

Near-optimal rate of consistency for linear prediction with missing values

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ICML 2022

Background

o Growing mass of data => **NA** (not attribut)/missing values

o Different sources:

1. Bugs
2. Cost

\$1	\$10	\$100	\$0
		NA	
	NA	NA	

3. Multiplication of sources (i.e. merge)

			NA	NA			
					NA	NA	NA
NA	NA	NA			NA	NA	NA

4. Sensitive data

Age	Job	Incomes
		NA
		NA
NA		NA
		NA

Background

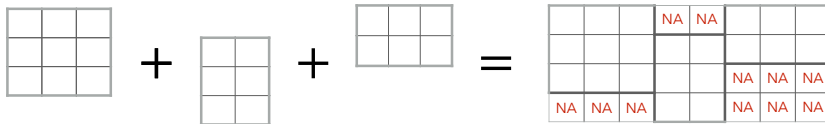
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○ Strategy 1: **complete** the dataset before the ML process (e.g. by collaborative filtering)

○ Strategy 2: **adapt** statistical analysis to handle missing values (e.g. EM algorithm to perform regression with NA)

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 +

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 =

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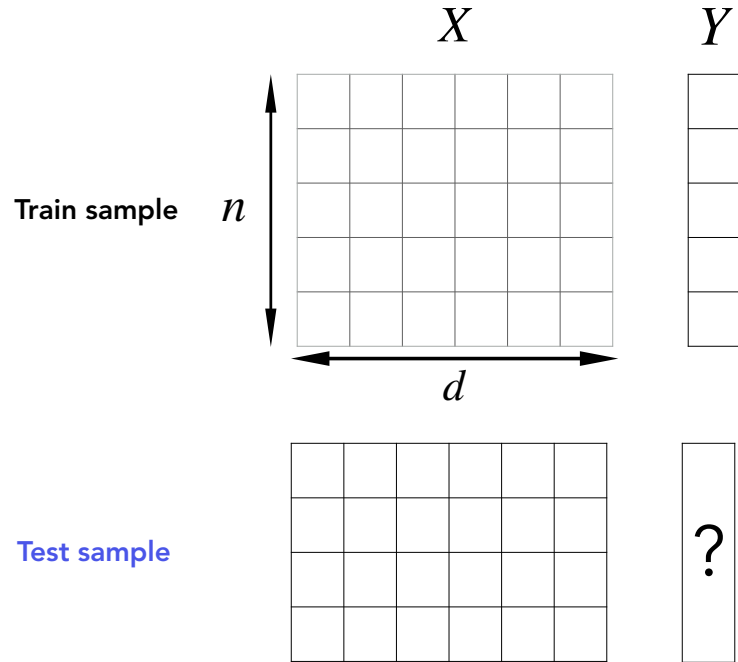
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○ Strategy 1: **complete** the dataset before the ML process (e.g. by collaborative filtering)

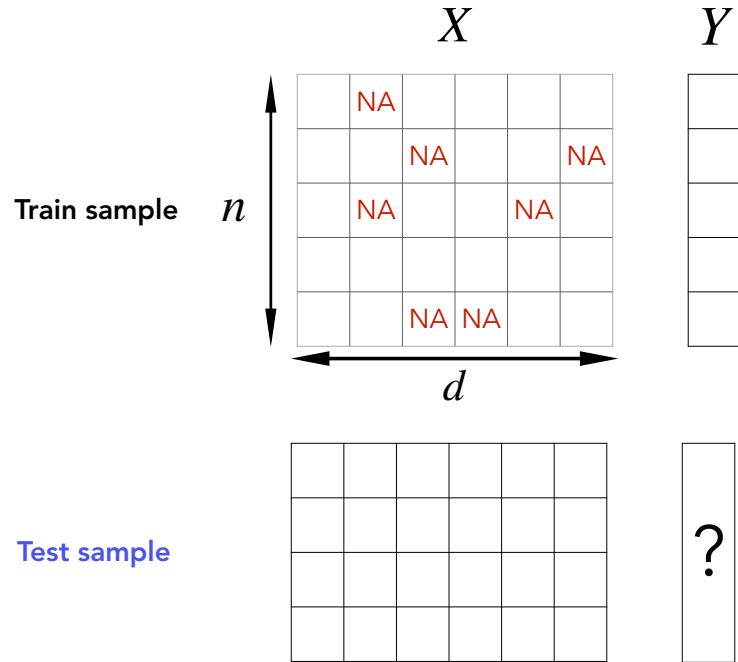
○ Strategy 2: **adapt** statistical analysis to handle missing values (e.g. EM algorithm to perform regression with NA)

What about **supervised learning**?
i.e. prediction with NAs

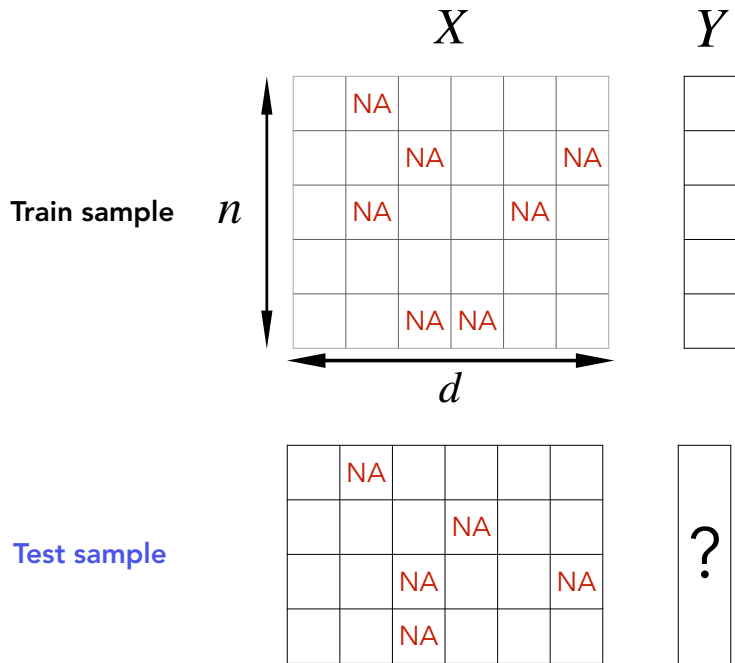
Supervised learning with missing values (NA)



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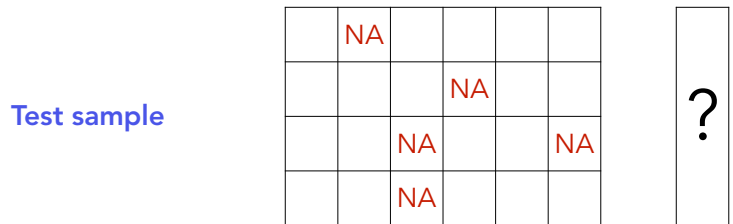
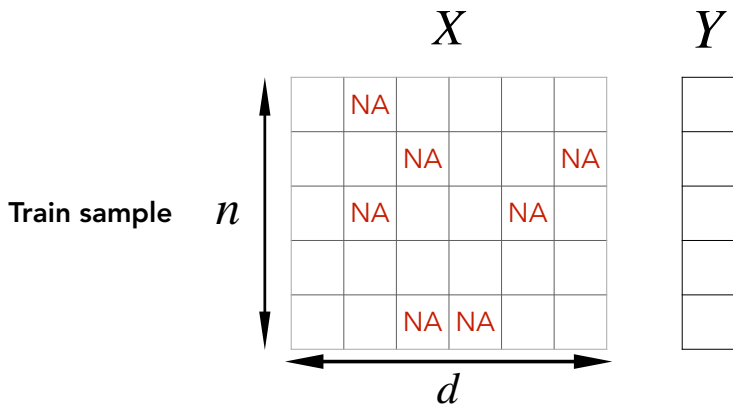


$$y = \beta^T x + \epsilon$$

$$\|\hat{\beta} - \beta\|_2^2$$

$$(\hat{\beta}^T x - \beta^T x)^2$$

Supervised learning with missing values (NA)



Missing pattern: $M_i \in \{0,1\}^d$

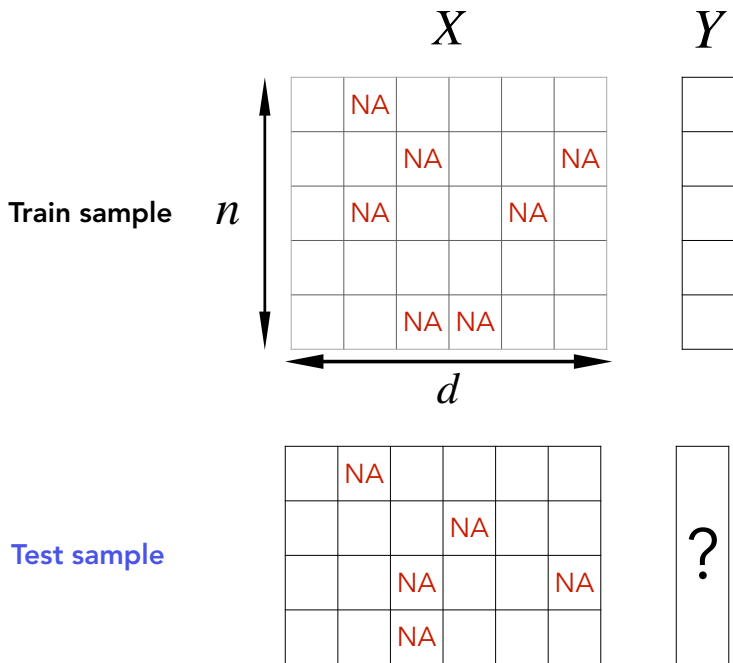
$$X_i = \begin{array}{|c|c|c|c|c|c|} \hline \text{NA} & 1 & -5 & \text{NA} & 0 & 2 \\ \hline \end{array}$$

$$M_i = (1, 0, 0, 1, 0, 0)$$

Input: $Z = (X_{\text{obs}}, M)$

Output: $Y \in \mathbb{R}$

Supervised learning with missing values (NA)



Missing pattern: $M_i \in \{0,1\}^d$

$$X_i = \begin{bmatrix} \text{NA} & 1 & -5 & \text{NA} & 0 & 2 \end{bmatrix}$$

$$M_i = (1, 0, 0, 1, 0, 0)$$

Input: $Z = (X_{\text{obs}}, M)$

Output: $Y \in \mathbb{R}$

Goal: Predict on **test sample** minimizing

$$R(f) = \mathbb{E}_{Z,Y} \left[(Y - f(Z))^2 \right]$$

Zoo of assumptions on NA

○ First point of view : $P(X, M) = P(M|X)P(X)$

- 1) Assumption on $P(X)$:
Example: X Gaussian Vector
- 2) Assumption on $P(M|X)$:

Bugs

Medical protocol

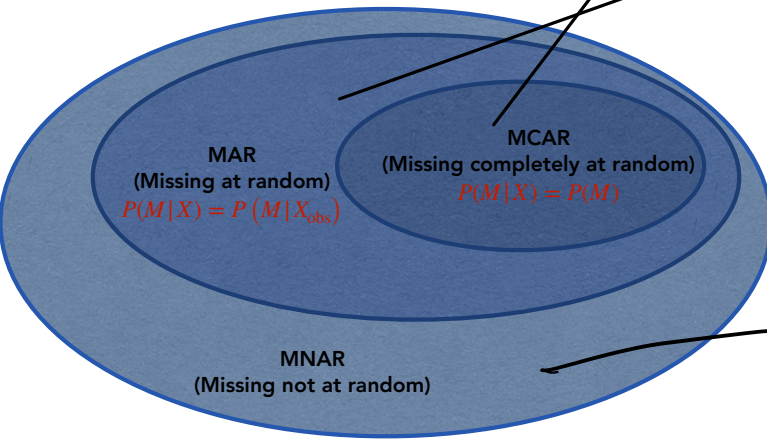
check up

Good

Bad

NA

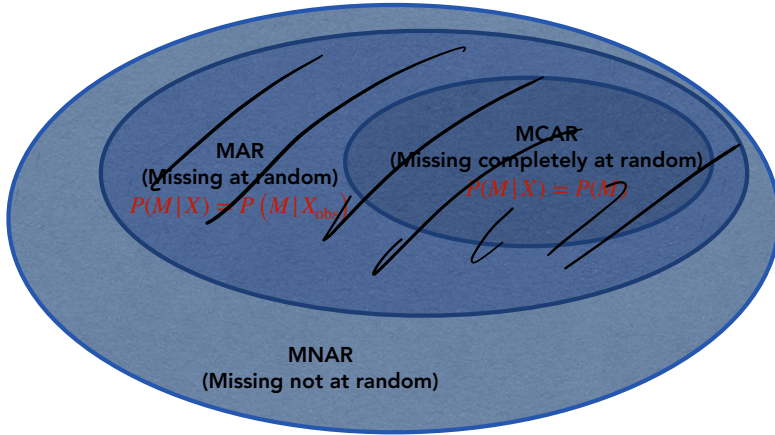
High value censorship



Zoo of assumptions on NA

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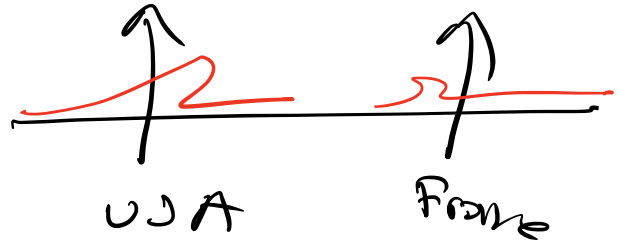
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○ Second point of view : $P(X, M) = P(M)P(X|M)$

- 1) Assumption on $P(M)$:
Example: $P(M = m) = p_m$
- 2) Assumption on $P(X|M)$:

GPM (Gaussian pattern mixture model):
 $X | (M = m)$ Gaussian Vector



Pattern-by-Pattern regression

- **Assumption:** linear model for complete inputs

$$y_i = \beta^\top X_i + \epsilon_i$$



! \ With NA, the Bayes predictor does *not* necessarily remain linear

- **Bayes predictor** (better prediction) decomposition

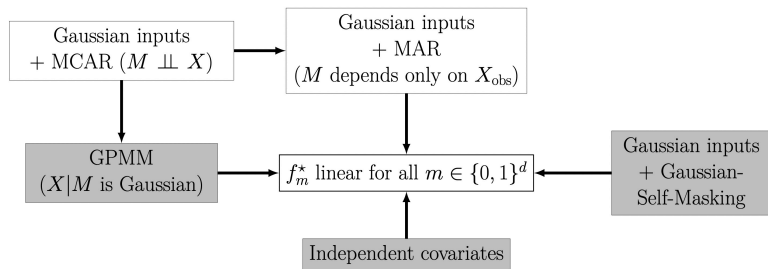
$$f^*(Z) = \sum_{m \in \{0,1\}^d} f_m^*(X_{obs(m)}) \mathbf{1}_{M=m}$$



Local **Bayes prediction** for the missing pattern ($M = m$)

Proposition: (Le Morvan et al. 2020)

Under **linear model** and several **missing data scenarios** (including MNAR), f_m^* are **linear**



Assumption to obtain linearity


Pattern-by-Pattern regression

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- **Bayes predictor** decomposition

$$f^\star(Z) = \sum_{m \in \{0,1\}^d} f_m^\star(X_{obs(m)}) \mathbf{1}_{M=m}$$


Local **Bayes prediction** for the missing pattern ($M = m$)

- **Pattern-by-pattern** predictor

$$\hat{f}(Z) = \sum_{m \in \{0,1\}^d} \hat{f}_m(X_{obs(m)}) \mathbf{1}_{M=m}$$


Local **Least-Square** regression on
 $\{(X_{i,obs}, Y_i), M_i = m\}$

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Theorem 1:

Under Lipschitz and sub-Gaussian assumptions,

$$\mathcal{E}(\hat{f}) := \mathbb{E} \left[\left(f^\star(Z) - \hat{f}(Z) \right)^2 \right] \leq A \log(n) 2^d \frac{d}{n} + \text{Approx}$$

- Optimal for equiprobable missing patterns $\left(p_m = \frac{1}{2^d} \right)$
- Tight for the worst case of pattern-by-pattern predictors
- Sub-optimal for other distributions?

Thresholded Pattern-by-Pattern regression

- Adaptivity to the missing pattern distribution to overcome the curse of dimensionality
- Overfitting reduction

via **Thresholded P-by-P** predictor:

$$\hat{f}(Z) = \sum_{m \in \{0,1\}^d} \hat{f}_m(X_{obs(m)}) \mathbf{1}_{M=m} \mathbf{1}_{\hat{p}_m > \frac{d}{n}}$$

$\hat{p}_m = \text{frequency of pattern } m$

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Local **Least-Square** regression on $\{(X_{i,obs}, Y_i), M_i = m\}$

- Definition: missing pattern **complexity**

$$\mathfrak{C}_p \left(\frac{d}{n} \right) = \sum_{m \in \{0,1\}^d} p_m \wedge \frac{d}{n}$$

Theorem 2: (Main result)

Under Lipschitz and Sub-Gaussian assumptions

$$\mathcal{E}(\hat{f}) \leq A \log(n) \mathfrak{C}_p \left(\frac{d}{n} \right) + \text{Approx}$$

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$$\mathcal{E}(\hat{f}) \leq A \log(n) \mathfrak{C}_p \left(\frac{d}{n} \right) + \text{Approx}$$

Examples:

1. **Uniform** distribution: $\mathfrak{C}_p \left(\frac{d}{n} \right) = 2^d \frac{d}{n}$
2. **Bernoulli** distribution: $M_j \sim \mathcal{B}(\epsilon)$ and $\epsilon \leq \frac{d}{n}$
$$\mathfrak{C}_p \left(\frac{d}{n} \right) \leq \frac{d^2}{n}$$

The thresholded P-by-P predictor is near-optimal

- Minimax risk

Worst case on a class of problem \mathcal{P}_p

↓

$$\mathcal{E}_{\min}(p) = \inf_{\tilde{f}} \sup_{\mathbb{P} \in \mathcal{P}_p} \mathbb{E}_{\mathbb{P}} \left[(\tilde{f}(Z) - f^*(Z))^2 \right]$$

↑

Best algorithm

where \mathcal{P}_p represents a class of data distributions

- for which the missing pattern distribution is p
- under Lipschitz and Sub-Gaussian assumptions

The thresholded P-by-P predictor is near-optimal

○ Minimax risk

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Best algorithm

where \mathcal{P}_p represents a class of data distributions
○ for which the missing pattern distribution is p
○ under Lipschitz and Sub-Gaussian assumptions

Theorem 3:

$$\sigma^2 \mathfrak{G}_p \left(\frac{1}{n} \right) \lesssim \mathcal{E}_{\min}(p) \leq \underbrace{\mathcal{E}(\hat{f})}_{\text{Theorem 2}} \leq A \log(n) \mathfrak{G}_p \left(\frac{d}{n} \right)$$

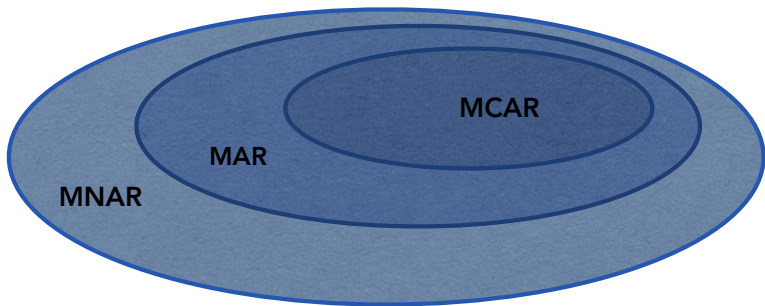
○ Lower bound still holds when \mathcal{P}_p includes **MAR** missing values

Examples

1. **Uniform** distribution: $\mathfrak{G}_p \left(\frac{1}{n} \right) = \frac{2^d}{n}$, $\mathfrak{G}_p \left(\frac{d}{n} \right) = 2^d \frac{d}{n}$
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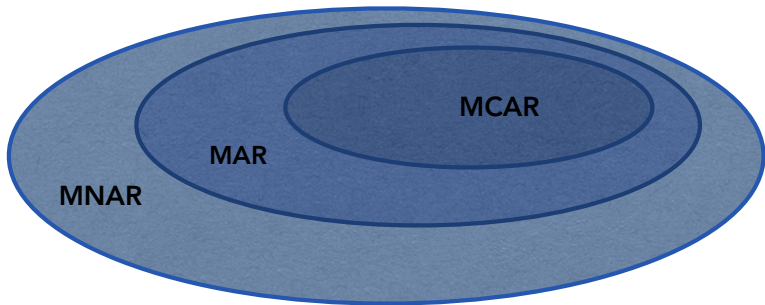
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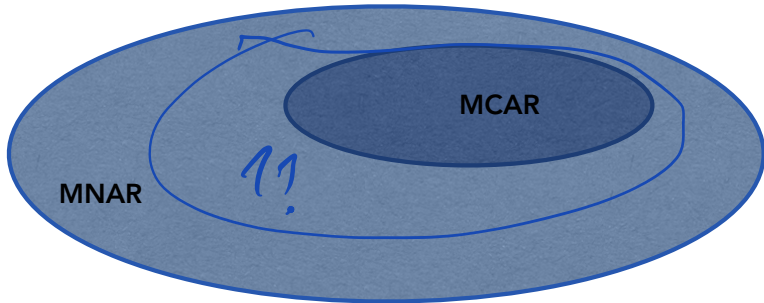
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○ Inference POV



○ Supervised learning POV



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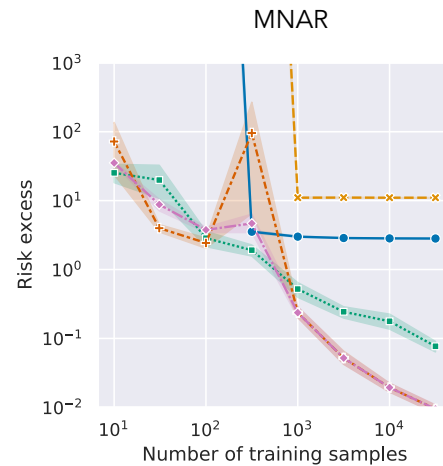
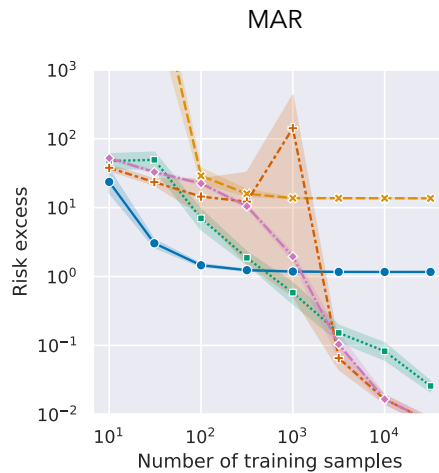
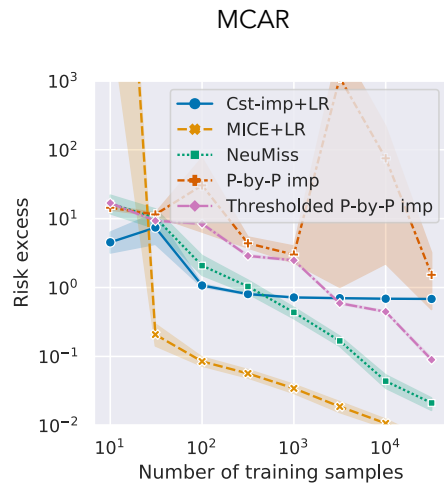
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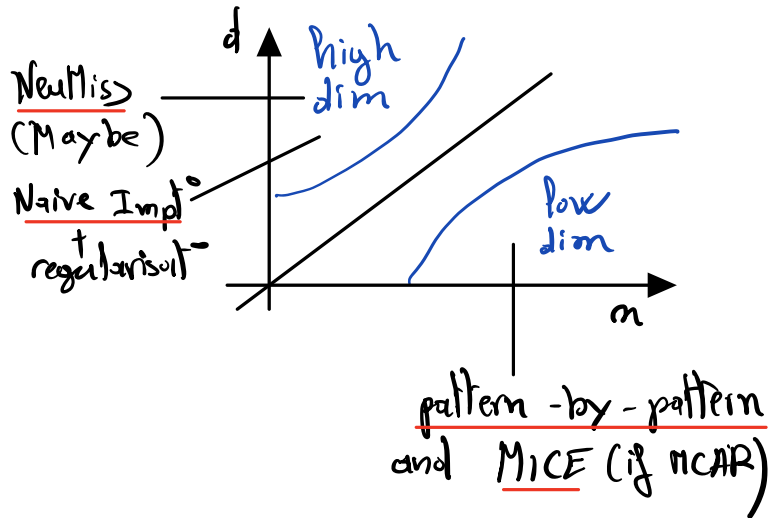
Numerical experiments



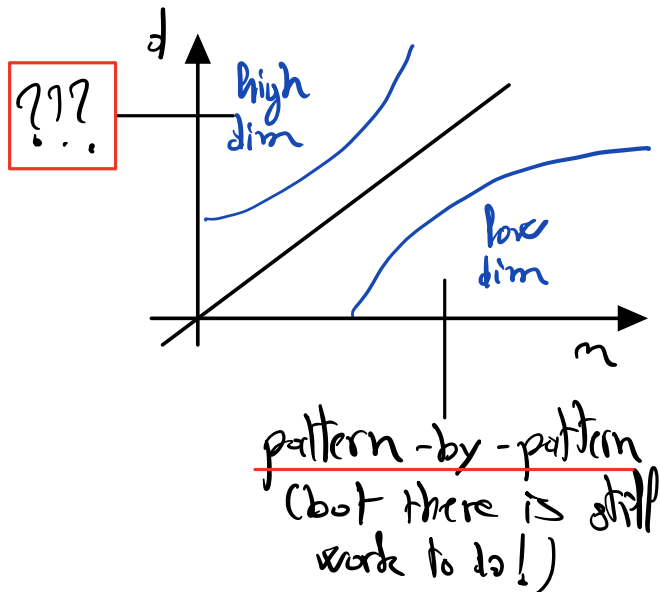
Excess risk w.r.t. n with $d = 8$

What's work?

In practice



In theory



Conclusion

Theoretical contributions

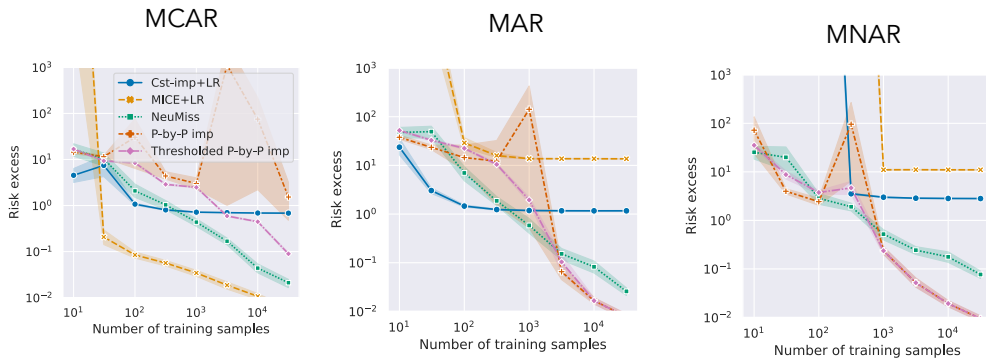
- New **thresholded** predictor
- **Adaptative** upper bound
- Near **optimal**

$$\hat{f}(Z) = \sum_{m \in \{0,1\}^d} \hat{f}_m(X_{obs(m)}) \mathbf{1}_{M=m} \mathbf{1}_{\hat{p}_m > \frac{d}{n}}$$

$$\sigma^2 \mathfrak{C}_p \left(\frac{1}{n} \right) \lesssim \mathcal{E}_{\min}(\mathcal{P}) \leq A \log(n) \mathfrak{C}_p \left(\frac{d}{n} \right)$$

Numerical experiments

- Thresholded P-by-P predictor:
 - reduced variance
 - **consistent** regardless of the missing scenario



Excess risk w.r.t. n with $d = 8$

Near-optimal rate of consistency for linear prediction with missing values

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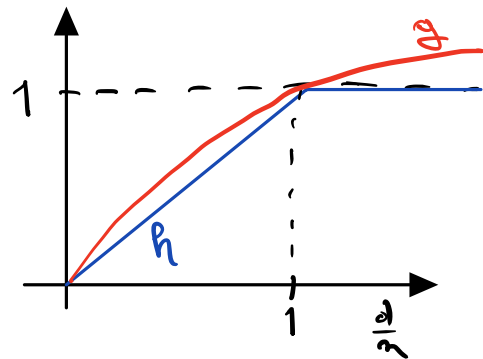


Special thanks to the Paris City Council for the financial support

Some theoretical details:

Why C_p is adaptive?

$$\begin{aligned} C_p\left(\frac{d}{n}\right) &= \sum_{m \in \{0,1\}^d} P_m \wedge \frac{d}{n} \\ &= \sum_m P_m \underbrace{\left(1 \wedge \frac{d}{n P_m}\right)}_{h\left(\frac{d}{n P_m}\right)} \\ &\leq \sum_m P_m g\left(\frac{d}{n P_m}\right) \end{aligned}$$



Two examples:

1) $g = \text{id} \implies C_p\left(\frac{d}{n}\right) \leq \sum_m P_m \frac{d}{n P_m} = 2 \frac{d}{n}$

2) $\left. \begin{array}{l} g(x) = x^\alpha \\ \alpha \in]0,1[\end{array} \right\} \implies C_p\left(\frac{d}{n}\right) \leq \sum_m P_m^\alpha \left(\frac{d}{n}\right)^\alpha = \left(\frac{d}{n}\right)^\alpha e^{\alpha H_{1-\alpha}(P)}$

where $H_{1-\alpha}(P) = \text{Renyi Entropy}$

