

Clustered Covariate Regression

Abdul-Nasah Soale & Emmanuel Selorm Tsyawo

Temple University

Presenter: Emmanuel Selorm Tsyawo (Job Market Candidate)
estsyawo@temple.edu

Oct. 11, 2019

Introduction

- ▶ Identification is necessary for consistency
- ▶ A crucial aspect to identification in single-index models is the non-singularity of the Gram matrix $E(\mathbf{x}'\mathbf{x})$
 - ▶ Singularity of $E(\mathbf{x}'\mathbf{x}) \implies$ non-identification in e.g. linear and probit regressions
- ▶ Singularity of the Gram matrix caused by
 - ▶ High dimensionality - many/more covariates relative to observations
 - ▶ Multicollinearity
 - ▶ Both

Motivation

- ▶ Singularities of $E(\mathbf{x}'\mathbf{x})$ in the limit theory engenders
 - ▶ inconsistency
 - ▶ differing rates of convergence

(Phillips 2016)

- ▶ Estimation impossible without adjustments when non-singularity holds in finite samples

Contribution

A novel approach to identification, inference, and estimation,
namely, *Clustered Covariate Regression* (CCR)

- ▶ in the presence of rank deficiency, weak identification due to
 - ▶ high-dimensionality
 - ▶ multicollinearity
 - ▶ both

How does the CCR work?

- ▶ Project an $n \times p$ rank-deficient \mathbf{x} using a $p \times k$ projection matrix \mathbf{m} such that $\mathbf{m}'\mathbf{x}'\mathbf{x}\mathbf{m}$ is non-singular.
- ▶ $p > k, n \gg k$

Outline

Introduction

Related Literature

The CCR

Estimation

Asymptotic Theory

Monte Carlo Experiments

Empirical Model

Conclusion

Strands of related literature

- ▶ Variable selection, e.g.,
 - ▶ the Lasso (Belloni, Chernozhukov, and Hansen (2014b))
 - ▶ sequential hypothesis tests (Chudik, Kapetanios, and Pesaran (2018))
- ▶ Dimension reduction (using pre-constructed projection matrices) e.g., Wold, Esbensen, and Geladi, 1987
 - ▶ principal component regression (PCR)
 - ▶ partial least squares (PLS)

What is gained by the CCR?

- ▶ The CCR spans the class of single-index models, e.g., linear, logit, and quantile regressions
- ▶ CCR obviates sparsity assumption in the high-dimensional parameter β
 - ▶ but assumes (approximate) reducibility of a high-dimensional $\beta \in \mathbb{R}^p$ vector to a smaller identifiable $\delta \in \mathbb{R}^k$, $p > k$
- ▶ CCR obviates the choice of tuning parameter and penalty function
 - ▶ but requires a criterion (e.g., BIC) for model selection
- ▶ The CCR projection matrix \mathbf{m} is model, outcome, covariate dependent; it is not pre-constructed
- ▶ Immune to multicollinearity via projection

The conditional functional

Suppose a conditional functional (e.g., conditional expectation, conditional quantile)

$$\nu(y_i|\mathbf{x}_i) = g(\mathbf{x}_i\boldsymbol{\beta}) = g(\mathbf{x}_i\mathbf{m}\boldsymbol{\delta})$$

- ▶ $\boldsymbol{\beta} = \mathbf{m}\boldsymbol{\delta}$ (or approximately)
- ▶ unknown parameters $\boldsymbol{\beta}$ is $p \times 1$, $\boldsymbol{\delta}$ is $k \times 1$
- ▶ \mathbf{m} is an unknown $p \times k$ (clustering) projection matrix

The projection matrix \mathbf{I}

Characteristics of the projection matrix \mathbf{m} :

- ▶ \mathbf{m} belongs to a set \mathcal{M} of $p \times k$ matrices
- ▶ has exactly p non-zero elements
- ▶ the columns of \mathbf{m} correspond to clusters
- ▶ each row comprises only one non-zero element
- ▶ the vector of non-zero elements in \mathbf{m} are researcher-specified, e.g. standard deviations or a vector of ones
- ▶ Cluster assignments unknown a priori

The projection matrix II

An example

- ▶ $\boldsymbol{\delta} = [\delta_1, \delta_2]'$, $\mathbf{x} = [x_1, x_2, x_3, x_4]$, and $\mathbf{m}' = \begin{bmatrix} \eta_1 & 0 & 0 & \eta_4 \\ 0 & \eta_2 & \eta_3 & 0 \end{bmatrix}$
- ▶ The linear predictor function
$$\mathbf{xm}\boldsymbol{\delta} = x_1\eta_1\delta_1 + x_2\eta_2\delta_2 + x_3\eta_3\delta_2 + x_4\eta_4\delta_1 = x_1\beta_1 + x_2\beta_2 + x_3\beta_3 + x_4\beta_4 = \mathbf{x}\boldsymbol{\beta}$$
- ▶ factorises $\boldsymbol{\beta}$ into $\mathbf{m}\boldsymbol{\delta}$
- ▶ if scaling, η_j specified by researcher else $\eta_j = 1$ for all $j = 1, \dots, p$

The CCR estimation problem

- ▶ An objective, e.g. linear regression

$$Q_n(\mathbf{x}m\delta) \equiv n^{-1} \sum_{i=1}^n q(\mathbf{x}_i m \delta), \quad q(\mathbf{x}_i m \delta) = (y_i - \mathbf{x}_i m \delta)^2$$

- ▶ $Q_n(\delta) \equiv \frac{1}{|\mathcal{A}_n|} \int_{\mathcal{M}} Q_n(\mathbf{x}m\delta) \mathbf{1}_{\mathcal{A}_n}(\mathbf{m}) d\mu(\mathbf{m}) =$
 $\frac{1}{|\mathcal{A}_n|} \sum_{\mathbf{m} \in \mathcal{A}_n} Q_n(\mathbf{x}m\delta)$

- ▶ μ is the counting measure

- ▶ set $\mathcal{A}_n \equiv \{\mathbf{m} \in \mathcal{M} : Q_n(\mathbf{x}m\delta) \leq Q_n(\mathbf{x}s\delta) \ \forall \mathbf{s} \in \mathcal{M} \setminus \mathbf{m}\}$

- ▶ $\mathbf{1}_{\mathcal{A}_n}(\mathbf{m}) \equiv \mathbf{1}\{\mathbf{m} \in \mathcal{A}_n\}$ indicator for $\mathbf{m} \in \mathcal{A}_n$

- ▶ $|\mathcal{A}_n| \equiv \mu(\mathcal{A}_n)$ denotes the cardinality of \mathcal{A}_n

The Estimation Problem

Minimisation

$$\hat{\delta} = \arg \min_{\delta \in \Delta} Q_n(\delta)$$

- ▶ If an $\mathbf{m} \in \mathcal{A}_n$ is known, δ can be estimated
- ▶ But, $\mathbf{m} \in \mathcal{A}_n$ is unknown, $\implies \delta$ is unknown
- ▶ Approach: use a sequential scheme to estimate $\mathbf{m} \in \mathcal{A}_n$ and δ

The Sequential CCR Algorithm

fix k (number of clusters)

1. Initialise counter $l = 0$, parameter vector $\beta^{(l)} = \mathbf{m}^{(l)}\delta^{(l)}$
2. Update $l \leftarrow l + 1$, for each $j = 1, \dots, p$,
 - ▶ update $\hat{\beta}_j^{(l)} = \arg \min_{\beta_j} Q_n(\mathbf{x}_{-j}\beta_{-j}^{(l)} + \mathbf{x}_j\beta_j)$
 - ▶ assign $\hat{\beta}_j^{(l)}$ to a cluster and update $\mathbf{m}^{(l)}$
 - ▶ update $\delta^{(l)} = \arg \min_{\delta \in \Delta} Q_n(\mathbf{x}\mathbf{m}^{(l)}\delta)$
 - ▶ update $\beta^{(l)} \leftarrow \mathbf{m}^{(l)}\delta^{(l)}$
3. Check convergence for $Q_n(\mathbf{x}\mathbf{m}^{(l-1)}\delta^{(l-1)}) - Q_n(\mathbf{x}\mathbf{m}^{(l)}\delta^{(l)}) < \epsilon$
else return to step 2

Without clustering, the algorithm is similar to the (block)-coordinate descent algorithm (used for e.g. Lasso)

Optimal k is determined using a model selection criterion, e.g., BIC

Assumptions

1. $\mu(\mathcal{A}_n) \rightarrow 1$ as $n \rightarrow \infty$, i.e., $\mathcal{A}_n \rightarrow \{\mathbf{m}_o\}$ as $n \rightarrow \infty$
2. $\sqrt{p/n} \rightarrow 0$ as $p, n \rightarrow \infty$

Theorem - Consistency

- ▶ $\hat{\delta}_n \xrightarrow{p} \delta_o$ under standard assumptions
- ▶ $\hat{\beta}_n \xrightarrow{p} \beta_o$ where $\hat{\beta}_n \equiv \mathbf{m}_n \hat{\delta}_n$ and $\beta_o \equiv \mathbf{m}_o \delta_o$

Theorem - Asymptotic Normality

$\sqrt{n}(\hat{\delta}_n - \delta_o) \xrightarrow{d} \mathcal{N}(\mathbf{0}, \mathbf{V}_o)$ where $\mathbf{V}_o = \mathbf{A}_o^{-1} \mathbf{B}_o \mathbf{A}_o^{-1}$,
 $\mathbf{A}_o \equiv E[\mathbf{H}(\mathbf{x}_i \mathbf{m}_o \delta_o)]$, and $\mathbf{B}_o \equiv E[\mathbf{s}(\mathbf{x}_i \mathbf{m}_o \delta_o) \mathbf{s}(\mathbf{x}_i \mathbf{m}_o \delta_o)']$

CCR in a baseline specification

Table: Baseline model: $\mathbf{x}_i \stackrel{iid}{\sim} \mathcal{N}(0, \mathbf{I}_p)$, $\boldsymbol{\delta} = [-2, -1, 0, 1, 2, 3]'$

| n | p | CCR | OLS | LASSO | PCR | PLS |
|----|----|----------------------|-----------------------|----------------------|--------------|-----------------------|
| 30 | 12 | 0.201(0.086) | 0.199(0.054) | 0.197 (0.055) | 0.395(0.215) | 0.325(0.112) |
| | 24 | 0.365 (0.136) | 0.375(0.145) | 0.372(0.140) | 0.515(0.173) | 0.400(0.119) |
| | 36 | 0.821(0.181) | - | 1.400(6.244) | 0.769(0.144) | 0.766(0.133) |
| 90 | 12 | 0.069 (0.023) | 0.091(0.021) | 0.091(0.022) | 0.357(0.231) | 0.188(0.055) |
| | 24 | 0.060 (0.023) | 0.010(0.018) | 0.099(0.018) | 1.061(0.158) | 0.353(0.078) |
| | 36 | 0.065 (0.027) | 0.110(0.018) | 0.109(0.018) | 1.225(0.117) | 0.490(0.083) |

Note: (1) Results: average $d_{\beta} = \|\hat{\beta} - \beta_o\|_1/p$. (2) 1000 simulations each (3) $\sigma(d_{\beta})$ in parentheses. (4) Optimal k - BIC. (5) Two-step Lasso, PCR, and PLS - 10-fold CV (6) $k^ = 6$.*

CCR under Multicollinearity

Table: Multicollinearity: $\mathbf{x}_i \stackrel{iid}{\sim} \mathcal{N}(0, \Sigma)$, $\Sigma_{jj'} = 0.5^{|j-j'|}$

| n | p | CCR | OLS | LASSO | PCR | PLS |
|----|----|----------------------|-----------------------|----------------------|----------------------|-----------------------|
| 30 | 12 | 0.283(0.106) | 0.250(0.075) | 0.246 (0.076) | 0.336(0.143) | 0.554(0.126) |
| | 24 | 0.468(0.168) | 0.481(0.194) | 0.448 (0.164) | 0.479(0.137) | 0.533(0.118) |
| | 36 | 1.118(0.241) | - | 1.197(2.346) | 0.714 (0.132) | 0.757(0.118) |
| 90 | 12 | 0.094 (0.041) | 0.116(0.030) | 0.114(0.031) | 0.250(0.127) | 0.389(0.067) |
| | 24 | 0.076 (0.036) | 0.128(0.026) | 0.126(0.026) | 0.300(0.118) | 0.285(0.061) |
| | 36 | 0.083 (0.043) | 0.142(0.026) | 0.140(0.026) | 0.327(0.111) | 0.389(0.065) |

Note: (1) Results: average $d_{\beta} = \|\hat{\beta} - \beta_o\|_1/p$. (2) 1000 simulations each (3) $\sigma(d_{\beta})$ in parentheses. (4) Optimal k - BIC. (5) Two-step Lasso, PCR, and PLS - 10-fold CV (6) $k^* = 6$.

The empirical model

Estimating private and spillover effects of R&D on productivity

$$E[y_{it} | \mathbf{w}_{it}, \mathbf{x}_t] = \alpha_0 + \mathbf{w}_{it}\boldsymbol{\theta} + x_{it}\gamma_{ii} + \sum_{j \neq i} x_{jt}\gamma_{ij} + \alpha_t + \delta_i$$

- ▶ $k_w + T + N + N^2$ parameters, NT observations
- ▶ e.g. $T = 27, N = 50, p = 2577$ parameters, $n = NT = 1350$ firm-year observations
- ▶ $p > n$

Conclusion

In a nutshell, we propose Clustered Covariate Regression. It is a novel approach to

- ▶ handling rank-deficiency in single-index models
 - ▶ multicollinearity, high-dimensionality, or both
- ▶ offers advantages: e.g. obviates sparsity in β and increases precision
- ▶ interesting extensions (left for future work)
 - ▶ high-dimensional causal inference
 - ▶ multicollinearity in non-linear models
 - ▶ estimating latent network structures from panel data