Efficient Inference on High-Dimensional Linear Models With Missing Outcomes

Yikun Zhang

Joint Work with Alexander Giessing and Yen-Chi Chen

Department of Statistics, University of Washington

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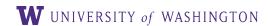




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Introduction



Problem of Interest

Consider a random sample $\{(Y_i, R_i, X_i)\}_{i=1}^n$ drawn from the joint distribution of (Y, R, X), where

- $Y \in \mathbb{R}$ is the outcome variable that could potentially be missing;
- $R \in \{0,1\}$ is the indicator of Y being observed;
- $X \in \mathbb{R}^d$ is the high-dimensional covariate vector with $d \gg n$.

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► Central Question of Interest:

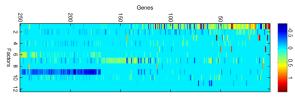
How can we conduct statistically and computationally efficient inference on $m_0(x) = E(Y|X=x)$ despite missing outcomes?



1 The covariates are easier to obtain within some population.



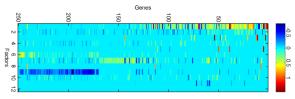
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- Micro-array gene expression data in biology (Carvalho et al., 2008).
- Home-price data with cross-sectional effects (Fan et al., 2011).



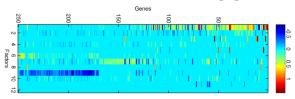
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- Home-price data with cross-sectional effects (Fan et al., 2011).
- Incorporating as many covariates as possible can control for potential confounders in causal inference (Wyss et al., 2022).
- Generating high-dimensional covariates with interaction terms or spline features enables the simple parametric (e.g., linear) model to capture complex patterns (Belloni et al., 2019).



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- Participants may drop out from the study in clinical trials (Higgins et al., 2008).
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- ▶ More Concrete Example: Some (estimated) stellar masses of the observed galaxies in the Sloan Digital Sky Survey (SDSS-IV) are missing in the Firefly value-added catalog (Comparat et al., 2017).



Motivations: Stellar Mass Inference Problem

The missingness of (estimated) stellar masses is due to

- Limiting usage of the observational run in SDSS-IV for galaxy targets;
- Potential data contamination;
- Misclassification of galaxies as stars.

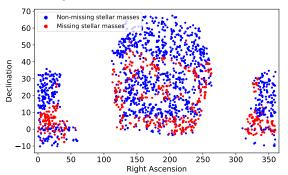


Figure 1: Galaxy distribution at a high redshift slice $0.4 \sim 0.401$.



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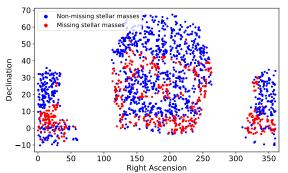


Figure 1: Galaxy distribution at a high redshift slice $0.4 \sim 0.401$.

► **Scientific Question:** How can we conduct valid inference on the (estimated) stellar mass based on the spectroscopic and photometric properties?



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● (*Linearity*) The data $\{(Y_i, R_i, X_i)\}_{i=1}^n \subset \mathbb{R} \times \{0, 1\} \times \mathbb{R}^d$ are i.i.d. observations from a sparse linear model

$$Y = X^T \beta_0 + \epsilon$$
 with $E(\epsilon | X) = 0$ and $E(\epsilon^2 | X) = \sigma_{\epsilon}^2$,

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 - Sparse additive model (Ravikumar et al., 2009);
 - Partially linear model (Müller and van de Geer, 2015);
 - Approximately/weakly sparse linear model (Belloni et al., 2019).
- ② (Missing At Random; MAR) $Y_i \perp \!\!\! \perp R_i | X_i$ for i = 1, ..., n.



Existing Works on High-Dimensional Inference

The existing works focus mainly on the statistical inference on $\beta_0 \in \mathbb{R}^d$.



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 (Fully observed outcomes) Debiased Lasso is applicable (Zhang and Zhang, 2014; van de Geer et al., 2014; Javanmard and Montanari, 2014):

$$\widehat{\beta}^{\text{debias}} = \widehat{\beta}_{\lambda} + \frac{1}{n} \widehat{\Theta} \sum_{i=1}^{n} X_{i} (Y_{i} - X_{i}^{T} \widehat{\beta}_{\lambda}),$$

- $\widehat{\beta}_{\lambda} = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^d} \left[\frac{1}{2n} \sum_{i=1}^n (Y_i X_i^T \beta)^2 + \lambda \left| \left| \beta \right| \right|_1 \right]$ is a Lasso solution with the regularization parameter $\lambda > 0$;
- $\widehat{\Theta} \in \mathbb{R}^{d \times d}$ is an approximation to the matrix inverse $\left(\frac{1}{n} \sum_{i=1}^{n} X_i X_i^T\right)^{-1}$.



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- $\widehat{\Theta} \in \mathbb{R}^{d \times d}$ is an approximation to the matrix inverse $\left(\frac{1}{n} \sum_{i=1}^{n} X_i X_i^T\right)^{-1}$.
- (MAR outcomes) Chakrabortty et al. (2019) proposed an M-estimation framework with a Lasso-type debiased and doubly robust estimator.



Drawback of Existing Works and Our Contributions

- **▶** Drawbacks of the Existing Approaches:
- (*Computational issue*) They require a good approximation to the $d \times d$ debiasing matrix $\widehat{\Theta}$.
- (Loss of statistical efficiency) Sample splitting or cross-fitting is necessary for the M-estimation framework.

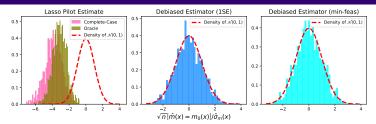


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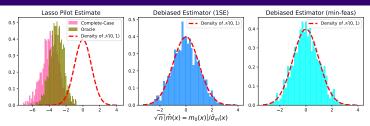
- (*Computational issue*) They require a good approximation to the $d \times d$ debiasing matrix $\widehat{\Theta}$.
- (Loss of statistical efficiency) Sample splitting or cross-fitting is necessary for the M-estimation framework.
- ▶ Our Contributions: Focus on the inference of $m_0(x) = x^T \beta_0$ instead.
- (*Computational efficiency*) Our core debiasing program is convex and only needs to solve for a *n*-dimensional weight vector.
- (*Statistical efficiency*) Our debiased estimator is semi-parametrically efficient among all asymptotically linear estimators.





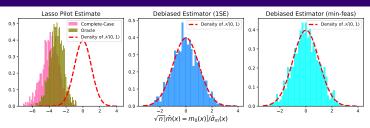
• Introduce our efficient debiasing method for inferring $m_0(x) = x^T \beta_0$.





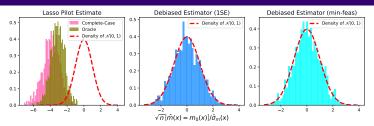
- Introduce our efficient debiasing method for inferring $m_0(x) = x^T \beta_0$.
 - Estimate $\pi(X) = P(R = 1|X)$ via any machine learning methods.
 - Design our debiasing program based on bias-variance trade-offs.
 - Fine-tune the program from its dual so as to debias the Lasso solution.





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 - Fine-tune the program from its dual so as to debias the Lasso solution.
- Discuss the asymptotic normality and semi-parametric efficiency of our final debiased estimator.
- Demonstrate the finite-sample performances via simulations and present an application to the stellar mass inference problem.

Methodology





For any fixed $\lambda > 0$, the Lasso solution (on the complete-case data) is a biased estimator of $\beta_0 \in \mathbb{R}^d$:

$$\widehat{\beta}_{\lambda} = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^d} \left[\frac{1}{2n} \sum_{i=1}^n R_i (Y_i - X_i^T \beta)^2 + \lambda ||\beta||_1 \right].$$

▶ **Question:** How can we correct for the bias in $\widehat{\beta}_{\lambda}$ or $\widehat{m}(x) = x^T \widehat{\beta}_{\lambda}$?



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- Optimality/KKT condition reads

$$\frac{1}{n} \sum_{i=1}^{n} R_{i} X_{i} \left(Y_{i} - X_{i}^{T} \widehat{\beta}_{\lambda} \right) = \lambda \widehat{z} \quad \text{with} \quad \widehat{z} \in \partial \left\| \widehat{\beta}_{\lambda} \right\|_{1} \in \mathbb{R}^{d}.$$
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• Linearity assumption $Y_i = X_i^T \beta_0 + \epsilon_i$ for i = 1, ..., n implies that

$$\frac{1}{n}\sum_{i=1}^{n}R_{i}X_{i}\epsilon_{i}+\widehat{\Sigma}\left(\beta_{0}-\widehat{\beta}_{\lambda}\right)=\lambda\widehat{z} \quad \text{with} \quad \widehat{\Sigma}=\frac{1}{n}\sum_{i=1}^{n}R_{i}X_{i}X_{i}^{T}.$$

• Given an approximation $\widehat{\Theta} \in \mathbb{R}^{d \times d}$ to the gram matrix $\widehat{\Sigma}$, it becomes

$$\widehat{\beta}_{\lambda} - \beta_0 + \widehat{\Theta}\lambda \widehat{z} = \underbrace{\frac{1}{n} \sum_{i=1}^{n} R_i \widehat{\Theta} X_i \epsilon_i}_{\text{Stochastic error } \sim \mathcal{N}_d(0, \widetilde{\Sigma})} + \underbrace{\left(\widehat{\Theta}\widehat{\Sigma} - I_d\right) \left(\beta_0 - \widehat{\beta}_{\lambda}\right)}_{\text{Asymptotically negligible bias}}.$$



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By KKT condition (1), the debiased Lasso estimate is thus given by

$$\begin{split} \widehat{\beta}^{\text{debias}} &= \widehat{\beta}_{\lambda} + \widehat{\Theta} \lambda \widehat{z} \\ &= \widehat{\beta}_{\lambda} + \frac{1}{n} \sum_{i=1}^{n} R_{i} \widehat{\Theta} X_{i} \left(Y_{i} - X_{i}^{T} \widehat{\beta}_{\lambda} \right). \end{split}$$



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• A candidate debiased estimator for $m_0(x) = x^T \beta_0$ is

$$\widehat{m}^{\text{debias}}(x) = x^T \widehat{\beta}^{\text{debias}} = x^T \widehat{\beta}_{\lambda} + \frac{1}{n} x^T \widehat{\Theta} \sum_{i=1}^n R_i X_i \left(Y_i - X_i^T \widehat{\beta}_{\lambda} \right).$$



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▶ **Issue:** Fitting the debiasing matrix $\widehat{\Theta} \in \mathbb{R}^{d \times d}$ is computationally inefficient; see, *e.g.*, the nodewise regression (Meinshausen and Bühlmann, 2006; van de Geer et al., 2014).



Heuristics From Debiased Lasso

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- ▶ **Solution:** Introduce the weight vector $\widehat{\boldsymbol{w}} = (\widehat{w}_1, ..., \widehat{w}_n)^T \in \mathbb{R}^n$ with (Giessing and Wang, 2023)

$$\widehat{w}_i = \begin{cases} \frac{1}{\sqrt{n}} x^T \widehat{\Theta} X_i & R_i = 1, \\ 0 & R_i = 0, \end{cases}$$

for i = 1, ..., n so that our final debiased estimator becomes

$$\widehat{m}^{\text{debias}}(x;\widehat{\boldsymbol{w}}) = x^T \widehat{\beta} + \frac{1}{\sqrt{n}} \sum_{i=1}^n \widehat{w}_i R_i \left(Y_i - X_i^T \widehat{\beta} \right). \tag{2}$$



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▶ Question: How do we estimate the weight vector $\hat{\boldsymbol{w}} = (\hat{w}_1, ..., \hat{w}_n)^T$?

Consider the generic debiased estimator $m^{\text{debias}}(x; w)$ from (2) as:

$$m^{\text{debias}}(x; \boldsymbol{w}) = x^T \beta + \frac{1}{\sqrt{n}} \sum_{i=1}^n w_i R_i \left(Y_i - X_i^T \beta \right). \tag{3}$$



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The conditional mean squared error of $\sqrt{n} m^{\text{debias}}(x; w)$ is given by

$$E\left[\left(\sqrt{n}\,m^{\text{debias}}(x;\boldsymbol{w})-\sqrt{n}\,m_0(x)\right)^2\,\Big|X_1,...,X_n\right]$$

$$= \sigma_{\epsilon}^{2} \sum_{i=1}^{n} w_{i}^{2} \pi(X_{i}) + \left[\left(\frac{1}{\sqrt{n}} \sum_{i=1}^{n} w_{i} \pi(X_{i}) X_{i} - x \right)^{T} \sqrt{n} \left(\beta_{0} - \beta \right) \right]^{2}$$
Main Conditional Variance

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$$\mathbf{x} \mathbf{x}^T$$

+
$$(\beta_0 - \beta)^T \left[\sum_{i=1}^n w_i^2 \pi(X_i) (1 - \pi(X_i)) X_i X_i^T \right] (\beta_0 - \beta),$$

Asymptotically Negligible Conditional Variance

where $\pi(X) = P(R = 1|X)$ is the propensity score under MAR condition.



$$E\left[\left(\sqrt{n}\,m^{\text{debias}}(x;\boldsymbol{w})-\sqrt{n}\,m_0(x)\right)^2\Big|X_1,...,X_n\right]$$

$$\approx \underbrace{\sigma_{\epsilon}^2\sum_{i=1}^nw_i^2\pi(X_i)}_{\text{Main Conditional Variance}} + \underbrace{\left[\left(\frac{1}{\sqrt{n}}\sum_{i=1}^nw_i\pi(X_i)X_i-x\right)^T\sqrt{n}\left(\beta_0-\beta\right)\right]^2}_{\text{Conditional Bias}}.$$

By Hölder's inequality, the "Conditional Bias" is upper bounded by

$$\left[\left| \left| \frac{1}{\sqrt{n}} \sum_{i=1}^{n} w_i \pi(X_i) X_i - x \right| \right| \quad \sqrt{n} \left| \left| \beta_0 - \beta \right| \right|_1 \right]^2.$$



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$$\approx \sigma_{\epsilon}^2 \sum_{i=1}^n w_i^2 \pi(X_i) + \left[\left(\frac{1}{\sqrt{n}} \sum_{i=1}^n w_i \pi(X_i) X_i - x\right)^T \sqrt{n} \left(\beta_0 - \beta\right)\right]^2.$$
Main Conditional Variance

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• We design our core debiasing program as:

$$\min_{\boldsymbol{w} \in \mathbb{R}^n} \sum_{i=1}^n \widehat{\pi}_i w_i^2 \quad \text{subject to} \quad \left\| x - \frac{1}{\sqrt{n}} \sum_{i=1}^n w_i \cdot \widehat{\pi}_i \cdot X_i \right\| \leq \frac{\gamma}{n},$$

where $\gamma > 0$ is a tuning parameter and $\widehat{\pi}_i$ is a consistent estimate of the propensity score $\pi(X_i)$ for i = 1, ..., n.

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$$\widehat{\beta}_{\lambda} = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^d} \left[\frac{1}{2n} \sum_{i=1}^n R_i (Y_i - X_i^T \beta)^2 + \lambda ||\beta||_1 \right].$$



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② Obtain consistent propensity score estimates $\widehat{\pi}_i$, i = 1, ..., n by any machine learning method based on $\{(X_i, R_i)\}_{i=1}^n \subset \mathbb{R}^d \times \{0, 1\}$.



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- Obtain consistent propensity score estimates $\widehat{\pi}_i$, i = 1, ..., n by any machine learning method based on $\{(X_i, R_i)\}_{i=1}^n \subset \mathbb{R}^d \times \{0, 1\}$.
- 3 Solve the debiasing program defined as:

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$$\widehat{\beta}_{\lambda} = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^d} \left[\frac{1}{2n} \sum_{i=1}^n R_i (Y_i - X_i^T \beta)^2 + \lambda ||\beta||_1 \right].$$

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6 Construct the asymptotic $(1 - \tau)$ -level confidence interval for $m_0(x)$ as:

$$\left[\widehat{m}^{\text{debias}}(x;\widehat{\boldsymbol{w}}) \pm \Phi^{-1}\left(1 - \frac{\tau}{2}\right) \cdot \widehat{\sigma}_{\epsilon} \cdot \sqrt{\frac{1}{n} \sum_{i=1}^{n} \widehat{\pi}_{i} \widehat{\boldsymbol{w}}_{i}^{2}}\right] \quad \text{ with } \Phi(\cdot) \text{ being the CDF of } \mathcal{N}(0,1).$$



Theory and Practice of Our Debiasing Program

There are two unanswered questions in our proposed debiasing inference procedure:

• How can we select the tuning parameter $\gamma > 0$ for our debiasing program?

$$\min_{w \in \mathbb{R}^n} \left\{ \sum_{i=1}^n \widehat{\pi}_i w_i^2 : \left| \left| x - \frac{1}{\sqrt{n}} \sum_{i=1}^n w_i \cdot \widehat{\pi}_i \cdot X_i \right| \right|_{\infty} \le \frac{\gamma}{n} \right\}.$$

o Why is the asymptotic $(1 - \tau)$ -level confidence interval for $m_0(x)$ valid?

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► **Answer:** The above two questions can be addressed by the *dual formulation/solution* of our debiasing program!



Dual Formulation of Our Debiasing Program

The primal form of our debiasing program is a quadratic programming problem with a box constraint:

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Proposition (Proposition 1 in Zhang et al. 2023)

The dual form of our debiasing program is given by

$$\min_{\ell \in \mathbb{R}^d} \left\{ \frac{1}{4n} \sum_{i=1}^n \widehat{\pi}_i \left[X_i^T \ell \right]^2 + x^T \ell + \frac{\gamma}{n} \left| \left| \ell \right| \right|_1 \right\}.$$

If the strong duality holds, we further have that

$$\widehat{w}_i = -\frac{1}{2\sqrt{n}} \cdot X_i^T \widehat{\ell} \quad \textit{for} \quad i = 1, ..., n,$$

where $\hat{w} \in \mathbb{R}^n$ and $\hat{\ell} \in \mathbb{R}^d$ are the solutions to the primal and dual debiasing program, respectively.



Practical Implication of Our Dual Debiasing Program

The dual form of our debiasing program is an *unconstrained* quadratic programming problem:

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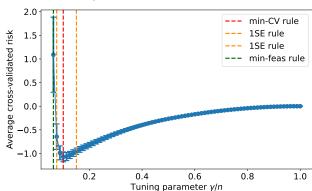


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We can fine-tune $\gamma > 0$ by cross-validation.





- Consider the regression function $m \equiv m(x) \in \mathbb{R}$ as the main parameter to be inferred and $\beta \in \mathbb{R}^d$ as the high-dimensional nuisance parameter.
- Our generic debiased estimator $m^{\text{debias}}(x, w)$ solves the sample-based estimating equation

$$\frac{1}{n}\sum_{i=1}^{n}\Xi_{x}(Y_{i},R_{i},X_{i};m^{\text{debias}},\beta)=m^{\text{debias}}(x;\boldsymbol{w})-x^{T}\beta-\frac{1}{\sqrt{n}}\sum_{i=1}^{n}w_{i}\cdot R_{i}\left(Y_{i}-X_{i}^{T}\beta\right)=0.$$



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• The Neyman near-orthogonalization condition (Chernozhukov et al., 2018) given $X = (X_1, ..., X_n)^T \in \mathbb{R}^{n \times d}$ at $(m_0, \beta_0) = (x^T \beta_0, \beta_0)$ requires

$$E\left[\frac{1}{n}\sum_{i=1}^{n}\Xi_{x}(Y_{i},R_{i},X_{i};m_{0},\beta_{0})\middle|\mathbf{X}\right]=0,$$

$$\sup_{\beta\in\mathcal{T}_{n}}\left|\left\{\frac{\partial}{\partial\beta}E\left[\frac{1}{n}\sum_{i=1}^{n}\Xi_{x}(Y_{i},R_{i},X_{i};m,\beta)\middle|\mathbf{X}\right]\middle|_{(m_{0},\beta_{0})}\right\}^{T}(\beta-\beta_{0})\right|\leq\frac{\delta_{n}}{\sqrt{n}},$$
(4)

where \mathcal{T}_n is a properly shrinking neighborhood of β_0 and $\delta_n = o(1)$.



Both conditions in (4) hold true, because for any $\beta \in \mathcal{T}_n$ and some convex set \mathcal{B} containing β_0 , we have that

$$\left| \left\{ \frac{1}{n} \sum_{i=1}^{n} \frac{\partial}{\partial \beta} \mathbb{E} \left[\Xi_{x}(Y_{i}, R_{i}, X_{i}; m, \beta) | X \right] \Big|_{(m_{0}, \beta_{0})} \right\}^{T} (\beta - \beta_{0}) \right|$$

$$= \left| \left[x - \frac{1}{\sqrt{n}} \sum_{i=1}^{n} w_{i} \cdot \pi(X_{i}) X_{i} \right]^{T} (\beta_{0} - \beta) \right|$$

$$" \leq " \left| \left| x - \frac{1}{\sqrt{n}} \sum_{i=1}^{n} w_{i} \cdot \widehat{\pi}_{i} \cdot X_{i} \right| \right|_{\infty} ||\beta - \beta_{0}||_{1} \quad \text{by H\"older's inequality}$$

$$\leq \frac{\gamma}{n} ||\beta - \beta_{0}||_{1} \quad \text{by the box constraint in our debiasing program}$$

$$\leq \frac{\delta_{n}}{\sqrt{n}} \quad \text{by setting } \mathcal{T}_{n} = \left\{ \beta \in \mathcal{B} \subset \mathbb{R}^{d} : ||\beta - \beta_{0}||_{1} \leq \frac{\sqrt{n}\delta_{n}}{\gamma} \right\}.$$



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- Our debiasing program optimizes the (estimated) variance among all the estimators satisfying Neyman near-orthogonalization (4).
- (4) also allows our debiasing program to *de-correlate* the Lasso pilot regression from propensity score estimation and weight optimization.

Asymptotic Theory





Theoretical Implication of Our Dual Debiasing Program

► Goal: Establish the asymptotic normality of our debiased estimator

$$\widehat{m}^{\text{debias}}(x; \widehat{\boldsymbol{w}}) = x^T \widehat{\beta} + \frac{1}{\sqrt{n}} \sum_{i=1}^n \widehat{w}_i R_i \left(Y_i - X_i^T \widehat{\beta} \right).$$



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▶ Naive Attempt: Linearity assumption $Y_i = X_i^T \beta_0 + \epsilon_i$ for i = 1, ..., n implies that

$$\sqrt{n} \left[\widehat{m}^{\text{debias}}(x; \widehat{\boldsymbol{w}}) - m_0(x) \right] = \underbrace{\sum_{i=1}^n \widehat{w}_i R_i \epsilon_i}_{\text{Not an i.i.d. sum!}} + \left[x - \frac{1}{\sqrt{n}} \sum_{i=1}^n \widehat{w}_i R_i X_i \right]^T \sqrt{n} \left(\widehat{\beta} - \beta_0 \right),$$



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▶ **Solution:** With the dual relation $\widehat{w}_i = -\frac{1}{2\sqrt{n}} \cdot X_i^T \widehat{\ell}, i = 1, ..., n$, we obtain

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= -\frac{1}{2\sqrt{n}} \sum_{i=1}^n R_i \epsilon_i X_i^T \ell_0(x) + \underbrace{\text{"Bias terms"}}_{o_P(1)}.$$
i.i.d. sum!



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whose exact solution is $\ell_0(x) = -2 \left[\mathbb{E} \left(RXX^T \right) \right]^{-1} x$. We assume that the r_ℓ -approximation $\widetilde{\ell}(x)$ to $\ell_0(x)$ is sparse with $r_\ell \in \left[0, \frac{1}{2}\right]$, *i.e.*,

$$s_{\ell}(x) = \left| \left| \widetilde{\ell}(x) \right| \right|_{0} \ll \min\{n,d\} \text{ with } \widetilde{\ell}(x) = \operatorname*{arg\,min}_{u \in \mathbb{R}^{d}} \left\{ \left| \left| u \right| \right|_{0} : \left| \left| u - \ell_{0}(x) \right| \right|_{2} \le r_{\ell} \left| \left| \ell_{0}(x) \right| \right|_{2} \right\}.$$



Consistency and Asymptotic Normality

- ① Consistency of Lasso pilot estimate: If $\lambda \asymp \sigma_{\epsilon} \sqrt{\frac{\log d}{n}}$ with $\log d = o(n)$, then $\left| \left| \widehat{\beta} \beta_0 \right| \right|_2 = O_P \left(\frac{1}{\kappa_p^2} \sqrt{\frac{s_\beta \log d}{n}} \right)$.
- **2** Consistency of the solution to the dual debiasing program: If r_{ℓ} shrinks to 0 in a certain rate and $\frac{\gamma}{n} \approx \frac{||x||_2}{\kappa_R} \sqrt{\frac{\log d}{n}} + \frac{||x||_2}{\kappa_R^2} \cdot r_{\pi}$, then

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Note: Under the same choice of $\gamma > 0$, the strong duality holds.

Theorem (Theorem 7 in Zhang et al. 2023)

If
$$\frac{(1+\kappa_R^2)s_{\max}\log(nd)}{\kappa_R^4} = o\left(\sqrt{n}\right)$$
, $\frac{(1+\kappa_R^4)\sqrt{s_{\max}\log(nd)}}{\kappa_R^6}$ $(r_\ell+r_\pi) = o(1)$, and $||x||_2 = O(1)$ with $s_{\max} = \{s_\beta, s_\ell(x)\}$, then
$$\frac{\sqrt{n}\left[\widehat{m}^{\text{debias}}(x; \widehat{w}) - m_0(x)\right]}{\sigma_m(x)} \stackrel{d}{\to} \mathcal{N}\left(0, 1\right).$$



Remarks on Our Theoretical Results

Our growth requirement $s_{\text{max}} = o\left(\frac{\sqrt{n}}{\log d}\right)$ on the sparsity level is a standard and *essentially necessary* condition for asymptotic normality; see Section 8.6 of Jankova and van de Geer (2018).



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Proposition (Proposition 8 in Zhang et al. 2023)

If
$$\frac{(1+\kappa_R^3)}{\kappa_R^5}\sqrt{\frac{s_\ell(x)\log(nd)}{n}} = o(1)$$
, $\frac{(1+\kappa_R^4)}{\kappa_R^6}\left[r_\ell + r_\pi\sqrt{s_\ell(x)}\right] = o(1)$, and $||x||_2 = O(1)$, then

$$\left| \sum_{i=1}^{n} \widehat{\pi}_{i} \widehat{w}_{i}^{2} - x^{T} \left[\mathbb{E} \left(RXX^{T} \right) \right]^{-1} x \right| = o_{P}(1).$$



Overfitting the Propensity Scores

Our theoretical results also provide insightful answers to the following two questions:

- Why don't we need sample splitting or cross fitting?
- Why can we estimate the propensity score by any machine learning methods without worrying about the overfitting issue?



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$$P\left(\max_{1\leq i\leq n}|\widehat{\pi}_i-\pi_i|>r_\pi\right)<\delta\quad\text{ with }\quad \pi_i=\pi(X_i), i=1,...,n.$$

• In other words, our debiased estimator performs even better when we overfit the propensity scores $\pi(X_i) = P(R_i = 1|X_i), i = 1,...,n$.



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- Why can we estimate the propensity score by any machine learning methods without worrying about the overfitting issue?
- ► **Answer:** Our asymptotic normality result depends on the *in-sample* estimation error r_{π} of the propensity score; recall that

$$P\left(\max_{1\leq i\leq n}|\widehat{\pi}_i-\pi_i|>r_\pi\right)<\delta\quad\text{ with }\quad \pi_i=\pi(X_i), i=1,...,n.$$

- In other words, our debiased estimator performs even better when we overfit the propensity scores $\pi(X_i) = P(R_i = 1|X_i), i = 1,...,n$.
- This coincides with "benign overfitting" in linear regression or neural networks (Bartlett et al., 2020; Li et al., 2021; Cao et al., 2022).

Comparative Simulations





Experimental Setups and Evaluation Metrics

We compare our debiasing method with L_1 -penalized logistic regression for the propensity score estimation with several existing methods:

- "DL-Jav": The debiased Lasso by Javanmard and Montanari (2014).
- "DL-vdG": The debiased Lasso by van de Geer et al. (2014).
- "Refit": Run the regular least-square regression on the support set of the Lasso pilot estimate (Belloni and Chernozhukov, 2013).



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These methods to be compared are implemented on

- Complete-case (CC) data $\{(X_i, Y_i, R_i = 1)\}_{i=1}^n$;
- Inverse probability weighted (IPW) data $\left\{ \left(\frac{X_i}{\sqrt{\widehat{\pi}_i}}, \frac{Y_i}{\sqrt{\widehat{\pi}_i}}, R_i = 1 \right) \right\}_{i=1}^n$;
- Oracle fully observed data (X_i, Y_i) for i = 1, ..., n.



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Evaluation metrics on 1000 Monte Carlo experiments include

- Average absolute bias $|\widehat{m}^{\text{debias}}(x) m_0(x)|$;
- Average coverage of the yielded 95% confidence intervals;
- Average length of the yielded 95% confidence intervals.

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Simulation Results Under Gaussian Noises (I)

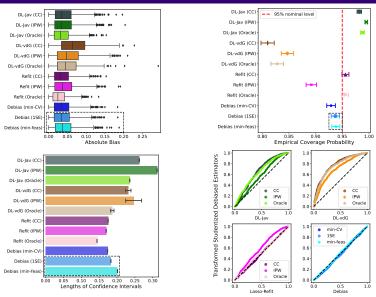


Figure 2: Sparse β_0^{sp} and sparse $x^{(2)}$ with $X_i \sim \mathcal{N}_d(\mathbf{0}, \Sigma^{cs}), i = 1, ..., n$.

Yikun Zhang



Simulation Results Under Gaussian Noises (II)

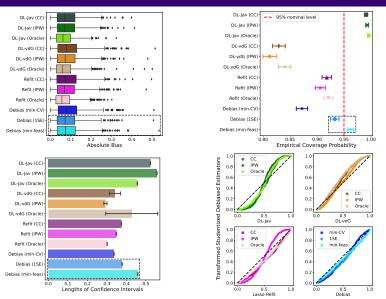


Figure 3: Pseudo-dense β_0^{pd} and sparse $x^{(2)}$ with $X_i \sim \mathcal{N}_d(\mathbf{0}, \Sigma^{ar}), i = 1, ..., n$.



Simulation Results Under Laplace $(0, 1/\sqrt{2})$ Noises

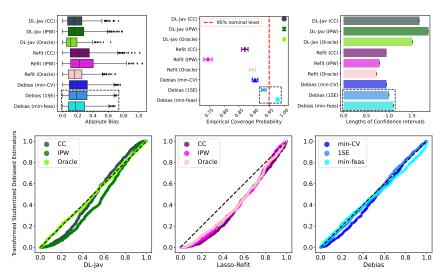


Figure 4: Dense β_0^{de} and sparse $x^{(4)}$ with $X_i \sim \mathcal{N}_d(\mathbf{0}, \Sigma^{cs}), i = 1, ..., n$.



Simulation Results Under t₂-Distributed Noises

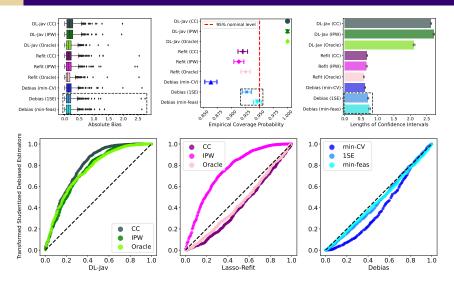


Figure 5: Pseudo-dense β_0^{pd} and dense $x^{(4)}$ with $X_i \sim \mathcal{N}_d(\mathbf{0}, \Sigma^{ar}), i = 1, ..., n$. Note that the mean-zero t_2 distribution has *infinite* variance.



Proposed Method With Nonparametric Propensity Scores

• True propensity score model: $P(R_i = 1|X_i) = \Phi\left(-4 + \sum_{k=1}^K Z_{ik}\right)$, where $(Z_{i1}, ..., Z_{iK})$ contains all polynomial combinations of the first eight components $X_{i1}, ..., X_{i8}$ of $X_i \in \mathbb{R}^{1000}$ with degrees ≤ 2 .



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- 2) Estimate the propensity scores $\pi(X_i)$, i = 1, ..., n by the following nonlinear/nonparametric machine learning methods:
 - Gaussian Naive Bayes ("NB").
 - **Random Forest ("RF"):** 100 trees, bootstrapping samples, and the Gini impurity.
 - **Support Vector Machine ("SVM"):** Gaussian radial basis function.
 - **Neural Network ("NN"):** Two hidden layers of size 80×50 and ReLU $h(x) = \max\{x, 0\}$ as the activation function.



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- Include an extra evaluation metric as the average mean absolute error ("Avg-MAE") for the estimated propensity scores.



Simulation Results With Nonparametric Propensity Scores

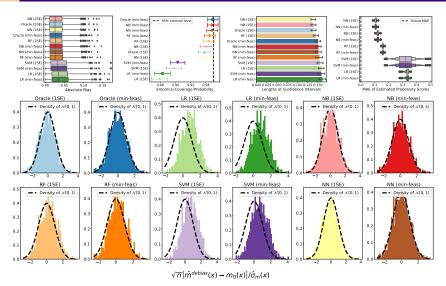


Figure 6: Sparse β_0^{sp} and (weakly) dense $x^{(4)}$.

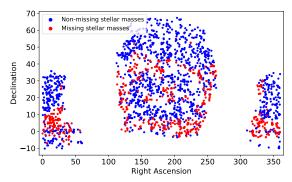
Real-World Applications





Background on Stellar Mass Inference

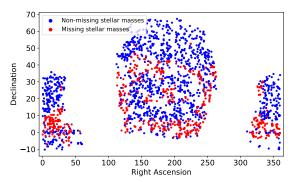
Recall that some estimated stellar masses of the observed galaxies in SDSS-IV are missing in the most recent Firefly value-added catalog.





Background on Stellar Mass Inference

Recall that some estimated stellar masses of the observed galaxies in SDSS-IV are missing in the most recent Firefly value-added catalog.



▶ Scientific Questions:

- How can we conduct valid inference on the (estimated) stellar mass based on the spectroscopic and photometric properties?
- ② Is it statistically significant that the stellar mass of a galaxy is negatively correlated with its distance to the nearby cosmic filament structures?



Ocnsider all the observed galaxies by SDSS-IV within a thin redshift slice $0.4 \sim 0.4005$, among which 30.2% of their stellar masses are missing in the Firefly value-added catalog.



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- Solution Apply feature transformation, remove highly linearly correlated covariates, and generate univariate B-spline base covariates of polynomial order 3 with 40 knots.



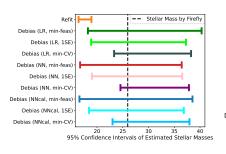
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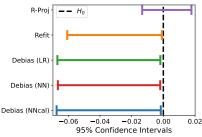


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- © Control for the confounding effects by including the distances from galaxies to candidate galaxy clusters.
- ▶ Final Dataset: n = 1185 and d = 1409.



Results on Stellar Mass Inference





- *Left Panel*: 95% confidence intervals by different debiasing methods for the estimated stellar mass of a new galaxy.
- Right Panel: 95% confidence intervals by different debiasing methods for the estimated regression coefficient associated with the distance to nearby cosmic filaments.

Conclusions and Future Works



W Conclusions

We develop an efficient debiasing method for conducting valid inference on high-dimensional linear models with MAR outcomes.

We develop an efficient debiasing method for conducting valid inference on high-dimensional linear models with MAR outcomes.

- Its computational and statistical efficiencies follow from the dual formulation.
- Sample splitting and cross fitting are not required, and the nuisance propensity score can be estimated by any machine learning method.
- We provide interpretations to our debiasing method from the viewpoints of bias-variance trade-off and Neyman near-orthogonalization.
- Comprehensive simulation studies and motivating applications demonstrate the potential of our proposed debiasing method.



Potential Application to Causal Inference (I)

The observable data in causal inference are

$$\{(\mathbb{Y}_i, T_i, X_i)\}_{i=1}^n \subset \mathbb{R} \times \{0, 1\} \times \mathbb{R}^d.$$

- $T_i \in \{0,1\}$ is a binary treatment assignment indicator;
- $\mathbb{Y}_i = T_i \cdot Y(1)_i + (1 T_i) \cdot Y(0)_i$ with Y(0), Y(1) as potential outcomes.
- ▶ **Objective:** Conduct valid inference on the regression function (or conditional mean outcome) of the treatment group.

	X_1^T	$Y(1)_1$	
Treatment Group	:	:	
	$X_{rac{n}{2}}^{T}$	$Y(1)_{\frac{n}{2}}$	$\mathrm{E}\left(Y X,T=1\right)$
Control Group		$Y(0)_{\frac{n}{2}+1}$:	based on
	X_n^T	$Y(0)_n$	

Figure 8: Traditional approaches for inferring E(Y|X,T=1).



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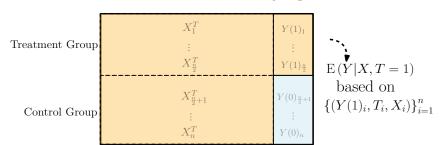


Figure 8: Our approach for inferring E(Y|X, T = 1).



Potential Application to Causal Inference (II)

Our debiasing method can be extended to valid inference on the linear average conditional treatment effect (ACTE)

$$\mathrm{E}[Y(1) - Y(0)|X]$$

with no unmeasured confounding and high-dimensional covariates.



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Our debiasing method can be extended to valid inference on the linear average conditional treatment effect (ACTE)

$$E[Y(1) - Y(0)|X]$$

with no unmeasured confounding and high-dimensional covariates.

• The modified debiasing program with tuning parameters $\gamma_1, \gamma_2 > 0$ is

$$\arg \min_{\boldsymbol{w}_{(0)}, \boldsymbol{w}_{(1)} \in \mathbb{R}^{n}} \sum_{i=1}^{n} \left[\widehat{\pi}_{i} w_{i(1)}^{2} + (1 - \widehat{\pi}_{i}) w_{i(0)}^{2} \right] \\
\text{s.t.} \quad \left\| x - \frac{1}{\sqrt{n}} \sum_{i=1}^{n} w_{i(1)} \cdot \widehat{\pi}_{i} \cdot X_{i} \right\| \leq \frac{\gamma_{1}}{n} \quad \text{and} \quad \left\| x - \frac{1}{\sqrt{n}} \sum_{i=1}^{n} w_{i(0)} \left(1 - \widehat{\pi}_{i} \right) X_{i} \right\| \leq \frac{\gamma_{2}}{n}.$$

The extended debiased estimator becomes

$$\begin{split} \widehat{m}^{\text{debias}}(x; \widehat{\boldsymbol{w}}_{(1)}, \widehat{\boldsymbol{w}}_{(0)}) \\ &= x^T \left(\widehat{\beta}_{(1)} - \widehat{\beta}_{(0)} \right) + \frac{1}{\sqrt{n}} \sum_{i=1}^{n} \left[\widehat{\boldsymbol{w}}_{i(1)} \cdot T_i \left(\mathbb{Y}_i - \boldsymbol{X}_i^T \widehat{\boldsymbol{\beta}}_{(1)} \right) - \widehat{\boldsymbol{w}}_{i(0)} \cdot (1 - T_i) \left(\mathbb{Y}_i - \boldsymbol{X}_i^T \widehat{\boldsymbol{\beta}}_{(0)} \right) \right]. \end{split}$$

The efficiency theory for this modified procedure is worth studying!

Thank you!

More details can be found in

[1] Y. Zhang, A. Giessing, and Y.-C. Chen. Efficient Inference on High-Dimensional Linear Models with Missing Outcomes. *arXiv* preprint, 2023. https://arxiv.org/abs/2309.06429.

Python Package: Debias-Infer and R Package: DebiasInfer.





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Implementation Details of the Proposed Debiasing Method

Lasso pilot estimate: We adopt the scaled Lasso (Sun and Zhang, 2012) with its universal regularization parameter $\lambda_0 = \sqrt{\frac{2 \log d}{n}}$ as the initialization. Specifically, it iteratively updates $\widehat{\beta}(\widetilde{\lambda})$, $\widehat{\sigma}_{\epsilon}(\widetilde{\lambda})$, $\widetilde{\lambda}$ via the jointly convex optimization program:

$$\left(\widehat{\beta}(\widetilde{\lambda}), \widehat{\sigma}_{\epsilon}(\widetilde{\lambda})\right) = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^{d}, \sigma_{\epsilon} > 0} \left[\frac{1}{2n\sigma_{\epsilon}} \sum_{i=1}^{n} R_{i} \left(Y_{i} - X_{i}^{T}\beta\right)^{2} + \frac{\sigma_{\epsilon}}{2} + \widetilde{\lambda} \left|\left|\beta\right|\right|_{1}\right].$$

Debiasing program: We solve the primal program by Python package "CVXPY" (Diamond and Boyd, 2016; Agrawal et al., 2018) or R package "CVXR" (Fu et al., 2020). For the dual program, we formulate a coordinate descent algorithm (Wright, 2015) as:

$$\left[\widehat{\ell}(x)\right]_{j} \leftarrow \frac{\mathcal{S}_{\frac{\gamma}{n}}\left(-\frac{1}{2n}\sum_{i=1}^{n}\widehat{\pi}_{i}\left(\sum_{k\neq j}X_{ik}X_{jk}\left[\widehat{\ell}(x)\right]_{k}\right) - x_{j}\right)}{\frac{1}{2n}\sum_{i=1}^{n}\widehat{\pi}_{i}X_{ij}^{2}} \text{ for } j = 1, ..., d,$$

where $S_{\frac{\gamma}{n}}(u) = \text{sign}(u) \cdot \left(u - \frac{\gamma}{n}\right)_{+}$ is the soft-thresholding operator.



One Standard Error (1SE) Rule For Model Selection

- Suppose that we conduct a K-fold cross-validation on a candidate set $\Gamma = \{\gamma_1, ..., \gamma_m\}$ of the tuning parameter.
- For each $\gamma_i \in \Gamma$, we compute the cross-validated risk or error on each fold of the data as:

$$CV_k(\gamma_i), k = 1, ..., K.$$

• For each $\gamma_i \in \Gamma$, we calculate the standard error of $CV_1(\gamma_i), ..., CV_K(\gamma_i)$ as:

$$SD(\gamma_i) = \sqrt{\text{Var}(CV_1(\gamma_i), ..., CV_K(\gamma_i))}, \quad SE(\gamma_i) = SD(\gamma_i)/\sqrt{K}.$$

Let

$$CV(\gamma) = \frac{1}{K} \sum_{k=1}^{K} CV_k(\gamma)$$
 and $\widehat{\gamma} = \operatorname*{arg\,min}_{\gamma \in \Gamma} CV(\gamma)$.

The 1SE rule (Breiman et al., 1984; Chen and Yang, 2021) selects $\gamma_{1SE} \in \Gamma$ with as the one with the smallest $CV(\gamma)$ such that

$$CV(\gamma_{1SE}) > CV(\widehat{\gamma}) + SE(\widehat{\gamma}).$$



One Standard Error (1SE) Rule For Model Selection

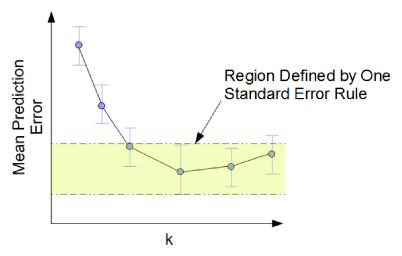


Figure 9: Illustration of the 1SE rule for selecting the model parameter.



Finger-of-God and Kaiser Effects

The galaxy distribution is distorted along the line of sight due to the peculiar velocities of galaxies, *i.e.*, the so-called *finger-of-god* (Jackson, 1972) and *Kaiser* (Kaiser, 1987) effects.

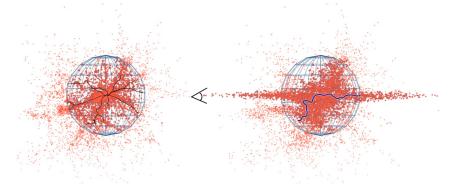


Figure 10: Redshift distortions along the line of sight (Kuchner et al., 2021).