

Active Learning in Dynamical Systems

Learning the behavior of an unknown dynamical system is a central problem in many applications

- RL: Exploration required for good performance, but how to explore efficiently is unclear
- Robotics: Dynamics of robots can be hard to model
- Airplane control: Different modes can be (de-) activated, and the current configuration needs to be detected

Problem: Obtaining **data** for learning often is **expensive** or induces a loss in performance

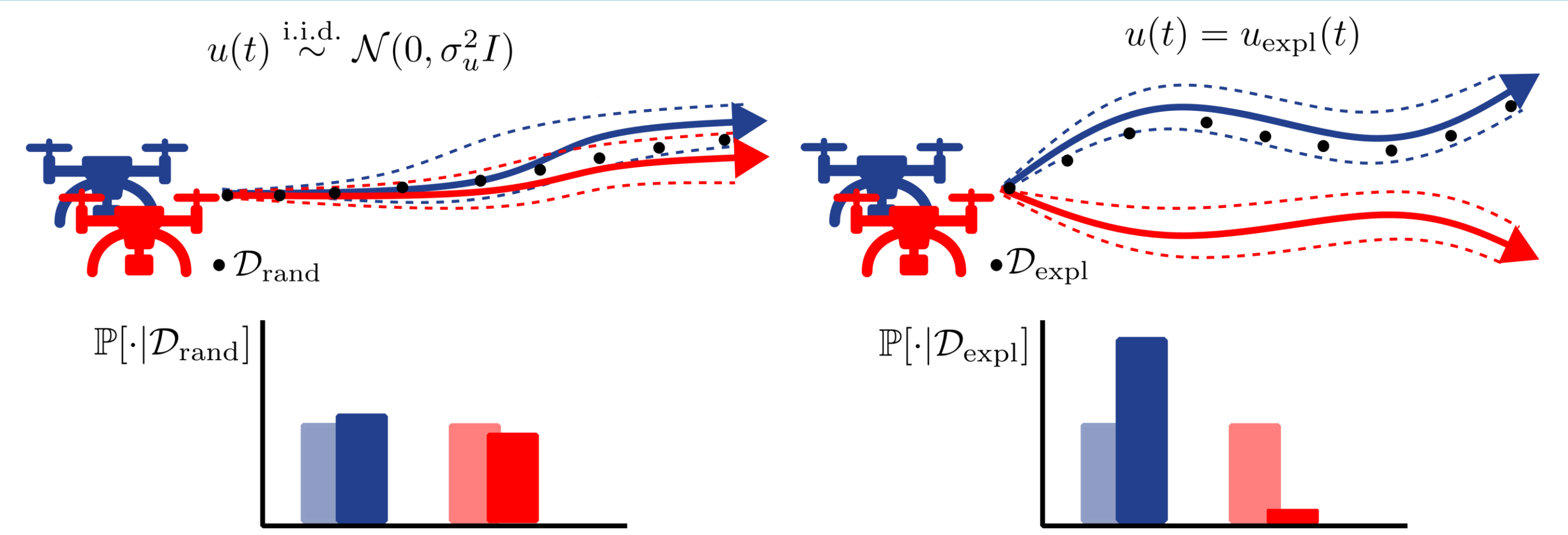
➔ Active Learning methods select the excitation input to **accelerate** the learning process

Limitations

- Isotropic Gaussian excitations are default exploration strategy, benefit of active learning methods often unclear
- Sample complexity lower bounds are restricted to specific input classes
- No modular framework to establish sample complexity upper bounds for general inputs
- Multi-model perspective has not been analyzed

Contributions

- Instance-dependent sample complexity lower bounds uncovering the potential **benefit of active learning**
- Tailored persistency of excitation (PE) notion, characterizing inputs yielding informative data
- **Modular framework** to provide sample complexity upper bounds for any PE input sequence
- Provably optimal sequential active learning algorithm with **close to oracle performance**



Setup:

- Unknown LTI system $x(t+1) = A_*x(t) + B_*u(t) + w(t)$, $\theta_* = (A_*, B_*)$
- Process noise $w(t) \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \Sigma_w)$
- Finite hypothesis class $(A_*, B_*) \in \mathcal{S} = \{(A_0, B_0), \dots, (A_N, B_N)\}$
- Bounded input energy $\mathbb{E} \left[\sum_{t=0}^{T-1} \|u(t)\|^2 \right] \leq \gamma_u^2 T$
- **Goal:** Identify the true system from data as fast as possible
- **Key Fact:** For simplicity, assume $x(0) = 0$. Define $U = [u(0)^\top \dots u(T)^\top]^\top$, then

$$\mathbb{E} \left[\sum_{s=0}^t \|\Delta A_i x(s) + \Delta B_i u(s)\|_{\Sigma_w^{-1}}^2 \right] = U^\top W_i(t) U + \text{Excitation due to noise.}$$
 - W_i captures the difficulty of distinguishing between θ_i and θ_* when applying input sequence U
 - $\Delta A_i = A_* - A_i$, $\Delta B_i = B_* - B_i$

Sample Complexity Lower Bound

- **δ -correct algorithm:** Any algorithm that yields $\mathbb{P}[\hat{\theta}_t \neq \theta_*] \leq \delta$, for $t \geq \bar{T}$.
- **Sample complexity lower bound:** For any excitation input sequence U and any δ -correct algorithm, it holds that

$$\min_{i \in [1, N]} \mathbb{E}[U^\top W_i(\bar{T})U] + \text{Excitation due to noise} \geq 2 \log \left(\frac{1}{2.4\delta} \right)$$

Further the lower bound is minimized by

$$U^* \in \arg \max_{U^\top U \leq \gamma_u^2 \bar{T}} \min_{i \in [1, N]} U^\top W_i U + \text{Excitation due to noise}$$

and when applying U^* any δ -correct algorithm satisfies

$$\min_{p \in \Delta_p} \gamma_u^2 \bar{T} \lambda_{\max} \left(\sum_{i=1}^N p_i W_i(\bar{T}) \right) + \text{Excitation due to noise} \geq 2 \log \left(\frac{1}{2.4\delta} \right)$$

Finally, for $u(t) \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \frac{\gamma_u^2}{n_u} I)$ and any δ -correct algorithm it holds that

$$\min_{p \in \Delta_N} \gamma_u^2 \bar{T} \lambda_{\text{mean}} \left(\sum_{i=1}^N p_i W_i(\bar{T}) \right) + \text{Excitation due to noise} \geq 2 \log \left(\frac{1}{2.4\delta} \right)$$

- ➔ The matrices W_i are the key quantities influencing the sample complexity lower bound
- ➔ Condition number of W_i determines the potential benefit of active learning

Sequential Active Learning Algorithm

- Optimal oracle excitation requires exact model knowledge ➔ Sequential algorithm

Algorithm 2 Input design subroutine

Require: \mathcal{S} , epoch length τ , prediction errors $\varepsilon_{\theta_i}((k-1)\tau)$, state x , scaling $\rho_k \in [0, 1]$

- 1: Compute weights using exponential weighting $w_{k+1}(i) = \exp(-\eta \varepsilon_{\theta_i}((k-1)\tau))$
- 2: Sample $\hat{i}_k \sim p_k(i)$, where $p_k(i) = w_k(i) / \sum_{j=0}^N w_k(j)$
- 3: Set $\hat{\theta}_k = \theta_{\hat{i}_k}$ and compute optimal input sequence $U_{\hat{\theta}_k}^*(x)$ by solving (28)
- 4: Define excitation according to ρ_k : $u_k^*(t) = \sqrt{1 - \rho_k} u_{\hat{\theta}_k}^*(t) + \sqrt{\rho_k} u_\eta(t)$, $u_\eta(t) \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \frac{\gamma_u^2}{n_u} I_{n_u})$
return excitation sequence u_k^*

- **Optimality of Alg. 2:** Let $\rho_k \xrightarrow{k \rightarrow \infty} 0$ slow enough. Then, for any $\delta \in (0, 1)$ the inputs derived by Alg. 2 yield $\mathbb{P}[\hat{\theta}_k \neq \theta_*] \leq \delta$ after at most k_δ epochs with

$$\lim_{\delta \rightarrow 0} \frac{c' \log \left(\frac{N}{\delta} \right)}{k_\delta} \leq \gamma_u^2 \min_{p \in \Delta_N} \lambda_{\max} \left(\sum_{i=1}^N p_i W_i(\tau) \right) + \text{Excitation due to noise}$$

- Input derived by Alg. 2 are optimal as $\delta \rightarrow 0$
- Sample complexity upper bound follows directly from PE

PE and Sample Complexity Upper Bound

- **Persistency of Excitation (PE):** An input sequence is PE for (θ_*, \mathcal{S}) if

$$\frac{1}{\tau} \sum_{t=0}^{\tau-1} \mathbb{E} \left[\|\Delta A_i x(t) + \Delta B_i u(t)\|_{\Sigma_w^{-1}}^2 \right] \geq c_u(\tau) \gamma_u^2 + \text{Excitation due to noise}$$

for a constant $c_u(\tau)$.

- Isotropic Gaussian excitation is PE with $c_u^{\text{rand}} = \min_{p \in \Delta_N} \lambda_{\text{mean}} \left(\sum_{i=1}^N p_i W_i(\tau) \right)$
- U^* is PE with $c_u^{\text{opt}} = \min_{p \in \Delta_N} \lambda_{\max} \left(\sum_{i=1}^N p_i W_i(\tau) \right)$

Algorithm 1 Sequential identification algorithm

Require: \mathcal{S} , epoch length τ , desired confidence δ

- 1: **for** $k = 1, 2, \dots$ **do**
- 2: Collect data using PE excitation input $\{u(t)\}_{t=(k-1)\tau}^{k\tau-1}$ with coefficients $c_{u_k}(\tau)$, $c_w(\tau)$
- 3: Compute $\varepsilon_{\theta_i}(k\tau)$ for all $i \in [0, N]$
- 4: **if** $\exists \hat{\theta} \in \mathcal{S} : \varepsilon_{\theta_i}(k\tau) - \varepsilon_{\hat{\theta}}(k\tau) > 2 \log \left(\frac{N}{\delta} \right)$ for all $\theta_i \in \mathcal{S} \setminus \hat{\theta}$ **then**
- 5: Stop and **return** estimate $\hat{\theta}$

- Prediction error $\varepsilon_{\theta_i}(t) = \sum_{s=0}^{t-1} \|x(s+1) - A_i x(s) - B_i u(s)\|_{\Sigma_w^{-1}}^2$

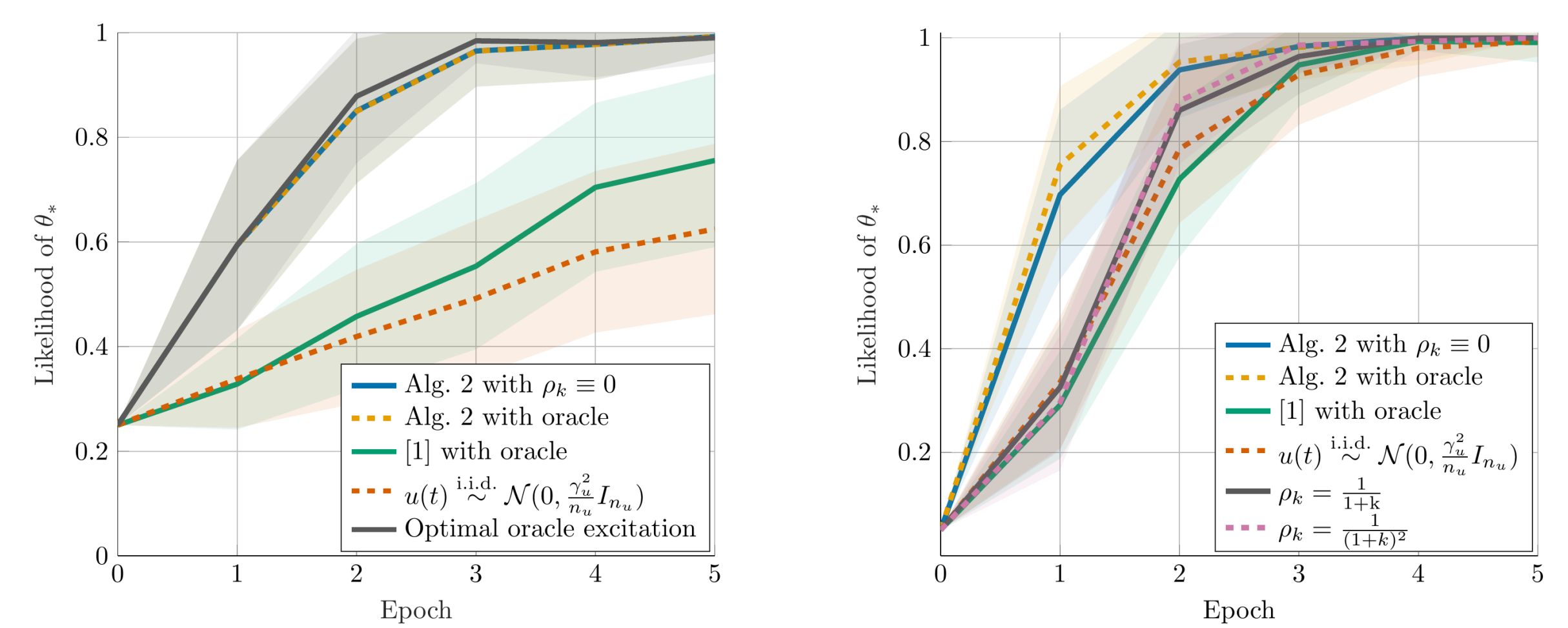
- **Sample complexity upper bound:** Algorithm 1 yields an estimate $\hat{\theta}$ satisfying $\mathbb{P}[\hat{\theta} \neq \theta_*] \leq \delta$ and terminates no later than when k satisfies

$$\sum_{j=1}^k c_{u_j}(\tau) \gamma_u^2 + \text{Excitation due to noise} \geq c \log \left(\frac{N}{\delta} \right).$$

Numerical Experiments

- Comparison of two problem setups with $\Sigma_w = I$, $\gamma_u = 1$, $n_x = 3$, $n_u = 2$, $\eta = 0.01$

- (a) Structured hypothesis class: Experiment design yields large benefits (W_i ill-conditioned)
- (b) Randomly generated hypothesis class: Isotropic random excitation is closer to optimal



(a) Structured set \mathcal{S}

(b) Random set \mathcal{S}

- Proposed algorithm achieves close to optimal performance without oracle knowledge
- Approach in [1] neglects finite hypothesis class \mathcal{S} , yielding significantly suboptimal performance
- CE ($\rho_k = 0$) empirically performs best
- Gap between isotropic Gaussian and optimal input increases when W_i is ill-conditioned

[1] Wagenmaker, A., Jamieson, K. (2020). Active Learning for Identification of Linear Dynamical Systems. *Proceedings of Thirty Third Conference on Learning Theory, 2020*

Conclusion & Outlook

Summary

- Benefit of experiment design is problem-dependent and is captured in matrices W_i
- Sequential identification algorithm that allows for modular analysis for any PE input
- Sequential active learning algorithm with

- sample complexity guarantees
- empirically close to oracle performance using CE

Outlook

- Analysis of computational complexity
- Extensions to task-optimal active learning
- Connecting to online settings, e.g., RL

