

Faster and more robust maximum likelihood estimation for random utility models

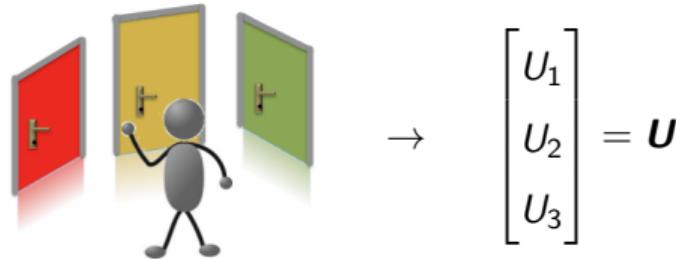
Lennart Oelschläger Dietmar Bauer

Bielefeld University, Empirical Methods Department, Econometrics Group

ICMC April 1-3 2024 in Puerto Varas, Chile

Faster and more robust MLE for RUMs

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- $\boldsymbol{\mu}$ are alternative-specific constants

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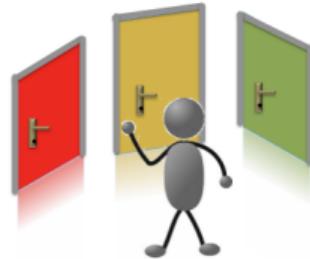
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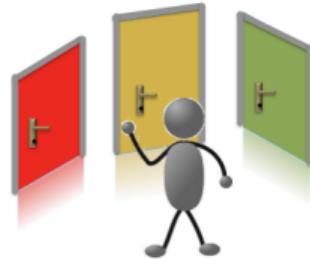
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Let θ denote the set of model parameters and assume a distribution for ε is fixed.

1. calculate choice probabilities $P_{nj}(\mathbf{X}_n, \theta)$ for each decider n and alternative j
2. build log-likelihood function $\log \mathcal{L}(\theta | \mathbf{X}, y) = \sum_{n,j} 1(y_n = j) \log P_{nj}(\mathbf{X}_n, \theta)$
3. maximize $\log \mathcal{L}(\theta | \mathbf{X}, y)$ with respect to θ

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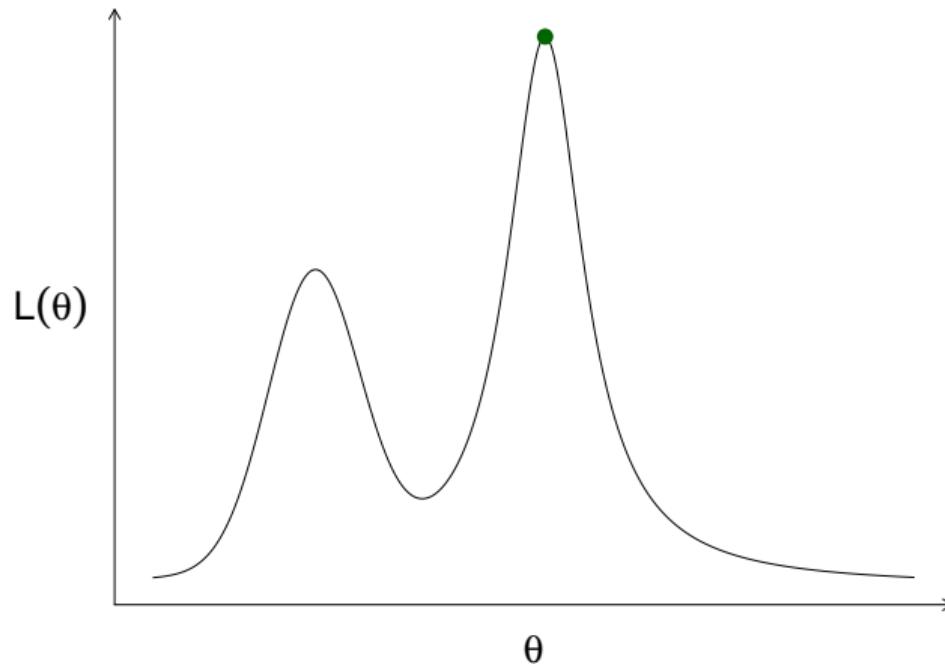


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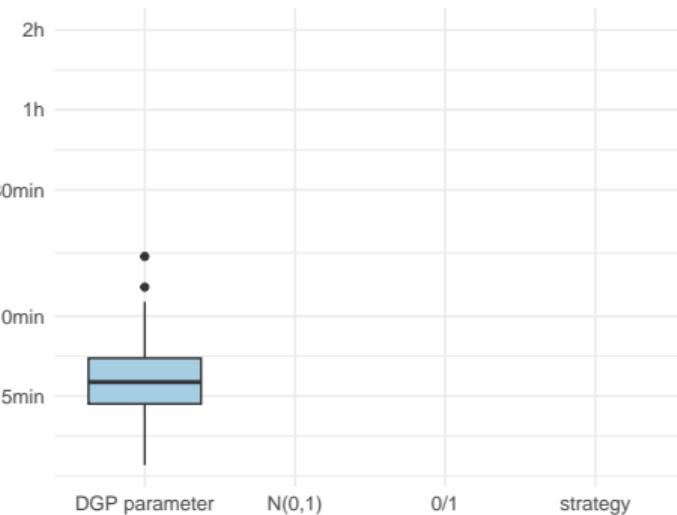
Outline of this talk

- 1** How slow and unreliable can it be?
- 2** Initialization strategies and proof of concept
- 3** Putting them together
- 4** Let's try with empirical data
- 5** Takeaways

100 data sets simulated from a probit model with $\theta_i \sim N(0, 1)$:

$N = 200$ deciders, $T = 10$ choices each, $J = 5$ alternatives, 5 covariates with normal random effects

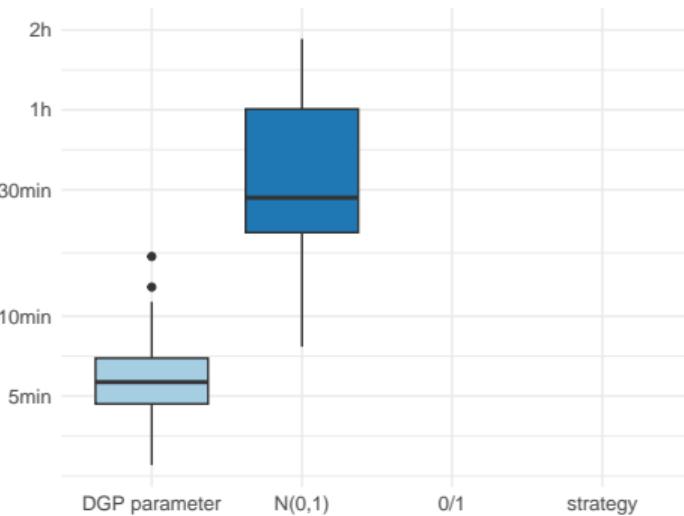
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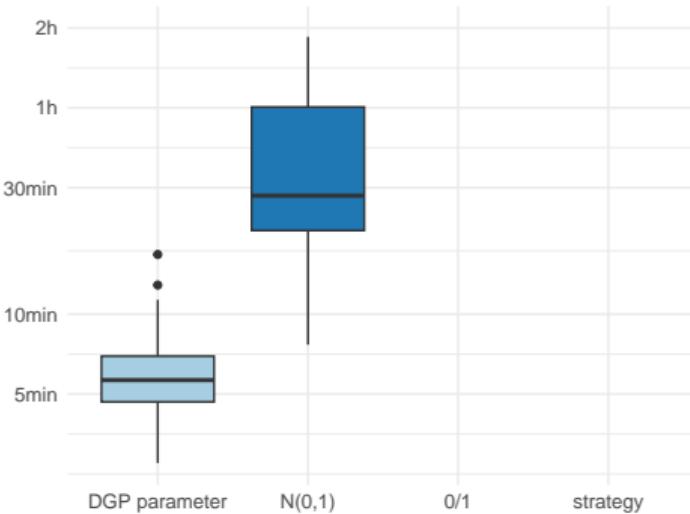
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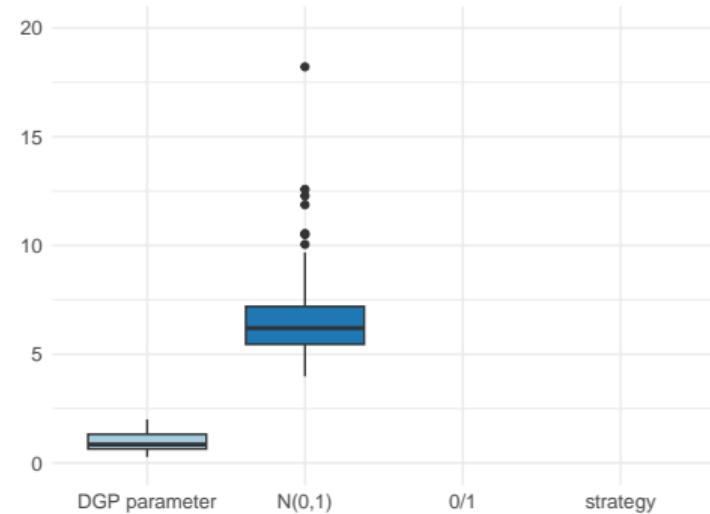
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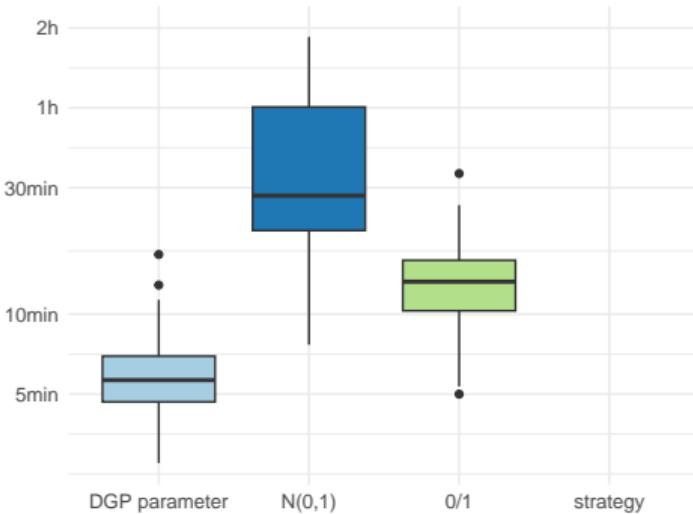
Euclidean distance from initials to MLE



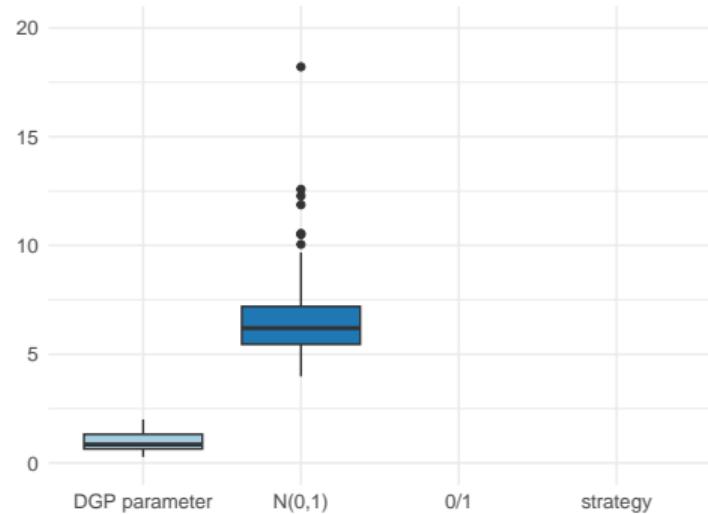
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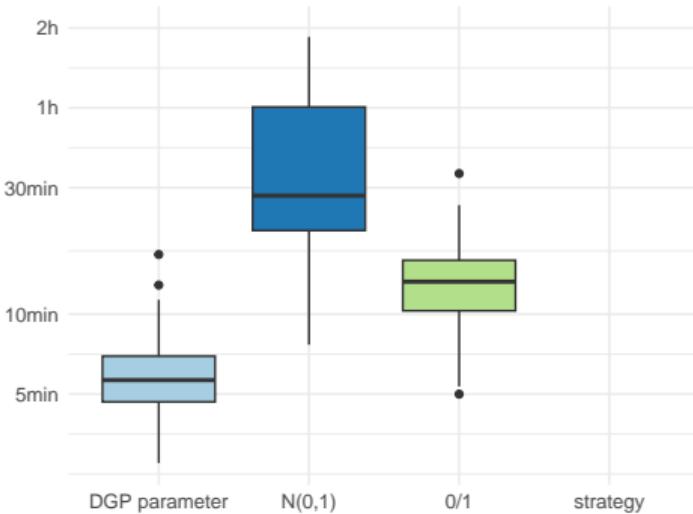
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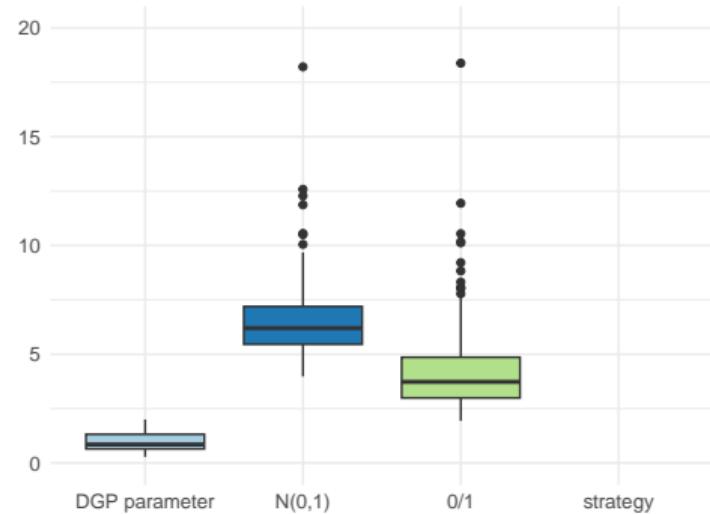
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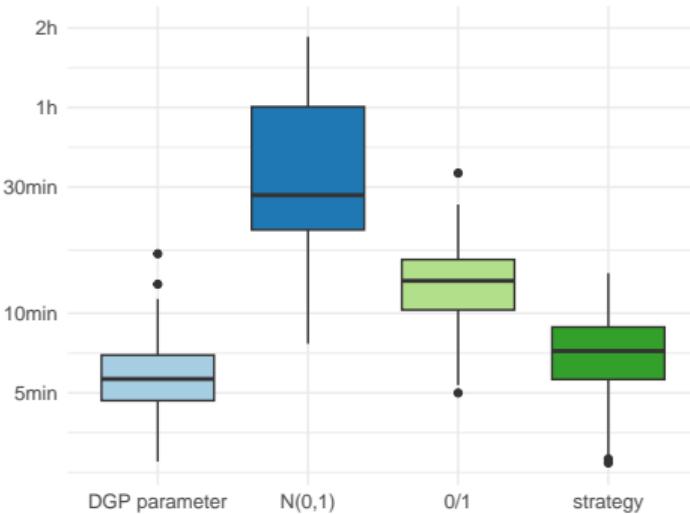
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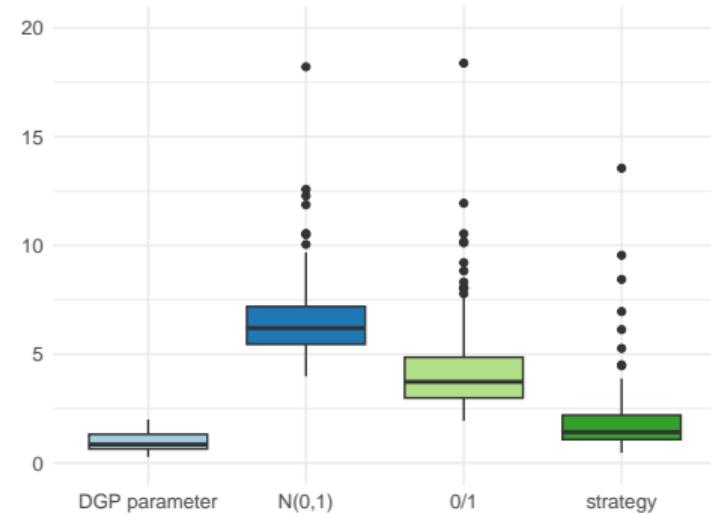
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$$U = \mu + X(\beta + \gamma) + \varepsilon$$

Parameter	Initialization strategy
β is alternative-varying and	
X is alternative-varying	constant utility direction
X is alternative-constant (ASCs μ is special case)	minimize choice frequency prediction error
β is alternative-constant	linear probability model
covariance Σ for ε	MCMC
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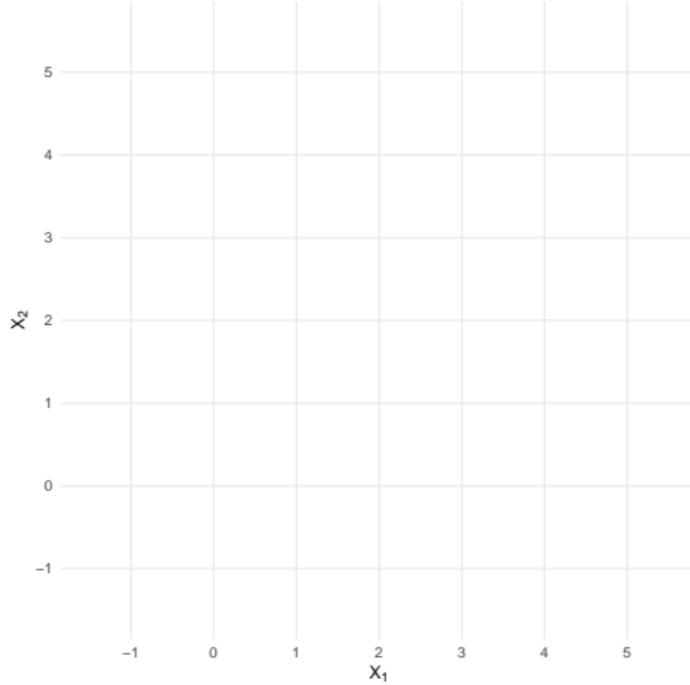
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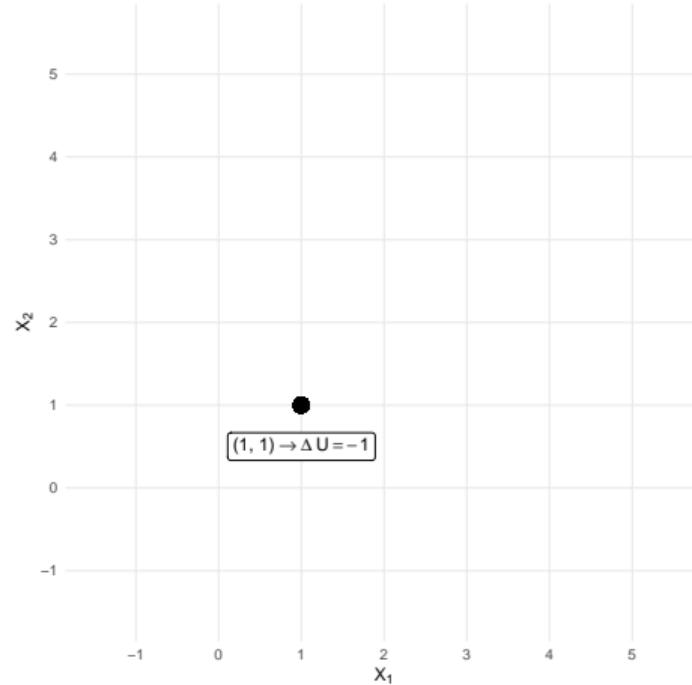
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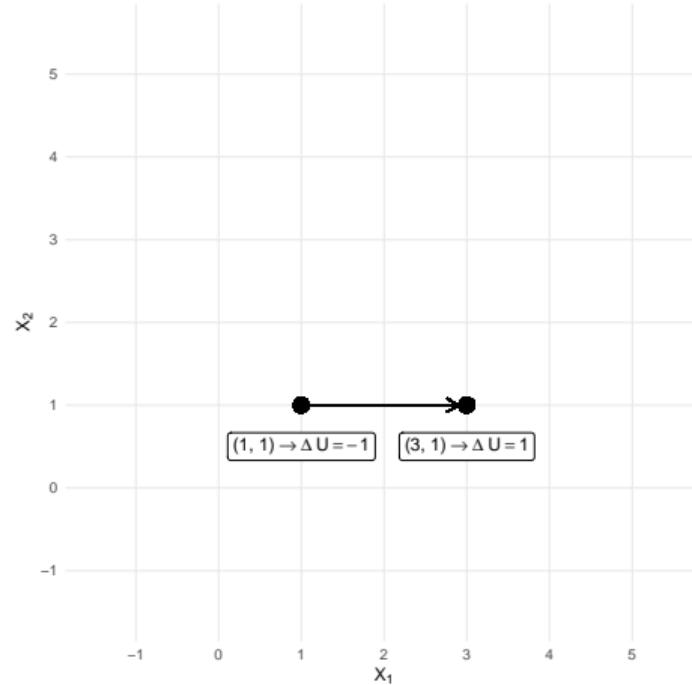
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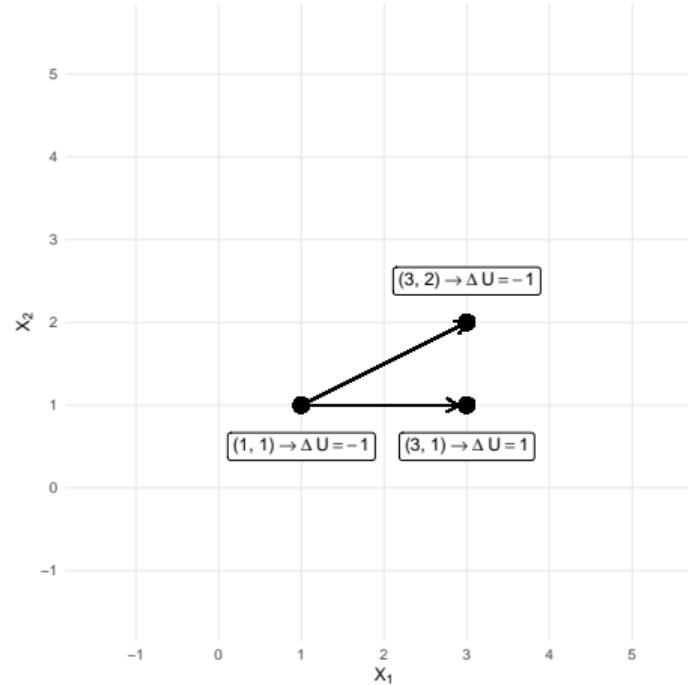
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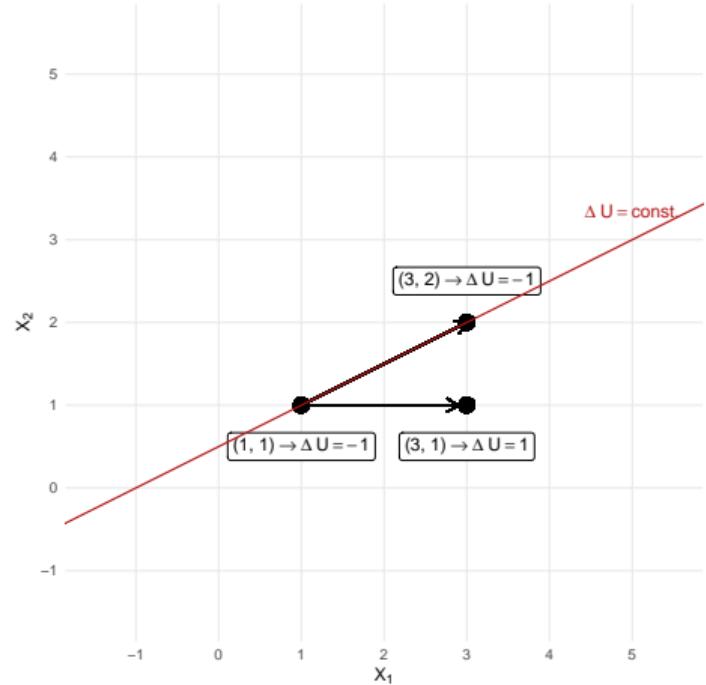
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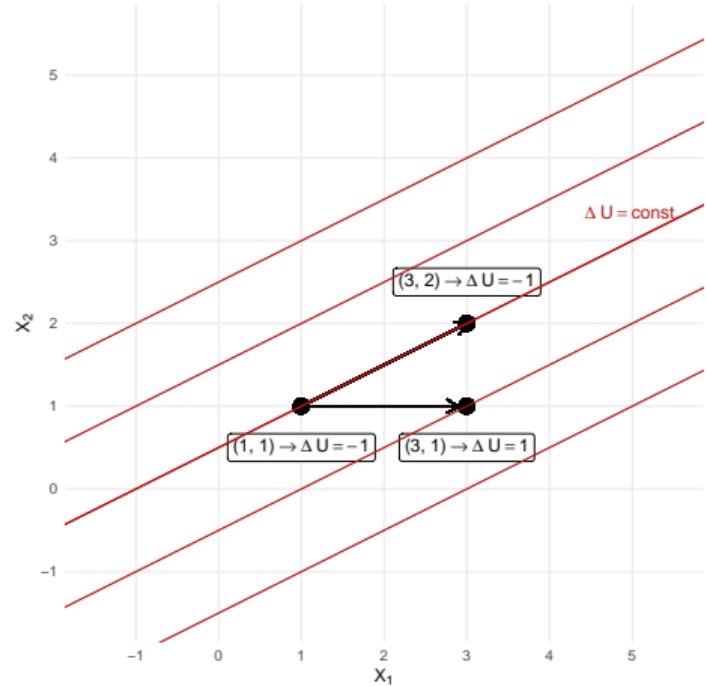


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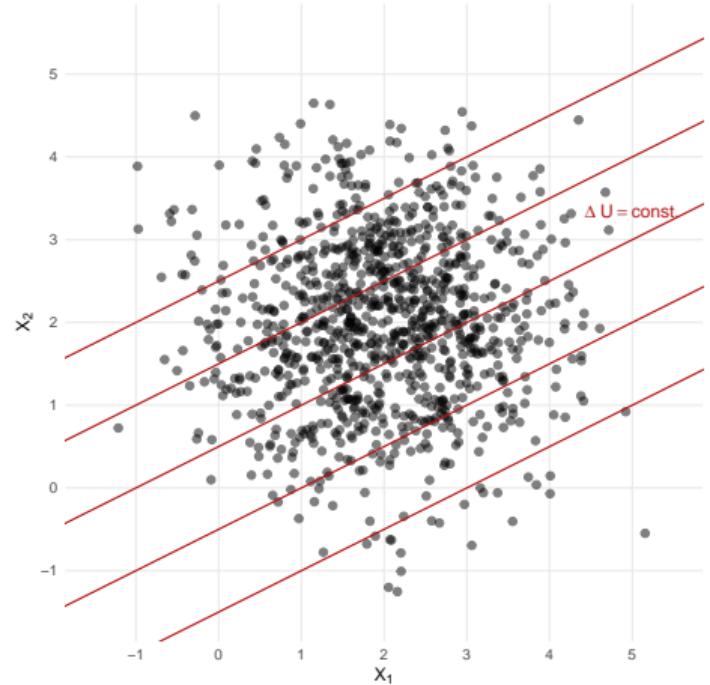


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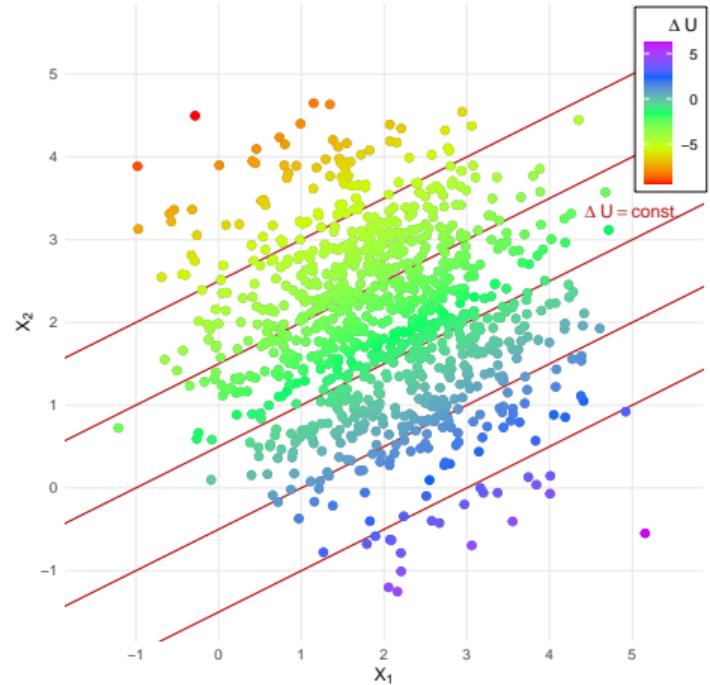


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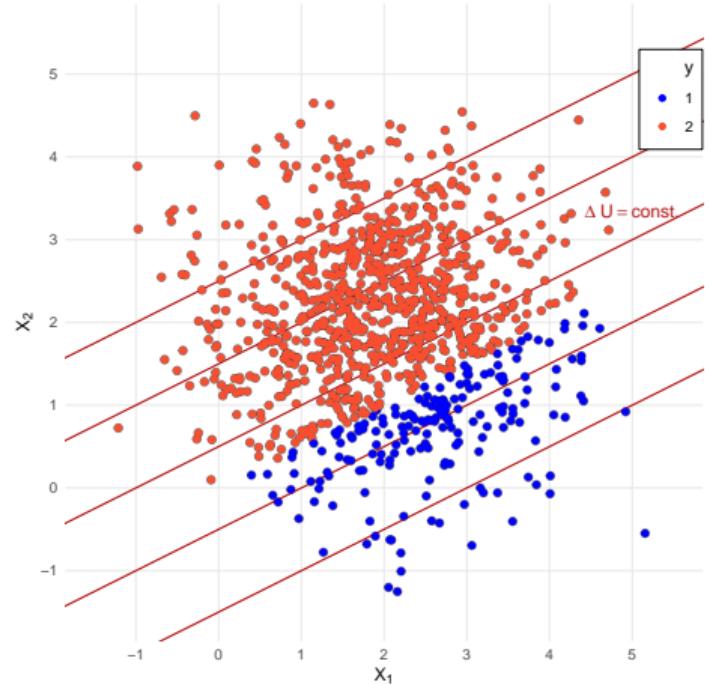


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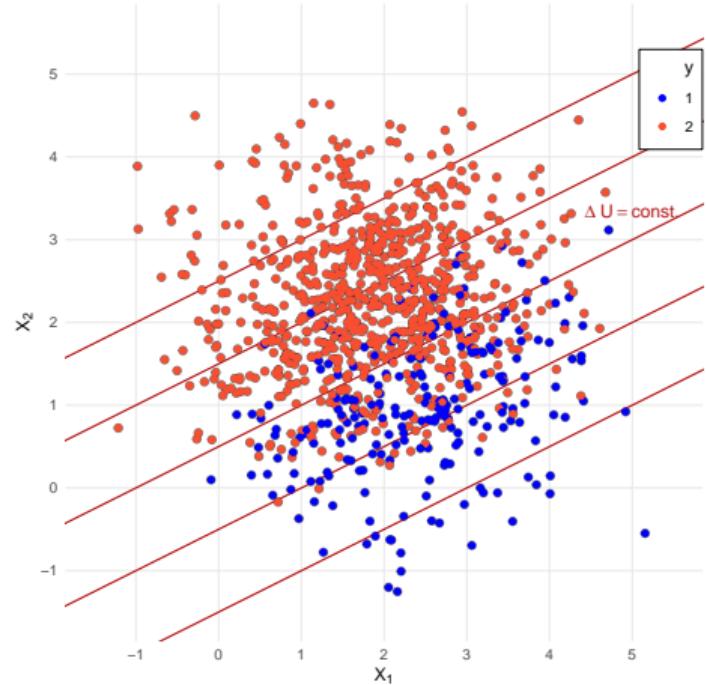


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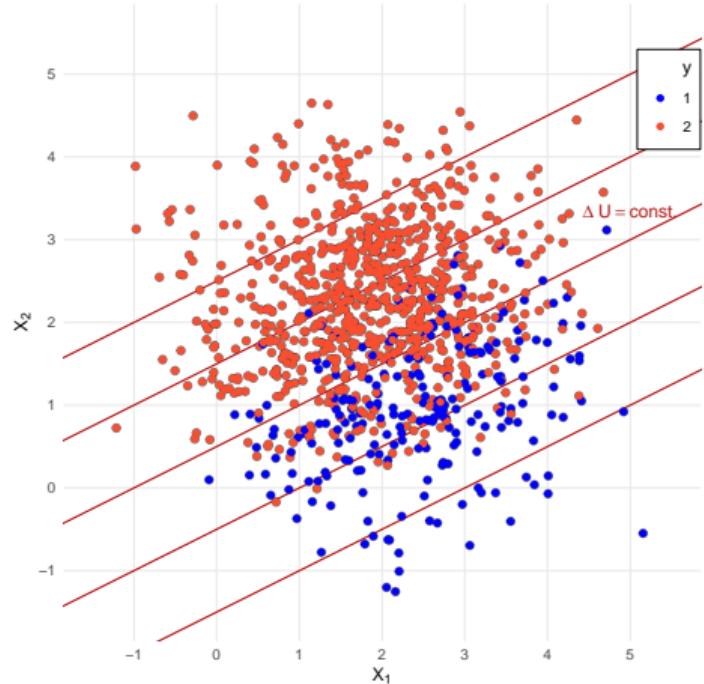
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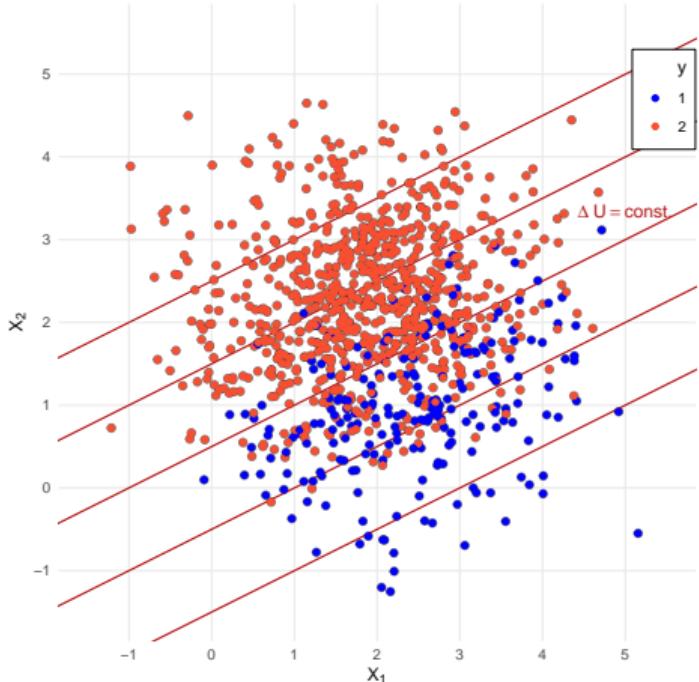
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this yields a consistent initial estimator $\hat{\boldsymbol{\beta}}$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, 1.000 replications

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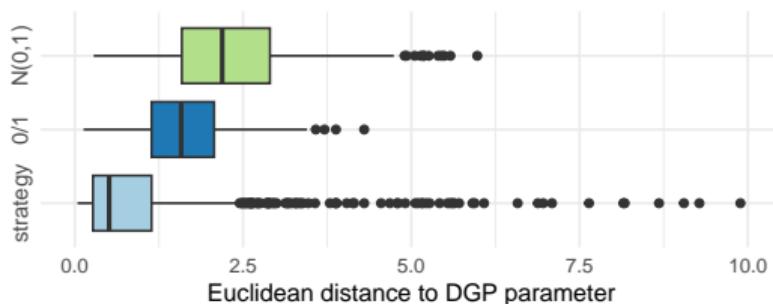
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N	1.000	■	10.000	□
J	4	■	8	□
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□

Average computation time: < 1 second



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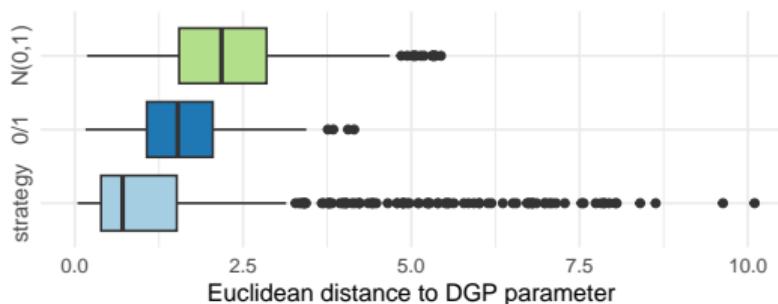
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$$\begin{pmatrix} U_1 \\ U_2 \end{pmatrix} = \begin{pmatrix} X_1 & 0 \\ 0 & X_2 \end{pmatrix} \begin{pmatrix} 1 \\ 2 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$y = \begin{cases} 1, & \text{if } \Delta U = U_1 - U_2 \geq 0 \\ 2, & \text{otherwise} \end{cases}$$

$$\Delta U = \text{const. in direction } \overrightarrow{\begin{pmatrix} 1 \\ 0.5 \end{pmatrix}} = \overrightarrow{\begin{pmatrix} 1/\beta_1 \\ 1/\beta_2 \end{pmatrix}}$$

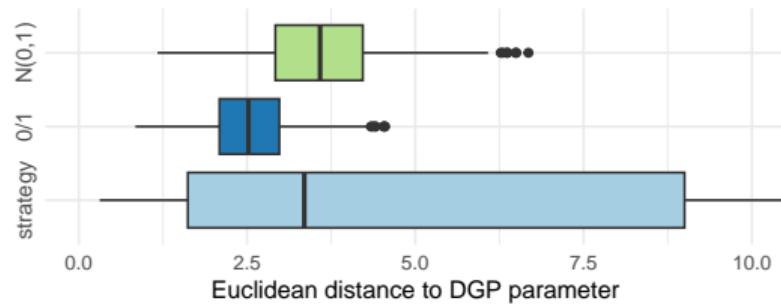
identify direction as the kernel of $\text{Cov}(y, \mathbf{X})$

this yields a consistent initial estimator $\hat{\boldsymbol{\beta}}$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

N	1.000	■	10.000	□
J	4	□	8	■
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, 1.000 replications

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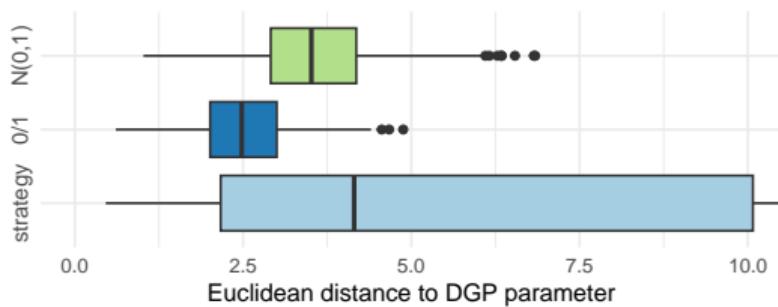
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J	4	□	8	■
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■

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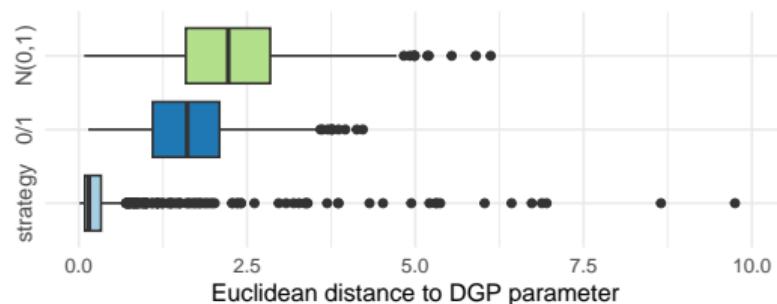
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N	1.000	<input type="checkbox"/>	10.000	<input checked="" type="checkbox"/>
J	4	<input checked="" type="checkbox"/>	8	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>

Average computation time: < 1 second



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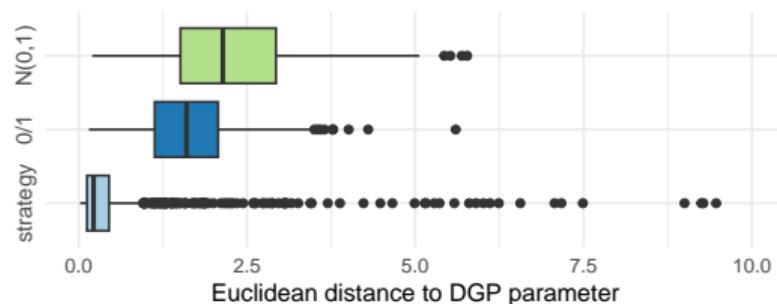
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J	4	■	8	□
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■

Average computation time: < 1 second



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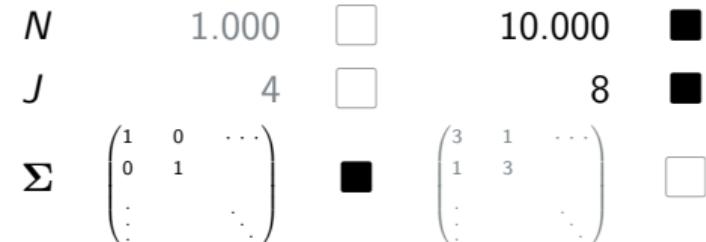
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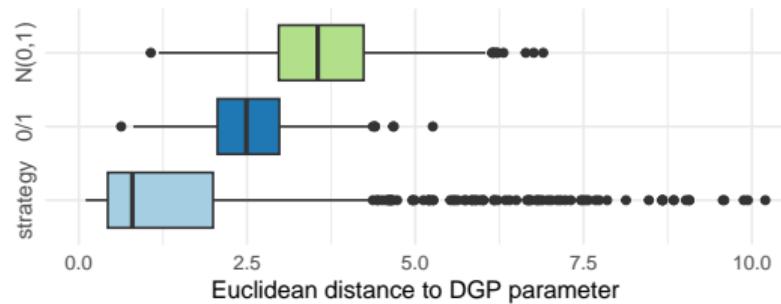
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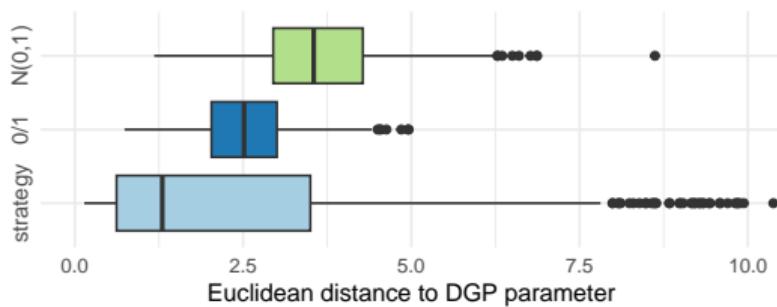
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N	1.000	□	10.000	■
J	4	□	8	■
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■

Average computation time: < 1 second



$$U = \mu + X(\beta + \gamma) + \varepsilon$$

Parameter

Initialization strategy

β is alternative-varying and

constant utility direction

X is alternative-varying

minimize choice frequency prediction error

► X is alternative-constant (ASCs μ is special case)

linear probability model

β is alternative-constant

MCMC

covariance Σ for ε

MCMC

covariance Ω for γ

$$\boldsymbol{U} = \boldsymbol{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma})$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} X & 0 & 0 \\ 0 & X & 0 \\ 0 & 0 & X \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

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$$y = \begin{cases} 1, & \text{if } U_1 = \max \mathbf{U} \\ 2, & \text{if } U_2 = \max \mathbf{U} \\ 3, & \text{if } U_3 = \max \mathbf{U} \end{cases}$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

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$$\mathbf{y}^d = \begin{cases} \begin{pmatrix} 1 \\ 0 \\ 0 \end{pmatrix}, & \text{if } U_1 = \max \mathbf{U} \\ \begin{pmatrix} 0 \\ 1 \\ 0 \end{pmatrix}, & \text{if } U_2 = \max \mathbf{U} \\ \begin{pmatrix} 0 \\ 0 \\ 1 \end{pmatrix}, & \text{if } U_3 = \max \mathbf{U} \end{cases}$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

Initialization strategy:

1. assume Σ is known, or choose $\Sigma = I_J$ else
2. let $\bar{\mathbf{y}}^d$ be the average of \mathbf{y}^d
3. find $\hat{\boldsymbol{\beta}} = \arg \min_{\boldsymbol{\beta}} \|P(\boldsymbol{\beta}, \Sigma) - \bar{\mathbf{y}}^d\|_2$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} \textcolor{red}{X} & 0 & 0 \\ 0 & \textcolor{red}{X} & 0 \\ 0 & 0 & \textcolor{red}{X} \end{pmatrix} \begin{pmatrix} 0 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

Initialization strategy:

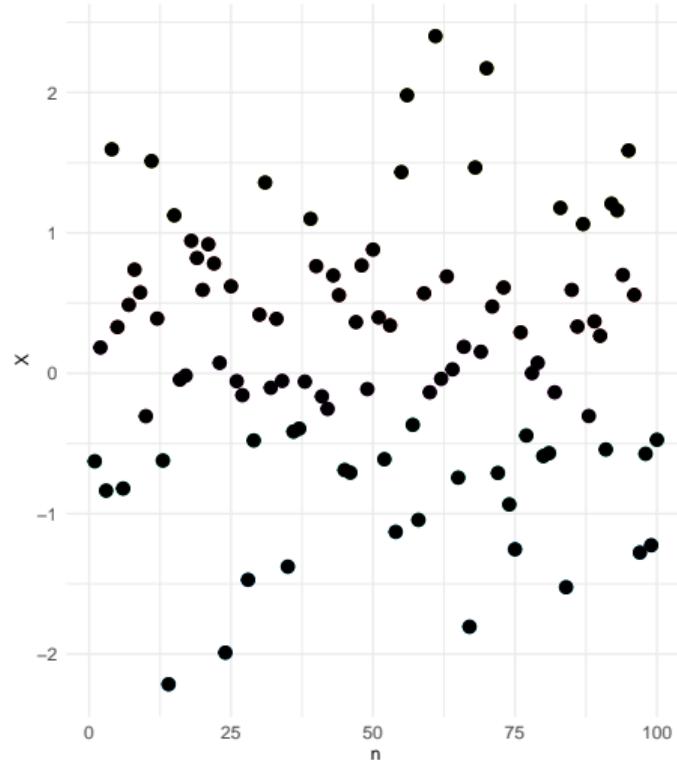
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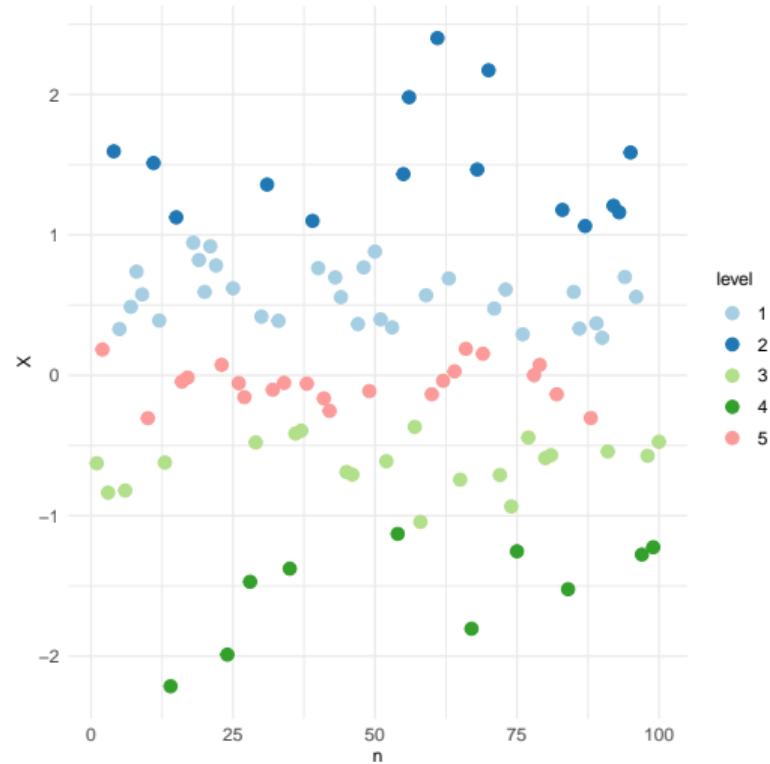


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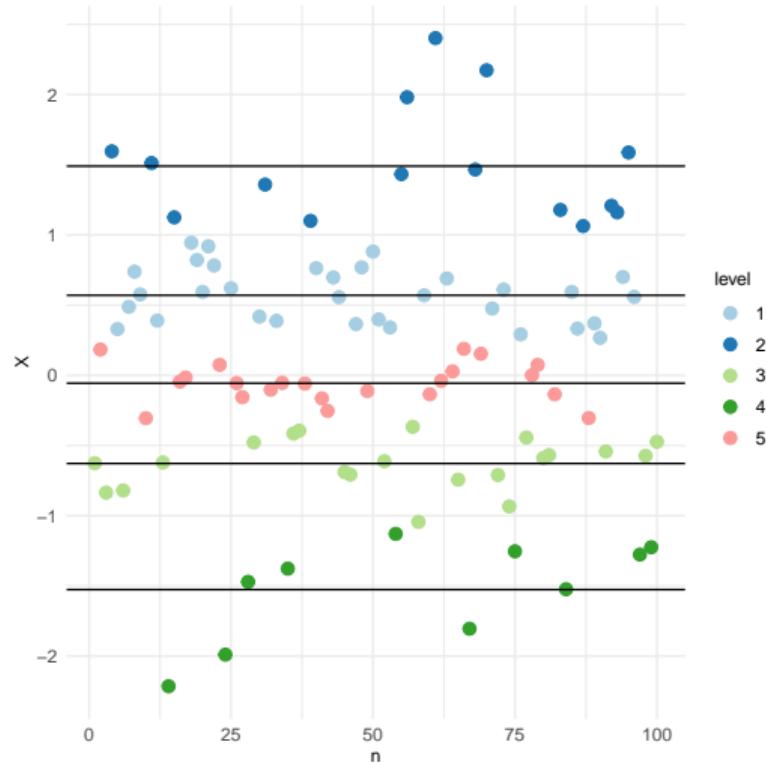


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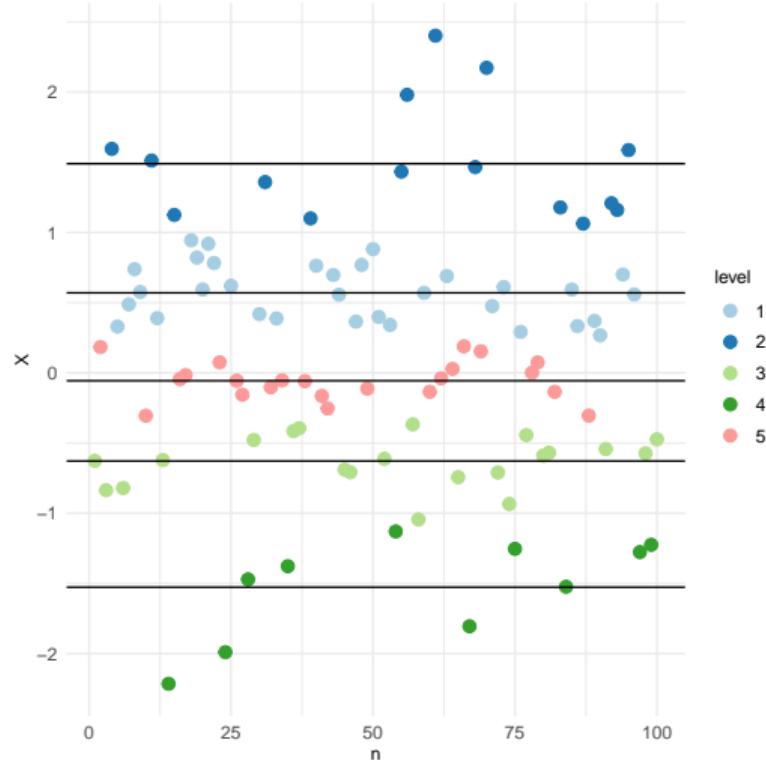


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Initialization strategy:

1. assume $\boldsymbol{\Sigma}$ is known, or choose $\boldsymbol{\Sigma} = I_J$ else
2. localize X , for each level $L_i \neq 0$:
 - 2.1 let $\bar{\mathbf{y}}_i^d$ be the average of \mathbf{y}_i^d
 - 2.2 find $\hat{\boldsymbol{\beta}}_i = \arg \min_{\boldsymbol{\beta}} \|P(\boldsymbol{\beta}, \boldsymbol{\Sigma}) - \bar{\mathbf{y}}_i^d\|_2$
 - 2.3 $\hat{\boldsymbol{\beta}}_i \leftarrow \hat{\boldsymbol{\beta}}_i / L_i$
3. $\hat{\boldsymbol{\beta}} = \bar{\hat{\boldsymbol{\beta}}}_i$



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

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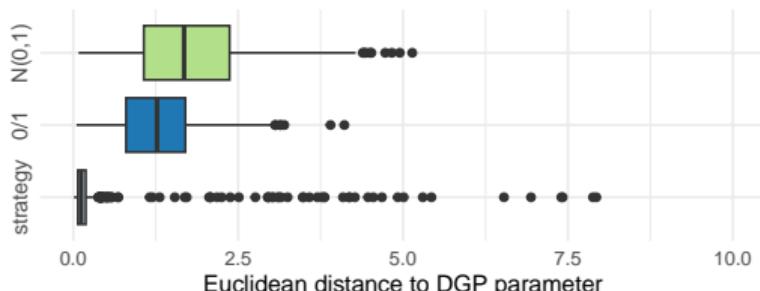
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Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $N = 1.000$, 1.000 rep.

J	4	<input checked="" type="checkbox"/>	8	<input type="checkbox"/>
levels	1	<input checked="" type="checkbox"/>	N	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>

Average computation time: < 1 second



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

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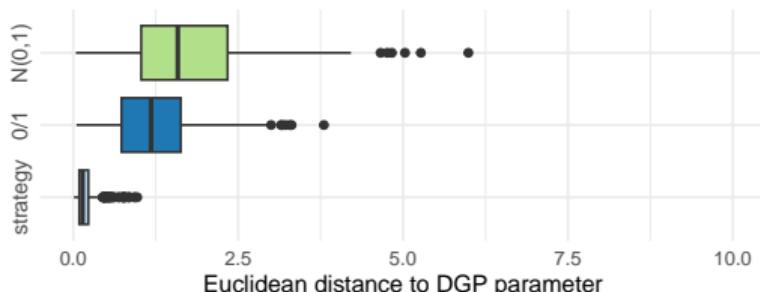
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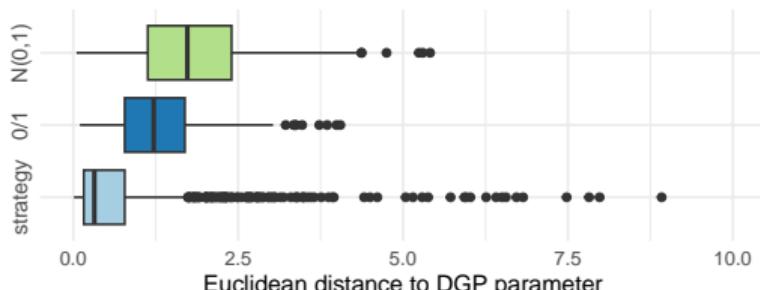
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J	4	<input checked="" type="checkbox"/>	8	<input type="checkbox"/>
levels	1	<input type="checkbox"/>	N	<input checked="" type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>

Average computation time: < 1 second



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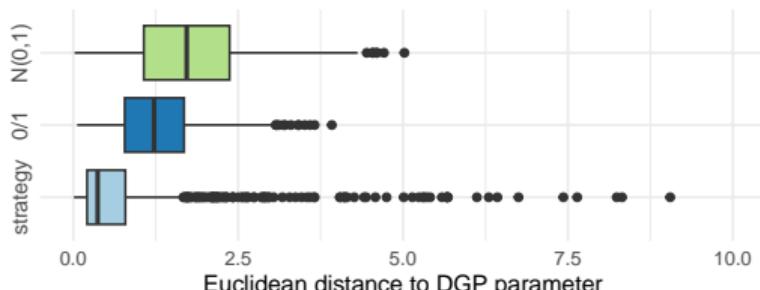
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Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $N = 1.000$, 1.000 rep.

J	4	■	8	□
levels	1	□	N	■
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■

Average computation time: < 1 second



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} X & 0 & 0 \\ 0 & X & 0 \\ 0 & 0 & X \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

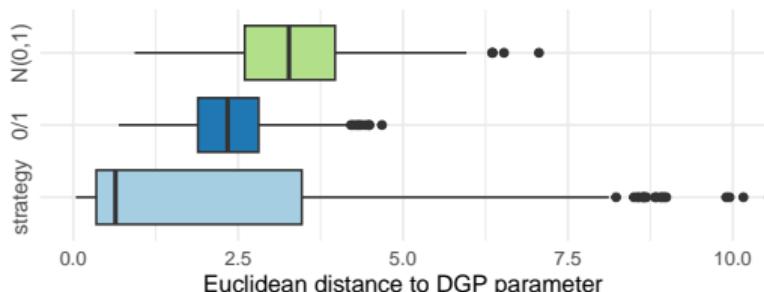
Initialization strategy:

1. assume Σ is known, or choose $\Sigma = I_J$ else
2. localize X , for each level $L_i \neq 0$:
 - 2.1 let $\bar{\mathbf{y}}_i^d$ be the average of \mathbf{y}_i^d
 - 2.2 find $\hat{\boldsymbol{\beta}}_i = \arg \min_{\boldsymbol{\beta}} \|P(\boldsymbol{\beta}, \Sigma) - \bar{\mathbf{y}}_i^d\|_2$
 - 2.3 $\hat{\boldsymbol{\beta}}_i \leftarrow \hat{\boldsymbol{\beta}}_i / L_i$
3. $\hat{\boldsymbol{\beta}} = \bar{\hat{\boldsymbol{\beta}}}$

Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $N = 1.000$, 1.000 rep.

J	4	<input type="checkbox"/>	8	<input checked="" type="checkbox"/>
levels	1	<input checked="" type="checkbox"/>	N	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>

Average computation time: 2 seconds



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} X & 0 & 0 \\ 0 & X & 0 \\ 0 & 0 & X \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

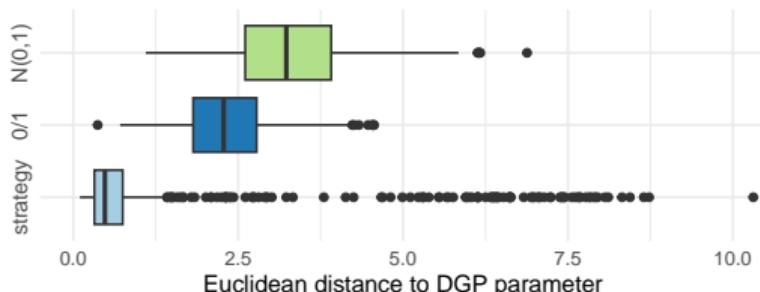
Initialization strategy:

1. assume Σ is known, or choose $\Sigma = I_J$ else
2. localize X , for each level $L_i \neq 0$:
 - 2.1 let $\bar{\mathbf{y}}_i^d$ be the average of \mathbf{y}_i^d
 - 2.2 find $\hat{\boldsymbol{\beta}}_i = \arg \min_{\boldsymbol{\beta}} \|P(\boldsymbol{\beta}, \Sigma) - \bar{\mathbf{y}}_i^d\|_2$
 - 2.3 $\hat{\boldsymbol{\beta}}_i \leftarrow \hat{\boldsymbol{\beta}}_i / L_i$
3. $\hat{\boldsymbol{\beta}} = \bar{\hat{\boldsymbol{\beta}}}$

Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $N = 1.000$, 1.000 rep.

J	4	<input type="checkbox"/>	8	<input checked="" type="checkbox"/>
levels	1	<input checked="" type="checkbox"/>	N	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>

Average computation time: 2 seconds



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} X & 0 & 0 \\ 0 & X & 0 \\ 0 & 0 & X \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

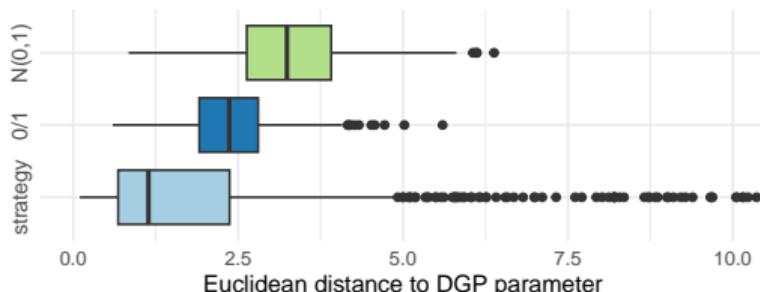
Initialization strategy:

1. assume Σ is known, or choose $\Sigma = I_J$ else
2. localize X , for each level $L_i \neq 0$:
 - 2.1 let $\bar{\mathbf{y}}_i^d$ be the average of \mathbf{y}_i^d
 - 2.2 find $\hat{\boldsymbol{\beta}}_i = \arg \min_{\boldsymbol{\beta}} \|P(\boldsymbol{\beta}, \Sigma) - \bar{\mathbf{y}}_i^d\|_2$
 - 2.3 $\hat{\boldsymbol{\beta}}_i \leftarrow \hat{\boldsymbol{\beta}}_i / L_i$
3. $\hat{\boldsymbol{\beta}} = \bar{\hat{\boldsymbol{\beta}}}$

Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $N = 1.000$, 1.000 rep.

J	4	<input type="checkbox"/>	8	<input checked="" type="checkbox"/>
levels	1	<input type="checkbox"/>	N	<input checked="" type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>

Average computation time: 2 seconds



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\begin{pmatrix} U_1 \\ U_2 \\ U_3 \end{pmatrix} = \begin{pmatrix} X & 0 & 0 \\ 0 & X & 0 \\ 0 & 0 & X \end{pmatrix} \begin{pmatrix} \beta_1 \\ \beta_2 \\ \beta_3 \end{pmatrix} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \varepsilon_3 \end{pmatrix}$$

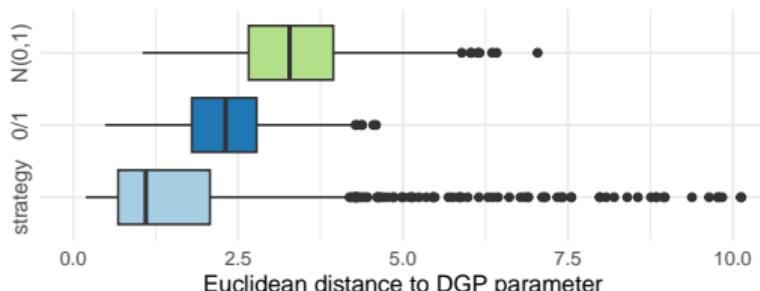
Initialization strategy:

1. assume Σ is known, or choose $\Sigma = I_J$ else
2. localize X , for each level $L_i \neq 0$:
 - 2.1 let $\bar{\mathbf{y}}_i^d$ be the average of \mathbf{y}_i^d
 - 2.2 find $\hat{\boldsymbol{\beta}}_i = \arg \min_{\boldsymbol{\beta}} \|P(\boldsymbol{\beta}, \Sigma) - \bar{\mathbf{y}}_i^d\|_2$
 - 2.3 $\hat{\boldsymbol{\beta}}_i \leftarrow \hat{\boldsymbol{\beta}}_i / L_i$
3. $\hat{\boldsymbol{\beta}} = \bar{\hat{\boldsymbol{\beta}}}$

Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $N = 1.000$, 1.000 rep.

J	4	<input type="checkbox"/>	8	<input checked="" type="checkbox"/>
levels	1	<input type="checkbox"/>	N	<input checked="" type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \dots \\ 0 & 1 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \dots \\ 1 & 3 & \dots \\ \vdots & \ddots & \ddots \end{pmatrix}$	<input checked="" type="checkbox"/>

Average computation time: 2 seconds



$$U = \mu + X(\beta + \gamma) + \varepsilon$$

Parameter	Initialization strategy
β is alternative-varying and	constant utility direction
X is alternative-varying	minimize choice frequency prediction error
X is alternative-constant (ASCs μ is special case)	linear probability model
► β is alternative-constant	MCMC
covariance Σ for ε	MCMC
covariance Ω for γ	

$$\boldsymbol{U} = \boldsymbol{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma})$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_1 \\ X_2 \end{pmatrix} \underbrace{\boldsymbol{\beta}_1}_{P=1} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$\boldsymbol{U} = \boldsymbol{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma})$$

$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_{1,1} & X_{1,2} \\ X_{2,1} & X_{2,2} \end{pmatrix} \underbrace{\begin{pmatrix} \beta_1 \\ \beta_2 \end{pmatrix}}_{P=2} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_{1,1} & X_{1,2} \\ X_{2,1} & X_{2,2} \end{pmatrix} \underbrace{\begin{pmatrix} \beta_1 \\ \beta_2 \end{pmatrix}}_{P=2} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$y = \begin{cases} 1, & \text{if } U_1 = \max \mathbf{U} \\ 2, & \text{if } U_2 = \max \mathbf{U} \end{cases}$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

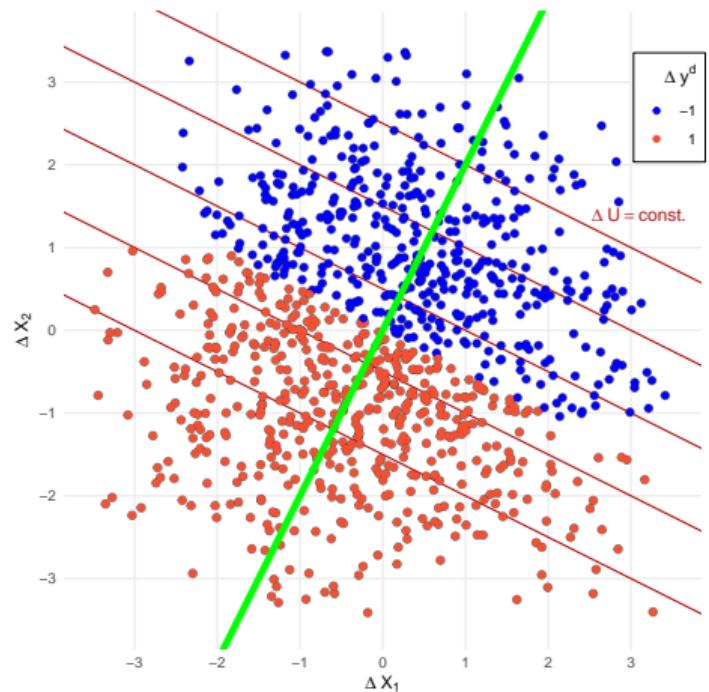
$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_{1,1} & X_{1,2} \\ X_{2,1} & X_{2,2} \end{pmatrix} \underbrace{\begin{pmatrix} \beta_1 \\ \beta_2 \end{pmatrix}}_{P=2} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$\mathbf{y}^d = \begin{cases} \begin{pmatrix} 1 \\ 0 \end{pmatrix}, & \text{if } U_1 = \max \mathbf{U} \\ \begin{pmatrix} 0 \\ 1 \end{pmatrix}, & \text{if } U_2 = \max \mathbf{U} \end{cases}$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_{1,1} & X_{1,2} \\ X_{2,1} & X_{2,2} \end{pmatrix} \underbrace{\begin{pmatrix} \beta_1 \\ \beta_2 \end{pmatrix}}_{P=2} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$\mathbf{y}^d = \begin{cases} \begin{pmatrix} 1 \\ 0 \end{pmatrix}, & \text{if } U_1 = \max \mathbf{U} \\ \begin{pmatrix} 0 \\ 1 \end{pmatrix}, & \text{if } U_2 = \max \mathbf{U} \end{cases}$$



$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma}) \qquad \mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_{1,1} & X_{1,2} \\ X_{2,1} & X_{2,2} \end{pmatrix} \underbrace{\begin{pmatrix} \beta_1 \\ \beta_2 \end{pmatrix}}_{P=2} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$\mathbf{y}^d = \begin{cases} \begin{pmatrix} 1 \\ 0 \end{pmatrix}, & \text{if } U_1 = \max \mathbf{U} \\ \begin{pmatrix} 0 \\ 1 \end{pmatrix}, & \text{if } U_2 = \max \mathbf{U} \end{cases}$$

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma})$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

$$\underbrace{\begin{pmatrix} U_1 \\ U_2 \end{pmatrix}}_{J=2} = \begin{pmatrix} X_{1,1} & X_{1,2} \\ X_{2,1} & X_{2,2} \end{pmatrix} \underbrace{\begin{pmatrix} \beta_1 \\ \beta_2 \end{pmatrix}}_{P=2} + \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \end{pmatrix}$$

$$\mathbf{y}^d = \begin{cases} \begin{pmatrix} 1 \\ 0 \end{pmatrix}, & \text{if } U_1 = \max \mathbf{U} \\ \begin{pmatrix} 0 \\ 1 \end{pmatrix}, & \text{if } U_2 = \max \mathbf{U} \end{cases}$$

- find $\hat{\boldsymbol{\alpha}}$ via OLS
- $\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$
 - 💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)
- robust to $\boldsymbol{\Sigma}$ and the existence of additional independent regressors

Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	■	10.000	□
P	3	■	10	□
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	■	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	□

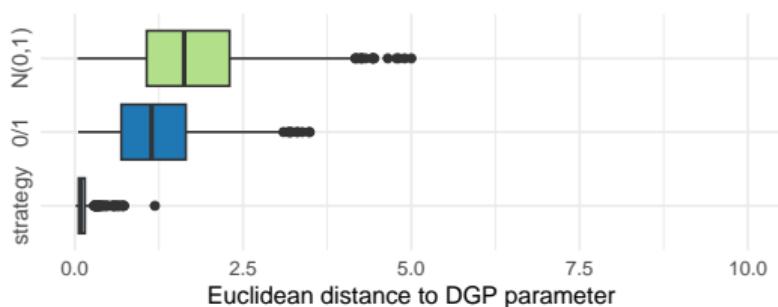
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	■	10.000	□
P	3	■	10	□
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	■

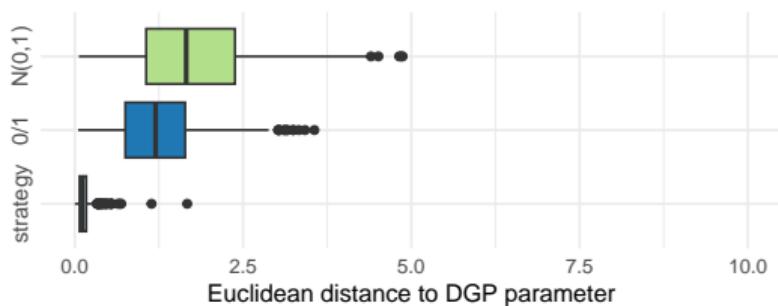
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	<input checked="" type="checkbox"/>	10.000	<input type="checkbox"/>
P	3	<input type="checkbox"/>	10	<input checked="" type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	<input type="checkbox"/>

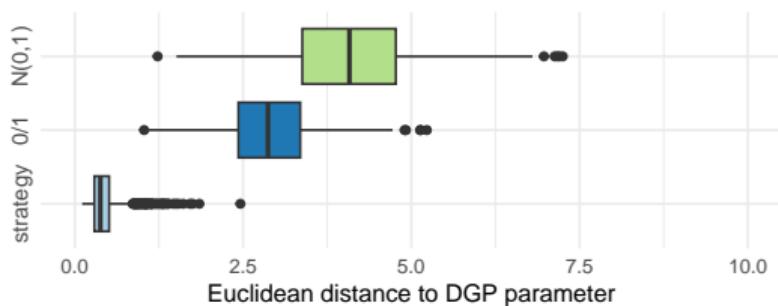
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	■	10.000	□
P	3	□	10	■
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	■

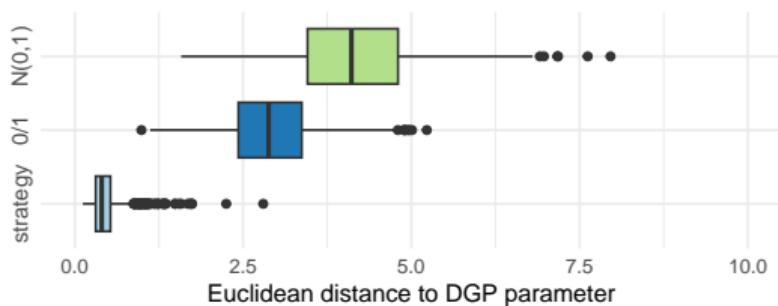
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	<input type="checkbox"/>	10.000	<input checked="" type="checkbox"/>
P	3	<input checked="" type="checkbox"/>	10	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	<input type="checkbox"/>

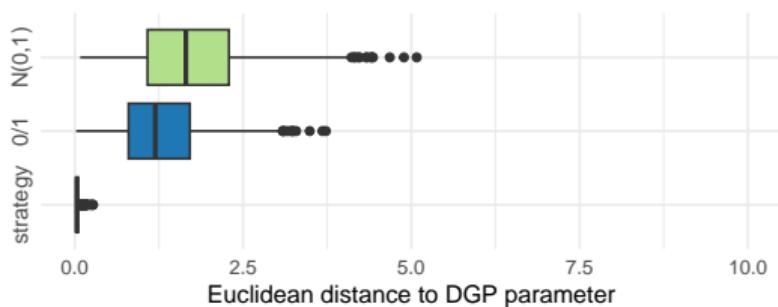
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	<input type="checkbox"/>	10.000	<input checked="" type="checkbox"/>
P	3	<input checked="" type="checkbox"/>	10	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	<input type="checkbox"/>	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	<input checked="" type="checkbox"/>

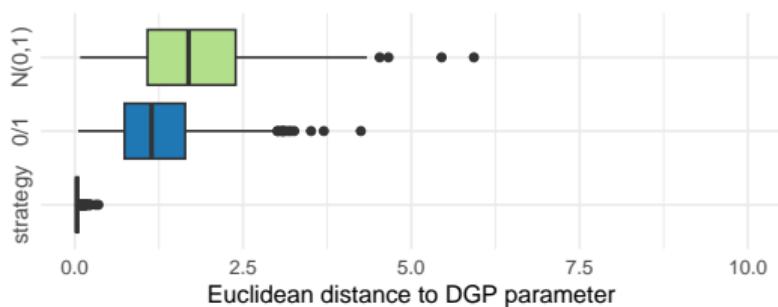
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	<input type="checkbox"/>	10.000	<input checked="" type="checkbox"/>
P	3	<input type="checkbox"/>	10	<input checked="" type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	<input type="checkbox"/>

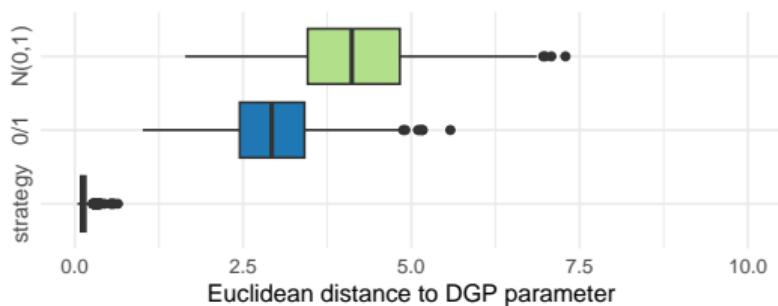
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



Simulation: $\beta_i \sim \mathcal{N}(0, 1)$, $J = 3$, 1.000 replications

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

$$\mathbf{y}^d = \mathbf{X}\boldsymbol{\alpha} + \mathbf{u}, \quad \mathbf{u} \sim \mathcal{N}(0, \mathbf{I})$$

N	1.000	<input type="checkbox"/>	10.000	<input checked="" type="checkbox"/>
P	3	<input type="checkbox"/>	10	<input checked="" type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$	<input type="checkbox"/>	$\begin{pmatrix} 3 & 1 & 1 \\ 1 & 3 & 1 \\ 1 & 1 & 3 \end{pmatrix}$	<input checked="" type="checkbox"/>

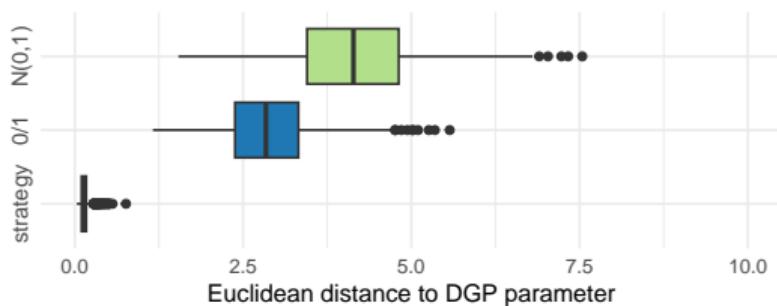
find $\hat{\boldsymbol{\alpha}}$ via OLS

$\hat{\boldsymbol{\alpha}}/\|\hat{\boldsymbol{\alpha}}\|$ is consistent for $\boldsymbol{\beta}/\|\boldsymbol{\beta}\|$

💡 Consistency as $N \rightarrow \infty$ and under a technical assumption on \mathbf{X} (normality is sufficient)

robust to Σ and the existence of additional independent regressors

Average computation time: < 1 second



$$U = \mu + X(\beta + \gamma) + \varepsilon$$

Parameter	Initialization strategy
β is alternative-varying and	
X is alternative-varying	constant utility direction
X is alternative-constant (ASCs μ is special case)	minimize choice frequency prediction error
β is alternative-constant	linear probability model
► covariance Σ for ε	MCMC
covariance Ω for γ	MCMC

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

1. assume $\boldsymbol{\beta}$ is known
2. for ≈ 100 iterations:
 - 2.1 $\mathbf{U} | \Sigma, \boldsymbol{\beta}, \mathbf{X}, y \sim \text{truncated } \mathcal{N}$
 - 2.2 $\Sigma | \boldsymbol{\beta}, \mathbf{U} \sim \mathcal{W}^{-1}$
3. derive $\hat{\Sigma}$ as sample average of Σ

Simulation: $\Sigma \sim \mathcal{W}^{-1}(J \cdot I, J + 1)$, $P = 5$, 1.000 rep.

$$\mathbf{U} = \mathbf{X}\beta + \epsilon, \quad \epsilon \sim \mathcal{N}(0, \Sigma)$$

N	1.000	<input checked="" type="checkbox"/>	10.000	<input type="checkbox"/>
J	4	<input checked="" type="checkbox"/>	8	<input type="checkbox"/>

1. assume β is known

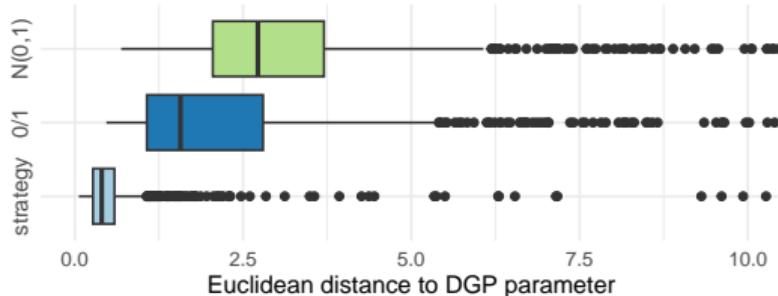
2. for ≈ 100 iterations:

2.1 $\mathbf{U} | \Sigma, \beta, \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\Sigma | \beta, \mathbf{U} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}$ as sample average of Σ

Average computation time: 3 seconds



Simulation: $\Sigma \sim \mathcal{W}^{-1}(J \cdot I, J + 1)$, $P = 5$, 1.000 rep.

$$\mathbf{U} = \mathbf{X}\beta + \epsilon, \quad \epsilon \sim \mathcal{N}(0, \Sigma)$$

N	1.000	■	10.000	□
J	4	□	8	■

1. assume β is known

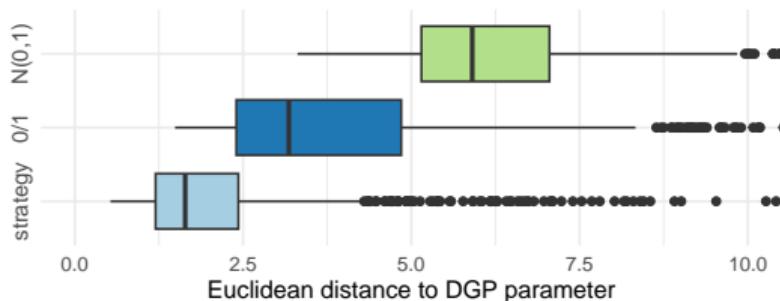
2. for ≈ 100 iterations:

2.1 $\mathbf{U} | \Sigma, \beta, \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\Sigma | \beta, \mathbf{U} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}$ as sample average of Σ

Average computation time: 3 seconds



Simulation: $\Sigma \sim \mathcal{W}^{-1}(J \cdot I, J + 1)$, $P = 5$, 1.000 rep.

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

N	1.000	<input type="checkbox"/>	10.000	<input checked="" type="checkbox"/>
J	4	<input checked="" type="checkbox"/>	8	<input type="checkbox"/>

1. assume $\boldsymbol{\beta}$ is known

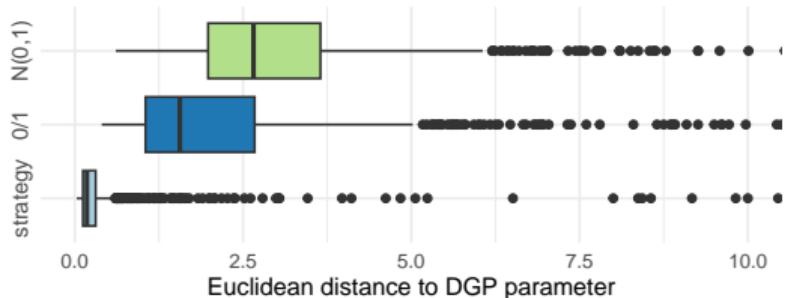
2. for ≈ 100 iterations:

2.1 $\mathbf{U} | \Sigma, \boldsymbol{\beta}, \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\Sigma | \boldsymbol{\beta}, \mathbf{U} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}$ as sample average of Σ

Average computation time: 30 seconds



Simulation: $\Sigma \sim \mathcal{W}^{-1}(J \cdot I, J + 1)$, $P = 5$, 1.000 rep.

$$\mathbf{U} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma)$$

N	1.000	□	10.000	■
J	4	□	8	■

1. assume $\boldsymbol{\beta}$ is known

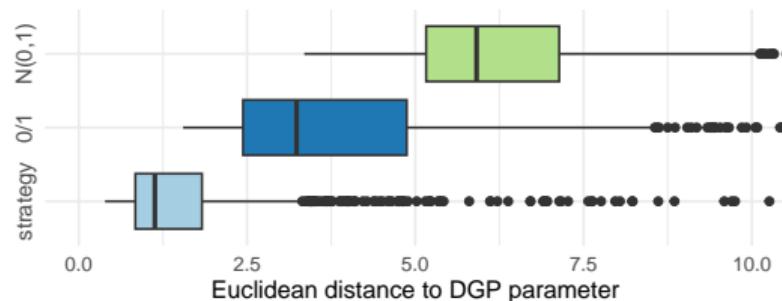
2. for ≈ 100 iterations:

2.1 $\mathbf{U} | \Sigma, \boldsymbol{\beta}, \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\Sigma | \boldsymbol{\beta}, \mathbf{U} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}$ as sample average of Σ

Average computation time: 30 seconds



$$U = \mu + X(\beta + \gamma) + \varepsilon$$

Parameter	Initialization strategy
β is alternative-varying and	
X is alternative-varying	constant utility direction
X is alternative-constant (ASCs μ is special case)	minimize choice frequency prediction error
β is alternative-constant	linear probability model
covariance Σ for ε	MCMC
► covariance Ω for γ	MCMC

$$\begin{aligned} \boldsymbol{U} &= \boldsymbol{X}(\boldsymbol{\beta} + \boldsymbol{\gamma}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma}), \\ \boldsymbol{\gamma} &\sim \mathcal{N}(0, \boldsymbol{\Omega}) \end{aligned}$$

$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma}), \\ \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \boldsymbol{\Omega})$$

$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \\ \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

1. assume $\boldsymbol{\beta}$ is known
2. for ≈ 200 iterations:
 - 2.1 $\mathbf{U} | \Sigma, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$
 - 2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \Omega, \mathbf{U} \sim \mathcal{N}$ for each n
 - 2.3 $\Sigma | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$
 - 2.4 $\Omega | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$
3. derive $\hat{\Sigma}, \hat{\Omega}$ as sample average of Σ, Ω

Simulation: $\Omega \sim \mathcal{W}^{-1}(P \cdot I, P + 1)$, $J = 3$, $T = 10$

$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \quad \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

2.1 $\mathbf{U} | \Sigma, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \Omega, \mathbf{U} \sim \mathcal{N}$ for each n

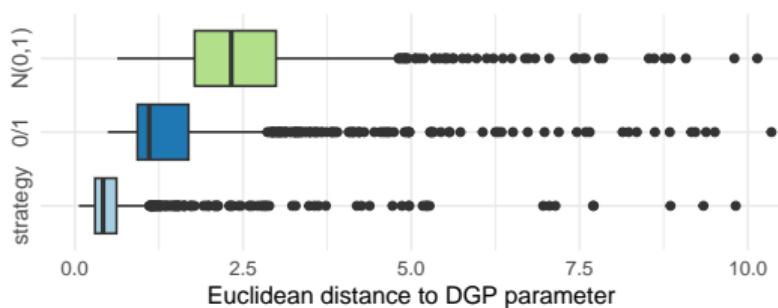
2.3 $\Sigma | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$

2.4 $\Omega | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}, \hat{\Omega}$ as sample average of Σ, Ω

N	100	<input checked="" type="checkbox"/>	500	<input type="checkbox"/>
P	3	<input checked="" type="checkbox"/>	5	<input type="checkbox"/>
Σ	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \end{pmatrix}$	<input checked="" type="checkbox"/>	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \end{pmatrix}$	<input type="checkbox"/>

Average computation time: 3 seconds



Simulation: $\Omega \sim \mathcal{W}^{-1}(P \cdot I, P + 1)$, $J = 3$, $T = 10$

$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \\ \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

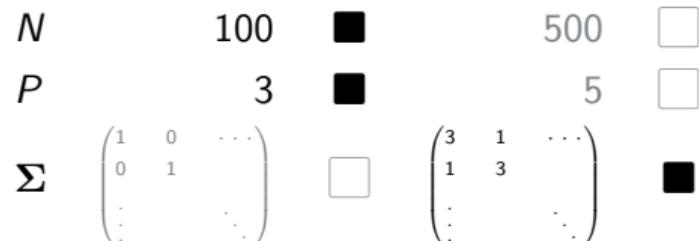
2.1 $\mathbf{U} | \Sigma, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \Omega, \mathbf{U} \sim \mathcal{N}$ for each n

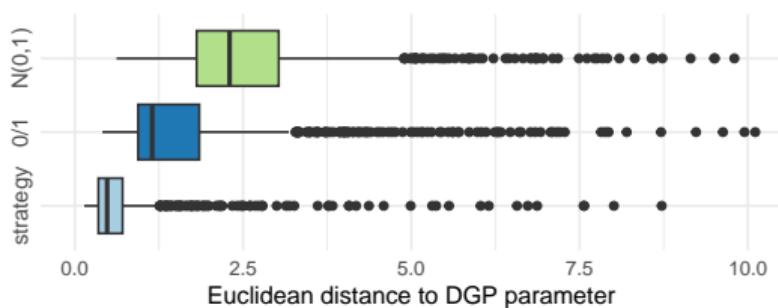
2.3 $\Sigma | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$

2.4 $\Omega | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}, \hat{\Omega}$ as sample average of Σ, Ω



Average computation time: 3 seconds



$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \quad \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

2.1 $\mathbf{U} | \Sigma, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

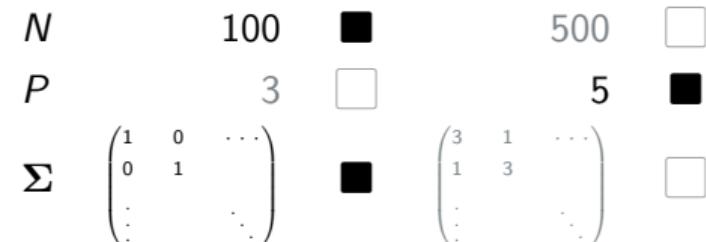
2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \Omega, \mathbf{U} \sim \mathcal{N}$ for each n

2.3 $\Sigma | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$

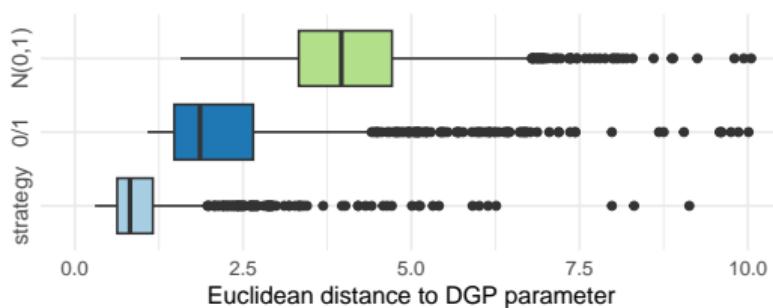
2.4 $\Omega | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}, \hat{\Omega}$ as sample average of Σ, Ω

Simulation: $\Omega \sim \mathcal{W}^{-1}(P \cdot \mathbf{I}, P + 1)$, $J = 3$, $T = 10$



Average computation time: 3 seconds



$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma}), \\ \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \boldsymbol{\Omega})$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

2.1 $\mathbf{U} | \boldsymbol{\Sigma}, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \boldsymbol{\Omega}, \mathbf{U} \sim \mathcal{N}$ for each n

2.3 $\boldsymbol{\Sigma} | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$

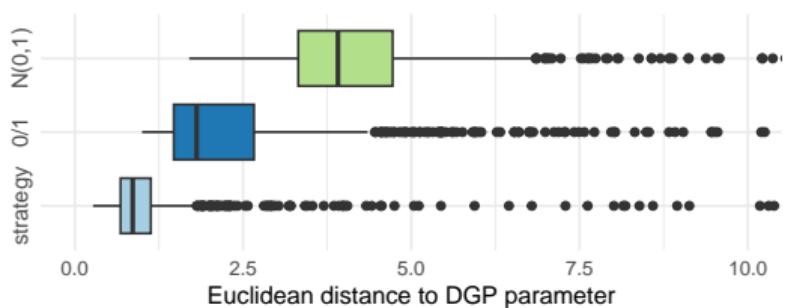
2.4 $\boldsymbol{\Omega} | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$

3. derive $\hat{\boldsymbol{\Sigma}}, \hat{\boldsymbol{\Omega}}$ as sample average of $\boldsymbol{\Sigma}, \boldsymbol{\Omega}$

Simulation: $\boldsymbol{\Omega} \sim \mathcal{W}^{-1}(P \cdot \mathbf{I}, P + 1)$, $J = 3$, $T = 10$

N	100	■	500	□
P	3	□	5	■
$\boldsymbol{\Sigma}$	$\begin{pmatrix} 1 & 0 & \cdots \\ 0 & 1 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	□	$\begin{pmatrix} 3 & 1 & \cdots \\ 1 & 3 & \cdots \\ \vdots & \ddots & \ddots \end{pmatrix}$	■

Average computation time: 3 seconds



$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \quad \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

2.1 $\mathbf{U} | \Sigma, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

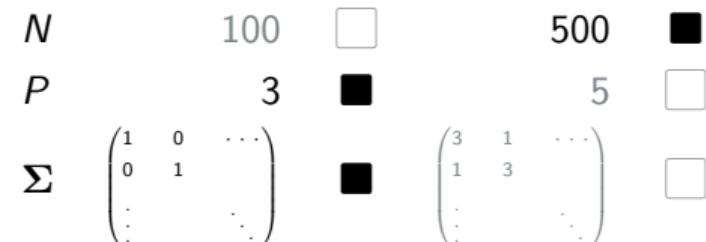
2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \Omega, \mathbf{U} \sim \mathcal{N}$ for each n

2.3 $\Sigma | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$

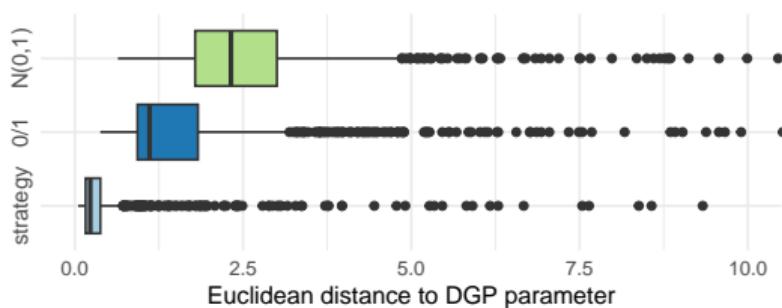
2.4 $\Omega | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$

3. derive $\hat{\Sigma}, \hat{\Omega}$ as sample average of Σ, Ω

Simulation: $\Omega \sim \mathcal{W}^{-1}(P \cdot \mathbf{I}, P + 1)$, $J = 3$, $T = 10$



Average computation time: 12 seconds



$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \boldsymbol{\Sigma}), \\ \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \boldsymbol{\Omega})$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

2.1 $\mathbf{U} | \boldsymbol{\Sigma}, (\boldsymbol{\beta}_n), \mathbf{X}, y \sim \text{truncated } \mathcal{N}$

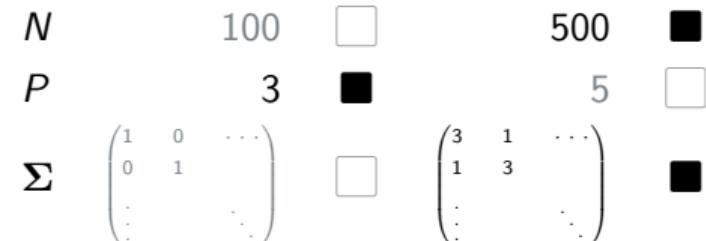
2.2 $\boldsymbol{\beta}_n | \boldsymbol{\beta}, \boldsymbol{\Omega}, \mathbf{U} \sim \mathcal{N}$ for each n

2.3 $\boldsymbol{\Sigma} | (\boldsymbol{\beta}_n), \mathbf{U} \sim \mathcal{W}^{-1}$

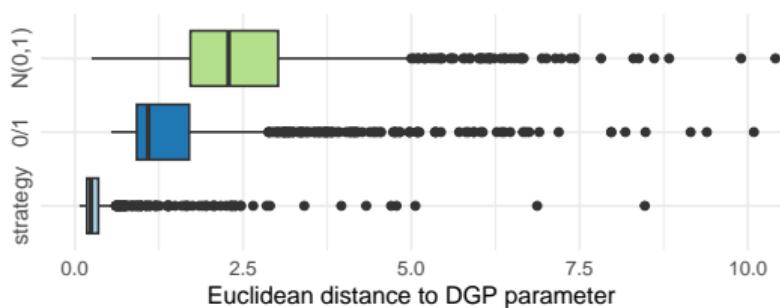
2.4 $\boldsymbol{\Omega} | (\boldsymbol{\beta}_n), \boldsymbol{\beta} \sim \mathcal{W}^{-1}$

3. derive $\hat{\boldsymbol{\Sigma}}, \hat{\boldsymbol{\Omega}}$ as sample average of $\boldsymbol{\Sigma}, \boldsymbol{\Omega}$

Simulation: $\boldsymbol{\Omega} \sim \mathcal{W}^{-1}(P \cdot \mathbf{I}, P + 1)$, $J = 3$, $T = 10$



Average computation time: 12 seconds



$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \quad \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

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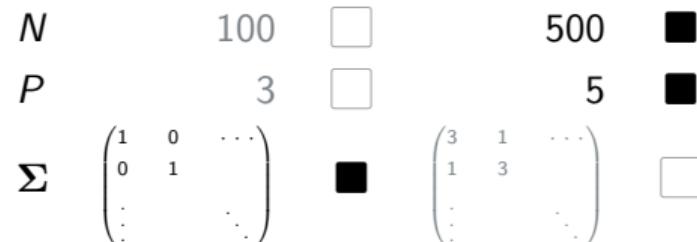
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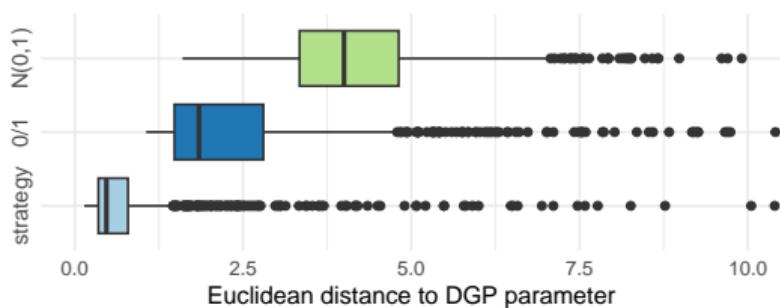
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3. derive $\hat{\Sigma}, \hat{\Omega}$ as sample average of Σ, Ω

Simulation: $\Omega \sim \mathcal{W}^{-1}(P \cdot \mathbf{I}, P + 1)$, $J = 3$, $T = 10$



Average computation time: 12 seconds



Simulation: $\Omega \sim \mathcal{W}^{-1}(P \cdot I, P + 1)$, $J = 3$, $T = 10$

$$\mathbf{U} = \mathbf{X}(\underbrace{\boldsymbol{\beta} + \boldsymbol{\gamma}}_{\boldsymbol{\beta}_n}) + \boldsymbol{\varepsilon}, \quad \boldsymbol{\varepsilon} \sim \mathcal{N}(0, \Sigma), \\ \boldsymbol{\beta}_n \sim \mathcal{N}(\boldsymbol{\beta}, \Omega)$$

1. assume $\boldsymbol{\beta}$ is known

2. for ≈ 200 iterations:

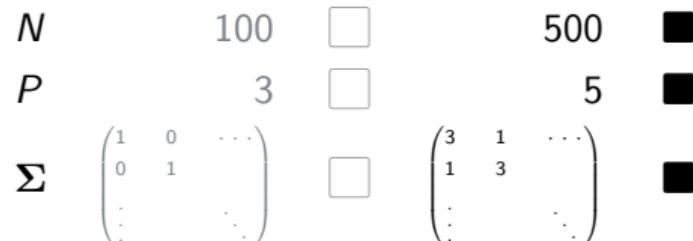
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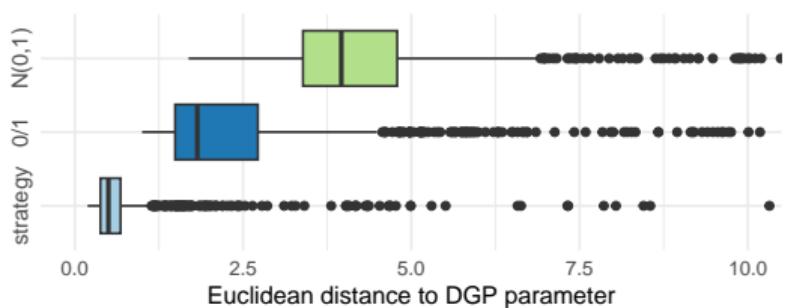
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Average computation time: 12 seconds



Outline

- 1 How slow and unreliable can it be?
- 2 Initialization strategies and proof of concept
- 3 Putting them together
- 4 Let's try with empirical data
- 5 Takeaways

Putting them together

We can combine the strategies in the general “ $\mathbf{U} = \mu + \mathbf{X}(\beta + \gamma) + \varepsilon$ ” case:

1. orthogonalize the regressors using the Frisch–Waugh–Lovell (FWL) theorem
2. get an initial estimate for μ and β :
 - 2.1 apply the presented strategies separately
 - 2.2 estimate a common utility scale via a linear probability model
3. perform MCMC jointly for Ω and Σ

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Empirical application

Stated-choice experiment by Helveston et al. (2015) on car purchase decision in USA:

<u>Attribute*</u>	<u>Option 1</u>	<u>Option 2</u>	<u>Option 3</u>
<u>Vehicle Type</u> ⓘ	Conventional  300 mile range on 1 tank	Plug-In Hybrid  300 mile range on 1 tank (first 40 miles electric)	Electric  75 mile range on full charge
<u>Brand</u> ⓘ	German	American	Japanese
<u>Purchase Price</u> ⓘ	\$18,000	\$32,000	\$24,000
<u>Fast Charging Capability</u> ⓘ	--	Not Available	Available
<u>Operating Cost (Equivalent Gasoline Fuel Efficiency)</u> ⓘ	19 cents per mile (20 MPG equivalent)	12 cents per mile (30 MPG equivalent)	6 cents per mile (60 MPG equivalent)
<u>0 to 60 mph Acceleration Time** ⓘ</u>	8.5 seconds (Medium-Slow)	8.5 seconds (Medium-Slow)	7 seconds (Medium-Fast)
			

Empirical application

Stated-choice experiment by Helveston et al. (2015) on car purchase choice (USA):

- $N = 384$ deciders
- a total of 5760 decisions
- $J = 3$ alternatives
- $P = 16$ alternative-specific regressors
- we assumed $\varepsilon \sim \text{iid } \mathcal{N}(0, \sqrt{1.6})$

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- With random initial values, only 15% of runs converged. This took a total of 3 hours.
- Informed initial values using the presented strategy led to a single run that converged in 2 minutes to the estimates reported by Helveston et al. (2015).

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Outline

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- 5 Takeaways

Takeaways

- RUMs are widely used in discrete choice applications, but MLE quickly becomes computational challenging (with uninformed initialization)
- We can improve optimization time and convergence rate by using consistent and numerically fast initial estimators:
 1. constant utility direction
 2. minimize choice frequency prediction error
 3. linear probability model
 4. MCMC

Takeaways

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- We can improve optimization time and convergence rate by using consistent and numerically fast initial estimators:
 1. constant utility direction
 2. minimize choice frequency prediction error
 3. linear probability model
 4. MCMC

Thanks for your attention! Do you have any questions or comments?

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-  loelschlaeger.de/talks