## Stability Selection: Theorem 10.1 in book

Assume:

- exchangeability condition:  $\{l(j \in \hat{S}_{\lambda}), j \in S_0^c\}$  is exchangeable for all  $\lambda \in \Lambda$
- $\hat{S}$  is not worse than random guessing

$$\frac{\mathbb{E}|S_0 \cap \hat{S}_{\Lambda}|)}{\mathbb{E}(|S_0^c \cap \hat{S}_{\Lambda}|)} \geq \frac{|S_0|}{|S_0^c|}$$

Then, for  $\pi_{\text{thr}} \in (1/2, 1)$ :

$$\mathbb{E}[V] \leq rac{1}{2\pi_{ ext{thr}}-1} rac{q_{\Lambda}^2}{
ho}$$

suppose we know  $q_{\Lambda}$  (see later) strategy: specify  $\mathbb{E}[V] = v_0$  (e.g. = 5)  $\sim$  for  $\pi_{\text{thr}} := \frac{1}{2} + \frac{q_{\Lambda}^2}{2\rho v_0}$ :  $\mathbb{E}[V] \le v_0$  example: regression model with p = 1000 variables

 $\hat{S}_{\lambda}$  = the top 10 variables from Lasso (e.g. the different  $\lambda$  from Lasso by CV and choose the top 10 variables with the largest absolute values of the corresponding estimated coefficients; if less than 10 variables are selected, take the selected variables) the value  $\lambda$  corresponds to the "top 10";  $\Lambda$  is a singleton

we then know that  $q_{\Lambda} = \mathbb{E}[|\hat{S}_{\lambda}(I)|] \leq 10$ 

For  $\mathbb{E}[V] = v_0 := 5$  we then obtain

$$\pi_{\rm thr} = \frac{1}{2} + \frac{q_{\Lambda}^2}{2\rho v_0} = 0.5 + \frac{10^2}{2*1000*5} = 0.51$$

# there is room to play around recommendation: take $|\hat{S}_{\lambda}|$ rather large and stability selection will reduce again to reasonable size

when taking the "top 30", the threshold becomes

$$\pi_{\rm thr} = \frac{1}{2} + \frac{q_{\Lambda}^2}{2\rho v_0} = 0.5 + \frac{30^2}{2*1000*5} = 0.59$$

adding noise... can always add (e.g. independent  $\mathcal{N}(0, 1)$ ) noise covariates enlarged dimension  $p_{\text{enlarged}}$ 

error control becomes better (for the same threshold)

$$\mathbb{E}[V] \leq \frac{1}{2\pi_{\rm thr} - 1} \frac{q_{\Lambda}^2}{p_{\rm enlarged}}$$

this sometimes helps indeed in practice – at the cost of loss in power

The assumptions for mathematical guarantees

not worse than random guessing

$$rac{\mathbb{E}|S_0 \cap \hat{S}_{\Lambda}|)}{\mathbb{E}(|S_0^c \cap \hat{S}_{\Lambda}|)} \ \geq \ rac{|S_0|}{|S_0^c|}$$

perhaps hard to check but very reasonable...

for Lasso in linear models it holds assuming the variable screening property asymptotically: if beta-min and compatibility condition hold

exchangeability condition:  $\{l(j \in \hat{S}_{\lambda}), j \in S_0^c\}$  is exchangeable for all  $\lambda \in \Lambda$ 

a restrictive assumption but the theorem is very general, for any algorithm  $\hat{S}$ 

a very special case where exchangeability condition holds: random equi-correlation design linear model

$$Y = X\beta^0 + \varepsilon$$
,  $\operatorname{Cov}(X)_{i,j} \equiv \rho \ (i \neq j)$ ,  $\operatorname{Var}(X_j) \equiv 1 \forall j$ 

distributions of  $(Y, X^{(S_0)}, \{X^{(j)}; j \in S_0^c\})$  and of  $(Y, X^{(S_0)}, \{X^{(\pi(j))}; j \in S_0^c\})$  are the same for any permutation  $\pi : S_0^c \to S_0^c$ 

- ► distribution of X<sup>(S<sub>0</sub>)</sup>, {X<sup>(π(j))</sup>; j ∈ S<sub>0</sub><sup>c</sup>} is the same for all π (because of equi-correlation)
- ► distribution of Y|X<sup>(S<sub>0</sub>)</sup>, {X<sup>(π(j))</sup>; j ∈ S<sup>c</sup><sub>0</sub>} is the same for all π (because it depends only on X<sup>(S<sub>0</sub>)</sup>)
- therefore: distribution of Y, X<sup>(S<sub>0</sub>)</sup>, {X<sup>(π(j))</sup>; j ∈ S<sub>0</sub><sup>c</sup>} is the same for all π and hence exchangeability condition holds for any (measurable) function Ŝ<sub>λ</sub>

### An illustration for graphical modeling

p = 160 gene expressions, n = 115GLasso estimator, selecting among the  $\binom{p}{2} = 12'720$  features stability selection with  $\mathbb{E}[V] \le v_0 = 30$ 



## with permutation (empty graph is correct)



Stability Selection is extremely easy to use and super-generic

the sufficient assumptions (far from necessary) for mathematical guarantees are restrictive but the method seems to work very well in practice

## P-values based on multi sample splitting (Ch. 11 in Bühlmann and van de Geer (2011))

Stability Selection

- uses subsampling many times a good thing!
- provides control of the expected number of false positives rather than e.g. the familywise error rate ~> we will "address" this with multi sample splitting and aggregation of P-values

familywise error rate (FWER):

 $FWER = \mathbb{P}[V > 0], V$  number of false positives

Fixed design linear model

$$Y = X\beta^0 + \varepsilon$$

instead of de-biased/de-sparsified method, consider the "older" technique (which is not statistically optimal but more generic and more in the spirit of stability selection)

split the sample into two parts  $I_1$  and  $I_2$  of equal size  $\lfloor n/2 \rfloor$ 

- use (e.g.) Lasso to select variables based on  $I_1$ :  $\hat{S}(I_1)$
- perform low-dimensional statistical inference on *l*<sub>2</sub> based on data (*X*<sup>(Ŝ(l<sub>1</sub>))</sup><sub>*l*<sub>2</sub></sub>); for example using the *t*-test for single coefficients β<sup>0</sup><sub>j</sub> (if *j* ∉ Ŝ(*l*<sub>1</sub>), assign the p-value 1 to the hypothesis *H*<sub>0,j</sub> : β<sup>0</sup><sub>j</sub> = 0)

due to independence of  $I_1$  and  $I_2$ , this is a "valid" strategy (see later)

validity of the (single) data splitting procedure consider testing  $H_{0,j}$ :  $\beta_j^0 = 0$  versus  $H_{A,j}$ :  $\beta_j^0 \neq 0$ assume Gaussian errors for the fixed design linear model : thus, use the *t*-test on the second half of the sample  $I_2$  to get a p-value

 $P_{\text{raw},j}$  from *t*-test based on  $X_{l_2}^{(\hat{S}(l_1))}, Y_{l_2}$ 

 $P_{\text{raw},j}$  is a valid p-value (controlling type I error) for testing  $H_{0,j}$  if  $\hat{S}(I_1) \supseteq S_0$  (i.e., the screening property holds)

if the screening property does not hold:  $P_{\text{raw},j}$  is still valid for  $H_{0,j}(M) : \beta_j(M) = 0$  where  $M = \hat{S}(I_1)$  is a selected sub-model and  $\beta(M) = ((X^{(M)})^T X^{(M)})^{-1} (X^{(M)})^T Y$ 

a p-value lottery depending on the random split of the data



→ should aggregate/average over multiple splits!

## Multiple testing and aggregation of p-values

the issue of multiple testing:

$$\tilde{P}_{j} = \begin{cases} P_{\text{raw},j} \text{ based on } Y_{l_{2}}, X_{l_{2}}^{(\hat{S}(l_{1}))} &, \text{if } j \in \hat{S}(l_{1}), \\ 1 &, \text{if } j \notin \hat{S}(l_{1}) \end{cases}$$

thus, we can have at most  $|\hat{S}(I_1)|$  false positives  $\sim$  can correct with Bonferroni with factor  $|\hat{S}(I_1)|$  (instead of factor *p*) to control the familywise error rate

$$\tilde{P}_{\operatorname{corr},j} = \min(\tilde{P}_j \cdot |\hat{S}(I_1)|, 1) \ (j = 1, \dots, p)$$

decision rule: reject  $H_{0,j}$  if and only if  $\tilde{P}_{\text{corr},j} \leq \alpha$  $\rightsquigarrow$  FWER  $\leq \alpha$  the issue with P-value aggregation:

if we run sample splitting B times, we obtain P-values

$$\tilde{P}^{[1]}_{\mathrm{corr},j},\ldots,\tilde{P}^{[B]}_{\mathrm{corr},j}$$

how to aggregate these dependent p-values to a single one? for  $\gamma \in (0, 1)$  define

$$Q_j(\gamma) = \min \left\{ q_{\gamma} \left( \{ \tilde{P}_{\text{corr},j}^{[b]} / \gamma; b = 1, \dots, B \} \right), 1 \right\},$$

where  $q_{\gamma}(\cdot)$  is the (empirical)  $\gamma$ -quantile function

Proposition 11.1 (Bühlmann and van de Geer, 2011) For any  $\gamma \in (0, 1)$ ,  $Q_i(\gamma)$  are P-values which control the FWER

example:  $\gamma=1/2$  aggregate the p-values with the sample median and multiply by the factor 2

avoid choosing  $\gamma$ :

$$P_{j} = \min \left\{ \underbrace{\underbrace{(1 - \log \gamma_{\min})}_{\text{price to optimize over } \gamma} \inf_{\substack{\gamma \in (\gamma_{\min}, 1)}} Q_{j}(\gamma), 1 \right\} (j = 1, \dots, p).$$

Theorem 11.1 (Bühlmann and van de Geer (2011)) For any  $\gamma_{\min} \in (0, 1)$ ,  $P_j$  are P-values which control the FWER

the entire framework for p-value aggregation holds whenever the single p-values are valid ( $\mathbb{P}[P_{\text{raw},j} \leq \alpha] \leq \alpha$  under  $H_{0,j}$ ) has nothing to do with high-dimensional regression and sample splitting









one can also adapt the method to control the False Discovery Rate (FDR)

multi sample splitting and p-value construction:

- ► is very generic, also for "any other" model class
- is powerful in terms of multiple testing correction: we only correct for multiplicity from |Ŝ(I<sub>1</sub>)| variables
- it relies in theory on the screening property of the selector in practice: it is a quite competitive method!
- Schultheiss et al. (2021): can improve multi sample splitting by multi carve methods, based on "technology" from selective inference

# Undirected graphical models

(Ch. 13 in Bühlmann and van de Geer (2011))

- In the graph G: set of vertices/nodes V = {1,...,p} set of edges E ⊆ V × V
- random variables X = X<sup>(1)</sup>,..., X<sup>(p)</sup> with distribution P identify nodes in V with components of X

graphical model: (*G*, *P*)

pairwise Markov property:

P satisfies the pairwise Markov property (w.r.t. G) if

$$(j,k) \notin E \Longrightarrow X^{(j)} \perp X^{(k)} | X^{(V \setminus \{j,k\})}$$

Global Markov property (stronger property than pairwise Markov prop): consider disjoint subsets  $A, B, C \subseteq V$ P satisfies the global Markov property (w.r.t. G) if

A and B are separated by  $C \implies X^{(A)} \perp X^{(B)}$ 

only condition on subset C



global Markov property  $\Longrightarrow$  pairwise Markov property

Proof: consider  $(j, k) \notin E$ denote by  $A = \{j\}, B = \{k\}, C = V \setminus \{j, k\};$ since  $(j, k) \notin E, A = \{j\}$  and  $B = \{k\}$  are separated by *C* by the global Markov property:  $X^{(j)} \perp X^{(k)} | X^{(V \setminus \{j, k\})}$ 

→ global Markov property is more "interesting"

consider graphical model (G, P)

if *P* has a positive and continuous density w.r.t. Lebesgue measure:

the global and pairwise Markov properties (w.r.t. *G*) coincide/are equivalent (Lauritzen, 1996)

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prime example: P is Gaussian
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the Markov properties imply some conditional independencies from graphical separation

for example with pairwise Markov property:

$$(j,k) \notin E \Longrightarrow X^{(j)} \perp X^{(k)} | X^{(V \setminus \{j,k\})}$$

how about reverse relation ?

$$(j,k) \in E \implies X^{(j)} \not\perp X^{(k)} | X^{(V \setminus \{j,k\})}$$

can we interpret existing edges?

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in general: no! (unfortunately)
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in some special cases:

$$(j,k) \in E \implies X^{(j)} \not\perp X^{(k)} | X^{(V \setminus \{j,k\})}$$

prime example: P is Gaussian

$$(j,k) \in E \iff X^{(j)} \not\perp X^{(k)} | X^{(V \setminus \{j,k\})}$$

for A and B not separated by C: in general not true that

$$X^{(A)} \not\perp X^{(B)} | X^{(C)}$$

... due to possible strange cancellations of "edge weights"

## Gaussian "counterexample"



$$\begin{aligned} \boldsymbol{X}^{(1)} &\leftarrow \boldsymbol{\varepsilon}^{(1)}, \\ \boldsymbol{X}^{(2)} &\leftarrow \boldsymbol{\alpha} \boldsymbol{X}^{(1)} + \boldsymbol{\varepsilon}^{(2)}, \\ \boldsymbol{X}^{(3)} &\leftarrow \boldsymbol{\beta} \boldsymbol{X}^{(1)} + \boldsymbol{\gamma} \boldsymbol{X}^{(2)} + \boldsymbol{\varepsilon}^{(3)}, \\ \boldsymbol{\varepsilon}^{(1)}, \boldsymbol{\varepsilon}^{(2)}, \boldsymbol{\varepsilon}^{(3)} \text{ i.i.d. } \mathcal{N}(0, 1) \end{aligned}$$

 $\rightsquigarrow$  a Gaussian distribution *P* for  $\beta + \alpha \gamma = 0$ : Corr(*X*<sub>1</sub>, *X*<sub>3</sub>) = 0 that is: *X*<sup>(1)</sup>  $\perp$  *X*<sup>(3)</sup> it is a Gaussian Graphical Model where P is Markov w.r.t. the following graph



we know that  $X^{(1)} \perp X^{(3)}$  (for special constellations of  $\alpha, \beta, \gamma$ )

take  $A = \{1\}, B = \{3\}, C = \emptyset$ although A and B are not separated (by the emptyset) since there is a direct edge it does not hold that  $X^{(1)} \not\perp X^{(3)}$  (conditional on  $\emptyset$ , i.e., marginal)

# Gaussian Graphical Model

conditional independence graph (CIG): (G, P) satisfies the pairwise Markov property

Gaussian Graphical Model (GGM): a conditional independence graph with *P* being Gaussian for simplicity, assume mean zero:  $P \sim N_p(0, \Sigma)$ 

we know already that edges are equivalent to conditional dependence given all other variables

for a GGM:

$$(j,k)\in E \iff (\Sigma^{-1})_{jk} \neq 0$$

Neighborhood selection: nodewise regression

$$X^{(j)} = \beta_k^{(j)} X^{(k)} + \sum_{r \neq j,k} \beta_r^{(j)} X^{(r)} + \varepsilon^{(j)}, \ j = 1 \dots, p$$
$$X^{(k)} = \beta_j^{(k)} X^{(j)} + \sum_{r \neq k,j} \beta_r^{(k)} X^{(r)} + \varepsilon^{(k)}$$

for GGM:

$$(j,k)\in \pmb{E} \Longleftrightarrow eta_k^{(j)}
eq \pmb{0} \iff eta_j^{(k)}
eq \pmb{0}$$