fast cube implementation

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Abstract

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1 Introduction

2 Notation

Consider a finite population U of size N whose units can be defined by labels $k \in \{1, 2, ..., N\}$. Let $S = \{s | s \subset U\}$ be the set of all possible samples. A sampling design is defined by a probability distribution p(.) on S such that

$$p(s) \ge 0$$
 for all $s \in \mathcal{S}$ and $\sum_{s \in \mathcal{S}} p(s) = 1$.

A random sample S is a random vector that maps elements of S to an N vector of 0 or 1 such that P(S = s) = p(s). Define $a_k(S)$, for k = 1, ..., N:

$$a_k = \begin{cases} 1 & \text{if } k \in S \\ 0 & \text{otherwise.} \end{cases}$$

Then a sample can be denoted by means of a vector notation: $\mathbf{a}^{\top} = (a_1, a_2, \dots, a_N)$. For each unit of the population, the inclusion probability $0 \le \pi_k \le 1$ is defined as the probability that unit k is selected into sample S:

$$\pi_k = P(k \in S) = E(a_k) = \sum_{s \in S | k \in S} p(s), \text{ for all } k \in U.$$

Let $\boldsymbol{\pi}^{\top} = (\pi_1, \dots, \pi_N)$ be the vector of inclusion probabilities. Then, $E(\mathbf{a}) = \boldsymbol{\pi}$. Let also $\pi_{k\ell}$ be the probability of selecting the units k and ℓ together in the sample, with $\pi_{kk} = \pi_k$. The matrix of second-order inclusion probabilities is given by $\Pi = E(\mathbf{a}\mathbf{a}^{\top})$. The sample is generally selected with the aim of estimating some population parameters. Let y_k denote a real number associated with unit $k \in U$, usually called the variable of interest. For example, the total

$$Y = \sum_{k \in U} y_k$$

can be estimated by using the classical Horvitz-Thompson estimator of the total defined by

$$\widehat{Y}_{HT} = \sum_{k \in U} \frac{y_k a_k}{\pi_k}.$$
 (1)

3 Balanced Sampling

Usually, some auxiliary information $\mathbf{x}_k^{\top} = (x_{k1}, x_{k2}, \dots, x_{kq}) \in \mathbb{R}^q$ regarding the population units is available. A sampling design is said to be balanced on the

auxiliary variables x_k if and only if it satisfies the balancing equations

$$\widehat{\mathbf{X}} = \sum_{k \in S} \frac{\mathbf{x}_k}{\pi_k} = \sum_{k \in U} \frac{\mathbf{x}_k a_k}{\pi_k} = \sum_{k \in U} \mathbf{x}_k = \mathbf{X}.$$

Sometimes it is not possible to select a sample that satisfies exactly the constraint. We write $\hat{\mathbf{X}} \approx \mathbf{X}$ to notice that the sample is approximately balanced. In many applications, inclusion probabilities are such that samples have a fixed size n. A sampling design of fixed size can be viewed as balanced on only one auxiliary variable $x_k = \pi_k$. Indeed, we have mathematically,

$$\sum_{k \in S} \frac{x_k}{\pi_k} = \sum_{k \in S} \frac{\pi_k}{\pi_k} = n_S.$$

Let denote the set of all samples that have fixed size equal to n by

$$S_n = \left\{ \mathbf{a} \in \{0, 1\}^N \mid \sum_{k=1}^N a_k = n \right\}.$$

More generally, we write the problem of selecting a balanced sample by the following linear system:

$$\begin{cases} \sum_{k \in U} \frac{\mathbf{x}_k a_k}{\pi_k} = \sum_{k \in U} \frac{\mathbf{x}_k}{\pi_k} \pi_k \\ a_k \in \{0, 1\}, \ k \in U. \end{cases}$$

Or also written in matrix form,

$$\mathbf{A}\mathbf{a} = \mathbf{A}\boldsymbol{\pi},\tag{2}$$

where $\mathbf{A} = \left(\frac{\mathbf{x}_1}{\pi_1}, \dots, \frac{\mathbf{x}_N}{\pi_N}\right)$. The aim consist then of obtaining a sample \mathbf{a} that satisfies the constraints.

4 Cube Method

The cube method select a sample that is balanced and respect the inclusion probabilities. The method can take equal or unequal inclusion probabilities. A each step, vector $\boldsymbol{\pi}$ is randomly modified. The subspace induced by the linear system (2) is defined by the following,

$$A = \{ \mathbf{a} \in \mathbb{R}^N | \mathbf{A}\mathbf{a} = \mathbf{A}\boldsymbol{\pi} \}$$

= $\boldsymbol{\pi} + \text{Null}(\mathbf{A}),$

where $\text{Null}(\mathbf{A}) = \{u \in \mathbb{R}^N | \mathbf{A}\mathbf{u} = 0\}$. The idea is then to use a vector of the null space of \mathbf{A} such that we ensure to have martingale property of the updated inclusion probabilities. More specifically we have the following equation,

$$E_p(\boldsymbol{\pi}^t | \boldsymbol{\pi}^{t-1}) = E_p(\boldsymbol{\pi}^{t-1}), \text{ for all } t = 1, ..., N.$$

At each step, at least one component is set to 0 or 1. Matrix A is updated from the new inclusion probabilities. This step is repeated until there is only one component that is not equal to 0 or 1. Algorithm 1 present the full picture of the method.

Algorithm 1 flight phase of the cube Method

Initializing π^0 by π , For t = 1, ..., N, we repeat the three next steps:

- 1. Find $\mathbf{u}^t \neq 0$ such that u^t is inside of the null space of \mathbf{Z}^{\top} and $u_k^t = 0$ if π_k^{t-1} is equal to 0 or 1.
- 2. Calculate $\widetilde{\lambda_1^t}$ and $\widetilde{\lambda_2^t}$ the two greater value such that

$$0 \leqslant \pi_k^{t-1} + \lambda_1^t u_k^t \leqslant 1,$$

$$0 \leqslant \pi_k^{t-1} - \lambda_2^t u_k^t \leqslant 1,$$
 for all $k \in U$

Observe that λ_1^t and λ_2^t are both greater than 0.

3. Update the inclusion probabilities using the rules:

$$oldsymbol{\pi}^t = \left\{ egin{array}{ll} oldsymbol{\pi}^{t-1} + \widetilde{\lambda_1^t} \mathbf{u}^t & ext{with probability} & q_1^t \ oldsymbol{\pi}^{t-1} - \widetilde{\lambda_2^t} \mathbf{u}^t & ext{with probability} & q_2^t \end{array}
ight.$$

where
$$q_1^t = \widetilde{\lambda_2^t}/(\widetilde{\lambda_1^t} + \widetilde{\lambda_2^t})$$
 and $q_2^t = \widetilde{\lambda_1^t}/(\widetilde{\lambda_1^t} + \widetilde{\lambda_2^t})$

We repeat these 3 steps until it is no more possible to find a vector \mathbf{u} that is inside of the null space.

- 4.1
- 5 Reduction
- 6 Simulation
- 7 Discussion