r^2(X)}
\]

and the step size satisfies

\[
\eta \leq C_1 \cdot \frac{1}{L \|A\|_2 \|B\|_2 \|X\|_2} \cdot \left( \sqrt{L(W^{(0)}) + \|A\|_2 \|B\|_2 \|X\|_2} \right),
\]

then for all iterates of GD in Algorithm 1, it holds that

\[
L(W^{(t)}) - L(W^*) \leq \left( 1 - \frac{\eta \sigma_{\min}^2(A) \sigma_{\min}^2(B) \sigma_r^2(X)}{L \|A\|_2 \|B\|_2 \|X\|_2} \right)^t \cdot (L(W^{(0)}) - L(W^*)).
\]

**Remark 3.2.** Theorem 3.1 can imply the convergence result in Bartlett et al. (2019). Specifically, in order to turn into the setting considered in Bartlett et al. (2019), we choose \( m = d = k, A = I, B = I \), \( L(W^*) = 0 \) and \( XX^T = I \). Then it can be easily observed that the condition in Theorem 3.1 becomes \( L(W^{(0)}) - L(W^*) \leq C^{-2} \). This implies that the global convergence can be established as long as \( L(W^{(0)}) - L(W^*) \) is smaller than some constant, which is equivalent to the condition proved in Bartlett et al. (2019).

In general, \( L(W^{(0)}) - L(W^*) \) can be large and thus the setting considered in Bartlett et al. (2019) may not be able to guarantee global convergence. Therefore, it is natural to ask in which setting

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**Algorithm 1 (Stochastic) Gradient descent with zero initialization**

1: **input:** Training data \( \{x_i, y_i\}_{i=1}^n \), step size \( \eta \), total number of iterations \( T \), minibatch size \( B \), input and output weight matrices \( A \) and \( B \).
2: **initialization:** For all \( l \in [L] \), each entry of weight matrix \( W_l^{(0)} \) is initialized as 0.
3: for \( t = 0, \ldots, T - 1 \) do
4: \( W_l^{(t+1)} = W_l^{(t)} - \eta \nabla_{W_l} L(W^{(t)}) \) for all \( l \in [L] \)
5: end for
6: **output:** \( W^{(T)} \)
7: for \( t = 0, \ldots, T - 1 \) do
8: Uniformly sample a subset \( B^{(t)} \) of size \( B \) from training data without replacement.
9: For all \( l \in [L] \), compute the stochastic gradient \( G_l^{(t)} = \frac{1}{B} \sum_{i \in B^{(t)}} \nabla_{W_l} \ell(W^{(t)}; x_i, y_i) \)
10: for all \( l \in [L] \), \( W_l^{(t+1)} = W_l^{(t)} - \eta G_l^{(t)} \)
11: end for
12: **output:** \( \{W^{(t)}\}_{t=0, \ldots, T} \)
the condition on $A$ and $B$ in Theorem 3.1 can be satisfied. Here we provide one possible choice which is commonly used in practice (another viable choices can be found in Section 4). We use Gaussian random input and output transformations, i.e., each entry in $A$ is independently generated from $N(0,\alpha^2)$ and each entry in $B$ is generated from $N(0,\beta^2)$. Based on this choice of transformations, we have the following proposition that characterizes the quantity of the largest and smallest singular values of $A$ and $B$, and the training loss at the initialization (i.e., $L(W^{(0)})$). The following proposition is proved in Section A.2.

**Proposition 3.3.** In Algorithm 1, if each entry in $A$ is independently generated from $N(0,\alpha^2)$ and each entry in $B$ is independently generated from $N(0,\beta^2)$, then if $m \geq C \cdot (d+k+\log(1/\delta))$ for some absolute constant $C$, with probability at least $1-\delta$, it holds that

$$
\sigma_{\text{min}}(A) = \Omega(\sqrt{m}), \quad \sigma_{\text{max}}(A) = O(\sqrt{m}), \quad \sigma_{\text{min}}(B) = \Omega(\sqrt{m}), \quad \sigma_{\text{max}}(B) = O(\sqrt{m}),
$$

and

$$
L(W^{(0)}) \leq O(\alpha^2 \beta^2 km \log(n/\delta) \|X\|^2_F + \|Y\|^2_F).
$$

Then based on Theorem 3.1 and Proposition 3.3, we provide the following corollary, proved in Section 3.4, which shows that GD is able to achieve global convergence if the neural network is wide enough.

**Corollary 3.4.** Suppose $\|Y\|_F = O(\|X\|_F)$. Then using Gaussian random input and output transformations in Proposition 3.3 with $\alpha = \beta = 1$, if the neural network width satisfies $m = \Omega(\max\{kr^2 \log(n/\delta), k + d + \log(1/\delta)\})$ then, with probability at least $1 - \delta$, the output of GD in Algorithm 1 achieves training loss at most $L(W^*) + \epsilon$ within $T = O(\kappa \log(1/\epsilon))$ iterations, where $\kappa = \frac{\|X\|^2_F}{\sigma^2(X)}$ denotes the condition number of the covariance matrix of training data.

**Remark 3.5.** For standard deep linear networks, Du & Hu (2019) proved that GD with Gaussian random initialization can converge to a $\epsilon$-suboptimal global minima within $T = O(\kappa \log(1/\epsilon))$ iterations if the neural network width satisfies $m = \tilde{O}(Lr^3 + d)$. In stark contrast, training deep linear ResNets achieves the same convergence rate as training deep linear networks and linear regression, while the condition on the neural network width is strictly milder than that for training standard deep linear networks by a factor of $O(L\kappa)$. This improvement may in part validate the empirical advantage of deep ResNets.

### 3.2 Convergence Guarantee of Stochastic Gradient Descent

The following theorem establishes the global convergence of SGD for training deep linear ResNets.

**Theorem 3.6.** There are absolute constants $C, C_1$ and $C_2$, such for any $0 < \delta \leq 1/6$ and $\epsilon > 0$, if the input and output weight matrices satisfy

$$
\frac{\sigma_{\text{min}}^2(A)\sigma_{\text{min}}^2(B)}{\|A\|_2^2\|B\|_2} \geq C \cdot \frac{n\|X\|_2 \cdot \log(L(W^{(0)})/\epsilon)}{B\sigma^2(X)} \cdot \sqrt{L(W^{(0)})},
$$

and the step size and maximum iteration number are set as

$$
\eta \leq C_1 \cdot \frac{B\sigma_{\text{min}}^2(A)\sigma_{\text{min}}^2(B)\sigma^2(X)}{Ln\|A\|_2^2\|B\|_2\|X\|_2^2} \cdot \min \left\{ \frac{\epsilon}{\|X\|_F^2 \cdot L(W^*)}, \frac{n\|X\|_2^2 \cdot \log(T/\delta) \log(L(W^{(0)})/\epsilon)}{B} \right\},
$$

$$
T = C_2 \cdot \frac{1}{\eta L\sigma_{\text{min}}^2(A)\sigma_{\text{min}}^2(B)\sigma^2(X)} \cdot \log \left( \frac{L(W^{(0)}) - L(W^*)}{\epsilon} \right),
$$

then with probability $\delta$ at least $1/2$ (with respect to the random choices of mini batches), SGD in Algorithm 1 can find a network that achieves training loss at most $L(W^*) + \epsilon$.

By combining Theorem 3.6 and Proposition 3.3, we can show that as long as the neural network is wide enough, SGD can achieve global convergence. Specifically, we provide the condition on the neural network width and the iteration complexity of SGD in the following corollary.

**Corollary 3.7.** Suppose $\|Y\|_F = O(\|X\|_F)$. Then using Gaussian random input and output transformations in Proposition 3.3 with $\alpha = \beta = 1$, for sufficiently small $\epsilon > 0$, if the neural network width satisfies $m = \tilde{O}(kr^2 \log^2(1/\epsilon) \cdot n^2 / B^2 + d)$, with constant probability, SGD in Algorithm 1 can find a point that achieves training loss at most $L(W^*) + \epsilon$ within $T = O(\kappa^2 \epsilon^{-1} \log(1/\epsilon) \cdot n / B)$ iterations.

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One can boost this probability to $1 - \delta$ by independently running $\log(1/\delta)$ copies of SGD in Algorithm 1.
From Corollaries 3.7 and 3.4, we can see that compared with the convergence guarantee of GD, the condition on the neural network width for SGD is worse by a factor of $\tilde{O}(n^2 \log^2(1/\epsilon)/B^2)$ and the iteration complexity is higher by a factor of $\tilde{O}(\kappa \epsilon^{-1} \cdot n/B)$. This is because for SGD, its trajectory length contains high uncertainty, and thus we need stronger conditions on the neural network in order to fully control it.

We further consider the special case that $L(W^*) = 0$, which implies that there exists a ground truth matrix $\Phi$ such that for each training data point $(x_i, y_i)$ we have $y_i = \Phi x_i$. In this case, we have the following theorem, which shows that SGD can attain a linear rate to converge to the global minimum.

**Theorem 3.8.** There are absolute constants $C$, and $C_1$ such that for any $0 < \delta < 1$, if the input and output weight matrices satisfy

$$\frac{\sigma_{\min}^2(A) \sigma_{\min}^2(B)}{||A||_2 \cdot ||B||_2} \geq C \cdot \frac{n ||X||_2}{B \sigma_p^2(X)} \cdot \sqrt{L(W(0))},$$

and the step size is set as

$$\eta \leq C_1 \cdot \frac{B^2 \sigma_{\min}^2(A) \sigma_{\min}^2(B) \sigma_p^2(X)}{L n^2 ||A||_2^2 \cdot ||B||_2^2 \cdot ||X||_2^2 \cdot \log(T/\delta)},$$

for some maximum iteration number $T$, then with probability at least $1 - \delta$, the following holds for all $t \leq T$,

$$L(W(t)) \leq 2L(W(0)) \cdot \left(1 - \eta L \sigma_{\min}^2(A) \sigma_{\min}^2(B) \sigma_p^2(X) \frac{1}{\epsilon} \right)^t.$$  

Similarly, using Gaussian random transformations in Proposition 3.3, we show that SGD can achieve global convergence for wide enough deep linear ResNets in the following corollary.

**Corollary 3.9.** Suppose $||Y||_F = O(||X||_F)$. Then using Gaussian random transformations in Proposition 3.3 with $\alpha = \beta = 1$, for any $\epsilon \leq \tilde{O}(B \|X\|_2 \cdot \sqrt{n} \|X\|_F^2)$, if the neural network width satisfies $m = \tilde{O}(kr n^2 \cdot n^2/B^2 + d)$, with high probability, SGD in Algorithm 1 can find a network that achieves training loss at most $\epsilon$ within $T = \tilde{O}(\kappa n^2 \log(1/\epsilon) \cdot n^2/B^2)$ iterations.

## 4 Discussion on Different Input and Output Linear Transformations

In this section, we will discuss several different choices of linear transformations at input and output layers and their effects to the convergence performance. For simplicity, we will only consider the condition for GD.

As we stated in Subsection 3.1, GD converges if the input and output weight matrices $A$ and $B$

$$\frac{\sigma_{\min}^2(A) \sigma_{\min}^2(B)}{||A||_2 \cdot ||B||_2} \geq C \cdot \frac{||X||_2}{\sigma_p^2(X)} \cdot (L(W(0)) - L(W^*))^{1/2}. \quad (4.1)$$

Then it is interesting to figure out what kind of choice of $A$ and $B$ can satisfy this condition. In Proposition 3.3, we showed that Gaussian random transformations (i.e., each entry of $A$ and $B$ is generated from certain Gaussian distribution) satisfy this condition with high probability, so that GD converges. Here we will discuss the following two other transformations.

**Identity transformations.** We first consider the transformations that $A = [I_d \times d, 0_{d \times (m-d)}]^\top$ and $B = \sqrt{m/k} \cdot [I_{k \times k}, 0_{k \times (m-k)}]$. which is equivalent to the setting in Bartlett et al. (2019) when $m = k = d$. Then it is clear that

$$\sigma_{\min}(B) = \sigma_{\max}(B) = \sqrt{m/k} \quad \text{and} \quad \sigma_{\min}(A) = \sigma_{\max}(A) = 1.$$  

Now let us consider $L(W(0))$. By our choices of $B$ and $A$ and zero initialization on weight matrices in hidden layers, in the case that $d = k$, we have

$$L(W(0)) = \frac{1}{2} ||BAX - Y||_F^2 = \frac{1}{2} \sqrt{m/k} X - Y ||_F^2.$$  

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We remark that \( \sqrt{m/k} \|X - Y\|_F^2 / 2 \) could be as big as \( \frac{1}{2} \left( \frac{m}{k} \|X\|_F^2 / k + \|Y\|_F^2 \right) \) (for example, when \( X \) and \( Y \) are orthogonal). Then plugging these results into (4.1), the condition on \( A \) and \( B \) becomes
\[
\sqrt{m/k} \geq C \cdot \frac{\|X\|_2}{\sigma^2(X)} \cdot \left( \frac{1}{2} \left( \frac{m}{k} \|X\|_F^2 / k + \|Y\|_F^2 \right) - L(W^*) \right)^{1/2} \geq C \cdot \frac{\|X\|_2}{\sigma^2(X)} \cdot \sqrt{\frac{m \|X\|_F^2}{2k}},
\]
where the second inequality is due to the fact that \( L(W^*) \leq \|Y\|_F^2 / 2 \). Then it is clear if \( \|X\|_F \geq \sqrt{2/C} \), the above inequality cannot be satisfied for any choice of \( m \), since it will be cancelled out on both sides of the inequality. Therefore, in such cases, our bound does not guarantee that GD achieves global convergence. Thus, it is consistent with the non-convergence results in (Bartlett et al., 2019). Note that replacing the scaling factor \( \sqrt{m/k} \) in the definition of \( B \) with any other function of \( d, k \) and \( m \) would not help.

**Modified identity transformations.** In fact, we show that a different type of identity transformations of \( A \) and \( B \) can satisfy the condition (4.1). Here we provide one such example. Assuming \( m \geq d + k \), we can construct two sets \( S_1, S_2 \subset [m] \) satisfying \( |S_1| = d \), \( |S_2| = k \) and \( S_1 \cap S_2 = \emptyset \). Let \( S_1 = \{i_1, \ldots, i_d\} \) and \( S_2 = \{j_1, \ldots, j_k\} \). Then we construct matrices \( A \) and \( B \) as follows:
\[
A_{ij} = \begin{cases} 
1 & \text{if } (i, j) = (i_j, j) \\
0 & \text{otherwise}
\end{cases} \quad B_{ij} = \begin{cases} 
\alpha & \text{if } (i, j) = (i_j, j) \\
0 & \text{otherwise}
\end{cases}
\]
where \( \alpha \) is a parameter which will be specified later. In this way, it can be verified that \( BA = 0 \), \( \sigma_{\min}(A) = \sigma_{\max}(A) = 1 \), and \( \sigma_{\min}(B) = \sigma_{\max}(B) = \alpha \). Thus it is clear that the initial training loss satisfies \( L(W^{(0)}) = \|Y\|_F^2 / 2 \). Then plugging these results into (4.1), the condition on \( A \) and \( B \) can be rewritten as
\[
\alpha \geq C \cdot \frac{\|X\|_2}{\sigma^2(X)} \cdot \left( \frac{\|Y\|_F^2}{2} - L(W^*) \right)^{1/2}.
\]
The R.H.S. of the above inequality does not depend on \( \alpha \), which implies that we can choose sufficiently large \( \alpha \) to make this inequality hold. Thus, GD can be guaranteed to achieve the global convergence. Moreover, it is worth noting that using modified identity transformation, a neural network with \( m = d + k \) suffices to guarantee the global convergence of GD. We further remark that similar analysis can be extended to SGD.

5 Experiments

In this section, we conduct various experiments to verify our theory on synthetic data, including i) comparison between different input and output transformations and ii) comparison between training deep linear ResNets and standard linear networks.

5.1 Different Input and Output Transformations

To validate our theory, we performed simple experiment on 10-d synthetic data. Specifically, we randomly generate \( X \in \mathbb{R}^{10 \times 1000} \) from a standard normal distribution and set \( Y = -X + 0.1 \cdot E \), where each entry in \( E \) is independently generated from standard normal distribution. Consider 10-hidden-layer linear ResNets, we apply three input and output transformations including identity transformations, modified identity transformations and random transformations. We evaluate the convergence performances for these three choices of transformations and report the results in Figures 1(a)-1(b), where we consider two cases \( m = 40 \) and \( m = 200 \). It can be clearly observed that gradient descent with identity initialization gets stuck, but gradient descent with modified identity initialization or random initialization converges well. This verifies our theory. It can be also observed that modified identity initialization can lead to slightly faster convergence rate as its initial training loss can be smaller. In fact, with identity transformations in this setting, only the first 10 entries of the \( m \) hidden variables in each layer ever take a non-zero value, so that, no matter how large \( m \) is, effectively, \( m = 10 \), and the lower bound of Bartlett et al. (2019) applies.

5.2 Comparison with Standard Deep Linear Networks

Then we compare the convergence performances with that of training standard deep linear networks. Specifically, we adopt the same training data generated in Section 5.1 and consider training \( L \)-hidden-layer neural network with fixed width \( m \). The convergence results are displayed in Figures
Sanjeev Arora, Nadav Cohen, Noah Golowich, and Wei Hu. A convergence analysis of gradient descent for deep linear neural networks. In International Conference on Learning Representations (ICLR 2019a).

Sanjeev Arora, Nadav Cohen, and Elad E Hazan. On the optimization of deep networks: Implicit acceleration by overparameterization. In 35th International Conference on Machine Learning, ICML 2018, pp. 372–389, 2018.

Sanjeev Arora, Nadav Cohen, Noah Golowich, and Wei Hu. A convergence analysis of gradient descent for deep linear neural networks. In International Conference on Learning Representations, 2019a.

1(c)-1(d), where we consider different choices of $L$. For training linear ResNets, we found that the convergence performances are quite similar for different $L$, thus we only plot the convergence result for the largest one (e.g., $L=20$ for $m=40$ and $L=100$ for $m=200$). However, it can be observed that for training standard linear networks, the convergence performance becomes worse as the depth increases. This is consistent with the theory as our condition on the neural network width is $m = O(kr^2)$ (please refer to Corollary 3.4), which has no dependency in $L$, while the condition for training standard linear network is $m = O(Lkr^3)$ (Du & Hu, 2019), which is linear in $L$.

6 CONCLUSION

In this paper, we proved the global convergence of GD and SGD for training deep linear ResNets with square loss. More specifically, we considered fixed linear transformations at both input and output layers, and proved that under certain conditions on the transformations, GD and SGD with zero initialization on all hidden weights can converge to the global minimum. In addition, we further proved that when specializing to appropriate Gaussian random linear transformations, GD and SGD can converge as long as the neural network is wide enough. Compared with the convergence results of GD for training standard deep linear networks, our condition on the neural network width is strictly milder. Our analysis can be generalized to prove similar results for different loss functions such as cross-entropy loss, and can potentially provide meaningful insights to the convergence analysis of deep non-linear ResNets.

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REFERENCES

Zeyuan Allen-Zhu, Yuanzhi Li, and Zhao Song. A convergence theory for deep learning via over-parameterization. In International Conference on Machine Learning, pp. 242–252, 2019.

Alexandr Andoni, Rina Panigrahy, Gregory Valiant, and Li Zhang. Learning polynomials with neural networks. In International Conference on Machine Learning, pp. 1908–1916, 2014.

Sanjeev Arora, Nadav Cohen, and Elad E Hazan. On the optimization of deep networks: Implicit acceleration by overparameterization. In 35th International Conference on Machine Learning, ICML 2018, pp. 372–389, 2018.

Sanjeev Arora, Nadav Cohen, Noah Golowich, and Wei Hu. A convergence analysis of gradient descent for deep linear neural networks. In International Conference on Learning Representations, 2019a.
Sanjeev Arora, Simon S Du, Wei Hu, Zhiyuan Li, Ruslan Salakhutdinov, and Ruosong Wang. On exact computation with an infinitely wide neural net. In Advances in Neural Information Processing Systems, 2019b.

Peter L Bartlett, David P Helmbold, and Philip M Long. Gradient descent with identity initialization efficiently learns positive-definite linear transformations by deep residual networks. Neural computation, 31(3):477–502, 2019.

Alon Brutzkus, Amir Globerson, Eran Malach, and Shai Shalev-Shwartz. SGD learns over-parameterized networks that provably generalize on linearly separable data. In International Conference on Learning Representations, 2018.

Lenaic Chizat, Edouard Oyallon, and Francis Bach. On lazy training in differentiable programming. In Advances in Neural Information Processing Systems, 2019.

Amit Daniely. SGD learns the conjugate kernel class of the network. In Advances in Neural Information Processing Systems, pp. 2422–2430, 2017.

Simon Du and Wei Hu. Width provably matters in optimization for deep linear neural networks. In International Conference on Machine Learning, pp. 1655–1664, 2019.

Simon Du, Jason Lee, Haochuan Li, Liwei Wang, and Xiyu Zhai. Gradient descent finds global minima of deep neural networks. In International Conference on Machine Learning, pp. 1675–1685, 2019.

Simon S Du, Wei Hu, and Jason D Lee. Algorithmic regularization in learning deep homogeneous models: Layers are automatically balanced. In Advances in Neural Information Processing Systems, pp. 384–395, 2018a.

Simon S Du, Jason D Lee, and Yuandong Tian. When is a convolutional filter easy to learn? In International Conference on Learning Representations, 2018b.

Simon S Du, Xiyu Zhai, Barnabas Poczos, and Aarti Singh. Gradient descent provably optimizes over-parameterized neural networks. arXiv preprint arXiv:1810.02054, 2018c.

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

Daniel C Freeman and Joan Bruna. Topology and geometry of half-rectified network optimization. In International Conference on Learning Representations, 2016.

Spencer Frei, Yuan Cao, and Quanquan Gu. Algorithm-dependent generalization bounds for over-parameterized deep residual networks. In Advances in Neural Information Processing Systems, 2019.

Suriya Gunasekar, Jason D Lee, Daniel Soudry, and Nati Srebro. Implicit bias of gradient descent on linear convolutional networks. In Advances in Neural Information Processing Systems, pp. 9461–9471, 2018.

Moritz Hardt and Tengyu Ma. Identity matters in deep learning. arXiv preprint arXiv:1611.04231, 2016.

Wei Hu, Lechao Xiao, and Jeffrey Pennington. Provable benefit of orthogonal initialization in optimizing deep linear networks. In International Conference on Learning Representations, 2020. URL https://openreview.net/forum?id=rkgqN1SYvr.

Arthur Jacot, Franck Gabriel, and Clément Hongler. Neural tangent kernel: Convergence and generalization in neural networks. In Advances in neural information processing systems, pp. 8571–8580, 2018.

Ziwei Ji and Matus Telgarsky. Gradient descent aligns the layers of deep linear networks. In ICLR, 2019.

Kenji Kawaguchi. Deep learning without poor local minima. In Advances in Neural Information Processing Systems, pp. 586–594, 2016.
Kenji Kawaguchi and Jiaoyang Huang. Gradient descent finds global minima for generalizable deep neural networks of practical sizes. *arXiv preprint arXiv:1908.02419*, 2019.

Yuanzhi Li and Yingyu Liang. Learning overparameterized neural networks via stochastic gradient descent on structured data. In *Proceedings of the 32nd International Conference on Neural Information Processing Systems*, pp. 8168–8177, 2018.

Yuanzhi Li and Yang Yuan. Convergence analysis of two-layer neural networks with ReLU activation. In *Proceedings of the 31st International Conference on Neural Information Processing Systems*, pp. 597–607. Curran Associates Inc., 2017.

Haihao Lu and Kenji Kawaguchi. Depth creates no bad local minima. *arXiv preprint arXiv:1702.08580*, 2017.

Samet Oymak and Mahdi Soltanolkotabi. Towards moderate overparameterization: global convergence guarantees for training shallow neural networks. *arXiv preprint arXiv:1902.04674*, 2019.

Ohad Shamir. Exponential convergence time of gradient descent for one-dimensional deep linear neural networks. *arXiv preprint arXiv:1809.08587*, 2018.

Lili Su and Pengkun Yang. On learning over-parameterized neural networks: A functional approximation prospective. *arXiv preprint arXiv:1905.10826*, 2019.

Yuandong Tian. An analytical formula of population gradient for two-layered ReLU network and its applications in convergence and critical point analysis. In *Proceedings of the 34th International Conference on Machine Learning-Volume 70*, pp. 3404–3413. JMLR. org, 2017.

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

Lei Wu, Qingcan Wang, and Chao Ma. Global convergence of gradient descent for deep linear residual networks. *arXiv preprint arXiv:1911.00645*, 2019.

Chulhee Yun, Suvrit Sra, and Ali Jadbabaie. Global optimality conditions for deep neural networks. In *International Conference on Learning Representations*, 2018.

Huishuai Zhang, Da Yu, Wei Chen, and Tie-Yan Liu. Training over-parameterized deep resnet is almost as easy as training a two-layer network. *arXiv preprint arXiv:1903.07120*, 2019.

Xiao Zhang, Yaodong Yu, Lingxiao Wang, and Quanquan Gu. Learning one-hidden-layer ReLU networks via gradient descent. *arXiv preprint arXiv:1806.07808*, 2018.

Kai Zhong, Zhao Song, Prateek Jain, Peter L Bartlett, and Inderjit S Dhillon. Recovery guarantees for one-hidden-layer neural networks. In *Proceedings of the 34th International Conference on Machine Learning-Volume 70*, pp. 4140–4149. JMLR. org, 2017.

Yi Zhou and Yingbin Liang. Critical points of linear neural networks: Analytical forms and landscape properties. 2018.

Difan Zou and Quanquan Gu. An improved analysis of training over-parameterized deep neural networks. In *Advances in Neural Information Processing Systems*, 2019.

Difan Zou, Yuan Cao, Dongruo Zhou, and Quanquan Gu. Stochastic gradient descent optimizes over-parameterized deep ReLU networks. *Machine Learning Journal*, 2019.

## A Proof of Main Theorems

We first provide the following lemma which proves upper and lower bounds on $\|\nabla_W L(W)\|_F^2$ when $W$ is staying inside a certain region. Its proof is in Section B.1.
Lemma A.1. For any weight matrices satisfying \( \max_{i \in [L]} \| W_i \|_2 \leq 0.5/L \), it holds that,

\[
\| \nabla_{W_i} L(W) \|_F^2 \geq \frac{2 \sigma_{\min}^2(A) \sigma_{\min}^2(B) \sigma_r^2(X)(L(W) - L(W^*))}{e},
\]

\[
\| \nabla_{W_i} L(W) \|_F^2 \leq 2e\| A \|_2^2 \| B \|_2^2 \| X \|_2^2 (L(W) - L(W^*))
\]

\[
\| \nabla_{W_i}(W; x_i, y_i) \|_F^2 \leq 2e\| A \|_2^2 \| B \|_2^2 \| x_i \|_2^2 \| (W; x_i, y_i) \|
\]

In addition, the stochastic gradient \( G_t \) in Algorithm 1 satisfies

\[
\| G_t \|_F^2 \leq \frac{2en^2\| A \|_2^2 \| B \|_2^2 \| X \|_2^2}{B^2} L(W),
\]

where \( B \) is the minibatch size.

The gradient lower bound can also be interpreted as the Polyak-Łojasiewicz condition, which is essential to the linear convergence rate. The gradient upper bound is crucial to bound the trajectory length, since this lemma requires that \( \max_{i \in [L]} \| W_i \| \leq 0.5/L \).

The following lemma proves the smoothness property of the training loss function \( L(W) \) when \( W \) is staying inside a certain region. Its proof is in Section B.2.

Lemma A.2. For any two collections of weight matrices, denoted by \( \tilde{W} = \{ \tilde{W}_1, \ldots, \tilde{W}_L \} \) and \( W = \{ W_1, \ldots, W_L \} \), satisfying \( \max_{i \in [L]} \| W_i \|_F, \max_{i \in [L]} \| \tilde{W}_i \|_F \leq 0.5/L \) that, it holds that

\[
L(\tilde{W}) - L(W) \leq \sum_{i=1}^{L} \langle \nabla_{W_i} L(W), \tilde{W}_i - W_i \rangle + L\| A \|_2 \| B \|_2 \| X \|_2 (2\sqrt{2e}L(W) + 0.5e\| A \|_2 \| B \|_2 \| X \|_2) \sum_{i=1}^{L} \| \tilde{W}_i - W_i \|_F^2.
\]

Based on these two lemmas, we are able to complete the proof of all theorems, which are provided as follows.

A.1 PROOF OF THEOREM 3.1

Proof of Theorem 3.1. In order to simplify the proof, we use the short-hand notations \( \lambda_A, \mu_A, \lambda_B \) and \( \mu_B \) to denote \( \| A \|_2, \sigma_{\min}(A), \| B \|_2 \) and \( \sigma_{\min}(B) \) respectively. Specifically, we rewrite the condition on \( A \) and \( B \) as follows

\[
\frac{\mu_A^2 \mu_B^2}{\lambda_A \lambda_B} \geq \frac{4\sqrt{2e}L(X)_{\| X \|_2}}{\sigma_r^2(X)} \cdot (L(W^{(0)}) - L(W^*))^{1/2}.
\]

We prove the theorem by induction on the update number \( s \), using the following two-part inductive hypothesis:

(i) \( \max_{i \in [L]} \| W_{i}^{(s)} \|_F \leq 0.5/L \),

(ii) \( L(W_{i}^{(s)}) - L(W^*) \leq \left( 1 - \frac{\eta L \mu_A^2 \mu_B^2 \sigma_r^2(X)}{e} \right)^s \cdot (L(W^{(0)}) - L(W^*)) \).

First, it can be easily verified that this holds for \( s = 0 \). Now, assume that the inductive hypothesis holds for \( s < t \).

Induction for Part (i): We first prove that \( \max_{i \in [L]} \| W_{i}^{(t)} \|_F \leq 0.5/L \). By triangle inequality and the update rule of gradient descent, we have

\[
\| W_{i}^{(t)} \|_F \leq \sum_{s=0}^{t-1} \eta \| \nabla_{W_i} L(W_{i}^{(s)}) \|_F
\]

\[
\leq \eta \sum_{s=0}^{t-1} \sqrt{2e} \lambda_A \lambda_B \| X \|_2 \cdot (L(W_{i}^{(s)}) - L(W^*))^{1/2}
\]

\[
\leq \sqrt{2e} \eta \lambda_A \lambda_B \| X \|_2 \cdot (L(W^{(0)}) - L(W^*))^{1/2} \cdot \sum_{s=0}^{t-1} \left( 1 - \frac{\eta L \mu_A^2 \mu_B^2 \sigma_r^2(X)}{e} \right)^{s/2}
\]
where the second inequality follows from Lemma A.1, and the third inequality follows from the inductive hypothesis. Since \( \sqrt{1 - x} \leq 1 - x/2 \) for any \( x \in [0, 1] \), we further have

\[
\|W_t^{(t)}\|_F \leq \sqrt{2}e\eta\lambda_A\lambda_B\|X\|_2 \cdot (L(W^{(0)}) - L(W^*))^{1/2} \leq \frac{\sqrt{8e^3\lambda_A\lambda_B\|X\|_2}}{L_{\mu_B}\mu_B\sigma_f^2(X)} \cdot (L(W^{(0)}) - L(W^*))^{1/2}.
\]

Under the condition that \( \mu_B^2\mu_B^2/(\lambda_A\lambda_B) \geq 2\sqrt{8e^3\|X\|_2(L(W^{(0)}) - L(W^*))^{1/2}/\sigma_f^2(X) \), it can be readily verified that \( \|W_t^{(t)}\|_F \leq 0.5/L \). Since this holds for all \( l \in [L] \), we have proved Part (i) of the inductive step, i.e., \( \max_{l \in [L]} \|W_l^{(t)}\|_F \leq 0.5/L \).

**Induction for Part (ii):** Now we prove Part (ii) of the inductive step, bounding the improvement in the objective function. Note that we have already shown that \( W^{(t)} \) satisfies \( \max_{l \in [L]} \|W_l^{(t)}\|_F \leq 0.5/L \), thus by Lemma A.2 we have

\[
L(W^{(t)}) \leq L(W^{(t-1)}) - \eta \sum_{l=1}^{L} \|\nabla W_l L(W^{(t-1)})\|_F^2 + \eta^2\lambda_A\lambda_B\|X\|_2 \cdot (\sqrt{eL(W^{(t-1)})} + 0.5e\lambda_A\lambda_B\|X\|_2) \cdot L(W^{(t-1)}) \|\nabla W_l L(W^{(t-1)})\|_F^2,
\]

where we use the fact that \( W_l^{(t)} - W_l^{(t-1)} = -\eta\nabla W_l L(W^{(t-1)}) \). Note that \( L(W^{(t-1)}) \leq L(W^{(0)}) \) and the step size is set to be

\[
\eta = \frac{1}{2L\lambda_A\lambda_B\|X\|_2 \cdot (\sqrt{eL(W^{(0)})} + 0.5e\lambda_A\lambda_B\|X\|_2)},
\]

so that we have

\[
L(W^{(t)}) - L(W^{(t-1)}) \leq -\eta \sum_{l=1}^{L} \|\nabla W_l L(W^{(t-1)})\|_F^2 \leq -\eta \mu_B^2\mu_B^2\sigma_f^2(X) \cdot (L(W^{(t-1)}) - L(W^*)),
\]

where the second inequality is by Lemma A.1. Applying the inductive hypothesis, we get

\[
L(W^{(t)}) - L(W^*) \leq \left( 1 - \frac{\eta \mu_B^2\mu_B^2\sigma_f^2(X)}{e} \right) \cdot (L(W^{(t-1)}) - L(W^*)) \leq \left( 1 - \frac{\eta \mu_B^2\mu_B^2\sigma_f^2(X)}{e} \right) \cdot (L(W^{(0)}) - L(W^*)),
\]

which completes the proof of the inductive step of Part (ii). Thus we are able to complete the proof.

\( \square \)

### A.2 Proof of Proposition 3.3

**Proof of Proposition 3.3.** We prove the bounds on the singular values and initial training loss separately.

**Bounds on the singular values:** Specifically, we set the neural network width as

\[
m \geq 100 \cdot \left( \sqrt{\max\{d, k\}} + \sqrt{2\log(12/d)} \right)^2
\]

By Corollary 5.35 in Vershynin (2010), we know that for a matrix \( U \in \mathbb{R}^{d_1 \times d_2} \) (\( d_1 \geq d_2 \)) with entries independently generated by standard normal distribution, with probability at least \( 1 - 2\exp(-t^2/2) \), its singular values satisfy

\[
\sqrt{d_1} - \sqrt{d_2} - t \leq \sigma_{\min}(U) \leq \sigma_{\max}(U) \leq \sqrt{d_1} + \sqrt{d_2} + t.
\]
Based on our constructions of $A$ and $B$, we know that each entry of $\frac{1}{\beta}B$ and $\frac{1}{\alpha}A$ follows standard Gaussian distribution. Therefore, set $t = 2\sqrt{\log(12/\delta)}$ and apply union bound, with probability at least $1 - \delta/3$, the following holds,
\[
\alpha(\sqrt{m} - \sqrt{d} - 2\sqrt{\log(12/\delta)}) \leq \sigma_{\min}(A) \leq \sigma_{\max}(A) \leq \alpha(\sqrt{m} + \sqrt{d} + 2\sqrt{\log(12/\delta)})
\]
\[
\beta(\sqrt{m} - \sqrt{k} - 2\sqrt{\log(12/\delta)}) \leq \sigma_{\min}(B) \leq \sigma_{\max}(B) \leq \beta(\sqrt{m} + \sqrt{k} + 2\sqrt{\log(12/\delta)}),
\]
where we use the facts that $\sigma_{\min}(\kappa U) = \kappa \sigma_{\min}(U)$ and $\sigma_{\max}(\kappa U) = \kappa \sigma_{\max}(U)$ for any scalar $\kappa$ and matrix $U$. Then applying our choice of $m$, we have with probability at least $1 - \delta/3$,\[0.9\alpha\sqrt{m} \leq \sigma_{\min}(A) \leq 1.1\alpha\sqrt{m} \quad \text{and} \quad 0.9\beta\sqrt{m} \leq \sigma_{\min}(B) \leq 1.1\beta\sqrt{m}.
\]
This completes the proof of the bounds on the singular values of $A$ and $B$.

**Bounds on the initial training loss:** The proof in this part is similar to the proof of Proposition 6.5 in Du & Hu (2019). Since we apply zero initialization on all hidden layers, by Young’s inequality, we have the following for any $(x, y)$,
\[
\ell(W^{(0)}; x, y) = \frac{1}{2} \|BAx - y\|_2^2 \leq \|BAx\|_F^2 + \|y\|_2^2. \tag{A.2}
\]
Since each entry of $B$ is generated from $\mathcal{N}(0, \beta^2)$, conditioned on $A$, each entry of $BAx$ is distributed according to $\mathcal{N}(0, \beta^2\|Ax\|_2^2)$, so $\frac{\|BAx\|_F^2}{\|Ax\|_2^2}$ follows a $\chi^2_k$ distribution. Applying a standard tail bound for $\chi^2_k$ distribution, we have, with probability at least $1 - \delta'$,
\[
\frac{\|BAx\|_F^2}{\|Ax\|_2^2} \leq \beta^2 k(1 + 2\sqrt{\log(1/\delta')}/k + 2\log(1/\delta')).
\]
Note that by our bounds of the singular values, if $m \geq 100 \cdot (\sqrt{\max[d,k]} + \sqrt{2\log(8/\delta)})^2$, we have with probability at least $1 - \delta/3$, $\|A\|_2 \leq 1.1\alpha\sqrt{m}$, thus, it follows that with probability at least $1 - \delta' - \delta$,
\[
\|BAx\|_F^2 \leq 1.21\alpha^2\beta^2 km\left[1 + 2\sqrt{\log(1/\delta')} + 2\log(1/\delta')\right]\|X\|_F^2.
\]
Then by union bound, it is evident that with probability $1 - n\delta' - \delta/3$,
\[
\|BAx\|_F^2 = \sum_{i=1}^n \|BAx_i\|_2^2 \leq 1.21\alpha^2\beta^2 km\left[1 + 2\sqrt{\log(1/\delta')} + 2\log(1/\delta')\right]\|X\|_F^2.
\]
Set $\delta' = \delta/(3n)$, suppose $\log(1/\delta') \geq 1$, we have with probability at least $1 - 2\delta/3$,
\[
L(W^{(0)}) = \frac{1}{2} \|BAx - y\|_F^2 \leq \|BAx\|_F^2 + \|y\|_F^2 \leq 6.05\alpha^2\beta^2 km \log(2n/\delta)\|X\|_F^2 + \|y\|_F^2.
\]
This completes the proof of the bounds on the initial training loss.

Applying a union bound on these two parts, we are able to complete the proof. \hfill \Box

### A.3 Proof of Corollary 3.4

**Proof of Corollary 3.4.** Recall the condition in Theorem 3.1:
\[
\frac{\sigma_{\min}^2(A)\sigma_{\min}^2(B)}{\|A\|_2\|B\|_2} \geq C \cdot \frac{\|X\|_2}{\sigma_r(X)} \cdot (L(W^{(0)}) - L(W^*))^{1/2}. \tag{A.3}
\]
By Proposition 3.3, we know that, with probability $1 - \delta$,
\[
\frac{\sigma_{\min}^2(A)\sigma_{\min}^2(B)}{\|A\|_2\|B\|_2} = \Theta(m),
\]
\[
\frac{\|X\|_2}{\sigma_r(X)} \cdot (L(W^{(0)}) - L(W^*))^{1/2} = O\left(\frac{(\sqrt{km\log(n/\delta)} + 1)\|X\|_F\|X\|_2}{\sigma_r(X)}\right).
\]
Note that \(|X|_F \leq \sqrt{r} |X|_2\), thus the condition (A.3) can be satisfied if \(m = \Omega(kr^{2/3} \log(n/\delta))\) where \(\kappa = |X|_2^2/\sigma^2(X)\).

Theorem 3.1 implies that \(L(W^{(i)}) - L(W^*) \leq \epsilon\) after \(T = O \left( \frac{1}{\eta L\sigma_{min}^2(A)\sigma_{min}^2(B)\sigma^2(X)} \log \frac{1}{\epsilon} \right)\) iterations. Plugging in the value of \(\eta\), we get

\[
T = O \left( \frac{\|A\|_2\|B\|_2\|X\|_2 \cdot (\sqrt{L(W^{(0)})} + \|A\|_2\|B\|_2\|X\|_2)}{\sigma_{min}^2(A)\sigma_{min}^2(B)\sigma^2(X)} \log \frac{1}{\epsilon} \right).
\]

By Proposition 3.3, we have

\[
T = O \left( \frac{\|X\|_2 \cdot (\sqrt{m \log(n/\delta)}\|X\|_F + \|A\|_2\|B\|_2\|X\|_2)}{m\sigma^2(X)} \log \frac{1}{\epsilon} \right)
\]

\[
= O \left( \frac{\|X\|_2 \cdot (\sqrt{kr \log(n/\delta)/m\|X\|_2 + \|X\|_2})}{\sigma^2(X)} \log \frac{1}{\epsilon} \right)
\]

\[
= O \left( \kappa \log \frac{1}{\epsilon} \right)
\]

for \(m = \Omega(kr \log(n/\delta))\), completing the proof.

**A.4 PROOF OF THEOREM 3.6**

**Proof of Theorem 3.6.** The guarantee is already achieved by \(W^{(0)}\) if \(\epsilon \geq L(W^{(0)}) - L(W^*)\), so we may assume without loss of generality that \(\epsilon < L(W^{(0)}) - L(W^*)\).

Similar to the proof of Theorem 3.1, we use the short-hand notations \(\lambda_A, \mu_A, \lambda_B\) and \(\mu_B\) to denote \(\|A\|_2, \sigma_{min}(A), \|B\|_2\) and \(\sigma_{min}(B)\) respectively. Then we rewrite the condition on \(A\) and \(B\), and our choices of \(\eta\) and \(T\) as follows

\[
\frac{\mu_A^2\mu_B^2}{\lambda_A \lambda_B} \geq \frac{\sqrt{8e^3n\|X\|_2 \cdot \log(L(W^{(0)})/\epsilon')}}{B\sigma^2(X)} \cdot \sqrt{2L(W^{(0)})}
\]

\[
\eta \leq \frac{B\mu_A^2\mu_B^2\sigma^2(X)}{6e^3Ln\lambda_A \lambda_B\|X\|_2} \cdot \min \left\{ \left( \frac{\epsilon'}{\log(2)L(W^{(0)})} \right)^2, \left( \frac{3n\|X\|_2 \cdot \log(T/\delta) \cdot \log(L(W^{(0)})/\epsilon')}{\epsilon'} \right)^2 \right\}
\]

\[
T = \frac{e}{\eta \mu_A^2\mu_B^2\sigma^2(X)} \cdot \log \left( \frac{L(W^{(0)}) - L(W^*)}{\epsilon'} \right)
\]

where we set \(\epsilon' = \epsilon/3\) for the purpose of the proof.

We first prove the convergence guarantees on expectation, and then apply the Markov inequality.

For SGD, our guarantee is not made on the last iterate but the best one. Define \(\mathcal{E}_t\) to be the event that there is no \(s \leq t\) such that \(L(W^{(t)}) - L(W^*) \leq \epsilon\). If \(I(\mathcal{E}_t) = 0\), then there is an iterate \(W_s\) with \(s \leq t\) that achieves training loss within \(\epsilon'\) of optimal.

Similar to the proof of Theorem 3.1, we prove the theorem by induction on the update number \(s\), using the following inductive hypothesis: either \(I(\mathcal{E}_s) = 0\) or the following three inequalities hold,

(i) \(\max_{i \leq s} \|W_i^{(s)}\|_F \leq \sqrt{2e\eta \sigma_{max}(A)\|X\|_2} \cdot \sqrt{2L(W^{(0)})}\)

(ii) \(E[(L(W^{(s)}) - L(W^*))] \leq \left( 1 - \frac{\eta L\sigma_{min}^2(A)\sigma_{min}^2(B)\sigma^2(X)}{\epsilon} \right)^s \cdot (L(W^{(0)}) - L(W^*))\)

(iii) \(L(W^{(s)}) \leq 2L(W^{(0)})\)

where the expectation in Part (ii) is with respect to all of the random choices of minibatches. Clearly, if \(I(\mathcal{E}_s) = 0\), we have already finished the proof since there is an iterate that achieves training loss
within $\epsilon'$ of optimal. Recalling that $\epsilon < L(W^{(0)}) - L(W^*)$, it is easy to verify that the inductive hypothesis holds when $s = 0$.

For the inductive step, we will prove that if the inductive hypothesis holds for $s < t$, then it holds for $s = t$. When $I(\mathcal{E}_{t-1}) = 0$, then $I(\mathcal{E}_t)$ is also 0 and we are done. Therefore, the remaining part is to prove the inductive hypothesis for $s = t$ under the assumption that $I(\mathcal{E}_{t-1}) = 1$, which implies that (i), (ii) and (iii) hold for all $s \leq t - 1$. For Parts (i) and (ii), we will directly prove that the corresponding two inequalities hold. For Part (iii), we will prove that either this inequality holds or $I(\mathcal{E}_t) = 0$.

**Induction for Part (i):** As we mentioned, this part will be proved under the assumption $I(\mathcal{E}_{t-1}) = 1$. Besides, combining Part (i) for $s = t - 1$ and our choice of $\eta$ and $T$ implies that $\max_{t[L]} \|W_t^{(t-1)}\|_F \leq 0.5/L$. Then by triangle inequality, we have the following for $\|W_t^{(t)}\|_F$:

$$\|W_t^{(t)}\|_F \leq \|W_t^{(t-1)}\|_F + \eta \|G_t^{(t-1)}\|_F.$$

By Lemma A.1, we have

$$\|G_t^{(t-1)}\|_F \leq \frac{\sqrt{2e}n\lambda_A\lambda_B \|X\|_2}{B} \cdot \sqrt{L(W_t^{(t-1)})}.$$

Then we have

$$\|W_t^{(t)}\|_F \leq (\|W_t^{(t-1)}\|_F + \eta \|G_t^{(t-1)}\|_F) \leq \|W_t^{(t-1)}\|_F + \frac{\sqrt{2e}n\lambda_A\lambda_B \|X\|_2}{B} \cdot \sqrt{L(W_t^{(t-1)})}. \quad (A.4)$$

By Part (iii) for $s = t - 1$, we know that $L(W_t^{(t-1)}) \leq 2L(W^{(0)})$. Then by Part (i) for $s = t - 1$, it is evident that

$$\|W_t^{(t)}\|_F \leq \frac{\sqrt{2e}n\lambda_A\lambda_B \|X\|_2}{B} \cdot \sqrt{2L(W^{(0)})}. \quad (A.5)$$

This completes the proof of the inductive step of Part (i).

**Induction for Part (ii):** As we previously mentioned, we will prove this part under the assumption $I(\mathcal{E}_{t-1}) = 1$. Thus, as mentioned earlier, the inductive hypothesis implies that $\max_{t[L]} \|W_t^{(t-1)}\|_F \leq 0.5/L$. By Part (i) for $s = t$, which has been verified in (A.5), it can be proved that $\max_{t[L]} \|W_t^{(t)}\|_F \leq 0.5/L$, then we have the following by Lemma A.2,

$$L(W_t^{(t)}) - L(W_t^{(t-1)}) \leq -\eta \sum_{l=1}^L \langle \nabla W_l L(W_t^{(t-1)}), G_t^{(t-1)} \rangle$$

$$+ \eta^2 L\lambda_A\lambda_B \|X\|_2 \cdot \left( \sqrt{eL(W_t^{(t-1)})} + 0.5\epsilon\lambda_A\lambda_B \|X\|_2 \right) \cdot \sum_{l=1}^L \|G_t^{(t-1)}\|_F^2. \quad (A.6)$$

By our condition on $A$ and $B$, it is easy to verify that

$$\lambda_A\lambda_B \geq \frac{\mu_A^2 \mu_B^2}{\lambda_A\lambda_B} \geq \frac{2\sqrt{2e}L(W^{(0)})}{\|X\|_2}.$$

Then by Part (iii) for $s = t - 1$ (A.6) yields

$$L(W_t^{(t)}) - L(W_t^{(t-1)}) \leq -\eta \sum_{l=1}^L \langle \nabla W_l L(W_t^{(t-1)}), G_t^{(t-1)} \rangle + \eta^2 L\lambda_A^2\lambda_B^2 \|X\|_2 \cdot \sum_{l=1}^L \|G_t^{(t-1)}\|_F^2. \quad (A.7)$$

Taking expectation conditioning on $W_t^{(t-1)}$ gives

$$\mathbb{E}[L(W_t^{(t)})W_t^{(t-1)}] - L(W_t^{(t-1)}) \leq -\eta \sum_{l=1}^L \|\nabla W_l L(W_t^{(t-1)})\|_F^2$$

$$+ \eta^2 L\lambda_A^2\lambda_B^2 \|X\|_2 \sum_{l=1}^L \mathbb{E}[[G_t^{(t-1)}]_F^2 | W_t^{(t-1)}]. \quad (A.8)$$
Note that, for \( i \) sampled uniformly from \( \{1, \ldots, n\} \), the expectation \( \mathbb{E}[\|G_i^{(t-1)}\|_F^2 | W^{(t-1)}] \) can be upper bounded by

\[
\mathbb{E}[\|G_i^{(t-1)}\|_F^2 | W^{(t-1)}] = \mathbb{E}[\|G_i^{(t-1)} - \nabla_{W_i} L(W^{(t-1)})\|_F^2 | W^{(t-1)}] + \|\nabla_{W_i} L(W^{(t-1)})\|_F^2 \\
\leq \frac{n^2}{B} \mathbb{E}\left[\|\nabla_{W_i} \ell(W^{(t-1)}; x_i, y_i)\|_F^2 | W^{(t-1)}\right] + \|\nabla_{W_i} L(W^{(t-1)})\|_F^2.
\]

(A.9)

By Lemma A.1, we have

\[
\mathbb{E}[\|\nabla_{W_i} \ell(W^{(t-1)}; x_i, y_i)\|_F^2 | W^{(t-1)}] \leq 2e^2 \lambda_A^2 \lambda_B^2 \mathbb{E}\left[\|x_i\|_2^2 \ell(W^{(t-1)}; x_i, y_i) | W^{(t-1)}\right] \\
\leq \frac{2e^2 \lambda_A^2 \lambda_B^2}{n} \sum_{i=1}^{n} \|x_i\|_2^2 \ell(W^{(t-1)}; x_i, y_i) \\
\leq \frac{2e^2 \lambda_A^2 \lambda_B^2 \|X\|_{2,\infty} L(W^{(t-1)})}{n}.
\]

Plugging the above inequality into (A.9) and (A.8), we get

\[
\mathbb{E}\left[L(W^{(t)}) | W^{(t-1)}\right] - L(W^{(t-1)}) \\
\leq - \eta \sum_{l=1}^{L} \left\|\nabla_{W_i} L(W^{(t-1)})\right\|_F^2 L
\]

\[+ e^2 \lambda_A^2 \lambda_B^2 \|X\|_{2,\infty} L(W^{(t-1)}) + \|\nabla_{W_i} L(W^{(t-1)})\|_F^2 \]

Recalling that \( \eta \leq 1/(6eL^2 \lambda_A^2 \lambda_B^2 \|X\|_2^2) \), we have

\[
\mathbb{E}\left[L(W^{(t)}) | W^{(t-1)}\right] - L(W^{(t-1)}) \leq - \frac{5\eta}{6} \sum_{l=1}^{L} \left\|\nabla_{W_i} L(W^{(t-1)})\right\|_F^2 L
\]

\[+ 2e^2 \lambda_A^2 \lambda_B^2 \|X\|_{2,\infty} L(W^{(t-1)}) \].

(A.10)

By Lemma A.1, we have

\[
\sum_{l=1}^{L} \left\|\nabla_{W_i} L(W^{(t-1)})\right\|_F^2 L \geq 2e^{-1} L \mu_A^2 \mu_B^2 \sigma_i^2(X) \left(L(W^{(t-1)}) - L(W^*)\right).
\]

If we set

\[
\eta \leq \frac{B \mu_A^2 \mu_B^2 \sigma_i^2(X)}{6e^3 L \mu_A^2 \lambda_B^2 \|X\|_{2,\infty}^2},
\]

then (A.10) yields

\[
\mathbb{E}\left[L(W^{(t)}) | W^{(t-1)}\right] - L(W^{(t-1)}) \\
\leq - \frac{5\eta L \mu_A^2 \mu_B^2 \sigma_i^2(X)}{3e} \left(L(W^{(t-1)}) - L(W^*)\right) \]

\[+ \frac{2e^2 \eta^2 L^2 n \lambda_A^4 \lambda_B^4 \|X\|_{2,\infty}^2 \left(L(W^{(t-1)}) - L(W^*)\right)}{B} \]

\[+ \frac{2e^2 \eta^2 L^2 n \lambda_A^4 \lambda_B^4 \|X\|_{2,\infty}^2 L(W^*)}{B} \]

\[\leq - \frac{4\eta L \mu_A^2 \mu_B^2 \sigma_i^2(X)}{3e} \left(L(W^{(t-1)}) - L(W^*)\right) + \frac{2e^2 \eta^2 L^2 n \lambda_A^4 \lambda_B^4 \|X\|_{2,\infty}^2 \|X\|_{2,\infty} L(W^*)}{BL^2}.
\]

(A.12)

Define

\[
\gamma_0 = \frac{4L \mu_A^2 \mu_B^2 \sigma_i^2(X)}{3e}, \quad \text{and} \quad \gamma_1 = \frac{2e^2 \eta^2 L^2 n \lambda_A^4 \lambda_B^4 \|X\|_{2,\infty}^2 \|X\|_{2,\infty} L(W^*)}{B}.
\]
rearranging (A.12) further gives
\[
E[L(W^{(t)})|W^{(t-1)}] - L(W^*) \leq (1 - \eta \gamma_0) \cdot (L(W^{(t)}) - L(W^*)) + \eta^2 \gamma_1. \tag{A.13}
\]
Therefore, setting the step size as
\[
\eta \leq \frac{\eta_0 c \gamma_1}{4},
\]
we further have
\[
E[L(W^{(t)}) - L(W^*)|W^{(t-1)}] \leq \frac{B \mu^2 \rho^2 \sigma_f^2(X)}{6e^3 L \lambda^2 \eta^2 \|X\|^2} \cdot \frac{c'}{L(W^*)},
\]
where the second inequality is by (A.13) and the last inequality is by the fact that we assume \(I(\mathcal{E}_{t-1}) = 1\), which implies that \(L(W^{(t-1)}) - L(W^*) \geq c' \geq 4\gamma_1 \eta / \gamma_0\). Further taking expectation over \(W^{(t-1)}\), we get
\[
E[L(W^{(t)}) - L(W^*)] \leq (1 - 3\gamma_0 / 4) \cdot E[L(W^{(t-1)}) - L(W^*)],
\]
where the second inequality follows from Part (ii) for \(s = t - 1\) and the assumption that \(I(\mathcal{E}_0) = 1\). Plugging the definition of \(\gamma_0\), we are able to complete the proof of the inductive step of Part (ii).

**Induction for Part (iii):** Recalling that for this part, we are going to prove that either \(L(W^{(t)}) \leq 2L(W^{(0)})\) or \(I(\mathcal{E}_t) = 0\), which is equivalent to \(L(W^{(t)}) \cdot I(\mathcal{E}_t) \leq 2L(W^{(0)})\) since \(L(W^{(0)})\) and \(L(W^{(t)})\) are both positive. We will prove this by martingale inequality. Let \(\mathcal{F}_t = \sigma(W^{(0)}, \ldots, W^{(t)})\) be a \(\sigma\)-algebra, and \(\mathcal{F} = \{\mathcal{F}_t\}_{t \geq 1}\) be a filtration. We first prove that \(E[L(W^{(t)}) | I(\mathcal{E}_t), \mathcal{F}_{t-1}] \leq L(W^{(t-1)}) I(\mathcal{E}_{t-1})\). Apparently, this inequality holds when \(I(\mathcal{E}_{t-1}) = 0\) since both sides will be zero. Then if \(I(\mathcal{E}_{t-1}) = 1\), by (A.14) we have
\[
E[L(W^{(t)}) | W^{(t-1)}, I(\mathcal{E}_{t-1}) = 1] \leq E[L(W^{(t)}) | \mathcal{F}_{t-1}, I(\mathcal{E}_{t-1}) = 1] \leq L(W^{(t-1)}).
\]
Combining these two cases, by Jensen’s inequality, we further have
\[
E[\log(L(W^{(t)}) | I(\mathcal{E}_t)) | \mathcal{F}_{t-1}] \leq \log(E[L(W^{(t)}) | I(\mathcal{E}_t), \mathcal{F}_{t-1}]),
\]
which implies that \(\{\log(L(W^{(t)}) | I(\mathcal{E}_t))\}_{t \geq 0}\) is a super-martingale. Then we will upper bound the martingale difference \(\log(L(W^{(t)}) | I(\mathcal{E}_t)) - \log(L(W^{(t-1)}) | I(\mathcal{E}_{t-1}))\). Clearly this quantity would be zero if \(I(\mathcal{E}_{t-1}) = 0\). Then if \(I(\mathcal{E}_{t-1}) = 1\), by (A.7) we have
\[
L(W^{(t)}) \leq L(W^{(t-1)}) + \eta \sum_{l=1}^T \|\nabla_w L(W^{(t-1)})\|_F \|G_l^{(t-1)}\|_F + c n^2 \rho^2 \|L \lambda A \lambda B\|_2 \sum_{l=1}^T \|G_l^{(t-1)}\|_2^2.
\]
By Part (i) for \(s = t - 1\), Lemma A.1, we further have
\[
L(W^{(t)}) \leq \left(1 + \frac{2c \eta \ln \lambda A \lambda B}{B} \|X\|_2^2 + \frac{2c^2 n^2 \rho^2}{B^2} L \lambda A \lambda B \|X\|_2^2 \right) L(W^{(t-1)}) \leq \left(1 + \frac{3c \eta \ln \lambda A \lambda B}{B} \|X\|_2^2 \right) L(W^{(t-1)}),
\]
where the second inequality follows from the choice of \(\eta\) that
\[
\eta \leq \frac{B}{2c n \lambda A \lambda B \|X\|_2^2}.
\]
Using the fact that \(I(\mathcal{E}_t) \leq 1\) and \(I(\mathcal{E}_{t-1}) = 1\), we further have
\[
\log(L(W^{(t)}) | I(\mathcal{E}_t)) \leq \log(L(W^{(t-1)}) | I(\mathcal{E}_{t-1})) + \frac{3c \ln \lambda A \lambda B \|X\|_2^2}{B}.
\]
which also holds for the case \( \mathbb{I}(\mathcal{E}_{t-1}) = 0 \). Recall that \( \{\log(L(W^{(t)}) \cdot \mathbb{I}(\mathcal{E}_t)\}_{t>0} \) is a super-martingale, thus by one-side Azuma’s inequality, we have with probability at least \( 1 - \delta' \),

\[
\log(L(W^{(t)}) \cdot \mathbb{I}(\mathcal{E}_t)) \leq \log(L(W^{(0)})) + \frac{3e\eta Ln\lambda^2_A \lambda^2_B \|X\|_2^2}{B} \cdot \sqrt{2t \log(1/\delta')},
\]

Setting \( \delta' = \delta/T \), using the fact that \( t \leq T \) and leveraging our choice of \( T \) and \( \eta \), we have with probability at least \( 1 - \delta/T \),

\[
\sqrt{T} \eta = \frac{\log(2)B}{3e\sqrt{2 \log(\delta/T) Ln\lambda^2_A \lambda^2_B \|X\|_2^2}},
\]

which implies that

\[
L(W^{(t)}) \mathbb{I}(\mathcal{E}_t) \leq \exp \left[ \log \left( L(W^{(0)}) \right) + \log(2) \right] \leq 2L(W^{(0)}). \tag{A.16}
\]

This completes the proof of the inductive step of Part (iii).

Note that this result holds with probability at least \( 1 - \delta/T \). Thus applying union bound over all iterates \( \{W^{(t)}\}_{t=0,\ldots,T} \) yields that all induction arguments hold for all \( t \leq T \) with probability at least \( 1 - \delta \).

Moreover, plugging our choice of \( T \) and \( \eta \) into Part (ii) gives

\[
\mathbb{E}[L(W^{(t)}) - L(W^*)] \leq \epsilon'.
\]

By Markov inequality, we further have with probability at least \( 2/3 \), it holds that \( [L(W^{(T)}) - L(W^*)] \cdot \mathbb{I}(\mathcal{E}_t) \leq 3\epsilon' = \epsilon \). Therefore, by union bound (together with the high probability arguments of (A.16)) and assuming \( \delta < 1/6 \), we have with probability at least \( 2/3 - \delta > 1/2 \), one of the iterates of SGD can achieve training loss within \( \epsilon' \) of optimal. This completes the proof.

\subsection{A.5 Proof of Corollary 3.7}

\textbf{Proof of Corollary 3.7.} Recall the condition in Theorem 3.6:

\[
\frac{\sigma^2_{\min}(A)\sigma^2_{\min}(B)}{\|A\|_2 \|B\|_2} \geq C \cdot \frac{n\|X\|_2 \cdot \log(L(W^{(0)})/\epsilon)}{B\sigma^2(X)} \cdot \sqrt{L(W^{(0)})}, \tag{A.17}
\]

Then plugging in the results in Proposition 3.3 and the fact that \( \|X\|_F \leq \sqrt{T} \|X\|_2 \), we obtain that condition (A.17) can be satisfied if \( m = O(kr n^2 \log^2(1/\epsilon) \cdot B/n) \).

In addition, consider sufficiently small \( \epsilon \) such that \( \epsilon \leq \tilde{O}(B \|X\|_2^2 \cdot \kappa \|X\|_2^2) \), then and use the fact that \( \|X\|_2 \leq \|X\|_2 \) we have \( \eta = O(kB\epsilon/(Lmnk\|X\|_2^2)) \) based on the results in Proposition 3.3. Then in order to achieve \( \epsilon \)-suboptimal training loss, the iteration complexity is

\[
T = \frac{e}{\eta L\sigma^2_{\min}(A)\sigma^2_{\min}(B) \sigma^2(X)} \log \left( \frac{L(W^{(0)}) - L(W^*)}{\epsilon} \right) = O(\kappa^2 \epsilon^{-1} \log(1/\epsilon) \cdot n/B).
\]

This completes the proof.

\subsection{A.6 Proof of Theorem 3.8}

\textbf{Proof of Theorem 3.8.} Similar to the proof of Theorem 3.6, we set the neural network width and step size as follows,

\[
\frac{\mu_A^2 \mu_B^2}{\lambda_A \lambda_B} \geq \frac{4\sqrt{2e^2 n} \|X\|_2}{B \sigma^2(X)} \cdot \sqrt{L(W^{(0)})}
\]

\[
\eta \leq \frac{\log(2)B^2 \mu_A^2 \mu_B^2 (A) \sigma^2(X)}{54e^3 Ln^3 \lambda_A \lambda_B \|X\|_2^4 \cdot \log(T/\delta)},
\]

where \( \lambda_A, \mu_A, \lambda_B \) and \( \mu_B \) denote \( \|A\|_2, \sigma_{\min}(A), \|B\|_2 \) and \( \sigma_{\min}(B) \) respectively.

Different from the proof of Theorem 3.6, the convergence guarantee established in this regime is made on the last iterate of SGD, rather than the best one. Besides, we will prove the theorem by induction on the update parameter \( t \), using the following two-part inductive hypothesis:
(i) \( \max_{i \in [L]} \| W^{(t)}_i \|_F \leq 0.5/L \)

(ii) \( L(W^{(t)}) \leq 2L(W^{(0)}) \cdot \left( 1 - \frac{\eta L \mu^2 \sigma^2(X)}{3\epsilon} \right)^s \).

**Induction for Part (i)** We first prove that \( \max_{i \in [L]} \| W^{(t)}_i \|_F \leq 0.5/L \). By triangle inequality and the update rule of SGD, we have

\[
\| W^{(t)}_i \|_F \leq \sum_{s=0}^{t-1} \eta \| G_i \|_F \\
\leq \eta \sum_{s=0}^{t-1} \frac{\sqrt{2} \epsilon n \lambda_A \lambda_B \| X \|_2}{B} (L(W^{(s)}) - L(W^*))^{1/2} \\
\leq \frac{\sqrt{2} \epsilon n \lambda_A \lambda_B \| X \|_2}{B} \cdot (L(W^{(0)}) - L(W^*))^{1/2} \cdot \sum_{s=0}^{t-1} \left( 1 - \frac{\eta L \mu^2 \sigma^2(X)}{2\epsilon} \right)^s \\
\leq \frac{\sqrt{2} \epsilon n \lambda_A \lambda_B \| X \|_2}{BL \mu^2 \sigma^2(X)} \cdot (L(W^{(0)}) - L(W^*))^{1/2}
\]

where the second inequality is by Lemma A.1, and the third inequality follows from Part (ii) for all \( s < t \) and the fact that \( (1 - x)^{1/2} \leq 1 - x/2 \) for all \( x \in [0, 1] \). Then applying our choice of \( m \) implies that \( \| W^{(t)}_i \|_F \leq 0.5/L \).

**Induction for Part (ii)** Similar to Part (ii) and (iii) of the induction step in the proof of Theorem 3.6, we first prove the convergence in expectation, and then use Azuma’s inequality to get the high-probability based results. It can be simply verified that

\[
\lambda_A \lambda_B \geq \frac{\mu^2_B}{\lambda_A \lambda_B} \geq 4\sqrt{2} 4n \lambda_A \lambda_B \| X \|_2 \cdot \log \left( \frac{L(W^{(0)})}{\epsilon} \right) \cdot \sqrt{2L(W^{(0)})} \geq 2\sqrt{2e^{-1}L(W^{(0)})} \\
\eta \leq \frac{\log(2) B \mu^2_B \sigma^2(X)}{96 \epsilon^3 L \lambda_A \lambda_B \| X \|_2^2 \cdot \log(\delta)} \leq \frac{B \mu^2_B \sigma^2(X)}{6\epsilon^3 L \lambda_A \lambda_B \| X \|_2^2 \cdot \log(\delta)}.
\]

Thus, we can leverage (A.12) and obtain

\[
E[L(W^{(t)}) | W^{(t-1)}] - L(W^{(t-1)}) \leq \frac{4\eta L \mu^2_B \sigma^2(X)}{3\epsilon} L(W^{(t-1)}),
\]

where we use the fact that \( L(W^*) = 0 \). Then by Jensen’s inequality, we have

\[
E[\log L(W^{(t)}) | W^{(t-1)}] \leq \log L(W^{(t-1)}) + \log \left( 1 - \frac{4\eta L \mu^2_B \sigma^2(X)}{3\epsilon} \right),
\]

\[
\leq \log L(W^{(t-1)}) - \frac{4\eta L \mu^2_B \sigma^2(X)}{3\epsilon},
\]

where the second inequality is by log(1 + x) ≤ x. Then similar to the proof of Theorem 3.6, we are going to apply martingale inequality to prove this part. Let \( \mathcal{F}_t = \sigma(W^{(0)}, \ldots, W^{(t)}) \) be a \( \sigma \)-algebra, and \( \mathcal{F} = \{ \mathcal{F}_t \}_{t \geq 1} \) be a filtration, the above inequality implies that

\[
E[\log L(W^{(t)}) | \mathcal{F}_{t-1}] \leq \frac{4\eta L \mu^2_B \sigma^2(X)}{3\epsilon},
\]

which implies that \( \{ \log L(W^{(t)}) + 4\eta L \mu^2_B \sigma^2(X)/(3\epsilon) \} \) is a super-martingale. Besides, by (A.15), we can obtain

\[
\log L(W^{(t)}) \leq \log L(W^{(t-1)}) + \frac{3\epsilon n \lambda_A \lambda_B \| X \|_2^2}{B},
\]

which implies that

\[
\log L(W^{(t)}) + \frac{4\eta L \mu^2_B \sigma^2(X)}{3\epsilon} \leq \log L(W^{(t-1)}) + \frac{4(\epsilon - 1) \eta L \mu^2_B \sigma^2(X)}{3\epsilon} + 4\epsilon n \lambda_A \lambda_B \| X \|_2^2.
\]
where we again use the fact that $\log(1 + x) \leq x$. Thus, by the one-sided Azuma’s inequality we have with probability at least $1 - \delta'$ that

$$
\log \left( L(W^{(t)}) \right) \leq \log \left( L(W^{(0)}) \right) - \frac{4\eta L \mu_A^2 \mu_B^2 \sigma_f^2(X)}{3e} + \frac{4\eta L n^2 \lambda_A^2 \lambda_B^2 \|X\|_2^2}{e} \cdot \sqrt{2t \log(1/\delta')}
$$

$$
\leq \log \left( L(W^{(0)}) \right) - \frac{t\eta L \mu_A^2 \mu_B^2 \sigma_f^2(X)}{e} + \frac{96e^3 \eta L n^2 \lambda_A^4 \lambda_B^4 \|X\|_2^2 \log(1/\delta')}{B^2 \mu_A^2 \mu_B^2 \sigma_f^2(X)}
$$

$$
\leq \log \left( L(W^{(0)}) \right) - \frac{t\eta L \mu_A^2 \mu_B^2 \sigma_f^2(X)}{e} + \log(2),
$$

where the second inequality follows from the fact that $-at + b\sqrt{t} \leq b^2/a$, and the last inequality is by our choice of $\eta$ that

$$
\eta \leq \frac{\log(2) B^2 \mu_A^2 \mu_B^2 \sigma_f^2(X)}{96e^3 L n^2 \lambda_A^4 \lambda_B^4 \|X\|_2^2 \cdot \log(1/\delta')}.
$$

Then it is clear that with probability at least $1 - \delta'$,

$$
L(W^{(t)}) \leq 2L(W^{(0)}) \cdot \exp \left( -\frac{t\eta L \mu_A^2 \mu_B^2 \sigma_f^2(X)}{e} \right),
$$

(A.19)

which completes the induction for Part (ii).

Similar to the proof of Theorem 3.6, (A.19) holds with probability at least $1 - \delta'$ for a given $t$. Then we can set $\delta' = \delta/T$ and apply union bound such that with probability at least $1 - \delta$, (A.19) holds for all $t \leq T$. This completes the proof.

A.7 PROOF OF COROLLARY 3.9

Proof of Corollary 3.9. Recall the condition in Theorem 3.8:

$$
\frac{\sigma_{\text{min}}^2(A) \sigma_{\text{min}}^2(B)}{\|A\|_2 \|B\|_2} \geq C \cdot \frac{n\|X\|_2}{B \sigma_f^2(X)} \cdot \sqrt{L(W^{(0)})},
$$

(A.20)

Then plugging in the results in Proposition 3.3 and the fact that $\|X\|_F \leq \sqrt{n} \|X\|_2$, we obtain that condition (A.17) can be satisfied if $m = O(kn^2 \cdot B/n)$.

In addition, it can be computed that $\eta = O(kB^2/(Lmn^2 \kappa \|X\|_2^2))$ based on the results in Proposition 3.3. Then in order to achieve $\epsilon$-suboptimal training loss, the iteration complexity is

$$
T = \frac{e}{\eta L \sigma_{\text{min}}^2 \sigma_{\text{min}}^2(B) \sigma_f^2(X)} \log \left( \frac{L(W^{(0)}) - L(W^*)}{\epsilon} \right) = O(k^2 \log(1/\epsilon) \cdot n^2 / B^2).
$$

This completes the proof.

B PROOFS OF TECHNICAL LEMMAS

B.1 PROOF OF LEMMA A.1

We first note the following useful lemmas.

Lemma B.1 (Claim B.1 in Du & Hu (2019)). Define $\Phi = \arg \min_{\Theta \in \mathbb{R}^{k \times d}} \|\Theta X - Y\|_F^2$, then for any $U \in \mathbb{R}^{k \times d}$ it holds that

$$
\|UX - Y\|_F^2 = \|UX - \Phi X\|_F^2 + \|\Phi X - Y\|_F^2.
$$

Lemma B.2 (Theorem 1 in Fang et al. (1994)). Let $U, V \in \mathbb{R}^{d \times d}$ be two positive definite matrices, then it holds that

$$
\lambda_{\text{min}}(U) \text{Tr}(V) \leq \text{Tr}(UV) \leq \lambda_{\text{max}}(U) \text{Tr}(V).
$$

The following lemma is proved in Section B.3.
Lemma B.3. Let $U \in \mathbb{R}^{d \times r}$ be a rank-$r$ matrix. Then for any $V \in \mathbb{R}^{r \times k}$, it holds that

$$\sigma_{\min}(U)\|V\|_F \leq \|UV\|_F \leq \sigma_{\max}(U)\|V\|_F.$$  

Proof of Lemma A.1. Proof of gradient lower bound: We first prove the gradient lower bound. Let $U = B(1 + W_L) \ldots (1 + W_i)A$, by Lemma B.1 and the definition of $L(W^*)$, we know that there exist a matrix $\Phi \in \mathbb{R}^{k \times d}$ such that

$$L(W) = \frac{1}{2}\|UX - \Phi X\|_F^2 + L(W^*).$$  

(B.1)

Therefore, based on the assumption that $\max_{l \in [L]} \|W_l\|_F \leq 0.5/L$, we have

$$\|\nabla_{W_l}L(W)\|_F^2 = \|B(1 + W_L) \ldots (1 + W_i)\|_F^2 (UX - \Phi X)(I + W_l) \ldots AX \|_F^2 \geq \sigma_{\min}^2((I + W_L) \ldots (I + W_i)) \cdot \sigma_{\min}^2((I + W_l) \ldots (I + W_i)) \cdot \|B^T(U - \Phi)XX^TA^\top\|_F^2 \geq (1 - 0.5/L)2L-2\|B^T(U - \Phi)XX^TA^\top\|_F^2,$$

where the last inequality follows from the fact that $\sigma_{\min}(I + W_i) \geq 1 - \|W_i\|_2 \geq 1 - \|W_i\|_F \geq 1 - 0.5/L$. Applying Lemma B.2, we get

$$\|B^T(U - \Phi)XX^TA^\top\|_F^2 = \text{Tr}(BB^\top(U - \Phi)XX^TA^\topXX^T(U - \Phi)^\top) \geq \lambda_{\min}(BB^\top) \cdot \text{Tr}(A^\topXX^T(U - \Phi)^\top(U - \Phi)XX^T) \geq \lambda_{\min}(BB^\top) \cdot \lambda_{\min}(A^\topA) \cdot \|U - \Phi\|_F^2.$$  

Note that $X$ is of $r$-rank, thus there exists a full-rank matrix $\hat{X} \in \mathbb{R}^{d \times r}$ such that $\hat{X}X^\top = XX^\top$. Thus we have

$$\|(U - \Phi)X\|_F^2 = \text{Tr}((U - \Phi)XX^T(U - \Phi)^\top) = \text{Tr}((U - \Phi)\hat{X}X^\top(U - \Phi)^\top) = \|(U - \Phi)\hat{X}\|_F^2.$$  

Therefore,

$$\|(U - \Phi)XX^\top\|_F^2 = \|(U - \Phi)\hat{X}X^\top\|_F^2 = \text{Tr}((U - \Phi)\hat{X}X^\top(U - \Phi)^\top) \geq \lambda_{\min}(\hat{X}^\top\hat{X}) \cdot \|(U - \Phi)\hat{X}\|_F^2.$$

(B.2)

where the inequality follows from Lemma B.2 and the last equality follows from (B.2), (B.1) and the fact that $\lambda_{\min}(\hat{X}^\top\hat{X}) = \lambda_r(XX^\top) = \sigma_r^2(X)$. Note that we assume $d, k \leq m$ and $d \leq n$. Thus it follows that $\lambda_{\min}(BB^\top) = \sigma_{\min}^2(B)$ and $\lambda_{\min}(A^\topA) = \sigma_{\min}^2(A)$. Then putting everything together, we can obtain

$$\|\nabla_{W_l}L(W)\|_F^2 \geq 2\sigma_{\min}^2(B)\sigma_{\min}^2(A)\sigma_r^2(X)(1 - 0.5/L)2L-2(L(W - L(W^*))$$

Then using the inequality $(1 - 0.5/L)^{2L-2} \geq e^{-1}$, we are able to complete the proof of gradient lower bound.

Proof of gradient upper bound: The gradient upper bound can be proved in a similar way. Specifically, Lemma B.3 implies

$$\|\nabla_{W_l}L(W)\|_F^2 = \|B(1 + W_L) \ldots (1 + W_i)\|_F^2 (UX - \Phi X)(I + W_l) \ldots AX \|_F^2 \leq \sigma_{\max}^2((I + W_L) \ldots (I + W_i)) \cdot \sigma_{\max}^2((I + W_l) \ldots (I + W_i)) \cdot \|B^T(U - \Phi)XX^TA^\top\|_F^2 \leq \|B\|_2^2 \|A\|_2^2 \cdot \|(U - \Phi)XX^\top\|_F^2 \leq (1 + 0.5/L)^{2L-2}\|B\|_2^2 \|A\|_2^2 \cdot \|(U - \Phi)XX^\top\|_F^2.$$
where the last inequality is by the assumption that \(\max_{i \in [L]} \|W_i\|_F \leq 0.5/L\). By (B.3), we have
\[
\|(U - \Phi)XX^T\|_F^2 = \|(U - \Phi)(XX^T)^{1/2}(XX^T)^{1/2}\|_F^2 \\
\leq \lambda_{\max}(XX^T) \cdot \|(U - \Phi)(XX^T)^{1/2}\|_F^2 \\
= \lambda_{\max}(XX^T) \cdot \|\nabla_xiw\|_F^2 \\
= 2\|X\|_2 \cdot (L(W) - L(W^*)),
\]
where the inequality is by Lemma B.3 and the second equality is by (B.2). Therefore, combining the above results yields
\[
\|\nabla_iw, L(W)\|_F^2 \leq 2\sigma_{\max}(B)\sigma_{\max}(A)\|X\|_2^2 (1 + 0.5/L)^{2L-2} (L(W) - L(W^*)).
\]
Using the inequality \((1 + 0.5/L)^{2L-2} \leq (1 + 0.5/L)^{2L} \leq e\), we are able to complete the proof of gradient upper bound.

**Proof of the upper bound of stochastic gradient:** Define by \(B\) the set of training data points used to compute the stochastic gradient, then define by \(X\) and \(Y\) the stacking of \(\{x_i\}_{i \in B}\) and \(\{y_i\}_{i \in B}\) respectively. Let \(U = B(I + W_L) \cdots (I + W_1)A\), the minibatch stochastic gradient takes form
\[
\nabla_{W_i} \ell(W; x_i, y_i) = \left[(B(I + W_L) \cdots (I + W_{l+1}))^\top (Ux_i - y_i) \right](I + W_{l-1}) \cdots A x_i]^\top.
\]
Therefore, by Lemma B.3, we have
\[
\|\nabla_{W_i} \ell(W; x_i, y_i)\|_F^2 \leq \sigma_{\max}^2 \left((I + W_L) \cdot (I + W_{l+1}) \cdots (I + W_1)\right) \cdot \sigma_{\max}^2 \left((I + W_{l-1}) \cdots (I + W_1)\right) \\
\cdot \|X\|_2 \cdot \|\nabla_xiw\|_F^2 \\
\leq (1 + 0.5/L)^{2L-2} \cdot \|B\|_2 \cdot \|A\|_2 \cdot \|X\|_2 \cdot \|Ux_i - y_i\|_F^2 \\
\leq 2e \cdot \|A\|_2 \cdot \|B\|_2 \cdot \|x_i\|_2 \cdot \ell(W; x_i, y_i),
\]
where the last inequality is by the fact that \((1 + 0.5/L)^{2L-2} \leq e\).

**Proof of the upper bound of stochastic gradient:** Define by \(B\) the set of training data points used to compute the stochastic gradient, then define by \(X\) and \(Y\) the stacking of \(\{x_i\}_{i \in B}\) and \(\{y_i\}_{i \in B}\) respectively. Let \(U = B(I + W_L) \cdots (I + W_1)A\), the minibatch stochastic gradient takes form
\[
G_l = \frac{n}{B} \sum_{i \in B} \nabla_{W_i} \ell(W; x_i, y_i) \\
= \frac{n}{B} \left[(B(I + W_L) \cdots (I + W_{l+1}))^\top (Ux_i - y_i) \right]
(I + W_{l-1}) \cdots A x_i]^\top.
\]
Then by Lemma B.3, we have
\[
\|G_l\|_F^2 \leq \frac{n^2}{B^2} \cdot \sigma_{\max}^2 \left((I + W_L) \cdot (I + W_{l+1}) \cdots (I + W_1)\right) \cdot \sigma_{\max}^2 \left((I + W_{l-1}) \cdots (I + W_1)\right) \\
\cdot \|X\|_2 \cdot \|\nabla_xiw\|_F^2 \\
\leq \frac{n^2}{B^2} \cdot \|B\|_2 \cdot (1 + 0.5/L)^{2L-2} \cdot \|B\|_2 \cdot \|A\|_2 \cdot \|X\|_2 \cdot \|Ux - Y\|_F^2 \\
\leq \frac{en^2}{B^2} \cdot \|B\|_2 \cdot \|A\|_2 \cdot \|X\|_2 \cdot \|Ux - Y\|_F^2.
\]
where the second inequality is by the assumptions that \(\max_{i \in [L]} \|W_i\|_F \leq 0.5/L\), and the last inequality follows from the the fact that \((1 + 0.5/L)^{2L-2} \leq (1 + 0.5/L)^{2L} \leq e\). Note that \(X\) and \(Y\) are constructed by stacking \(B\) columns from \(X\) and \(Y\) respectively, thus we have \(\|X\|_2^2 \leq \|X\|_2^2\) and \(\|Ux - Y\|_F^2 \leq \|Ux - Y\|_F^2 = 2L(W)\). Then it follows that
\[
\|G_l\|_F^2 \leq \frac{2en^2}{B^2} \cdot \|B\|_2 \cdot \|A\|_2 \cdot \|X\|_2 \cdot L(W).
\]
This completes the proof of the upper bound of stochastic gradient.
B.2 Proof of Lemma A.2

Proof of Lemma A.2. Let $U = B(I + W_L) \cdots (I + W_1)A$ and $\tilde{U} = B(I + \tilde{W}_L) \cdots (I + \tilde{W}_1)A$ and $\Delta = \tilde{U} - U$. We have

\[
L(\tilde{W}) - L(W) = \frac{1}{2}(\|\tilde{U}X - Y\|_F^2 - \|UX - Y\|_F^2) = \frac{1}{2}(\|U + \Delta\|X - Y\|_F^2 - \|UX - Y\|_F^2) = \frac{1}{2}(\|UX - Y + \Delta X\|_F^2 - \|UX - Y\|_F^2) = \frac{1}{2}(\|2(UX - Y, \Delta X) + \|\Delta X\|_F^2) = \langle UX - Y, (\tilde{U} - U)X \rangle + \frac{1}{2}\|U - \tilde{U}\|X\|_F^2. \tag{B.4}
\]

We begin by working on the first term. Let $V = (I + W_L) \cdots (I + W_1)$ and $\tilde{V} = (I + \tilde{W}_L) \cdots (I + \tilde{W}_1)$, so that $\tilde{U} - U = B(\tilde{V} - V)A$. Breaking down the effect of transforming $V = \prod_{j=L}^1 (I + W_j)$ into $\tilde{V} = \prod_{j=L}^1 (I + \tilde{W}_j)$ into the effects of replacing one layer at a time, we get

\[
\tilde{V} - V = \prod_{l=1}^L \left[ \left( \prod_{j=L}^{l+1} (I + W_j) \right) \left( \prod_{j=L}^{l} (I + \tilde{W}_j) \right) \right]
\]

and, for each $l$, pulling out a common factor of $\left( \prod_{j=L}^{l+1} (I + W_j) \right) \left( \prod_{j=L}^{l-1} (I + \tilde{W}_j) \right)$ gives

\[
\tilde{V} - V = \sum_{l=1}^L (I + W_L) \cdots (I + W_{l+1})(\tilde{W}_l - W_l)(I + \tilde{W}_{l-1}) \cdots (I + \tilde{W}_1) + \sum_{l=1}^L (I + W_L) \cdots (I + W_{l+1})(\tilde{W}_l - W_l)
\]

\[
\cdot \left[ (I + \tilde{W}_{l-1}) \cdots (I + \tilde{W}_1) - (I + W_{l-1}) \cdots (I + W_1) \right]. \tag{B.5}
\]
The first term $V_1$ satisfies

$$
\langle UX - Y, BV_1AX \rangle 
= \left\langle UX - Y, B \left( \sum_{l=1}^{L} (I + W_L) \cdots (I + W_{l+1})(\widehat{W}_l - W_l)(I + W_{l-1}) \cdots (I + W_1) \right) AX \right\rangle
$$

$$
= \sum_{l=1}^{L} \left\langle UX - Y, B(I + W_L) \cdots (I + W_{l+1})(\widehat{W}_l - W_l)(I + W_{l-1}) \cdots (I + W_1)AX \right\rangle
$$

$$
= \sum_{l=1}^{L} \text{Tr}((UX - Y)\top B(I + W_L) \cdots (I + W_{l+1})(\widehat{W}_l - W_l)(I + W_{l-1}) \cdots (I + W_1)AX)
$$

$$
= \sum_{l=1}^{L} \text{Tr}((I + W_{l-1}) \cdots (I + W_1)AX(UX - Y)\top B(I + W_L) \cdots (I + W_{l+1})(\widehat{W}_l - W_l))
$$

$$
= \sum_{l=1}^{L} \left\langle [B(I + W_L) \cdots (I + W_{l+1})]\top (UX - Y)[(I + W_{l-1}) \cdots AX]\top, \widehat{W}_l - W_l \right\rangle
$$

$$
= \sum_{l=1}^{L} \langle \nabla_w L(W), \widehat{W}_l - W_l \rangle,
$$

where the first equality is by the definition of $V_1$. Now we focus on the second term $V_2$ of (B.5),

$$
V_2 = \sum_{l=1}^{L} (I + W_L) \cdots (I + W_{l+1})(\widehat{W}_l - W_l)
$$

$$
\cdot \sum_{s=1}^{l-1} (I + W_{s-1}) \cdots (I + W_{s+1})(\widehat{W}_s - W_s)(I + \widehat{W}_{s+1}) \cdots (I + \widehat{W}_1).
$$

Recalling that $\|W_l\|_F, \|\widehat{W}_l\|_F \leq 0.5/L$ for all $l \in [L]$, by triangle inequality we have

$$
\|V_2\|_F \leq (1 + 0.5/L)^L \cdot \sum_{l,s \in [L]: l > s} |\widehat{W}_l - W_l|_F \cdot |\widehat{W}_s - W_s|_F
$$

$$
\leq (1 + 0.5/L)^L \cdot \left( \sum_{l=1}^{L} |\widehat{W}_l - W_l|_F \right)^2,
$$

where we use the fact that $\sum_{l,s \in [L]: l > s} a_l a_s \leq \sum_{l \in [L]} a_l a_s = (\sum_l a_l)^2$ holds for all $a_1, \ldots, a_L \geq 0$. Therefore, the following holds regarding $V_2$:

$$
\langle UX - Y, BV_2AX \rangle \leq \|UX - Y\|_F \|BV_2AX\|_F
$$

$$
\leq \sqrt{2L(W)} \|B\|_2 \|A\|_2 \|X\|_2 \|V_2\|_F
$$

$$
\leq \sqrt{2c} \sqrt{L(W)} \|B\|_2 \|A\|_2 \|X\|_2 \left( \sum_{l=1}^{L} |\widehat{W}_l - W_l|_F \right)^2
$$

(B.7)

where the third inequality follows from the fact that $(1 + 0.5/L)^L = (1 + 0.5/L)^L \leq \sqrt{e}$. Next, we are going to upper bound the second term of (B.4): $\frac{1}{c}(\langle U - \bar{U}, X \rangle)^2 \leq \|B(V - \nu)AX\|_F^2 \leq \|A\|_2 \|B\|_2 \|X\|_2 \|\bar{V} - X \|_F^2$, it suffices to bound the norm $\|\bar{V} - X \|_F^2$. By (B.5), we have

$$
\|ar{V} - V\|_F = \left\| \sum_{l=1}^{L} (I + W_L) \cdots (I + W_{l+1})(\widehat{W}_l - W_l)(I + \widehat{W}_{l-1}) \cdots (I + \widehat{W}_1) \right\|_F
$$

$$
\leq (1 + 0.5/L)^L \sum_{l=1}^{L} |\widehat{W}_l - W_l|_F.
$$

(B.8)
Plugging (B.6), (B.7) and (B.8) into (B.4), we have
\begin{equation}
\begin{aligned}
L(\hat{\mathbf{W}}) - L(\mathbf{W})
&= \langle \mathbf{U} \mathbf{X} - \mathbf{Y}, \mathbf{B} (\mathbf{V}_1 + \mathbf{V}_2) \mathbf{X} \rangle + \frac{1}{2} \mathbf{B} (\hat{\mathbf{V}} - \mathbf{V}) \mathbf{A} \mathbf{X} \|_F^2 \\
&\leq \sum_{l=1}^L \langle \nabla_{\mathbf{W}_l} L(\mathbf{W}), \hat{\mathbf{W}}_l - \mathbf{W}_l \rangle \\
&\quad + \| \mathbf{A} \|_2 \| \mathbf{B} \|_2 \| \mathbf{X} \|_2 \left( \sqrt{2e L(\mathbf{W})} + 0.5 e \mathbf{A} \|_2 \| \mathbf{B} \|_2 \| \mathbf{X} \|_2 \right) \left( \sum_{l=1}^L \| \hat{\mathbf{W}}_l - \mathbf{W}_l \|_F \right)^2 \\
&\leq \sum_{l=1}^L \langle \nabla_{\mathbf{W}_l} L(\mathbf{W}), \hat{\mathbf{W}}_l - \mathbf{W}_l \rangle \\
&\quad + L \| \mathbf{A} \|_2 \| \mathbf{B} \|_2 \| \mathbf{X} \|_2 \left( \sqrt{2e L(\mathbf{W})} + 0.5 e \mathbf{A} \|_2 \| \mathbf{B} \|_2 \| \mathbf{X} \|_2 \right) \sum_{l=1}^L \| \hat{\mathbf{W}}_l - \mathbf{W}_l \|_F^2,
\end{aligned}
\end{equation}

(B.9)

where the last inequality is by Jesen’s inequality. This completes the proof.

\hfill \Box

B.3 Proof of Lemma B.3

Proof of Lemma B.3. Note that we have
\begin{equation}
\| \mathbf{U} \mathbf{V} \|_F^2 = \text{Tr}(\mathbf{U} \mathbf{V} \mathbf{V}^\top \mathbf{U}^\top) = \text{Tr}(\mathbf{U}^\top \mathbf{U} \mathbf{V} \mathbf{V}^\top).
\end{equation}

By Lemma B.2, it is clear that
\begin{equation}
\lambda_{\text{min}}(\mathbf{U}^\top \mathbf{U}) \text{Tr}(\mathbf{V} \mathbf{V}^\top) \leq \text{Tr}(\mathbf{U}^\top \mathbf{U} \mathbf{V} \mathbf{V}^\top) \leq \lambda_{\text{max}}(\mathbf{U}^\top \mathbf{U}) \text{Tr}(\mathbf{V} \mathbf{V}^\top).
\end{equation}

Since \( \mathbf{U} \in \mathbb{R}^{d \times r} \) is of \( r \)-rank, thus we have \( \lambda_{\text{min}}(\mathbf{U}^\top \mathbf{U}) = \sigma_{\text{min}}^2(\mathbf{U}) \). Then applying the facts that \( \lambda_{\text{max}}(\mathbf{U}^\top \mathbf{U}) = \sigma_{\text{max}}^2(\mathbf{U}) \) and \( \text{Tr}(\mathbf{V} \mathbf{V}^\top) = \| \mathbf{V} \|_F^2 \), we are able to complete the proof.

\hfill \Box