APES: Approximated Exhaustive Search for GLM

Kevin Y.X. Wang

School of Mathematics and Statistics The University of Sydney

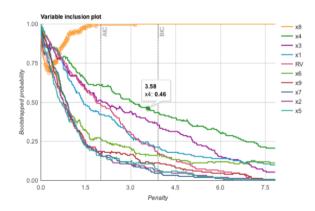
Peking University - Beijing 29 December 2019



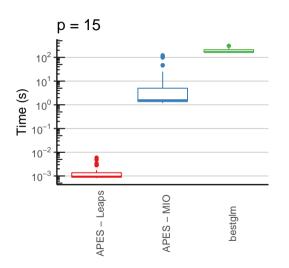


Acknowledgement

- This is joint work with Prof Samuel Müller, Dr Garth Tarr and Prof Jean Yang.
- mplot **Tarr2018** is a package to assess model stability and variable selection for linear models and generalised linear models.



My talk in one slide



- Exhaustive variable selection for GLM is slow.
- APES improves the speed of exhaustive search using approximations.
- APES is now published at Australian and New Zealand Journal of Statistics: https://doi.org/10.1111/ anzs.12276.

Background









Data and models

- $y = (y_1, y_2, \dots, y_n)^{\top}$, with independent $y_i \in \{0, 1\}$, $i = 1, \dots, n$.
- Design matrix $\boldsymbol{X} = (\boldsymbol{x}_1, \dots, \boldsymbol{x}_n) \in \mathbb{R}^{n \times p}$.
- Index the columns of X by $\{1, \ldots, p\}$.
- Let α denote any subset of p_{α} distinct elements from $\{1, \ldots, p\}$. Use \mathcal{A} to denote the collection of all α , so $|\mathcal{A}| = 2^p$.
- \boldsymbol{X}_{α} denote the $n \times p_{\alpha}$ matrix with columns given by the columns of \boldsymbol{X} whose indices appear in α .

Logistic regression

- We model the conditional response variable $Y_i|X$ as $Bernoulli(\pi_i)$, where $\pi_i = \mathbb{P}(Y_i = 1|X)$.
- We will use the **logistic function** as our link function, so $x_i^{\mathsf{T}} \boldsymbol{\beta} = \operatorname{logit}(\pi_i) = \ln{(\pi_i/(1-\pi_i))}.$
- Model fitting usually involves estimating $oldsymbol{eta}$ (or equivalently, $oldsymbol{\pi}$).

Iterative Reweighted Least Square (IRLS)

- 1. Denote weights $w_i = \pi_i(1 \pi_i)$ and other estimates at the t-th iteration with a superscript (t).
- 2. Construct

$$z_i^{(t)} = \underbrace{\log it \left(\pi_i^{(t)}\right)}_{\boldsymbol{x}_i^{\top} \boldsymbol{\beta}^{(t)}} + \frac{y_i - \pi_i^{(t)}}{\pi_i^{(t)} (1 - \pi_i^{(t)})}.$$

3. Update via

$$\widehat{oldsymbol{eta}}^{(t+1)} \leftarrow (oldsymbol{X}^ op oldsymbol{W}^{(t)} oldsymbol{X})^{-1} oldsymbol{X}^ op oldsymbol{W}^{(t)} oldsymbol{z}^{(t)},$$

with $\mathbf{W}^{(t)} = \operatorname{diag}\left(w_i^{(t)}\right)$.

4. At convergence, $\hat{\boldsymbol{\beta}} = (\boldsymbol{X}^{\top} \boldsymbol{W} \boldsymbol{X})^{-1} \boldsymbol{X}^{\top} \boldsymbol{W} \boldsymbol{z}$, which equals to the MLE of logistic regression model.

Challenges in exhaustive GLM variable selection

For large p, exhaustive variable selection in GLM is difficult:

1. The computational cost of IRLS is $\mathcal{O}(np^2)$ per iteration.



Challenges in exhaustive GLM variable selection

For large p, exhaustive variable selection in GLM is difficult:

- 1. The computational cost of IRLS is $\mathcal{O}(np^2)$ per iteration.
- 2. We need to explore all 2^p models.



Aim of APES

Can we perform linear exhaustive variable selection (which benefits from fast algorithms) and use the results to approximate exhaustive GLM variable selection?

GLM exhaustive Can this be done? LM exhaustive

1.	Turning	GLM	to	LM

Exhaustively computing modified MLE

- Hosmer1989 described an approximation to exhaustive variable selection for logistic regression without the need for numerical optimisation.
- Their method starts with the estimated probability $\widehat{\pi}(\alpha_f)$, from the **full** logistic model.
- Then, for each model $\alpha \in \mathcal{A}$, we calculate:

$$\widehat{\boldsymbol{\beta}}(\boldsymbol{\alpha};\widehat{\boldsymbol{\pi}}(\alpha_f)) = (\boldsymbol{X}_{\boldsymbol{\alpha}}^{\top}\boldsymbol{W}(\widehat{\boldsymbol{\pi}}(\alpha_f))\boldsymbol{X}_{\boldsymbol{\alpha}})^{-1}\boldsymbol{X}_{\boldsymbol{\alpha}}^{\top}\boldsymbol{W}(\widehat{\boldsymbol{\pi}}(\alpha_f))\boldsymbol{z}(\widehat{\boldsymbol{\pi}}(\alpha_f)).$$

• This is **NOT** the MLE for α , which should be $\widehat{\beta}(\alpha; \widehat{\pi}(\alpha))$.

Variable selection using the modified estimator

- Given $\widehat{\beta}(\alpha; \widehat{\pi}(\alpha_f))$, we could **approximate** RSS or BIC for all $\alpha \in A$.
- Upon selection of a small set of desired models, we can recompute the MLE and calculate other model fit statistics.

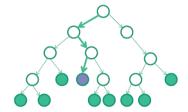
GLM exhaustive
Hosmer approx.

LM exhaustive

2. Reducing computational time

Best subsets search

- Application of this approximation method is limited by the number of LMs we can explore.
- For $p\approx 50$, ${\cal A}$ is approximately 1 quadrillion in size, which is too large to explore exhaustively.
- A <u>best subset</u> algorithm limits our search to a subset of A but it guarantees to contain the global RSS-optimal model.
- leaps (Furnival1974; Lumley2017) discard "branches" of models with insufficient fit.



Mixed Integer Optimisation

- Bertsimas2016 showed that it is feasible to perform best subset search for linear models with p in the hundreds.
- The most attractive component: guaranteed sub-optimality if algorithm is terminated before full convergence. Thus allowing a upper bound for real time limit.
- Current implementation in R is bestsubset, (Hastie2017), which outputs the RSS-best linear model for each model size.

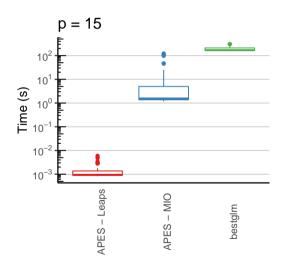
APES: Approximated \underline{E} xhaustive \underline{S} earch



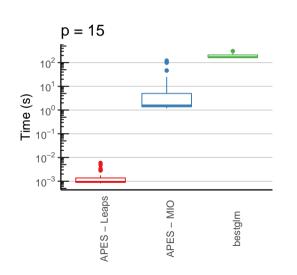
APES: Approximated \underline{E} xhaustive \underline{S} earch

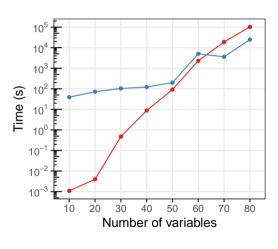


How fast is APES compare to genuine exhaustive search? Very.



How fast is APES compare to genuine exhaustive search? Very.





APES - leaps → APES - MIO

Simulation

Simulation set-up

- n = 500, p = 100, number of non-zero coefficient is k = 6.
- Intercept term is set to 0, then we tried 3 different choices of β :
 - 1. Equally spaced indices:

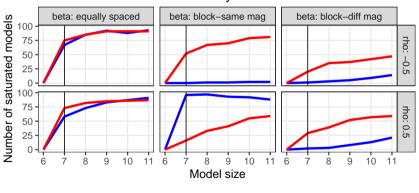
$$(\frac{1}{2},0,\ldots,0,\frac{1}{2},0,\ldots,0,\frac{1}{2},0,\ldots,0,\frac{1}{2},0,\ldots,0,\frac{1}{2},0,\ldots,\frac{1}{2}).$$

- 2. Block of indices, same magnitude/sign: $(\frac{1}{2}, \frac{1}{2}, \frac{1}{2}, \frac{1}{2}, \frac{1}{2}, \frac{1}{2}, \frac{1}{2}, \frac{1}{2}, 0, \dots, 0)$. 3. Block of indices, different magnitude/sign: $(\frac{1}{3}, -1, 1, \frac{2}{3}, -\frac{1}{3}, 0, \dots, 0)$.
- Generating $X \sim \mathcal{N}(0, \Sigma)$, with $\Sigma_{ij} = \rho^{|i-j|}$, $\rho = 0.5$ or -0.5. Then standardise.
- We repeated the simulation 100 times and compared APES against de-biased Lasso using various evaluation metrics.

Evaluation 1: saturated models

- For each model size, APES and Lasso outputs one model.
- In most cases, APES has less false exclusion of variables than Lasso.

Number of saturated models by each methods

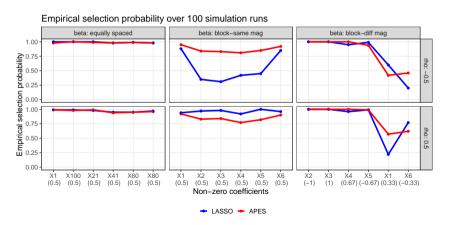


Lasso -

APES

Evaluation 2: variable selection

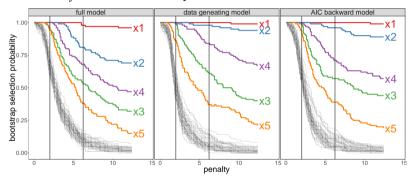
- BIC was used to select an optimal model.
- In most cases, APES has higher selection probability of active variables than Lasso.



Some extensions

The choice of $\widehat{\pi}(\alpha_f)$ can be relaxed in two different ways:

• The full model α_f is not necessarily the best for variable selection.



 We could replace the MLE by other estimators, e.g. the Lasso or robust quasi-likelihood estimator.

Final remarks

- 1. APES is a **fast approximation** method for **exhaustive** variable selection in GLM.
- 2. APES pushes model dimensions into the hundreds and serves as a standard of comparison like a true exhaustive search.
- APES is now published at ANZJS: https://doi.org/10.1111/anzs.12276.
 - https://github.com/kevinwang09/APES
 - https://github.com/garthtarr/mplot
 - Email: kevin.wang@sydney.edu.au.
 - Twitter: @KevinWang009

References