Stochastic Gradient Descent Learns State Equations with Nonlinear Activations

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Abstract

We study discrete time dynamical systems governed by the state equation $h_{t+1} = \phi(Ah_t + Bu_t)$. Here A, B are weight matrices, ϕ is an activation function, and u_t is the input data. This relation is the backbone of recurrent neural networks (e.g. LSTMs) which have broad applications in sequential learning tasks. We utilize stochastic gradient descent to learn the weight matrices from a finite input/state trajectory $\{u_t, h_t\}_{t=0}^N$. We prove that SGD estimate linearly converges to the ground truth weights while using near-optimal sample size. Our results apply to increasing activations whose derivatives are bounded away from zero. The analysis is based on i) a novel SGD convergence result with nonlinear activations and ii) careful statistical characterization of the state vector. Numerical experiments verify the fast convergence of SGD on ReLU and leaky ReLU in consistence with our

Keywords: state equation, dynamical systems, sample complexity, stochastic gradient descent

1. Introduction

A wide range of problems involve sequential data with a natural temporal ordering. Examples include natural language processing, time series prediction, system identification, and control design, among others. State-of-the-art algorithms for sequential problems often stem from dynamical systems theory and are tailored to learn from temporally dependent data. Linear models and algorithms; such as Kalman filter, PID controller, and linear dynamical systems, have a long history and are utilized in control theory since 1960's with great success (Brown et al. (1992); Ho and Kalman (1966); Aström and Hägglund (1995)). More recently, nonlinear models such as recurrent neural networks (RNN) found applications in complex tasks such as machine translation and speech recognition (Bahdanau et al. (2014); Graves et al. (2013); Hochreiter and Schmidhuber (1997)). Unlike feedforward neural networks, RNNs are dynamical systems that use their internal state to process inputs. The goal of this work is to shed light on the inner workings of RNNs from a theoretical point of view. In particular, we focus on the RNN state equation which is characterized by a nonlinear activation function ϕ , state weight matrix A, and input weight matrix B as follows

$$\boldsymbol{h}_{t+1} = \phi(\boldsymbol{A}\boldsymbol{h}_t + \boldsymbol{B}\boldsymbol{u}_t), \tag{1.1}$$

Here h_t is the state vector and u_t is the input data at timestamp t. This equation is the source of dynamic behavior of RNNs and distinguishes RNN from feedforward networks. The weight matrices A and B govern the dynamics of the state equation and are inferred from data. We will explore the statistical and computational efficiency of stochastic gradient descent (SGD) for learning these weight matrices.

Contributions: Suppose we are given a finite trajectory of input/state pairs $(u_t, h_t)_{t=0}^N$ generated from the state equation (1.1). We consider a least-squares regression obtained from N equations; with inputs $(u_t, h_t)_{t=1}^N$ and outputs $(h_{t+1})_{t=1}^N$. For a class of activation functions including leaky ReLU and for stable systems¹, we show that SGD linearly converges to the ground truth weight matrices while requiring near-optimal trajectory length N. In particular, the required sample size is $\mathcal{O}(n+p)$ where n and p are the dimensions of the state and input vectors respectively. The results are extended to unstable systems when the samples are collected from multiple independent RNN trajectories rather than a single trajectory. Our theory applies to increasing activation functions whose derivatives are bounded away from zero, which includes leaky ReLU, and (sub)gaussian input data. Numerical experiments on ReLU and leaky ReLU corroborate our theory and demonstrate that SGD converges faster as the activation slope increases. To obtain our results, we i) characterize the statistical properties of the state vector (e.g. well-conditioned covariance) and ii) derive a novel SGD convergence result with nonlinear activations; which may be of independent interest. As a whole, this paper provides a step towards foundational understanding of RNN training via SGD.

1.1. Related Work

Our work is related to the recent optimization and statistics literature on linear dynamical systems (LDS) and neural networks.

Linear dynamical systems: The state-equation (1.1) reduces to a LDS when ϕ is the linear activation $(\phi(x) = x)$. Identifying the weight matrices is a core problem in linear system identification and is related to the optimal control problem (e.g. linear quadratic regulator) with unknown system dynamics. While these problems are studied since 1950's (Ljung (1998, 1987); Åström and Eykhoff (1971)), our work is closer to the recent literature that provides data dependent bounds and characterize the non-asymptotic learning performance. Recht and coauthors have a series of papers exploring optimal control problem (Simchowitz et al. (2018); Tu et al. (2018, 2017)). In particular, Hardt et al. (2016) shows gradient descent learns single-input-single-output (SISO) LDS with polynomial guarantees. Oymak and Ozay (2018) and Faradonbeh et al. (2018) provide sample complexity bounds for learning LDS. Sanandaji et al. (2011b,a); Pereira et al. (2010) study the identification of sparse systems.

Neural networks: There is a growing literature on the theoretical aspects of deep learning and provable algorithms for training neural networks. Most of the existing results are concerned with feedforward networks. Ge et al. (2017); Li and Yuan (2017); Mei et al. (2018b); Soltanolkotabi (2017); Janzamin et al. (2015); Zhong et al. (2017b) consider learning fully-connected shallow networks with gradient descent. Mei et al. (2018a); Soltanolkotabi et al. (2017); Foster et al. (2018) analyze empirical landscape of related nonlinear learning problems. Brutzkus and Globerson (2017); Zhong et al. (2017a); Du et al. (2017); Goel et al. (2018) address convolutional neural networks; which is an efficient weight-sharing architecture. Several works (Ji and Telgarsky (2018); Oymak and Soltanolkotabi (2019); Allen-Zhu et al. (2018); Chizat and Bach (2018); Belkin et al. (2018); Li et al. (2019)) consider the training problem when the network is over-parametrized and study the over-fitting ability of such networks. Baldi and Vershynin (2019); Bartlett et al. (2017); Neyshabur et al. (2018, 2017) studies the capacity and generalization ability of deep networks. For recurrent networks, Sedghi and Anandkumar (2016) proposed tensor algorithms with polynomial guarantees and Khrulkov et al. (2017) studied their expressive power. More recently, Miller and Hardt (2018) showed that stable RNNs can be approximated by feed-forward networks.

¹Throughout this work, a system is called stable if the spectral norm of the state matrix A is less than 1.

Algorithm 1 Learning state equations with nonlinear activations

- 1: **Inputs:** $(y_t, h_t, u_t)_{t=1}^N$ sampled from a trajectory. Scaling μ , learning rate η . Initialization A_0, B_0 .
- 2: **Outputs:** Estimates \hat{A} , \hat{B} of the weight matrices A, B.
- 3: $\boldsymbol{x}_t \leftarrow [\mu \boldsymbol{h}_t^T \ \boldsymbol{u}_t^T]^T$ for $1 \leq t \leq N$.
- 4: $\Theta_0 \leftarrow [\mu^{-1} A_0 B_0]$
- 5: **For** τ from 1 to END
- 6: Pick γ_{τ} from $\{1, 2, \dots, N\}$ uniformly at random.
- 7: $\mathbf{\Theta}_{\tau} \leftarrow \mathbf{\Theta}_{\tau-1} \eta \nabla \mathcal{L}_{\gamma_{\tau}}(\mathbf{\Theta}_{\tau-1})$
- 8: EndFor
- 9: **return** $[\hat{A} \ \hat{B}] \leftarrow \Theta_{\text{END}} \begin{bmatrix} \mu I_n & 0 \\ 0 & I_p \end{bmatrix}$.

2. Problem Setup

We first introduce the notation. $\|\cdot\|$ returns the spectral norm of a matrix and $s_{\min}(\cdot)$ returns the minimum singular value. The activation $\phi:\mathbb{R}\to\mathbb{R}$ applies entry-wise if its input is a vector. Throughout, ϕ is assumed to be a 1-Lipschitz function. With proper scaling of its parameters, the system (1.1) with a Lipschitz activation can be transformed into a system with 1-Lipschitz activation. The functions $\Sigma[\cdot]$ and $\operatorname{var}[\cdot]$ return the covariance of a random vector and variance of a random variable respectively. I_n is the identity matrix of size $n \times n$. Normal distribution with mean μ and covariance Σ is denoted by $\mathcal{N}(\mu, \Sigma)$. Throughout, c, C, c_0, c_1, \ldots denote positive absolute constants.

Setup: We consider the dynamical system parametrized by an activation function $\phi(\cdot)$ and weight matrices $A \in \mathbb{R}^{n \times n}$, $B \in \mathbb{R}^{n \times p}$ as described in (1.1). Here, h_t is the n dimensional state-vector and u_t is the p dimensional input to the system at time t. As mentioned previously, (1.1) corresponds to the state equation of a recurrent neural network. For most RNNs of interest, the state h_t is hidden and we only get to interact with h_t via an additional output equation. For Elman networks Elman (1990), this equation is characterized by some output activation ϕ_y and output weights C, D as follows

$$y_t = \phi_y(Ch_t + Du_t). \tag{2.1}$$

In this work, our attention is restricted to the state equation (1.1); which corresponds to setting $y_t = h_{t+1}$ in the output equation. To analyze (1.1) in a non-asymptotic data-dependent setup, we assume a finite input/state trajectory of $\{u_t, h_t\}_{t=0}^N$ generated by some ground truth weight matrices (A, B). Our goal is learning the unknown weights A and B in a data and computationally efficient way. In essence, we will show that, if the trajectory length satisfies $N \gtrsim n + p$, SGD can quickly and provably accomplish this goal using a constant step size.

Appoach: Our approach is described in Algorithm 1. It takes two hyperparameters; the scaling factor μ and learning rate η . Using the RNN trajectory, we construct N triples of the form $\{u_t, h_t, h_{t+1}\}_{t=1}^N$. We formulate a regression problem by defining the output vector y_t , input vector x_t , and the target parameter C as follows

$$y_t = h_{t+1}$$
 , $x_t = \begin{bmatrix} \mu h_t \\ u_t \end{bmatrix} \in \mathbb{R}^{n+p}$, $C = [\mu^{-1} A B] \in \mathbb{R}^{n \times (n+p)}$. (2.2)

With this reparameterization, we find the input/output identity $y_t = \phi(Cx_t)$. We will consider the least-squares regression given by

$$\mathcal{L}(\mathbf{\Theta}) = \frac{1}{N} \sum_{t=1}^{N} \mathcal{L}_t(\mathbf{\Theta}) \quad \text{where} \quad \mathcal{L}_t(\mathbf{\Theta}) = \frac{1}{2} \| \mathbf{y}_t - \phi(\mathbf{\Theta} \mathbf{x}_t) \|_{\ell_2}^2. \tag{2.3}$$

For learning the ground truth parameter C, we utilize SGD on the loss function (2.3) with a constant learning rate η . Starting from an initial point Θ_0 , after END SGD iterations, Algorithm 1 returns an estimate $\hat{C} = \Theta_{\text{END}}$. Estimates of A and B are decoded from the left and right submatrices of \hat{C} respectively.

3. Main Results

3.1. Preliminaries

The analysis of the state equation naturally depends on the choice of the activation function; which is the source of nonlinearity. We first define a class of Lipschitz and increasing activation functions.

Definition 1 (β -increasing activation) Given $1 \ge \beta \ge 0$, the activation function ϕ satisfies $\phi(0) = 0$ and $1 \ge \phi'(x) \ge \beta$ for all $x \in \mathbb{R}$.

Our results will apply to strictly increasing activations where ϕ is β -increasing for some $\beta>0$. Observe that, this excludes ReLU activation which has zero derivative for negative values. However, it includes Leaky ReLU which is a generalization of ReLU and is frequently utilized in practical deep learning models He et al. (2015). Parameterized by $1 \geq \beta \geq 0$, Leaky ReLU is a β -increasing function given by

$$LReLU(x) = \max(\beta x, x). \tag{3.1}$$

In general, given an increasing and 1-Lipschitz activation ϕ , a β -increasing function ϕ_{β} can be obtained by blending ϕ with the linear activation, i.e. $\phi_{\beta}(x) = (1 - \beta)\phi(x) + \beta x$.

A critical property that enables SGD is that the state-vector covariance $\Sigma[h_t]$ is well-conditioned under proper assumptions. The lemma below provides upper and lower bounds on this covariance matrix in terms of problem variables.

Lemma 2 (State vector covariance) Consider the state equation (1.1) where $h_0 = 0$ and $u_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \mathbf{I}_p)$. Define the upper bound term B_t as

$$B_t = \|\mathbf{B}\| \sqrt{\frac{1 - \|\mathbf{A}\|^{2t}}{1 - \|\mathbf{A}\|^2}}.$$
 (3.2)

- Suppose ϕ is 1-Lipschitz and $\phi(0) = 0$. Then, for all $t \geq 0$, $\Sigma[\mathbf{h}_t] \leq B_t^2 \mathbf{I}_n$.
- Suppose ϕ is a β -increasing function and $p \geq n$. Then, $\Sigma[h_t] \succeq \beta^2 s_{\min}(B)^2 I_n$.

As a natural extension from linear dynamical systems, we will say the system is stable if ||A|| < 1 and unstable otherwise. For activations we consider, stability implies that if the input is set to 0, state vector h_t will exponentially converge to 0 i.e. the system forgets the past states quickly. This is also the reason $(B_t)_{t\geq 0}$ sequence converges for stable systems and diverges otherwise. The condition

number of the covariance will play a critical role in our analysis. Using Lemma 2, this number can be upper bounded by ρ defined as

$$\rho = \left(\frac{B_{\infty}}{\beta s_{\min}(\mathbf{B})}\right)^2 = \left(\frac{\|\mathbf{B}\|}{s_{\min}(\mathbf{B})}\right)^2 \frac{1}{\beta^2 (1 - \|\mathbf{A}\|^2)}.$$
 (3.3)

Observe that, the condition number of B appears inside the ρ term.

3.2. Learning from Single Trajectory

Our main result applies to stable systems ($\|A\| < 1$) and provides a non-asymptotic convergence guarantee for SGD in terms of the upper bound on the state vector covariance. This result characterizes the sample complexity and the rate of convergence of SGD; and also provides insights into the role of activation function and the spectral norm of A.

Theorem 3 (Main result) Let $\{u_t, h_{t+1}\}_{t=1}^N$ be a finite trajectory generated from the state equation (1.1). Suppose $\|A\| < 1$, ϕ is β -increasing, $h_0 = 0$, $p \ge n$, and $u_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, I_p)$. Let ρ be same as (3.3) and c, C, c_0 be properly chosen absolute constants. Pick the trajectory length N to satisfy

$$N \ge CL\rho^2(n+p),$$

where $L=1-\frac{\log(cn\rho)}{\log\|A\|}$. Pick scaling $\mu=1/B_{\infty}$, learning rate $\eta=c_0\frac{\beta^2}{\rho n(n+p)}$, and consider the loss function (2.3). With probability $1-4N\exp(-100n)-8L\exp(-\mathcal{O}(\frac{N}{L\rho^2}))$, starting from an initial point Θ_0 , for all $\tau\geq 0$, the SGD iterations described in Algorithm 1 satisfies

$$\mathbb{E}[\|\mathbf{\Theta}_{\tau} - \mathbf{C}\|_{F}^{2}] \le (1 - c_{0} \frac{\beta^{4}}{2\rho^{2} n(n+p)})^{\tau} \|\mathbf{\Theta}_{0} - \mathbf{C}\|_{F}^{2}.$$
(3.4)

Here the expectation is over the randomness of the SGD updates.

Sample complexity: Theorem 3 essentially requires $N \gtrsim (n+p)/\beta^4$ samples for learning. This can be seen by unpacking (3.3) and ignoring the logarithmic L term and the condition number of B. Observe that $\mathcal{O}(n+p)$ growth achieves near-optimal sample size for our problem. Each state equation (1.1) consists of n sub-equations (one for each entry of h_{t+1}). We collect N state equations to obtain a system of Nn equations. On the other hand, the total number of unknown parameters in A and B are n(n+p). This implies Theorem 3 is applicable as soon as the problem is mildly overdetermined i.e. $Nn \gtrsim n(n+p)$.

Input distribution: As shown in Theorem 6, the results can be extended to other distributions as long as one can show that the covariance matrix of the state vector is well-conditioned. For Gaussian input distribution, we are indeed able to prove this hence we opted to state a comprehensive result under this model. Finally, we believe well-conditioned covariance assumption can be guaranteed by adding a small Gaussian excitation to arbitrary input sequences which is left as a future direction.

Computational complexity: Theorem 3 requires $\mathcal{O}(n(n+p)\log\frac{1}{\varepsilon})$ iterations to reach ε -neighborhood of the ground truth. Our analysis reveals that, this rate can be accelerated if the state vector is zeromean. This happens for odd activation functions satisfying $\phi(-x) = -\phi(x)$ (e.g. linear activation). The result below is a corollary and requires $\times n$ less iterations.

Theorem 4 (Faster learning for odd activations) Consider the same setup provided in Theorem 3. Additionally, assume that ϕ is an odd function. Pick scaling $\mu = 1/B_{\infty}$, learning rate $\eta = c_0 \frac{\beta^2}{\rho(n+p)}$, and consider the loss function (2.3). With probability $1 - 4N \exp(-100n) - 8L \exp(-\mathcal{O}(\frac{N}{L\rho^2}))$, starting from an initial point Θ_0 , for all $\tau \geq 0$, the SGD iterations described in Algorithm 1 satisfies

$$\mathbb{E}[\|\mathbf{\Theta}_{\tau} - \mathbf{C}\|_{F}^{2}] \le (1 - c_{0} \frac{\beta^{4}}{2\rho^{2}(n+p)})^{\tau} \|\mathbf{\Theta}_{0} - \mathbf{C}\|_{F}^{2}, \tag{3.5}$$

where the expectation is over the randomness of the SGD updates.

Another aspect of the convergence rate is the dependence on β . In terms of β , the SGD error (3.4) decays as $(1 - \mathcal{O}(\beta^8))^{\tau}$. While it is not clear how optimal is the exponent 8, numerical experiments in Section 6 demonstrate that larger β indeed results in drastically faster convergence.

4. Main Ideas and Proof Strategy

We first outline our high-level proof strategy for Theorem 3; which brings together ideas from statistics and optimization.

- 1. We first show that input data is well-behaved by proving that state-vector h_t has a well-conditioned covariance as discussed in Lemma 2 and shown in Appendix B. The key idea is if ϕ is β -increasing, then the random input data u_t provides sufficient excitation for the output state h_{t+1} .
- 2. Even if individual samples are well-behaved, analyzing (2.3) is still challenging due to temporal dependencies between the samples. These dependencies prevent us from directly using statistical learning results that typically assume i.i.d. samples. We show that the dependency between samples at time t and t+T decay exponentially fast in separation T (for stable systems). This is outlined in Appendix C.
- 3. This observation allows us to obtain nearly independent data by subsampling the original trajectory to get $(h_{iT}, u_{iT})_{i \geq 0}$. Thanks to exponential decay, a logarithmically small T can be chosen to generate large subtrajectories of size N/T. Appendix D uses additional perturbation arguments to establish the well-behavedness of the overall data matrix.
- 4. To conclude, we obtain a deterministic result which establishes fast convergence result for β -increasing activations and well-behaved dataset. This is provided in Theorem 5 and proved in Appendix A.

The first three steps are related to the statistical nature of the problem which can be decoupled from the last step. Specifically, the last step derives a deterministic result that establishes the linear convergence of SGD for β -increasing functions. For linear convergence proofs, a typical strategy is showing the *strong convexity* of the loss function i.e. showing that, for some $\alpha > 0$ and all points $\boldsymbol{v}, \boldsymbol{u}$, the gradient satisfies

$$\langle \nabla \mathcal{L}(\boldsymbol{v}) - \nabla \mathcal{L}(\boldsymbol{u}), \boldsymbol{v} - \boldsymbol{u} \rangle \ge \alpha \|\boldsymbol{v} - \boldsymbol{u}\|_{\ell_2}^2.$$

The core idea of our convergence result is that the strong convexity parameter of the loss function with β -increasing activations can be connected to the loss function with *linear activations*. In particular, recalling (2.3), set $y_t^{\text{lin}} = Cx_t$ and define the linear loss to be

$$\mathcal{L}^{ ext{lin}}(oldsymbol{\Theta}) = rac{1}{2N} \sum_{i=1}^N \|oldsymbol{y}_t^{ ext{lin}} - oldsymbol{\Theta} oldsymbol{x}_t\|_{\ell_2}^2.$$

Denoting the strong convexity parameter of the original loss by α_{ϕ} and that of linear loss by α_{lin} , we argue that $\alpha_{\phi} \geq \beta^2 \alpha_{\text{lin}}$; which allows us to establish a convergence result as soon as α_{lin} is strictly positive. Next result is our SGD convergence theorem which follows from this discussion.

Theorem 5 (Deterministic convergence) Suppose a data set $\{x_i, y_i\}_{i=1}^N$ is given; where output y_i is related to input x_i via $y_i = \phi(\langle x_i, \theta \rangle)$ for some $\theta \in \mathbb{R}^n$. Suppose $\beta > 0$ and ϕ is a β -increasing. Let $\gamma_+ \geq \gamma_- > 0$ be scalars. Assume that input samples satisfy the bounds

$$\gamma_{+} \boldsymbol{I}_{n} \succeq \frac{1}{N} \sum_{i=1}^{N} \boldsymbol{x}_{i} \boldsymbol{x}_{i}^{T} \succeq \gamma_{-} \boldsymbol{I}_{n} \quad , \quad \|\boldsymbol{x}_{i}\|_{\ell_{2}}^{2} \leq B \, \textit{for all } i.$$

Let $\{r_{\tau}\}_{\tau=0}^{\infty}$ be a sequence of i.i.d. integers uniformly distributed between 1 to N. Then, starting from an arbitrary point θ_0 , setting learning rate $\eta=\frac{\beta^2\gamma_-}{\gamma_+B}$, for all $\tau\geq 0$, the SGD iterations for quadratic loss

$$\boldsymbol{\theta}_{\tau+1} = \boldsymbol{\theta}_{\tau} - \eta(\phi(\boldsymbol{x}_{r_{\tau}}^T \boldsymbol{\theta}_{\tau}) - \boldsymbol{y}_{r_{\tau}})\phi'(\boldsymbol{x}_{r_{\tau}}^T \boldsymbol{\theta}_{\tau})\boldsymbol{x}_{r_{\tau}}, \tag{4.1}$$

satisfies the error bound

$$\mathbb{E}[\|\boldsymbol{\theta}_{\tau} - \boldsymbol{\theta}\|_{\ell_{2}}^{2}] \leq \|\boldsymbol{\theta}_{0} - \boldsymbol{\theta}\|_{\ell_{2}}^{2} (1 - \frac{\beta^{4} \gamma_{-}^{2}}{\gamma_{+} B})^{\tau}, \tag{4.2}$$

where the expectation is over the random selection of the SGD iterations $\{r_{\tau}\}_{\tau=0}^{\infty}$.

This theorem provides a clean convergence rate for SGD for β -increasing activations and naturally generalizes standard results on linear regression which corresponds to $\beta=1$. We remark that related results appear in the literature on generalized linear models. Kakade et al. (2011); Foster et al. (2018); Mei et al. (2018a) provide learning theoretic loss/gradient/hessian convergence results for isotonic regression, robust regression, and β -increasing activations. Goel et al. (2018) establishes a similar result for leaky ReLU activations under the assumption of symmetric input distribution and infinitely many samples (i.e. in population limit). Compared to these, we establish a *deterministic* linear convergence guarantee for SGD that works whenever the data matrix is full rank. We believe extensions to proximal gradient methods might be beneficial for high-dimensional nonlinear problems (e.g. sparse/low-rank approximation, manifold constraints Cai et al. (2010); Beck and Teboulle (2009); Oymak et al. (2018); Agarwal et al. (2010); Pereira et al. (2010)) and is left as a future work.

To derive our main results in Section 3, we need to address the first three steps outlined earlier and determine the conditions under which Theorem 5 is applicable to the data obtained from RNN state equation with high probability. Below we provide desirable characteristics of the state vector; which enables our statistical results.

Assumption 1 (Well-behaved state vector) *Let* L > 1 *be an integer. There exists positive scalars* $\gamma_+, \gamma_-, \theta$ *and an absolute constant* C > 0 *such that* $\theta \le 3\sqrt{n}$ *and the following holds*

- Lower bound: $\Sigma[h_{L-1}] \succeq \gamma_- I_n$,
- **Upper bound:** for all t, the state vector satisfies

$$\Sigma[\mathbf{h}_t] \leq \gamma_+ \mathbf{I}_n$$
, $\|\mathbf{h}_t - \mathbb{E}[\mathbf{h}_t]\|_{\psi_2} \leq C\sqrt{\gamma_+}$ and $\|\mathbb{E}[\mathbf{h}_t]\|_{\ell_2} \leq \theta\sqrt{\gamma_+}$. (4.3)

Here $\|\cdot\|_{\psi_2}$ returns the subgaussian norm of a vector (see Def. 5.22 of Vershynin (2010)).

Assumption 1 ensures that covariance is well-conditioned, state vector is well-concentrated, and it has a reasonably small expectation. Our next theorem establishes statistical guarantees for learning the RNN state equation based on this assumption.

Theorem 6 (General result) Let $\{u_t, h_{t+1}\}_{t=1}^N$ be a length N trajectory of the state equation (1.1). Suppose $\|A\| < 1$, ϕ is β -increasing, $h_0 = 0$, and $u_t \stackrel{i.i.d.}{\sim} u$ where $u \in \mathbb{R}^p$ is zero-mean with identity covariance and subgaussian norm bounded above by a constant. Given scalars $\gamma_+ \geq \gamma_- > 0$, set the condition number as $\rho = \gamma_+/\gamma_-$. For absolute constants $C, c, c_0 > 0$, choose trajectory length N to satisfy

$$N \ge CL\rho^2(n+p)$$
 where $L = \lceil 1 - \frac{\log(cn\rho)}{\log\|\mathbf{A}\|} \rceil$.

Suppose Assumption 1 holds with $L, \gamma_+, \gamma_-, \theta$. Pick scaling to be $\mu = 1/\sqrt{\gamma_+}$ and learning rate to be $\eta = c_0 \frac{\beta^2}{\rho(\theta + \sqrt{2})^2(n+p)}$. With probability $1 - 4N \exp(-100n) - 8L \exp(-\mathcal{O}(\frac{N}{L\rho^2}))$, starting from Θ_0 , for all $\tau \geq 0$, the SGD iterations on loss (2.3) as described in Algorithm 1 satisfies

$$\mathbb{E}[\|\boldsymbol{\Theta}_{\tau} - \boldsymbol{C}\|_{F}^{2}] \leq (1 - c_{0} \frac{\beta^{4}}{2\rho^{2}(\theta + \sqrt{2})^{2}(n+p)})^{\tau} \|\boldsymbol{\Theta}_{0} - \boldsymbol{C}\|_{F}^{2}, \tag{4.4}$$

where the expectation is over the randomness of SGD updates.

The advantage of this theorem is that, it isolates the optimization problem from the statistical properties of state vector. If one can prove tighter bounds on achievable $(\gamma_+, \gamma_-, \theta)$, it will immediately imply improved performance for SGD. In particular, Theorems 3 and 4 are simple corollaries of Theorem 6 with proper choices when input is normally distributed i.e. $u_t \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}(0, I_p)$.

- Theorem 3 follows by setting $\gamma_+ = B_{\infty}^2$, $\gamma_- = \beta^2 s_{\min}(\mathbf{B})^2$, and $\theta = \sqrt{n}$.
- Theorem 4 follows by setting $\gamma_+=B_\infty^2, \gamma_-=\beta^2 s_{\min}({\pmb B})^2,$ and $\theta=0.$

5. Learning Unstable Systems

So far, we considered learning from a single RNN trajectory for stable systems (||A|| < 1). For such systems, as the time goes on, the impact of the earlier states disappear. In our analysis, this allows us to split a single trajectory into multiple nearly-independent trajectories. This approach will not work for unstable systems (A is arbitrary) where the impact of older states may be amplified over time. To address this, we consider a model where the data is sampled from multiple independent trajectories.

Suppose N independent trajectories of the state-equation (1.1) are available. Pick some integer $T_0 \geq 1$. Denoting the ith trajectory by the triple $(\boldsymbol{h}_{t+1}^{(i)}, \boldsymbol{h}_{t}^{(i)}, \boldsymbol{u}_{t}^{(i)})_{t \geq 0}$, we collect a single sample from each trajectory at time T_0 to obtain the triple $(\boldsymbol{h}_{T_0+1}^{(i)}, \boldsymbol{h}_{T_0}^{(i)}, \boldsymbol{u}_{T_0}^{(i)})$. To utilize the existing optimization framework (2.3); for $1 \leq i \leq N$, we set,

$$(\mathbf{y}_i, \mathbf{h}_i, \mathbf{u}_i) = (\mathbf{h}_{T_0+1}^{(i)}, \mathbf{h}_{T_0}^{(i)}, \mathbf{u}_{T_0}^{(i)}).$$
 (5.1)

With this setup, we can again use the SGD Algorithm 1 to learn the weights A and B. The crucial difference compared to Section 3 is that, the samples $(y_i, h_i, u_i)_{i=1}^N$ are now independent of each other; hence, the analysis is simplified. As previously, having an upper bound on the condition number of the state-vector covariance is critical. This upper bound can be shown to be ρ defined as

$$\rho = \begin{cases} \bar{\rho} & \text{if } n > 1\\ \bar{\rho} \frac{1 - \beta^2 |\mathbf{A}|^2}{1 - (\beta |\mathbf{A}|)^{2T_0}} & \text{if } n = 1 \end{cases} \quad \text{where} \quad \bar{\rho} = \frac{B_{T_0}^2}{\beta^2 s_{\min}(\mathbf{B})^2}. \tag{5.2}$$

The $\bar{\rho}$ term is similar to the earlier definition (3.3); however it involves B_{T_0} rather than B_{∞} . This modification is indeed necessary since $B_{\infty} = \infty$ when $\|A\| > 1$. On the other hand, note that, $B_{T_0}^2$ grows proportional to $\|A\|^{2T_0}$; which results in exponentially bad condition number in T_0 . Our ρ definition remedies this issue for single-output systems; where n=1 and A is a scalar. In particular, when $\beta=1$ (e.g. ϕ is linear) ρ becomes equal to the correct value 1^2 . The next theorem provides our result on unstable systems in terms of this condition number and other model parameters.

Theorem 7 (Unstable systems) Suppose we are given N independent trajectories $(\boldsymbol{h}_t^{(i)}, \boldsymbol{u}_t^{(i)})_{t \geq 0}$ for $1 \leq i \leq N$. Each trajectory is sampled at time T_0 to obtain N samples $(\boldsymbol{y}_i, \boldsymbol{h}_i, \boldsymbol{u}_i)_{i=1}^N$ where the ith sample is given by (5.1). Suppose the sample size satisfies

$$N \ge C\rho^2(n+p)$$

where ρ is given by (5.2). Assume the initial states are 0, ϕ is β -increasing, $p \geq n$, and $\mathbf{u}_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \mathbf{I}_p)$. Set scaling $\mu = 1/\sqrt{B_{T_0}}$, learning rate $\eta = c_0 \frac{\beta^2}{\rho n(n+p)}$, and run SGD over the equations described in (2.2) and (2.3). Starting from Θ_0 , with probability $1 - 2N \exp(-100(n+p)) - 4 \exp(-\mathcal{O}(\frac{N}{\sigma^2}))$, all SGD iterations satisfy

$$\mathbb{E}[\|\boldsymbol{\Theta}_{\tau} - \boldsymbol{C}\|_{F}^{2}] \leq (1 - c_{0} \frac{\beta^{4}}{2\rho^{2} n(n+p)})^{\tau} \|\boldsymbol{\Theta}_{0} - \boldsymbol{C}\|_{F}^{2},$$

where the expectation is over the randomness of the SGD updates.

6. Numerical Experiments

We conducted experiments on ReLU and Leaky ReLU activations. Let us first describe the experimental setup. We pick the state dimension n=50 and the input dimension p=100. We choose the ground truth matrix \boldsymbol{A} to be a scaled random unitary matrix; which ensures that all singular values of \boldsymbol{A} are equal. \boldsymbol{B} is generated with i.i.d. $\mathcal{N}(0,1)$ entries. Instead of using the theoretical

²Clearly, any nonzero 1×1 covariance matrix has condition number 1. However, due to subtleties in the proof strategy, we don't use $\rho = 1$ for $\beta < 1$. Obtaining tighter bounds on the subgaussian norm of the state-vector would help resolve this issue.

Algorithm 2 Empirical hyperparameter selection.

- 1: **Inputs:** $(\boldsymbol{h}_t, \boldsymbol{u}_t)_{t=1}^N$ sampled from a trajectory.
- 2: Outputs: Scaling μ .
- 3: Form the empirical covariance matrix Σ_h from $\{\boldsymbol{h}_t\}_{t=1}^N$. 4: Form the empirical covariance matrix Σ_u from $\{\boldsymbol{u}_t\}_{t=1}^N$.
- 5: **return** $\sqrt{\|\Sigma_u\|/\|\Sigma_h\|}$.

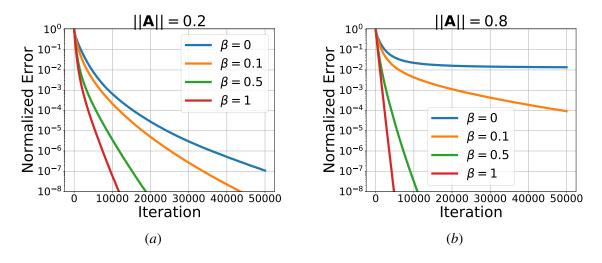


Figure 1: SGD convergence behavior for Leaky ReLUs with varying minimum slope β . Figures (a) and (b) repeat the same experiments. The difference is the spectral norm of the ground truth state matrix A.

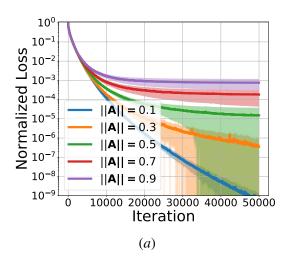
scaling choice, we determine the scaling μ from empirical covariance matrices outlined in Algorithm 2. Similar to our proof strategy, this algorithm equalizes the spectral norms of the input and state covariances to speed up convergence. We also empirically determined the learning rate and used $\eta = 1/100$ in all experiments.

Evaluation: We consider two performance measures in the experiments. Let \hat{C} be an estimate of the ground truth parameter $C = [\mu^{-1} A B]$. The first measure is the normalized error defined as $\|\hat{C} - C\|_F^2 / \|C\|_F^2$. The second measure is the normalized loss defined as

$$\frac{\sum_{i=1}^{N} \|\boldsymbol{y}_{t} - \phi(\hat{\boldsymbol{C}}\boldsymbol{x}_{t})\|_{\ell_{2}}^{2}}{\sum_{i=1}^{N} \|\boldsymbol{y}_{t}\|_{\ell_{2}}^{2}}.$$

In all experiments, we run Algorithm 1 for 50000 SGD iterations and plot these measures as a function of τ ; by using the estimate available at the end of the τ th SGD iteration for $0 \le \tau \le 50000$. Each curve is obtained by averaging the outcomes of 20 independent realizations. Our first experiments use N = 500; which is mildly larger than the total dimension n + p = 150. In Figure 1, we plot the Leaky ReLU errors with varying slopes as described in (3.1). Here $\beta = 0$ corresponds to ReLU and $\beta = 1$ is the linear model. In consistence with our theory, SGD achieves linear convergence and as β increases, the rate of convergence drastically improves³. The improvement is more visible for

³Note that convergence becomes faster for larger β under the realizable model i.e. there exists a ground truth state equation with activation slope β that can fit the observed trajectory. This is consistent with the technical setup our results are



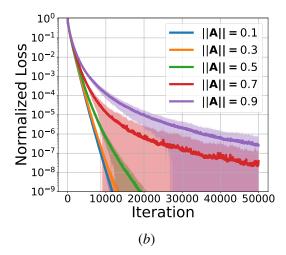


Figure 2: SGD convergence behavior for ReLU with varying spectral norm of the state matrix A. Figures a) and b) repeats the same experiments. The difference is that a) uses N=500 trajectory length whereas b) uses N=2500 (i.e. $\times 5$ more data). Shaded regions highlight the one standard deviation around the mean.

less stable systems driven by A with a larger spectral norm. In particular, while ReLU converges for small ||A||, SGD gets stuck before reaching the ground truth when ||A|| = 0.8.

To understand, how well SGD fits the training data, in Figure 2a, we plotted the normalized loss for ReLU activation. For more unstable system ($\|A\|=0.9$), training loss stagnates in a similar fashion to the parameter error. We also verified that the norm of the overall gradient $\|\nabla \mathcal{L}(\Theta_{\tau})\|_F$ continues to decay (where Θ_{τ} is the τ th SGD iterate); which implies that SGD converges before reaching a global minima. As A becomes more stable, rate of convergence improves and linear rate is visible. Finally, to better understand the population landscape of the quadratic loss with ReLU activations, Figure 2b repeats the same ReLU experiments while increasing the sample size five times to N=2500. For this more overdetermined problem, SGD converges even for $\|A\|=0.9$; indicating that

- population landscape of loss with ReLU activation is well-behaved,
- however ReLU problem requires more data compared to the Leaky ReLU for finding global minima (for our planted setup).

Overall, as predicted by our theory, experiments verify that SGD indeed quickly finds the optimal weight matrices of the state equation (1.1) and as the activation slope β increases, the convergence rate improves.

7. Conclusions

This work showed that SGD can learn the nonlinear dynamical system (1.1); which is characterized by weight matrices and an activation function. This problem is of interest for recurrent neural networks as well as nonlinear system identification. We showed that efficient learning is possible with optimal sample complexity and good computational performance. Our results apply to strictly

proven. Also note that the data distribution in the experiments changes with the activation slope β . If the dataset is fixed and not realizable, the results may be different as we vary the slope β .

increasing activations such as Leaky ReLU. We empirically showed that Leaky ReLU converges faster than ReLU and requires less samples; in consistence with our theory. We list a few unanswered problems that would provide further insights into recurrent neural networks.

- Covariance of the state-vector: Our results depend on the covariance of the state-vector and requires it to be positive definite. One might be able to improve the current bounds on the condition number and relax the assumptions on the activation function. Deriving similar performance bounds for ReLU is particularly interesting.
- **Hidden state:** For RNNs, the state vector is hidden and is observed through an additional equation (2.1); which further complicates the optimization landscape. Even for linear dynamical systems, learning the (A, B, C, D) system ((1.1), (2.1)) is a non-trivial task Ho and Kalman (1966); Hardt et al. (2016). What can be said when we add the nonlinear activations?
- Classification task: In this work, we used normally distributed input and least-squares regression for our theoretical guarantees. More realistic input distributions might provide better insight into contemporary problems, such as natural language processing; where the goal is closer to classification (e.g. finding the best translation from another language).

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Appendix A. Deterministic Convergence Result for SGD

Proof [Proof of Theorem 5] Given two distinct scalars a,b; define $\phi'(a,b) = \frac{\phi(a) - \phi(b)}{a - b}$. $\phi'(a,b) \ge \beta$ since ϕ is β -increasing. Define \boldsymbol{w}_{τ} to be the residual $\boldsymbol{w}_{\tau} = \boldsymbol{\theta}_{\tau} - \boldsymbol{\theta}$. Observing

$$\phi(\boldsymbol{x}_{r_{\tau}}^T \boldsymbol{\theta}_{\tau}) - \boldsymbol{y}_{r_{\tau}} = \phi'(\boldsymbol{x}_{r_{\tau}}^T \boldsymbol{\theta}_{\tau}, \boldsymbol{x}_{r_{\tau}}^T \boldsymbol{\theta}) \boldsymbol{x}_{r_{\tau}}^T \boldsymbol{w}_{\tau},$$

the SGD recursion obeys

$$\begin{aligned} \|\boldsymbol{w}_{\tau+1}\|_{\ell_{2}}^{2} &= \|\boldsymbol{w}_{\tau} - \eta(\phi(\boldsymbol{x}_{r_{\tau}}^{T}\boldsymbol{\theta}_{\tau}) - \boldsymbol{y}_{r_{\tau}})\phi'(\boldsymbol{x}_{r_{\tau}}^{T}\boldsymbol{\theta}_{\tau})\boldsymbol{x}_{r_{\tau}}\|_{\ell_{2}}^{2}. \\ &= \|\boldsymbol{w}_{\tau} - \eta\boldsymbol{x}_{r_{\tau}}\phi'(\boldsymbol{x}_{r_{\tau}}^{T}\boldsymbol{\theta}_{\tau})\phi'(\boldsymbol{x}_{r_{\tau}}^{T}\boldsymbol{\theta}_{\tau}, \boldsymbol{x}_{r_{\tau}}^{T}\boldsymbol{\theta})\boldsymbol{x}_{r_{\tau}}^{T}\boldsymbol{w}_{\tau}\|_{\ell_{2}}^{2} \\ &= \|(\boldsymbol{I} - \eta\boldsymbol{G}_{r_{\tau}})\boldsymbol{w}_{\tau}\|_{\ell_{2}}^{2}. \end{aligned}$$

where $G_{r_{\tau}} = x_{r_{\tau}} \phi'(x_{r_{\tau}}^T \theta_{\tau}) \phi'(x_{r_{\tau}}^T \theta_{\tau}, x_{r_{\tau}}^T \theta) x_{r_{\tau}}^T$. Since ϕ is 1-Lipschitz and β -increasing, $G_{r_{\tau}}$ is a positive-semidefinite matrix satisfying

$$\begin{aligned} \boldsymbol{x}_{r_{\tau}} \boldsymbol{x}_{r_{\tau}}^T \succeq \boldsymbol{G}_{r_{\tau}} \succeq \beta^2 \boldsymbol{x}_{r_{\tau}} \boldsymbol{x}_{r_{\tau}}^T, \\ \boldsymbol{G}_{r_{\tau}}^T \boldsymbol{G}_{r_{\tau}} \preceq \boldsymbol{x}_{r_{\tau}} \boldsymbol{x}_{r_{\tau}}^T \boldsymbol{x}_{r_{\tau}} \boldsymbol{x}_{r_{\tau}}^T \preceq B \boldsymbol{x}_{r_{\tau}} \boldsymbol{x}_{r_{\tau}}^T. \end{aligned}$$

Consequently, we find the following bounds in expectation

$$\gamma_{+} \mathbf{I}_{n} \succeq \mathbb{E}[\mathbf{G}_{r_{\tau}}] \succeq \beta^{2} \gamma_{-} \mathbf{I}_{n},$$

$$\mathbb{E}[\mathbf{G}_{r_{\tau}}^{T} \mathbf{G}_{r_{\tau}}] \preceq B \gamma_{+} \mathbf{I}_{n}.$$
(A.1)

Observe that (A.1) essentially lower bounds the *strong convexity* parameter of the problem with $\beta^2 \gamma_-$; which is the strong convexity of the identical problem with the linear activation (i.e. $\beta=1$). However, we only consider strong convexity around the ground truth parameter θ i.e. we restricted our attention to (θ, θ_τ) pairs. With this, $w_{\tau+1}$ can be controlled as,

$$\mathbb{E}[\|\boldsymbol{w}_{\tau+1}\|_{\ell_{2}}^{2}] = \mathbb{E}[\|(\boldsymbol{I} - \eta \boldsymbol{G}_{r_{\tau}})\boldsymbol{w}_{\tau}\|_{\ell_{2}}^{2}]$$

$$= \|\boldsymbol{w}_{\tau}\|_{\ell_{2}}^{2} - 2\eta \,\mathbb{E}[\boldsymbol{w}_{\tau}^{T}\boldsymbol{G}_{r_{\tau}}\boldsymbol{w}_{\tau}] + \eta^{2} \,\mathbb{E}[\boldsymbol{w}_{\tau}^{T}\boldsymbol{G}_{r_{\tau}}^{T}\boldsymbol{G}_{r_{\tau}}\boldsymbol{w}_{\tau}]$$

$$\leq \|\boldsymbol{w}_{\tau}\|_{\ell_{2}}^{2}(1 - 2\eta\beta^{2}\gamma_{-} + \eta^{2}B\gamma_{+}).$$

Setting $\eta = \frac{\beta^2 \gamma_-}{\gamma_+ B}$, we find the advertised bound

$$\mathbb{E}[\|\boldsymbol{w}_{\tau+1}\|_{\ell_2}^2] \leq \mathbb{E}[\|\boldsymbol{w}_{\tau}\|_{\ell_2}^2](1 - \frac{\beta^4 \gamma_{-}^2}{\gamma_{+} B}).$$

Applying induction over the iterations τ , we find the advertised bound (4.2)

$$\mathbb{E}[\|\boldsymbol{w}_{\tau}\|_{\ell_{2}}^{2}] \leq \|\boldsymbol{w}_{0}\|_{\ell_{2}}^{2} (1 - \frac{\beta^{4} \gamma_{-}^{2}}{\gamma_{+} B})^{\tau}.$$

Lemma 8 (Merging L **splits)** Assume matrices $\mathbf{X}^{(i)} \in \mathbb{R}^{N_i \times q}$ are given for $1 \leq i \leq L$. Suppose for all $1 \leq i \leq L$, rows of $\mathbf{X}^{(i)}$ has ℓ_2 norm at most \sqrt{B} and each $\mathbf{X}^{(i)}$ satisfies

$$\gamma_{+}\boldsymbol{I}_{n} \succeq \frac{\boldsymbol{X}^{(i)}^{T}\boldsymbol{X}^{(i)}}{N_{i}} \succeq \gamma_{-}\boldsymbol{I}_{n}.$$

Set
$$N = \sum_{i=1}^{L} N_i$$
 and form the concatenated matrix $\mathbf{X} = \begin{bmatrix} \mathbf{X}^{(1)} \\ \mathbf{X}^{(2)} \\ \vdots \\ \mathbf{X}^{(L)} \end{bmatrix}$. Denote ith row of \mathbf{X} by \mathbf{x}_i . Then, for

each i, $\|\boldsymbol{x}_i\|_{\ell_2}^2 \leq B$ and

$$\gamma_{+}\boldsymbol{I}_{n}\succeq \frac{\boldsymbol{X}^{T}\boldsymbol{X}}{N}=\frac{1}{N}\sum_{i=1}^{N}\boldsymbol{x}_{i}\boldsymbol{x}_{i}^{T}\succeq \gamma_{-}\boldsymbol{I}_{n}.$$

Proof The bound on the rows $\|\boldsymbol{x}_i\|_{\ell_2}$ directly follows by assumption. For the remaining result, first observe that $\boldsymbol{X}^T\boldsymbol{X} = \sum_{i=1}^L \boldsymbol{X}^{(i)} \boldsymbol{X}^{(i)}$. Next, we have

$$N\gamma_{+}\boldsymbol{I}_{n} = \sum_{i=1}^{L} N_{i}\gamma_{+}\boldsymbol{I}_{n} \succeq \sum_{i=1}^{L} \boldsymbol{X^{(i)}}^{T} \boldsymbol{X^{(i)}} \succeq \sum_{i=1}^{L} N_{i}\gamma_{-}\boldsymbol{I}_{n} = N\gamma_{-}\boldsymbol{I}_{n}.$$

Combining these two yields the desired upper/lower bounds on X^TX/N .

Appendix B. Properties of the nonlinear state equations

This section characterizes the properties of the state vector h_t when input sequence is normally distributed. These bounds will be crucial for obtaining upper and lower bounds for the singular values of the data matrix $X = [x_1 \dots x_N]^T$ described in (2.2). For probabilistic arguments, we will use the properties of subgaussian random variables. Orlicz norm provides a general framework that subsumes subgaussianity.

Definition 9 (Orlicz norms) For a scalar random variable Orlicz-a norm is defined as

$$||X||_{\psi_a} = \sup_{k>1} k^{-1/a} (\mathbb{E}[|X|^k])^{1/k}$$

Orlicz-a norm of a vector $\mathbf{x} \in \mathbb{R}^p$ is defined as $\|\mathbf{x}\|_{\psi_a} = \sup_{\mathbf{v} \in \mathcal{B}^p} \|\mathbf{v}^T \mathbf{x}\|_{\psi_a}$ where \mathcal{B}^p is the unit ℓ_2 ball. The subexponential norm is the Orlicz-1 norm $\|\cdot\|_{\psi_1}$ and the subgaussian norm is the Orlicz-2 norm $\|\cdot\|_{\psi_2}$.

Lemma 10 (Lipschitz properties of the state vector) Consider the state equation (1.1). Suppose activation ϕ is 1-Lipschitz. Observe that h_{t+1} is a deterministic function of the input sequence $\{u_{\tau}\}_{\tau=0}^{t}$. Fixing all vectors $\{u_{i}\}_{i\neq\tau}$ (i.e. all except u_{τ}), h_{t+1} is $\|A\|^{t-\tau}\|B\|$ Lipschitz function of u_{τ} for $0 \le \tau \le t$.

Proof Fixing $\{u_i\}_{i\neq\tau}$, denote h_{t+1} as a function of u_{τ} by $h_{t+1}(u_{\tau})$. Given a pair of vectors u_{τ}, u'_{τ} using 1-Lipschitzness of ϕ , for any $t > \tau$, we have

$$egin{aligned} \|m{h}_{t+1}(m{u}_{ au}) - m{h}_{t+1}(m{u}_{ au}')\|_{\ell_2} &\leq \|\phi(m{A}m{h}_t(m{u}_{ au}) + m{B}m{u}_t) - \phi(m{A}m{h}_t(m{u}_{ au}') + m{B}m{u}_t)\|_{\ell_2} \ &\leq \|m{A}(m{h}_t(m{u}_{ au}) - m{h}_t(m{u}_{ au}'))\|_{\ell_2} \ &\leq \|m{A}\|\|m{h}_t(m{u}_{ au}) - m{h}_t(m{u}_{ au}')\|_{\ell_2}. \end{aligned}$$

Proceeding with this recursion until $t = \tau$, we find

$$egin{aligned} \|m{h}_{t+1}(m{u}_{ au}) - m{h}_{t+1}(m{u}_{ au}')\|_{\ell_2} & \leq \|m{A}\|^{t- au}\|m{h}_{ au+1}(m{u}_{ au}) - m{h}_{ au+1}(m{u}_{ au}')\|_{\ell_2} \ & \leq \|m{A}\|^{t- au}\|\phi(m{A}m{h}_{ au} + m{B}m{u}_{ au}) - \phi(m{A}m{h}_{ au} + m{B}m{u}_{ au}')\|_{\ell_2} \ & \leq \|m{A}\|^{t- au}\|m{B}\|\|m{u}_{ au} - m{u}_{ au}'\|_{\ell_2}. \end{aligned}$$

This bound implies $h_{t+1}(u_{\tau})$ is $||A||^{t-\tau}||B||$ Lipschitz function of u_{τ} .

Lemma 11 (Upper bound) Consider the state equation governed by equation (1.1). Suppose $u_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \mathbf{I}_p)$, ϕ is 1-Lipschitz, $\phi(0) = 0$ and $\mathbf{h}_0 = 0$. Recall the definition (3.2) of B_t . We have the following properties

- h_t is a B_t -Lipschitz function of the vector $q_t = [u_0^T \dots u_{t-1}^T]^T \in \mathbb{R}^{tp}$.
- There exists an absolute constant c > 0 such that $\|\mathbf{h}_t \mathbb{E}[\mathbf{h}_t]\|_{\psi_2} \le cB_t$ and $\mathbf{\Sigma}[\mathbf{h}_t] \le B_t^2 \mathbf{I}_n$.
- h_t satisfies

$$\mathbb{E}[\|\boldsymbol{h}_t\|_{\ell_2}^2] \le tr(\boldsymbol{B}\boldsymbol{B}^T) \frac{1 - \|\boldsymbol{A}\|^{2t}}{1 - \|\boldsymbol{A}\|^2} \le \min\{n, p\} B_t^2.$$

Also, there exists an absolute constant c>0 such that for any $m\geq n$, with probability $1-2\exp(-100m)$, $\|\mathbf{h}_t\|_{\ell_2}\leq c\sqrt{m}B_t$.

Proof i) Bounding Lipschitz constant: Observe that h_t is a deterministic function of q_t i.e. $h_t = f(q_t)$ for some function f. To bound Lipschitz constant of f, for all (deterministic) vector pairs q_t and \hat{q}_t , we find a scalar L_f satisfying,

$$||f(q_t) - f(\hat{q}_t)||_{\ell_2} \le L_f ||q_t - \hat{q}_t||_{\ell_2}.$$
 (B.1)

Define the vectors, $\{a_i\}_{i=0}^t$, as follows

$$\boldsymbol{a}_i = [\hat{\boldsymbol{u}}_0^T \ \dots \ \hat{\boldsymbol{u}}_{i-1}^T \ \boldsymbol{u}_i^T \ \dots \ \boldsymbol{u}_{t-1}^T]^T.$$

Observing that $a_0 = q_t$, $a_t = \hat{q}_t$, we write the telescopic sum,

$$\|f(\boldsymbol{q}_t) - f(\hat{\boldsymbol{q}}_t)\|_{\ell_2} \leq \sum_{i=0}^{t-1} \|f(\boldsymbol{a}_{i+1}) - f(\boldsymbol{a}_i)\|_{\ell_2}.$$

Focusing on the individual terms $f(a_{i+1}) - f(a_i)$, observe that the only difference is the u_i , \hat{u}_i terms. Viewing h_t as a function of u_i and applying Lemma 10,

$$||f(\boldsymbol{a}_{i+1}) - f(\boldsymbol{a}_i)||_{\ell_2} \le ||\boldsymbol{A}||^{t-1-i} ||\boldsymbol{B}|| ||\boldsymbol{u}_i - \hat{\boldsymbol{u}}_i||_{\ell_2}.$$

To bound the sum, we apply the Cauchy-Schwarz inequality; which yields

$$|f(\mathbf{q}_{t}) - f(\hat{\mathbf{q}}_{t})| \leq \sum_{i=0}^{t-1} ||\mathbf{A}||^{t-1-i} ||\mathbf{B}|| ||\mathbf{u}_{i} - \hat{\mathbf{u}}_{i}||_{\ell_{2}}$$

$$\leq (\sum_{i=0}^{t-1} ||\mathbf{A}||^{2(t-1-i)} ||\mathbf{B}||^{2})^{1/2} \underbrace{(\sum_{i=0}^{t-1} ||\mathbf{u}_{i} - \hat{\mathbf{u}}_{i}||_{\ell_{2}}^{2})^{1/2}}_{||\mathbf{q}_{t} - \hat{\mathbf{q}}_{t}||_{\ell_{2}}}$$

$$\leq \sqrt{\frac{||\mathbf{B}||^{2} (1 - ||\mathbf{A}||^{2t})}{1 - ||\mathbf{A}||^{2}}} ||\mathbf{q}_{t} - \hat{\mathbf{q}}_{t}||_{\ell_{2}}}$$

$$= B_{t} ||\mathbf{q}_{t} - \hat{\mathbf{q}}_{t}||_{\ell_{2}}. \tag{B.2}$$

The final line achieves the inequality (B.1) with $L_f = B_t$ hence h_t is B_t Lipschitz function of q_t .

ii) Bounding subgaussian norm: When $u_t \overset{\text{i.i.d.}}{\sim} \mathcal{N}(0, I_p)$, the vector q_t is distributed as $\mathcal{N}(0, I_{tp})$. Since h_t a B_t Lipschitz function of q_t , for any fixed unit length vector v, $\alpha_v := v^T h_t = v^T f(q_t)$ is still B_t -Lipschitz function of q_t . Hence, using Gaussian concentration of Lipschitz functions, α_v satisfies

$$\mathbb{P}(|\alpha_{\boldsymbol{v}} - \mathbb{E}[\alpha_{\boldsymbol{v}}]| \ge t) \le 2\exp(-\frac{t^2}{2B_*^2}).$$

This implies that for any v, $\alpha_v - \mathbb{E}[\alpha_v]$ is $\mathcal{O}(B_t)$ subgaussian Vershynin (2010). This is true for all unit v, hence using Definition 9, the vector \mathbf{h}_t satisfies $\|\mathbf{h}_t - \mathbb{E}[\mathbf{h}_t]\|_{\psi_2} \leq \mathcal{O}(B_t)$ as well. Secondly, B_t -Lipschitz function of a Gaussian vector obeys the variance inequality $\mathbf{var}[\alpha_v] \leq B_t^2$ (page 49 of Ledoux (2001)), which implies the covariance bound

$$\Sigma[\boldsymbol{h}_t] \leq B_t^2 \boldsymbol{I}_n$$
.

iii) Bounding ℓ_2 -norm: To obtain this result, we first bound $\mathbb{E}[\|\boldsymbol{h}_t\|_{\ell_2}^2]$. Since ϕ is 1-Lipschitz and $\phi(0)=0$, we have the deterministic relation

$$\|\boldsymbol{h}_{t+1}\|_{\ell_2} \leq \|\boldsymbol{A}\boldsymbol{h}_t + \boldsymbol{B}\boldsymbol{u}_t\|_{\ell_2}.$$

Taking squares of both sides, expanding the right hand side, and using the independence of h_t , u_t and the covariance information of u_t , we obtain

$$\mathbb{E}[\|\boldsymbol{h}_{t+1}\|_{\ell_2}^2] \le \mathbb{E}[\|\boldsymbol{A}\boldsymbol{h}_t + \boldsymbol{B}\boldsymbol{u}_t\|_{\ell_2}^2] = \mathbb{E}[\|\boldsymbol{A}\boldsymbol{h}_t\|_{\ell_2}^2] + \mathbb{E}[\|\boldsymbol{B}\boldsymbol{u}_t\|_{\ell_2}^2]$$
(B.3)

$$\leq \|A\|^2 \mathbb{E}[\|h_t\|_{\ell_2}^2] + \operatorname{tr}(BB^T).$$
 (B.4)

Now that the recursion is established, expanding h_t on the right hand side until $h_0 = 0$, we obtain

$$\mathbb{E}[\|\boldsymbol{h}_{t+1}\|_{\ell_2}^2] \leq \sum_{i=0}^t \|\boldsymbol{A}\|^{2i} \mathrm{tr}(\boldsymbol{B}\boldsymbol{B}^T) \leq \mathrm{tr}(\boldsymbol{B}\boldsymbol{B}^T) \frac{1 - \|\boldsymbol{A}\|^{2(t+1)}}{1 - \|\boldsymbol{A}\|^2}.$$

Now using the fact that $tr(BB^T) \le rank(B)||B||^2 \le min\{n, p\}||B||^2$, we find

$$\mathbb{E}[\|\boldsymbol{h}_{t+1}\|_{\ell_2}]^2 \leq \mathbb{E}[\|\boldsymbol{h}_{t+1}\|_{\ell_2}^2] \leq \min\{n, p\}B_{t+1}^2.$$

Finally, using the fact that h_t is B_t -Lipschitz function and utilizing Gaussian concentration of $q_t \sim \mathcal{N}(0, I_{tp})$, we find

$$\mathbb{P}(\|\boldsymbol{h}_{t+1}\|_{\ell_2} - \mathbb{E}[\|\boldsymbol{h}_{t+1}\|_{\ell_2}] \ge t) \le \exp(-\frac{t^2}{2B_t^2}).$$

Setting $t = (c-1)\sqrt{m}B_t$ for sufficiently large c > 0, we find $\mathbb{P}(\|\mathbf{h}_t\|_{\ell_2} \ge \sqrt{n}B_t + (c-1)\sqrt{m}B_t) \le \exp(-100m)$.

Lemma 12 (Odd activations) Suppose ϕ is strictly increasing and obeys $\phi(x) = -\phi(-x)$ for all x and $h_0 = 0$. Consider the state equation (1.1) driven $u_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \mathbf{I}_p)$. We have that $\mathbb{E}[\mathbf{h}_t] = 0$.

Proof We will inductively show that $\{h_t\}_{t\geq 0}$ has a symmetric distribution around 0. Suppose the vector h_t satisfies this assumption. Let $S \subset \mathbb{R}^n$ be a set. We will argue that $\mathbb{P}(h_{t+1} \subset S) = \mathbb{P}(h_{t+1} \subset -S)$. Since ϕ is strictly increasing, it is bijective on vectors, and we can define the unique inverse set $S' = \phi^{-1}(S)$. Also since ϕ is odd, $\phi(-S') = -S$. Since h_t , u_t are independent and symmetric, we reach the desired conclusion as follows

$$\mathbb{P}(\mathbf{h}_{t+1} \subset S) = \mathbb{P}(\mathbf{A}\mathbf{h}_t + \mathbf{B}\mathbf{u}_t \subset S') = \mathbb{P}(\mathbf{A}(-\mathbf{h}_t) + \mathbf{B}(-\mathbf{u}_t) \subset S')$$
(B.5)

$$= \mathbb{P}(\mathbf{A}\mathbf{h}_t + \mathbf{B}\mathbf{u}_t \subset -S') = \mathbb{P}(\phi(\mathbf{A}\mathbf{h}_t + \mathbf{B}\mathbf{u}_t) \subset \phi(-S')) = \mathbb{P}(\mathbf{h}_{t+1} \subset -S). \tag{B.6}$$

Theorem 13 (State-vector lower bound) Consider the nonlinear state equation (1.1) with $\{u_t\}_{t\geq 0} \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \mathbf{I}_p)$. Suppose ϕ is a β -increasing function for some constant $\beta > 0$. For any $t \geq 1$, the state vector obeys

$$\Sigma[\boldsymbol{h}_t] \succeq \beta^2 s_{\min}(\boldsymbol{B}\boldsymbol{B}^T) \boldsymbol{I}_n$$

Proof The proof is an application of Lemma 15. The main idea is to write h_t as sum of two independent vectors, one of which has independent entries. Consider a multivariate Gaussian vector $\mathbf{g} \sim \mathcal{N}(0, \mathbf{\Sigma})$. \mathbf{g} is statistically identical to $\mathbf{g}_1 + \mathbf{g}_2$ where $\mathbf{g}_1 \sim \mathcal{N}(0, s_{\min}(\mathbf{\Sigma})\mathbf{I}_d)$ and $\mathbf{g}_2 \sim \mathcal{N}(0, \mathbf{\Sigma} - s_{\min}(\mathbf{\Sigma})\mathbf{I}_d)$ are independent multivariate Gaussians.

Since $Bu_t \sim \mathcal{N}(0, BB^T)$, setting $\Sigma = BB^T$ and $s_{\min} = s_{\min}(\Sigma)$, we have that $Bu_t \sim g_1 + g_2$ where g_1, g_2 are independent and $g_1 \sim \mathcal{N}(0, s_{\min}I_n)$ and $g_2 \sim \mathcal{N}(0, \Sigma - s_{\min}I_n)$. Consequently, we may write

$$Bu_t + Ah_t \sim g_1 + g_2 + Ah_t$$
.

For lower bound, the crucial component will be the g_1 term; which has i.i.d. entries. Applying Lemma 15 by setting $x = g_1$ and $y = g_2 + Ah_t$, and using the fact that h_t, g_1, g_2 are all independent of each other, we find the advertised bound, for all t > 0, via

$$\Sigma[\boldsymbol{h}_{t+1}] = \Sigma[\phi(\boldsymbol{g}_1 + \boldsymbol{g}_2 + \boldsymbol{A}\boldsymbol{h}_t)] \succeq \beta^2 s_{\min} \boldsymbol{I}_n.$$

The next theorem applies to multiple-input-single-output (MISO) systems where A is a scalar and B is a row vector. The goal is refining the lower bound of Theorem 13.

Theorem 14 (MISO lower bound) Consider the setup of Theorem 13 with single output i.e. n = 1. For any $t \ge 1$, the state vector obeys

$$\mathbf{var}[\boldsymbol{h}_t] \ge \beta^2 \|\boldsymbol{B}\|_{\ell_2}^2 \frac{1 - (\beta|\boldsymbol{A}|)^{2t}}{1 - \beta^2 |\boldsymbol{A}|^2}.$$

Proof For any random variable X, applying Lemma 15, we have $\mathbf{var}[\phi(X)] \ge \beta^2 \mathbf{var}[X]$. Recursively, this yields

$$\operatorname{var}[\boldsymbol{h}_{t+1}] = \operatorname{var}[\phi(\boldsymbol{A}\boldsymbol{h}_t + \boldsymbol{B}\boldsymbol{u}_t)] \ge \beta^2 \operatorname{var}[\boldsymbol{A}\boldsymbol{h}_t + \boldsymbol{B}\boldsymbol{u}_t] = \beta^2 (|\boldsymbol{A}|^2 \operatorname{var}[\boldsymbol{h}_t] + \|\boldsymbol{B}\|_{\ell_2}^2).$$

Expanding these inequalities till h_0 , we obtain the desired bound

$$extbf{var}[oldsymbol{h}_t] \geq \sum_{i=1}^t (eta^i |oldsymbol{A}|^{i-1} \|oldsymbol{B}\|_{\ell_2})^2.$$

Lemma 15 (Vector lower bound) Suppose ϕ is a β -increasing function. Let $\mathbf{x} = [\mathbf{x}_1 \dots \mathbf{x}_n]^T$ be a vector with i.i.d. entries distributed as $\mathbf{x}_i \sim X$. Let \mathbf{y} be a random vector independent of \mathbf{x} . Then,

$$\Sigma[\phi(x+y)] \succeq \beta^2 \text{var}[X] I_n$$
.

Proof We first apply law of total covariance (e.g. Lemma 16) to simplify the problem using the following lower bound based on the independence of x and y,

$$\Sigma[\phi(x+y)] \succeq \mathbb{E}_{y}[\Sigma[\phi(x+y) \mid y]]$$
 (B.7)

$$= \mathbb{E}_{\boldsymbol{y}}[\boldsymbol{\Sigma}_{\boldsymbol{x}}[\phi(\boldsymbol{x} + \boldsymbol{y})]]. \tag{B.8}$$

Now, focusing on the covariance $\Sigma_x[\phi(x+y)]$, fixing a realization of y, and using the fact that x has i.i.d. entries; $\phi(x+y)$ has independent entries as ϕ applies entry-wise. This implies that $\Sigma_x[\phi(x+y)]$ is a diagonal matrix. Consequently, its lowest eigenvalue is the minimum variance over all entries,

$$\mathbf{\Sigma}_{m{x}}[\phi(m{x}+m{y})] \succeq \min_{1 \leq i \leq n} \mathbf{var}[\phi(m{x}_i+m{y}_i)] m{I}_n.$$

Fortunately, Lemma 17 provides the lower bound $\mathbf{var}[\phi(\mathbf{x}_i + \mathbf{y}_i)] \ge \beta^2 \mathbf{var}[X]$. Since this lower bound holds for any fixed realization of \mathbf{y} , it still holds after taking expectation over \mathbf{y} ; which concludes the proof.

The next two lemmas are helper results for Lemma 15 and are provided for the sake of completeness.

Lemma 16 (Law of total covariance) Let x, y be two random vectors and assume y has finite covariance. Then

$$\mathbf{\Sigma}[oldsymbol{y}] = \mathbb{E}[\mathbf{\Sigma}[oldsymbol{y} \mid oldsymbol{x}]] + \mathbf{\Sigma}[\mathbb{E}[oldsymbol{y} \mid oldsymbol{x}]].$$

Proof First, write $\Sigma[y] = \mathbb{E}[yy^T] - \mathbb{E}[y]\mathbb{E}[y^T]$. Then, applying the law of total expectation to each term,

$$oldsymbol{\Sigma}[oldsymbol{y}] = \mathbb{E}[\mathbb{E}[oldsymbol{y}oldsymbol{y}^T \mid oldsymbol{x}]] - \mathbb{E}[\mathbb{E}[oldsymbol{y} \mid oldsymbol{x}]] \, \mathbb{E}[\mathbb{E}[oldsymbol{y}^T \mid oldsymbol{x}]].$$

Next, we can write the conditional expectation as $\mathbb{E}[\mathbb{E}[yy^T \mid x]] = \mathbb{E}[\Sigma[y \mid x]] + \mathbb{E}[\mathbb{E}[y \mid x]\mathbb{E}[y \mid x]]^T$. To conclude, we obtain the covariance of $\mathbb{E}[y \mid x]$ via the difference,

$$\mathbb{E}[\mathbb{E}[\boldsymbol{y} \mid \boldsymbol{x}] \, \mathbb{E}[\boldsymbol{y} \mid \boldsymbol{x}]]^T - \mathbb{E}[\mathbb{E}[\boldsymbol{y} \mid \boldsymbol{x}]] \, \mathbb{E}[\mathbb{E}[\boldsymbol{y}^T \mid \boldsymbol{x}]] = \boldsymbol{\Sigma}[\mathbb{E}[\boldsymbol{y} \mid \boldsymbol{x}]],$$

which yields the desired bound.

Lemma 17 (Scalar lower bound) Suppose ϕ is a β -increasing function with $\beta > 0$ as defined in Definition 1. Given a random variable X and a scalar y, we have

$$\operatorname{var}[\phi(X+y)] \ge \beta^2 \operatorname{var}[X].$$

Proof Since ϕ is β -increasing, it is invertible and ϕ^{-1} is strictly increasing. Additionally, ϕ^{-1} is $1/\beta$ Lipschitz since,

$$|\phi(a) - \phi(b)| \ge \beta |a - b| \implies |a - b| \ge \beta |\phi^{-1}(a) - \phi^{-1}(b)|.$$

Using this observation and the fact that $\mathbb{E}[X]$ minimizes $\mathbb{E}(X-\alpha)^2$ over α , $\operatorname{var}[\phi(X+y)]$ can be lower bounded as follows

$$\begin{aligned} \mathbf{var}[\phi(X+y)] &= \mathbb{E}(\phi(X+y) - \mathbb{E}[\phi(X+y)])^2 \\ &\geq \beta^2 \, \mathbb{E}((X+y) - \phi^{-1}(\mathbb{E}[\phi(X+y)]))^2 \\ &\geq \beta^2 \, \mathbb{E}(X+y - \mathbb{E}[X+y])^2 \\ &= \beta^2 \, \mathbb{E}(X - \mathbb{E}\,X)^2 = \beta^2 \mathbf{var}[X]. \end{aligned}$$

Note that, the final line is the desired conclusion.

Appendix C. Truncating Stable Systems

One of the challenges in analyzing dynamical systems is the fact that samples from the same trajectory have temporal dependence. This section shows that, for stable systems, the impact of the past states decay exponentially fast and the system can be approximated by using the recent inputs only. We first define the truncation of the state vector.

Definition 18 (Truncated state vector) Suppose $\phi(0) = 0$, initial condition $h_0 = 0$, and consider the state equation (1.1). Given a timestamp t, L-truncation of the state vector \mathbf{h}_t is denoted by $\bar{\mathbf{h}}_{t,L}$ and is equal to \mathbf{q}_t where

$$\mathbf{q}_{\tau+1} = \phi(\mathbf{A}\mathbf{q}_{\tau} + \mathbf{B}\mathbf{u}_{\tau}') \quad , \quad q_0 = 0 \tag{C.1}$$

is the state vector generated by the inputs \mathbf{u}'_{τ} satisfying

$$oldsymbol{u}_{ au}' = egin{cases} 0 & \textit{if } au < t - L \ oldsymbol{u}_{ au} & \textit{else} \end{cases}$$
 .

In words, L truncated state vector $\bar{h}_{t,L}$ is obtained by unrolling h_t until time t-L and setting the contribution of the state vector h_{t-L} to 0. This way, $\bar{h}_{t,L}$ depends only on the variables $\{u_t\}_{t=t-L}^{t-1}$.

The following lemma states that impact of truncation can be made fairly small for stable systems (||A|| < 1).

Lemma 19 (Truncation impact – deterministic) Consider the state vector \mathbf{h}_t and its L-truncation $\bar{\mathbf{h}}_{t,L}$ from Definition 18. Suppose ϕ is 1-Lipschitz. We have that

$$\|oldsymbol{h}_t - ar{oldsymbol{h}}_{t,L}\|_{\ell_2} \leq egin{cases} 0 & \textit{if } t \leq L \\ \|oldsymbol{A}\|^L \|oldsymbol{h}_{t-L}\|_{\ell_2} & \textit{else} \end{cases}.$$

Proof When $t \leq L$, Definition 18 implies $u'_{\tau} = u_{\tau}$ hence $h_t = q_t = \bar{h}_{t,L}$. When t > L, we again use Definition 18 and recall that $u'_{\tau} = 0$ until time $\tau = t - L - 1$. For all $t - L < \tau \leq t$, using 1-Lipschitzness of ϕ , we have that

$$egin{aligned} \|m{h}_{ au} - m{q}_{ au}\|_{\ell_2} &= \|\phi(m{A}m{h}_{ au-1} + m{B}m{u}_{ au-1}) - \phi(m{A}m{q}_{ au-1} + m{B}m{u}_{ au-1})\|_{\ell_2} \ &\leq \|(m{A}m{h}_{ au-1} + m{B}m{u}_{ au-1}) - (m{A}m{q}_{ au-1} + m{B}m{u}_{ au-1})\|_{\ell_2} \ &\leq \|m{A}(m{h}_{ au-1} - m{q}_{ au-1})\|_{\ell_2} \leq \|m{A}\|\|m{h}_{ au-1} - m{q}_{ au-1}\|_{\ell_2}. \end{aligned}$$

Applying this recursion between $t-L < \tau \le t$ and using the fact that $q_{t-L} = 0$ implies the advertised result

$$egin{aligned} \|m{h}_t - m{q}_t\|_{\ell_2} & \leq \|m{A}\|^L \|m{h}_{t-L} - m{q}_{t-L}\|_{\ell_2} \ & \leq \|m{A}\|^L \|m{h}_{t-L}\|_{\ell_2}. \end{aligned}$$

C.1. Near independence of sub-trajectories

We will now argue that, for stable systems, a single trajectory can be split into multiple nearly independent trajectories. First, we describe how the sub-trajectories are constructed.

Definition 20 (Sub-trajectory) Let sampling rate $L \ge 1$ and offset $1 \le \bar{\tau} \le L$ be two integers. Let $\bar{N} = \bar{N}_{\bar{\tau}}$ be the largest integer obeying $(\bar{N} - 1)L + \bar{\tau} \le N$. We sample the trajectory $\{h_t, u_t\}_{t=0}^N$ at the points $\bar{\tau}, \bar{\tau} + L, \dots, \bar{\tau} + (\bar{N} - 1)L + \bar{\tau}$ and define the $\bar{\tau}$ th sub-trajectory as

$$(\boldsymbol{h}^{(i)},\boldsymbol{u}^{(i)}) := (\boldsymbol{h}^{(i,\bar{\tau})},\boldsymbol{u}^{(i,\bar{\tau})}) = (\boldsymbol{h}_{(i-1)L+\bar{\tau}},\boldsymbol{u}_{(i-1)L+\bar{\tau}}).$$

Definition 21 (Truncated sub-trajectory) Consider the state equation (1.1) and recall Definition 18. Given offset $\bar{\tau}$ and sampling rate L, for $1 \leq i \leq \bar{N}$, the ith truncated sub-trajectory states are $\{\bar{\mathbf{h}}^{(i)}\}_{i=1}^{\bar{N}}$ where the ith state is defined as

$$ar{h}^{(i)} = ar{h}_{L(i-1)+ar{ au},L-1}.$$

The truncated samples are independent of each other as shown in the next lemma.

Lemma 22 Consider the truncated states of Definition 21. If (1.1) is generated by independent vectors $\{u_t\}_{t\geq 0}$, for any offset $\bar{\tau}$ and sampling rate L, the vectors $\{\bar{\boldsymbol{h}}^{(i)}\}_{i=1}^{\bar{N}}$, $\{\boldsymbol{u}^{(i)}\}_{i=1}^{\bar{N}}$ are all independent of each other.

Proof By construction $\bar{\boldsymbol{h}}^{(i)}$ only depends on the vectors $\{\boldsymbol{u}_{\tau}\}_{\tau=L(i-2)+\bar{\tau}+1}^{L(i-1)+\bar{\tau}-1}$. Note that the dependence ranges $[L(i-2)+\bar{\tau}+1,L(i-1)+\bar{\tau}-1]$ are disjoint intervals for different i's; hence $(\bar{\boldsymbol{h}}^{(i)})_{i=1}^{\bar{N}}$ are independent of each other. To show the independence of $\boldsymbol{u}^{(i)}$ and $\bar{\boldsymbol{h}}^{(i)}$; observe that inputs $\boldsymbol{u}^{(i)}=\boldsymbol{u}_{L(i-1)+\bar{\tau}}$ have timestamp $\bar{\tau}$ modulo L; which is not covered by the dependence range of $(\bar{\boldsymbol{h}}^{(i)})_{i=1}^{\bar{N}}$.

If the input is randomly generated, Lemma 19 can be combined with a probabilistic bound on h_t , to show that truncated states $\bar{h}^{(i)}$ are fairly close to the actual states $h^{(i)}$.

Lemma 23 (Truncation impact – random) Given offset $\bar{\tau}$ and sampling rate L, consider the state vectors of the sub-trajectory $\{\boldsymbol{h}^{(i)}\}_{i=1}^{\bar{N}}$ and L-1-truncations $(\bar{\boldsymbol{h}}^{(i)})_{i=1}^{\bar{N}}$. Suppose $\|\boldsymbol{A}\| < 1$, $\boldsymbol{h}_0 = 0$, ϕ is 1-Lipschitz, and $\phi(0) = 0$. Also suppose upper bound (4.3) of Assumption 1 holds for some $\theta \leq \sqrt{n}$, $\gamma_+ > 0$. There exists an absolute constant c > 0 such that with probability at least $1 - 2\bar{N} \exp(-100n)$, for all $1 \leq i \leq \bar{N}$, the following bound holds

$$\|\boldsymbol{h}^{(i)} - \bar{\boldsymbol{h}}^{(i)}\|_{\ell_2} \le c\sqrt{n}\|\boldsymbol{A}\|^{L-1}\sqrt{\gamma_+}.$$

In particular, we can always pick $\gamma_+ = B_{\infty}^2$ (via Lemma 11).

Proof Using Assumption 1, we can apply Lemma 31 on vectors $\{h_{(i-2)L+\bar{\tau}+1}\}_{i=1}^{\bar{N}}$. Using a union bound, with desired probability, all vectors obey

$$\|\boldsymbol{h}_{(i-2)L+\bar{\tau}+1} - \mathbb{E}[\boldsymbol{h}_{(i-2)L+\bar{\tau}+1}]\|_{\ell_2} \le (c-1)\sqrt{n\gamma_+},$$

for sufficiently large c. Since $\theta \leq \sqrt{n}$, triangle inequality implies $\|\mathbf{h}_{(i-2)L+\bar{\tau}+1}\|_{\ell_2} \leq c\sqrt{n\gamma_+}$. Now, applying Lemma 19, for all $1 \leq i \leq \bar{N}$, we find

$$\|\boldsymbol{h}^{(i)} - \bar{\boldsymbol{h}}^{(i)}\|_{\ell_{2}} = \|\boldsymbol{h}_{(i-1)L+\bar{\tau}} - \bar{\boldsymbol{h}}_{(i-1)L+\bar{\tau},L-1}\|_{\ell_{2}}$$

$$\leq \|\boldsymbol{A}\|^{L-1}\|\boldsymbol{h}_{(i-2)L+\bar{\tau}+1}\|_{\ell_{2}}$$

$$\leq c\|\boldsymbol{A}\|^{L-1}\sqrt{n\gamma_{+}}.$$

Appendix D. Properties of the data matrix

This section utilizes the probabilistic estimates from Section B to provide bounds on the condition number of data matrices obtained from the RNN trajectory (1.1). Following (2.2), these matrices H, U and X are defined as

$$\boldsymbol{H} = [\boldsymbol{h}_1 \dots \boldsymbol{h}_N]^T$$
 , $\boldsymbol{U} = [\boldsymbol{u}_1 \dots \boldsymbol{u}_N]^T$, $\boldsymbol{X} = [\boldsymbol{x}_1 \dots \boldsymbol{x}_N]^T$. (D.1)

The challenge is that, the state matrix H has dependent rows; which will be addressed by carefully splitting the trajectory $\{u_t, h_t\}_{t=0}^N$ into multiple sub-trajectories which are internally weakly dependent as discussed in Section C. We first define the matrices obtained from these sub-trajectories.

Definition 24 Given sampling rate L and offset $\bar{\tau}$, consider the L-subsampled trajectory $\{\boldsymbol{h}^{(i)}, \boldsymbol{u}^{(i)}\}_{i=1}^{\bar{N}}$ as described in Definitions 20 and 21. Define the matrices $\bar{\boldsymbol{H}} = \bar{\boldsymbol{H}}^{(\bar{\tau})} \in \mathbb{R}^{\bar{N} \times n}, \ \tilde{\boldsymbol{H}} = \tilde{\boldsymbol{H}}^{(\bar{\tau})} \in \mathbb{R}^{\bar{N} \times n}, \ \tilde{\boldsymbol{U}} = \tilde{\boldsymbol{U}}^{(\bar{\tau})} \in \mathbb{R}^{\bar{N} \times p}, \ \text{and} \ \tilde{\boldsymbol{X}} = \tilde{\boldsymbol{X}}^{(\bar{\tau})} \in \mathbb{R}^{\bar{N} \times (n+p)} \ \text{as}$

$$m{ar{H}} = [m{ar{h}}^{(1)} \ \dots \ m{ar{h}}^{(ar{N})}]^T, \ m{ ilde{H}} = [m{h}^{(1)} \ \dots \ m{h}^{(ar{N})}]^T, \ m{ ilde{U}} = [m{u}^{(1)} \ \dots \ m{u}^{(ar{N})}]^T, \ m{ ilde{X}} = [m{\mu} m{ ilde{H}} \ m{ ilde{U}}].$$

Lemma 25 (Handling perturbation) Consider the nonlinear state equation (1.1). Given sampling rate L>0 and offset $\bar{\tau}$, consider the matrices $\bar{\boldsymbol{H}}$, $\tilde{\boldsymbol{H}}$, $\tilde{\boldsymbol{X}}$ of Definition 24 and let $\boldsymbol{Q}=[\gamma_+^{-1/2}\bar{\boldsymbol{H}}\;\tilde{\boldsymbol{U}}]\in\mathbb{R}^{\bar{N}\times(n+p)}$. Suppose Assumption 1 holds, ϕ is β -increasing, and $\boldsymbol{u}_t\stackrel{i.i.d.}{\sim}\boldsymbol{u}$ where \boldsymbol{u} is zero-mean with identity covariance and subgaussian norm bounded above by a constant. There exists an absolute constant C>0 such that if $\bar{N}\geq C\frac{\gamma_+^2}{\gamma_-^2}(n+p)$, with probability $1-8\exp(-c\frac{\gamma_-^2}{\gamma_+^2}\bar{N})$, for all matrices \boldsymbol{M} obeying $\|\boldsymbol{M}-\bar{\boldsymbol{H}}\|\leq \frac{\sqrt{\gamma_-\bar{N}}}{10}$, the perturbed \boldsymbol{Q} matrices given by,

$$\tilde{\boldsymbol{Q}} = [\gamma_+^{-1/2} \boldsymbol{M} \; \tilde{\boldsymbol{U}}], \tag{D.2}$$

satisfy

$$(\Theta + \sqrt{2})^2 \succeq \frac{\tilde{\boldsymbol{Q}}^T \tilde{\boldsymbol{Q}}}{\bar{N}} \succeq \frac{\gamma_-}{2\gamma_+}.$$
 (D.3)

Proof This result is a direct application of Theorem 29 after determining minimum/maximum eigenvalues of population covariance. The cross covariance obeys $\mathbb{E}[\bar{\boldsymbol{H}}^T\tilde{\boldsymbol{U}}]=0$ due to independence. Also, for i>1, the truncated state vector $\bar{\boldsymbol{h}}^{(i)}$ is statistically identical to \boldsymbol{h}_{L-1} hence $\boldsymbol{\Sigma}[\bar{\boldsymbol{h}}^{(i)}] \succeq \gamma_- \boldsymbol{I}_n$. Consequently,

$$\boldsymbol{\Sigma}[\boldsymbol{u}^{(i)}] = \boldsymbol{I}_p, \frac{1}{\gamma_+} \boldsymbol{\Sigma}[\bar{\boldsymbol{h}}^{(i)}] \leq \boldsymbol{I}_n \text{ for all } i \text{ and } \frac{\gamma_-}{\gamma_+} \boldsymbol{I}_n \leq \frac{1}{\gamma_+} \boldsymbol{\Sigma}[\bar{\boldsymbol{h}}^{(i)}] \text{ for all } i > 1. \text{ Hence, setting } \boldsymbol{q}_i = \begin{bmatrix} \frac{1}{\sqrt{\gamma_+}} \bar{\boldsymbol{h}}^{(i)} \\ \boldsymbol{u}^{(i)} \end{bmatrix},$$
 for all $i > 1$

$$rac{\gamma_-}{\gamma_+} oldsymbol{I}_n \preceq oldsymbol{\Sigma}[oldsymbol{q}_i] \preceq oldsymbol{I}_n.$$

Set the matrix $\bar{Q} = [q_2 \dots q_{\bar{N}}]^T$ and note that $Q = [q_1 \ \bar{Q}^T]^T$. Applying Theorem 29 on \bar{Q} and Corollary 30 on Q, we find that, with the desired probability,

$$\theta + \sqrt{3/2} \ge \frac{1}{\sqrt{N}} \|\boldsymbol{Q}\| \ge \frac{1}{\sqrt{N}} s_{\min}(\boldsymbol{Q}) \ge \frac{1}{\sqrt{N}} s_{\min}(\bar{\boldsymbol{Q}}) \ge \sqrt{\frac{N-1}{N}} \sqrt{\frac{2\gamma_{-}}{3\gamma_{+}}} \ge 0.99 \times \sqrt{\frac{2\gamma_{-}}{3\gamma_{+}}}.$$

Setting $E = M - \bar{H}$ and observing $\tilde{Q} = Q + [\gamma_+^{-1/2} E \ 0]$, the impact of the perturbation E can be bounded naively via $s_{\min}(Q) - \gamma_+^{-1/2} \|E\| \le s_{\min}(\tilde{Q}) \le \|\tilde{Q}\| \le \|Q\| + \gamma_+^{-1/2} \|E\|$. Using the assumed bound on $\|E\|$, this yields

$$\theta + \sqrt{2} \geq \frac{1}{\sqrt{\bar{N}}} \|\tilde{\boldsymbol{Q}}\| \geq \frac{1}{\sqrt{\bar{N}}} s_{\min}(\tilde{\boldsymbol{Q}}) \geq \sqrt{\frac{\gamma_{-}}{2\gamma_{+}}}.$$

This final inequality is identical to the desired bound (D.3).

Theorem 26 (Data matrix condition) Consider the nonlinear state-equation (1.1). Given $\gamma_+ \geq \gamma_- > 0$, define the condition number $\rho = \frac{\gamma_+}{\gamma_-}$. For some absolute constants c, C > 0, pick a trajectory length N where

$$L = \lceil 1 - \frac{\log(cn\rho)}{\log \|\mathbf{A}\|} \rceil$$
 , $N_0 = \lfloor \frac{N}{L} \rfloor \ge C\rho^2(n+p)$,

and pick scaling $\mu = \frac{1}{\sqrt{\gamma_+}}$. Suppose $\|A\| < 1$, ϕ is β -increasing, $u_t \overset{i.i.d.}{\sim} u$ where u is zero-mean with identity covariance and subgaussian norm bounded above by a constant, and Assumption 1 holds with $\gamma_+, \gamma_-, \theta, L$. Matrix $\boldsymbol{X} = [\boldsymbol{x}_1 \ \dots \ \boldsymbol{x}_N]^T$ of (D.1) satisfies the following with probability $1 - 4N \exp(-100n) - 8L \exp(-\mathcal{O}(N_0/\rho^2))$.

- Each row of X has ℓ_2 norm at most $c_0\sqrt{p+n}$ where c_0 is an absolute constant.
- X^TX obeys the bound

$$(\Theta + \sqrt{2})^2 \boldsymbol{I}_{n+p} \succeq \frac{\boldsymbol{X}^T \boldsymbol{X}}{N} \succeq \rho^{-1} \boldsymbol{I}_{n+p} / 2.$$
 (D.4)

Proof The first statement on ℓ_2 -norm bound can be concluded from Lemma 27 and holds with probability $1-2N\exp(-100(n+p))$. To show the second statement, for a fixed offset $1\leq \bar{\tau}\leq L$, consider Definition 24 and the matrices $\tilde{\boldsymbol{H}}^{(\bar{\tau})}, \tilde{\boldsymbol{U}}^{(\bar{\tau})}, \tilde{\boldsymbol{X}}^{(\bar{\tau})}$. Observe that \boldsymbol{X} is obtained by merging multiple sub-trajectory matrices $\{\tilde{\boldsymbol{X}}^{(\bar{\tau})}\}_{\bar{\tau}=1}^L$. We will first show the advertised bound for an individual $\tilde{\boldsymbol{X}}^{(\bar{\tau})}$ by applying Lemma 25 and then apply Lemma 8 to obtain the bound on the combined matrix \boldsymbol{X} .

Recall that $\bar{N}_{\bar{\tau}}$ is the length of the $\bar{\tau}$ th sub-trajectory i.e. number of rows of $\tilde{\boldsymbol{X}}^{(\bar{\tau})}$. By construction $2N_0 \geq \bar{N}_{\bar{\tau}} \geq N_0$ for all $1 \leq \bar{\tau} \leq L$. Given $1 \leq \bar{\tau} \leq L$ and triple $\bar{\boldsymbol{H}}^{(\bar{\tau})}, \tilde{\boldsymbol{H}}^{(\bar{\tau})}, \tilde{\boldsymbol{U}}^{(\bar{\tau})}$, set $\boldsymbol{Q} = [\mu \bar{\boldsymbol{H}}^{(\bar{\tau})}, \tilde{\boldsymbol{U}}^{(\bar{\tau})}]$. Since N_0 is chosen to be large enough, applying Lemma 25 with $\mu = 1/\sqrt{\gamma_+}$ choice, and noting $\rho = \gamma_+/\gamma_-$, we find that, with probability $1 - 4 \exp(-c_1 N_0/\rho^2)$, all matrices \boldsymbol{M} satisfying $\|\boldsymbol{M} - \bar{\boldsymbol{H}}^{(\bar{\tau})}\| \leq \sqrt{\gamma_- N_0}/10$ and $\tilde{\boldsymbol{Q}}$ as in (D.2) obeys

$$(\Theta + \sqrt{2})^2 \succeq \frac{\tilde{\mathbf{Q}}^T \tilde{\mathbf{Q}}}{N} \succeq \rho^{-1}/2.$$
 (D.5)

Let us call this Event 1. To proceed, we will argue that with high probability $\|\tilde{\boldsymbol{H}}^{(\bar{\tau})} - \bar{\boldsymbol{H}}^{(\bar{\tau})}\|$ is small so that the bound above is applicable with $\boldsymbol{M} = \tilde{\boldsymbol{H}}^{(\bar{\tau})}$ choice; which sets $\tilde{\boldsymbol{Q}} = \tilde{\boldsymbol{X}}^{(\bar{\tau})}$ in (D.5). Applying Lemma 23, we find that, with probability $1 - 2\bar{N}_{\bar{\tau}} \exp(-100n)$,

$$\|\bar{\boldsymbol{H}}^{(\bar{\tau})} - \tilde{\boldsymbol{H}}^{(\bar{\tau})}\| \le \sqrt{2N_0} \max\{\|\boldsymbol{h}^{(i)} - \bar{\boldsymbol{h}}^{(i)}\|_{\ell_2}\} \le c_0 \sqrt{2N_0} \sqrt{n\gamma_+} \|\boldsymbol{A}\|^{L-1}.$$

Let us call this Event 2. We will show that our choice of L ensures right hand side is small enough and guarantees $\|\bar{\boldsymbol{H}}^{(\bar{\tau})} - \tilde{\boldsymbol{H}}^{(\bar{\tau})}\| \leq \sqrt{\gamma_- N_0}/10$. Set $c = \max\{200c_0^2, 1\}$. Desired claim follows by taking logarithms of upper/lower bounds and cancelling out $\sqrt{N_0}$ terms as follows

$$c_0\sqrt{n}\|\mathbf{A}\|^{L-1}\sqrt{\gamma_+} \le \sqrt{\gamma_-}/10\sqrt{2} \iff (L-1)\log\|\mathbf{A}\| + \log\sqrt{cn\rho} \le 0$$
 (D.6)

$$\iff -\frac{\log cn\rho}{2\log\|\boldsymbol{A}\|} \le L - 1 \tag{D.7}$$

$$\iff L = \lceil 1 - \frac{\log(cn\rho)}{\log\|\mathbf{A}\|} \rceil. \tag{D.8}$$

Here we use the fact that $\log \|\boldsymbol{A}\| < 0$ since $\|\boldsymbol{A}\| < 1$ and $cn\rho \geq 0$. Consequently, both Event 1 and Event 2 hold with probability $1 - 4\exp(-c_1N_0/\rho^2) - 2\bar{N}_{\bar{\tau}}\exp(-100n)$, implying (D.5) holds with $\tilde{\boldsymbol{Q}} = \tilde{\boldsymbol{X}}^{(\bar{\tau})}$. Union bounding this over $1 \leq \bar{\tau} \leq L$, (D.5) uniformly holds with $\tilde{\boldsymbol{Q}} = \tilde{\boldsymbol{X}}^{(\bar{\tau})}$ and all rows of \boldsymbol{X} are ℓ_2 -bounded with probability $1 - 4N\exp(-100n) - 8L\exp(-c_1N_0/\rho^2)$. Applying Lemma 8 on $(\tilde{\boldsymbol{X}}^{(\bar{\tau})})_{\bar{\tau}=1}^L$, we conclude with the bound (D.4) on the merged matrix \boldsymbol{X} .

Lemma 27 (ℓ_2 -bound on rows) Consider the setup of Theorem 26. With probability $1 - 2N \exp(-100(n + p))$, each row of X has ℓ_2 -norm at most $c\sqrt{p+n}$ for some constant c > 0.

Proof The tth row of \boldsymbol{X} is equal to $\boldsymbol{x}_t = [\frac{\boldsymbol{h}_t^T}{\sqrt{\gamma_+}} \ \boldsymbol{u}_t^T]^T$. Since $\|\boldsymbol{h}_t - \mathbb{E}[\boldsymbol{h}_t]\|_{\psi_2} \leq \mathcal{O}(\sqrt{\gamma_+})$ and $\|\boldsymbol{u}_t\|_{\psi_2} \leq \mathcal{O}(1)$, we have that $\|\boldsymbol{x}_t - \mathbb{E}[\boldsymbol{x}_t]\|_{\psi_2} \leq \mathcal{O}(1)$. Now, applying Lemma 31 on all rows $\{\boldsymbol{x}_t\}_{t=1}^N$, and using a union bound, with probability at least $1 - 2N \exp(-100(n+p))$, we have that $\|\boldsymbol{x}_t - \mathbb{E}[\boldsymbol{x}_t]\|_{\ell_2} \leq c\sqrt{n+p}$ for all t. To conclude, note that $\|\mathbb{E}[\boldsymbol{x}_t]\|_{\ell_2} = \|\mathbb{E}[\boldsymbol{h}_t]\|_{\ell_2}/\sqrt{\gamma_+} \leq \theta \leq 3\sqrt{n}$ via Assumption 1.

Appendix E. Proofs of Main Results

E.1. Proof of Lemma 2

The statement follows from upper bound Lemma 11 and lower bound Lemma 13.

E.2. Proof of Theorem 6

Proof To prove this theorem, we combine Theorem 26 with deterministic SGD convergence result of Theorem 5. Applying Theorem 26, with the desired probability, inequality (D.4) holds and for all i, input data satisfies the bound $\|\boldsymbol{x}_i\|_{\ell_2} \leq \sqrt{(n+p)/(2c_0)}$ for a sufficiently small constant $c_0 > 0$. As the next step, we will argue that these two events imply the convergence of SGD.

Let $\boldsymbol{\theta}^{(i)}, \boldsymbol{c}^{(i)} \in \mathbb{R}^{n+p}$ denote the ith rows of $\boldsymbol{\Theta}, \boldsymbol{C}$ respectively. Observe that the square-loss is separable along the rows of \boldsymbol{C} via $\|\boldsymbol{\Theta} - \boldsymbol{C}\|_F^2 = \sum_{i=1}^n \|\boldsymbol{\theta}^{(i)} - \boldsymbol{c}^{(i)}\|_{\ell_2}^2$. Hence, SGD updates each row $\boldsymbol{c}^{(i)}$ via its own state equation

$$\boldsymbol{y}_{t,i} = \phi(\left\langle \boldsymbol{c}^{(i)}, \boldsymbol{x}_t \right\rangle),$$

where $y_{t,i}$ is the *i*th entry of y_t . Consequently, we can establish the convergence result for an individual row of C. Convergence of all individual rows will imply the convergence of the overall matrix Θ_{τ} to the ground

truth C. Pick a row index i $(1 \le i \le n)$, set $c = c^{(i)}$ and denote ith row of Θ_{τ} by θ_{τ} . Also denote the label corresponding to ith row by $y_t = y_{t,i}$. With this notation, SGD over (2.3) runs SGD over the ith row with equations $y_t = \phi(\langle c, x_t \rangle)$ and with loss functions

$$\mathcal{L}(\boldsymbol{\theta}) = N^{-1} \sum_{t=1}^{N} \mathcal{L}_t(\boldsymbol{\theta}), \ \mathcal{L}_t(\boldsymbol{\theta}) = \frac{1}{2} (y_t - \phi(\langle \boldsymbol{\theta}, \boldsymbol{x}_t \rangle))^2.$$

Substituting our high-probability bounds on x_t (e.g. (D.4)) into Theorem 5, we can set $B = (n+p)/(2c_0)$, $\gamma_+ = (\theta + \sqrt{2})^2$, and $\gamma_- = \rho^{-1}/2$. Consequently, using the learning rate $\eta = c_0 \frac{\beta^2 \rho^{-1}}{(\theta + \sqrt{2})^2 (n+p)}$, for all $\tau \geq 0$, the τ th SGD iteration θ_τ obeys

$$\mathbb{E}[\|\boldsymbol{\theta}_{\tau} - \boldsymbol{c}\|_{\ell_{2}}^{2}] \leq \|\boldsymbol{\theta}_{0} - \boldsymbol{c}\|_{\ell_{2}}^{2} (1 - c_{0} \frac{\beta^{4} \rho^{-2}}{2(\theta + \sqrt{2})^{2}(n+p)})^{\tau},$$
(E.1)

where the expectation is over the random selection of SGD updates. This establishes the convergence for a particular row of C. Summing up these inequalities (E.1) over all rows $\theta_{\tau}^{(1)}, \ldots, \theta_{\tau}^{(n)}$ (which converge to $c^{(1)}, \ldots, c^{(n)}$ respectively) yields the targeted bound (4.4).

E.3. Proofs of main results on stable systems

E.3.1. PROOF OF THEOREM 3

Proof Applying Lemmas 11 and 2, independent of L, Assumption 1 holds with parameters

$$\gamma_{+} = B_{\infty}^{2}$$
 , $\gamma_{-} = \beta^{2} s_{\min}(\mathbf{B})^{2}$, $\theta = \sqrt{6n} - \sqrt{2} \ge \sqrt{n}$.

This yields $(\theta + \sqrt{2})^2 = 6n$. Hence, we can apply Theorem 6 with the learning rate $\eta = c_0 \frac{\beta^2}{6\rho n(n+p)}$ where

$$\rho = \frac{B_{\infty}^2}{\beta^2 s_{\min}(\mathbf{B})^2} = \frac{\gamma_+}{\gamma_-},\tag{E.2}$$

and convergence rate $1-\frac{\beta^2\eta}{2\rho}$. To conclude with the stated result, we use the change of variable $c_0/6\to c_0$.

E.3.2. PROOF OF THEOREM 4

Proof The proof is similar to that of Theorem 3. Applying Lemmas 11, 12, and 2, independent of L, Assumption 1 holds with parameters

$$\gamma_+ = B_\infty^2$$
 , $\gamma_- = s_{\min}(\boldsymbol{B})^2$, $\theta = 0$.

Hence, we again apply Theorem 6 with the learning rate $\eta = c_0 \frac{\beta^2}{2\rho(n+p)}$ where ρ is given by (E.2). Use the change of variable $c_0/2 \to c_0$ to conclude with the stated result.

E.4. Learning unstable systems

In a similar fashion to Section 4, we provide a more general result on unstable systems that makes a parametric assumption on the statistical properties of the state vector.

Assumption 2 (Well-behaved state vector – single timestamp) Given timestamp $T_0 > 0$, there exists positive scalars $\gamma_+, \gamma_-, \theta$ and an absolute constant C > 0 such that $\theta \le 3\sqrt{n}$ and the following holds

$$\gamma_+ I_n \succeq \Sigma[h_{T_0}] \succeq \gamma_- I_n$$
 , $\|h_{T_0} - \mathbb{E}[h_{T_0}]\|_{\psi_2} \le C\sqrt{\gamma_+}$ and $\|\mathbb{E}[h_t]\|_{\ell_2} \le \theta\sqrt{\gamma_+}$. (E.3)

The next theorem provides the parametrized result on unstable systems based on this assumption.

Theorem 28 (Unstable system - general) Suppose we are given N independent trajectories $(\boldsymbol{h}_t^{(i)}, \boldsymbol{u}_t^{(i)})_{t\geq 0}$ for $1\leq i\leq N$. Sample each trajectory at time T_0 to obtain N samples $(\boldsymbol{y}_i, \boldsymbol{h}_i, \boldsymbol{u}_i)_{i=1}^N$ where ith sample is

$$(\boldsymbol{y}_i, \boldsymbol{h}_i, \boldsymbol{u}_i) = (\boldsymbol{h}_{T_0+1}^{(i)}, \boldsymbol{h}_{T_0}^{(i)}, \boldsymbol{u}_{T_0}^{(i)}).$$

Let $C, c_0 > 0$ be absolute constants. Suppose Assumption 1 holds with T_0 and sample size satisfies $N \ge C\rho^2(n+p)$ where $\rho = \gamma_+/\gamma_-$. Assume ϕ is β -increasing, zero initial state conditions, and $\mathbf{u}_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \mathbf{I}_p)$. Set scaling to be $\mu = 1/\sqrt{\gamma_+}$ and learning rate to be $\eta = c_0 \frac{\beta^2}{\rho(\theta + \sqrt{2})^2(n+p)}$. Starting from $\mathbf{\Theta}_0$, we run SGD over the equations described in (2.2) and (2.3). With probability $1 - 2N \exp(-100(n+p)) - 4 \exp(-\mathcal{O}(\frac{N}{\rho^2}))$, all iterates satisfy

$$\mathbb{E}[\|\Theta_i - C\|_F^2] \le (1 - c_0 \frac{\beta^4}{2\rho^2(\theta + \sqrt{2})^2(n+p)})^{\tau} \|\Theta_0 - C\|_F^2,$$

where the expectation is over the randomness of the SGD updates.

Proof Set $x_i = [\gamma_+^{-1/2} h_i^T u_i^T]^T$ and $X = [x_1 \dots x_N]^T$. Since X has i.i.d. rows, we can apply Theorem 29 and Lemma 31 to find with the desired probability that

- Rows of x_i satisfy $\|x_i \mathbb{E}[x_i]\|_{\psi_2} \leq \mathcal{O}(1)$ and $\mathbb{E}[\|x_i\|_{\ell_2}] \leq 3\sqrt{n}$, hence all rows of X obeys $\|x_i\|_{\ell_2} \leq \sqrt{(n+p)/(2c_0)}$,
- X satisfies

$$(\theta + \sqrt{2})^2 \succeq \frac{\boldsymbol{X}^T \boldsymbol{X}}{N} \succeq \rho^{-1}/2.$$

To proceed, using $\gamma_- = \rho^{-1}/2$, $B = (n+p)/(2c_0)$, and $\gamma_+ = (\theta + \sqrt{2})^2$, we apply Theorem 5 on the loss function (2.3); which yields the desired result.

E.5. Proof of Theorem 7

Proof The proof is a corollary of Theorem 28. We need to substitute the proper values in Assumption 2. Applying Lemma 11, we can substitute $\gamma_+ = B_{T_0}^2$ and $\theta = \sqrt{6n} - \sqrt{2} \ge \sqrt{n}$. Next, we need to find a lower bound. Applying Lemma 2 for n > 1 and Lemma 14 for n = 1, we can substitute $\gamma_- = \gamma_+/\rho$ with the ρ definition of (5.2). With these, the result follows as an immediate corollary of Theorem 28.

Appendix F. Supplementary Statistical Results

The following theorem bounds the empirical covariance of matrices with independent subgaussian rows. Given a random vector x, define the de-biasing operation as $\mathbf{zm}(x) = x - \mathbb{E}[x]$.

Theorem 29 Let $A \in \mathbb{R}^{n \times d}$ be a matrix with independent subgaussian rows $\{a_i\}_{i=1}^n$ satisfying $\|\mathbf{zm}(a_i)\|_{\psi_2} \leq \mathcal{O}(K)$ and $\Sigma[a_i] \leq K^2 \mathbf{I}_d$ for some K > 0 and $\|\mathbb{E}[a_i]\|_{\ell_2} \leq \theta$. Suppose $\Sigma[a_i] \succeq \lambda \mathbf{I}_d$. Suppose $n \geq \mathcal{O}(K^4 d/\lambda^2)$. Then, each of the following happens with probability at least $1 - 2\exp(-cK^{-4}\lambda^2 n)$,

- $\theta + \sqrt{3/2}K \geq \frac{1}{\sqrt{n}} \|\boldsymbol{A}\|$.
- Suppose all rows of **A** have equal expectations. Then $\frac{1}{\sqrt{n}}s_{\min}(\mathbf{A}) \geq \sqrt{2\lambda/3}$.

Proof Let $E = \mathbb{E}[A]$, $\bar{A} = A - \mathbb{E}[A]$, $\bar{a}_i = \mathbf{zm}(a_i)$. We will decompose $A = \bar{A} + E$ hence we will first focus on bounding the upper and lower singular values of \bar{A} by studying the random processes $X_v = \|\bar{A}v\|_{\ell_2}^2$ and $Y_v = X_v - \mathbb{E}[X_v]$ over the unit sphere \mathcal{S}^{d-1} . First, we provide a deviation bound for the quantity $\sup_{v \in \mathcal{S}^{d-1}} |Y_v|$. To achieve this, we will utilize Talagrand's mixed tail bound and show that increments of Y_v are subexpoential. Pick two unit vectors $v, u \in \mathbb{R}^d$. Write x = u + v, y = u - v. We have that

$$X_{\boldsymbol{u}} - X_{\boldsymbol{v}} = \|\bar{\boldsymbol{A}}\boldsymbol{u}\|_{\ell_2}^2 - \|\bar{\boldsymbol{A}}\boldsymbol{v}\|_{\ell_2}^2 = \|\bar{\boldsymbol{A}}(\boldsymbol{x} + \boldsymbol{y})/2\|_{\ell_2}^2 - \|\bar{\boldsymbol{A}}(\boldsymbol{x} - \boldsymbol{y})/2\|_{\ell_2}^2 = \boldsymbol{x}^T\bar{\boldsymbol{A}}^T\bar{\boldsymbol{A}}\boldsymbol{y} = \sum_{i=1}^n (\bar{\boldsymbol{a}}_i^T\boldsymbol{x})(\bar{\boldsymbol{a}}_i^T\boldsymbol{y}).$$

Letting $\hat{x} = x/\|x\|_{\ell_2}$, $\hat{y} = y/\|y\|_{\ell_2}$, observe that, multiplication of subgaussians $x^T \bar{a}_i, y^T \bar{a}_i$ obey

$$\|(\boldsymbol{x}^T \bar{\boldsymbol{a}}_i)(\boldsymbol{y}^T \bar{\boldsymbol{a}}_i)\|_{\psi_1} \leq \mathcal{O}(\|\boldsymbol{x}\|_{\ell_2} \|\boldsymbol{y}\|_{\ell_2} K^2) \leq \mathcal{O}(K^2 \|\boldsymbol{y}\|_{\ell_2}).$$

Centering this subexponential variable around zero introduces a factor of 2 when bounding subexponential norm and yields $\|(\boldsymbol{x}^T\bar{\boldsymbol{a}}_i)(\boldsymbol{y}^T\bar{\boldsymbol{a}}_i) - \mathbb{E}[(\boldsymbol{x}^T\bar{\boldsymbol{a}}_i)(\boldsymbol{y}^T\bar{\boldsymbol{a}}_i)]\|_{\psi_1} \leq \mathcal{O}(K^2\|\boldsymbol{y}\|_{\ell_2})$. Now, using the fact that $Y_{\boldsymbol{u}} - Y_{\boldsymbol{v}}$ is sum of n independent zero-mean subexponential random variables, we have the tail bound

$$\mathbb{P}(n^{-1}|Y_{\boldsymbol{u}} - Y_{\boldsymbol{v}}| \ge t) \le 2\exp(-c'n\min\{\frac{t^2}{K^4\|\boldsymbol{y}\|_{\ell_2}^2}, \frac{t}{K^2\|\boldsymbol{y}\|_{\ell_2}}\}).$$

Applying Talagrand's chaining bound for mixed tail processes with distance metrics $\rho_2 = \frac{K^2\|\cdot\|_{\ell_2}}{\sqrt{n}}, \rho_1 = \frac{K^2\|\cdot\|_{\ell_2}}{n}$, (Theorem 3.5 of Dirksen (2013) or Theorem 2.2.23 of Talagrand (2014)) and using the fact that for unit sphere \mathcal{S}^{d-1} , Talagrand's γ functionals (see Talagrand (2014)) obey $\gamma_1(\mathcal{S}^{d-1}), \gamma_2^2(\mathcal{S}^{d-1}) \leq \mathcal{O}(d)$,

$$n^{-1} \sup_{v \in S^{d-1}} |Y_v| \le cK^2 (\sqrt{d/n} + d/n + t/\sqrt{n}), \tag{F.1}$$

with probability $1-2\exp(-\min\{t^2,\sqrt{n}t\})$. Since $n\geq C\lambda^{-2}K^4d$ for sufficiently large C>0, picking $t=\frac{1}{16c}K^{-2}\lambda\sqrt{n}$, with probability $1-2\exp(-\mathcal{O}(K^{-4}\lambda^2n))$, we ensure that right hand side of (F.1) is less than $\lambda/8$. This leads to the following inequalities

$$\frac{1}{n} \|\bar{\boldsymbol{A}}^T \bar{\boldsymbol{A}} - \mathbb{E}[\bar{\boldsymbol{A}}^T \bar{\boldsymbol{A}}]\| \le \frac{\lambda}{8} \implies \frac{9K^2}{8} \boldsymbol{I}_d \succeq \frac{1}{n} \bar{\boldsymbol{A}}^T \bar{\boldsymbol{A}} \succeq \frac{7\lambda}{8} \boldsymbol{I}_d. \tag{F.2}$$

$$\implies \frac{9}{8} K \ge \frac{1}{\sqrt{n}} \|\bar{\boldsymbol{A}}\| \ge s_{\min}(\bar{\boldsymbol{A}}) \ge \sqrt{\frac{7}{8}} \lambda.$$

Upper bound on spectral norm: For spectral norm of A, we use the triangle inequality

$$\frac{1}{\sqrt{n}} \|\mathbf{A}\| \le \frac{1}{\sqrt{n}} (\|\mathbf{E}\| + \|\bar{\mathbf{A}}\|) \le \max_{1 \le i \le n} \|\mathbb{E}[\mathbf{a}_i]\|_{\ell_2} + 9K/8 \le \theta + \sqrt{3/2}K.$$

Lower bound on minimum singular value: This part assumes that all row expectations are same. Denote the size n all ones vector by $\mathbf{1}_n$ and define the process $Z_{\boldsymbol{v}} = \frac{1}{\sqrt{n}} \mathbf{1}_n^T \bar{A} \boldsymbol{v}$. Observe that $\bar{A}^T \mathbf{1}_n = \sum_{i=1}^n \bar{a}_i \in \mathbb{R}^d$ is a vector satisfying $\|\bar{A}^T \mathbf{1}_n / \sqrt{n}\|_{\psi_2} \leq \mathcal{O}(K)$. Hence, again using $n \geq CK^4 \lambda^{-2}d$ for sufficiently large C > 0, applying Lemma 31 with $m = c_0 K^{-4} \lambda^2 n > d$ by picking a sufficiently small constant $c_0 > 1/C$, with probability at least $1 - 2 \exp(-100c_0 K^{-4} \lambda^2 n)$

$$\frac{1}{\sqrt{n}} \sup_{\|\mathbf{v}\|_{\ell_2} = 1} |Z_{\mathbf{v}}| = \frac{1}{n} \|\bar{\mathbf{A}}^T \mathbf{1}_n\|_{\ell_2} \le \frac{1}{12} K K^{-2} \lambda \le \frac{\sqrt{\lambda}}{12}.$$

Let $P = I_n - \frac{1}{n} \mathbf{1}_n \mathbf{1}_n^T$ be the projection onto the orthogonal complement of the all ones vector. Note that PEv = 0 as the rows of E are equal. With this observation, with desired probability, for any unit length v,

$$||Av||_{\ell_2} \ge ||PAv||_{\ell_2} = ||P\bar{A}v||_{\ell_2} \ge ||\bar{A}v||_{\ell_2} - |Z_v|$$
 (F.3)

$$\geq s_{\min}(\bar{\mathbf{A}}) - \sup_{\mathbf{v} \in \mathcal{S}^{d-1}} |Z_{\mathbf{v}}| \geq (\sqrt{7/8} - 1/12)\sqrt{\lambda n},\tag{F.4}$$

which implies $s_{\min}(\mathbf{A})/\sqrt{n} \geq \sqrt{2\lambda/3}$.

The corollary below is obtained by slightly modifying the proof above by using $\frac{1}{n} \|\bar{A}^T \bar{A} - \mathbb{E}[\bar{A}^T \bar{A}]\| \leq \frac{K^2}{8}$ in line (F.2) and only focusing on the spectral norm bound.

Corollary 30 Let $A \in \mathbb{R}^{n \times d}$ be a matrix with independent $\{a_i\}_{i=1}^n$ subgaussian rows satisfying $\|\mathbf{zm}(a_i)\|_{\psi_2} \leq \mathcal{O}(K)$ and $\mathbf{\Sigma}[a_i] \leq K^2 \mathbf{I}_d$ for some K > 0 and $\|\mathbb{E}[a_i]\|_{\ell_2} \leq \theta$. Suppose $\mathbf{\Sigma}[a_i] \succeq \lambda \mathbf{I}_d$. Suppose $n \geq \mathcal{O}(K^2 d)$. Then, with probability at least $1 - 4 \exp(-cK^{-2}n)$,

$$\theta + \sqrt{3/2}K \ge \frac{1}{\sqrt{n}} \|\boldsymbol{A}\|.$$

The following lemma is fairly standard and is proved for the sake of completeness.

Lemma 31 (Subgaussian vector length) Let $a \in \mathbb{R}^n$ be a zero-mean subgaussian vector with $||a||_{\psi_2} \leq L$. Then, for any $m \geq n$, there exists C > 0 such that

$$\mathbb{P}(\|\boldsymbol{a}\|_{\ell_2} \le CL\sqrt{m}) \ge 1 - 2\exp(-100m).$$

Proof We can pick a 1/2 cover $\mathcal C$ of the unit ℓ_2 -sphere with size $\log |\mathcal C| \leq 2n$. For any $v \in \mathcal C$, subgaussianity implies, $\mathbb P(|v^T a| \geq t) \leq 2 \exp(-\frac{ct^2}{2L^2})$. Setting $t = CL\sqrt{m}$ for sufficiently large constant C > 0, and union bounding over all $v \in \mathcal C$, we find

$$\mathbb{P}(\bigcap_{\boldsymbol{v} \in \mathcal{C}} \|\boldsymbol{v}\|_{\ell_2} \le CL\sqrt{m}) \ge 1 - 2\exp(2n - \frac{cC^2L^2m}{2L^2}) \le 1 - 2\exp(-100m).$$

To conclude, let $v(a) \in \mathcal{C}$ be a's neighbor satisfying $\|v - \frac{a}{\|a\|_{\ell_2}}\|_{\ell_2} \le 1/2$. Hence, we have

$$\|\boldsymbol{a}\|_{\ell_2} \le \|(\boldsymbol{a} - \boldsymbol{v}(\boldsymbol{a}))^T \boldsymbol{a}\|_{\ell_2} + \|\boldsymbol{v}^T \boldsymbol{a}\|_{\ell_2} \le \|\boldsymbol{a}\|_{\ell_2}/2 + CL\sqrt{m} \implies \|\boldsymbol{a}\|_{\ell_2} \le 2CL\sqrt{m}.$$

To conclude, use the change of variable $C \to C/2$.