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by

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SELECTING PROCEDURES FOR OPTIMAL SUBSET OF REGRESSION VARIABLES*

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Recently, a number of methods have been developed for selecting the "best" or at least a "good" subset of variables in regression analysis. For various reasons, we may be interested in including only a subset, say, of size $r < p$, the number of independent variables. Various authors have considered this problem and a variety of techniques are presently being used to construct such subsets.

Arvesen and McCabe (1975) proposed a procedure for selecting a subset within a class of subsets with t (fixed) independent variables, taking into account the statistical variation of the residual mean squares. Huang and Panchapakesan (1982) proposed a selection procedure based on the expected residual sums of squares. Hsu and Huang (1982) studied a sequential selection procedure for good regression models.

In this paper, we are interested in deriving an optimal decision procedure based on residual mean squares to select a subset excluding all "inferior" independent variables. This kind of optimality criterion is related to the approach of Gupta and Huang (1977).

Let $\pi_0, \pi_1, \ldots, \pi_k$ denote k+1 normal populations with unknown variances $\sigma_0^2, \sigma_1^2, \ldots, \sigma_k^2$ respectively. Assume that σ_0^2 is known. A population (model) is said to be superior (or good) if $\sigma_i^2 < \Delta\, \sigma_0^2$, to be inferior (or bad) if

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$\sigma_i^2 \geq \Delta\, \sigma_0^2$, where Δ is a specified constant greater than 1. Let Ω be the parameter space which is the collection of all possible parameters.

Let CD stand for a correct decision which is defined to be the selection of any subset which excludes all the inferior populations.

Assuming the following model

$$(1) \qquad \underset{\sim}{Y} = X\underset{\sim}{\beta} + \underset{\sim}{\epsilon}$$

where $X = [\underset{\sim}{1}, \underset{\sim}{X}_1, \ldots, \underset{\sim}{X}_{p-1}]$ is an nxp known matrix of rank $p \leq n$, $\underset{\sim}{\beta}' = (\beta_0, \beta_1, \ldots, \beta_{p-1})$ is a 1xp parameter vector, and $\underset{\sim}{\epsilon} \sim N(\underset{\sim}{0}, \sigma_0^2 I_n)$, and $\underset{\sim}{1}' = [1, \ldots, 1]_{1xN}$, I_n is an identity matrix with nxn.

In what follows, (1) which has p-1 independent variables, will be viewed as the true model. Without loss of generality we can assume that $\sigma_0^2 = 1$.

Consider the models for any r, $2 \leq r \leq p-1$,

$$(2) \qquad \underset{\sim}{Y} = X_{ri}\, \underset{\sim}{\beta}_{ri} + \underset{\sim}{\epsilon}_{ri}$$

where X_{ri} is an nxr matrix of rank r with $X_{11}' = [1, \ldots, 1]_{1xn}$, $\underset{\sim}{\beta}_{ri}$ is a rx1 parameter vector, and $\underset{\sim}{\epsilon}_{ri} \sim N(\underset{\sim}{0}, \sigma_{ri}^2 I_n)$, $i = 1, 2, \ldots, k_r = \binom{p-1}{r-1}$. Let $k = \sum_{r=2}^{p-1} k_r$. It should be noted that in stating the reduced model (2), our comparisons of models are made under the true model assumptions. The goal is to include all the designs X_{ri} (or sets of independent variables) associated with $\sigma_{[j]}^2$, $j = 1, \ldots, k-t$, where $\sigma_{[1]}^2 \leq \sigma_{[2]}^2 \leq \cdots \leq \sigma_{[k-t]}^2$ are ordered values from some of σ_{ri}'s, $i = 1, \ldots, k_r$, $r = 2, \ldots, p-1$.

Note that for any r, $2 \leq r \leq p-1$, if

$$SS_{ri} = \underset{\sim}{Y}'\{I - X_{ri}(X_{ri}'X_{ri})^{-1}X_{ri}'\}\underset{\sim}{Y} = \underset{\sim}{Y}'Q_i\underset{\sim}{Y},$$

then

$$SS_{ri} \sim \chi^2(\nu_r, (X\underset{\sim}{\beta})'Q_{ri}(X\underset{\sim}{\beta})/2)$$

(under the true model), where $\nu_r = n-r$, for $1 \leq i \leq k_r$. Note that the noncentrality parameter, in general, is not zero, and that

$$\sigma^2_{ri} = 1+(X\underset{\sim}{\beta})'Q_{ri}(X\underset{\sim}{\beta})/\nu_r.$$

Now we need some notation. Deleting a set of β_i's without specifying which ones are deleted, we use ri to denote the special subset that is not deleted. For example, if p = 3, r = 2 then there are three subsets with size 2; namely, $\{\beta_1,\beta_2\}$, $\{\beta_1,\beta_3\}$ and $\{\beta_2,\beta_3\}$. Then r1 denotes the set $\{\beta_1,\beta_2\}$, r2 denotes $\{\beta_2,\beta_3\}$ and r3 denotes $\{\beta_1,\beta_3\}$. Then, we use $\tilde{\beta}$ to denote the vector with the following subsets: $\{\beta_1,\beta_2\}$ with $(\beta_1,\beta_2,0)$, $\{\beta_1,\beta_3\}$ with $(\beta_1,0,\beta_3)$, and $\{\beta_2,\beta_3\}$ with $(0,\beta_2,\beta_3)$, where 0 is the parameter value which is omitted from the true model of the appropriate β_i's. Thus, in the following, we will use $\Omega_{0,ri}$ to denote those sets of $\underset{\sim}{\tilde{\beta}}$ as described above with the further condition that $\sigma^2_{ri} = \sigma^2_0 = 1$. Similarly, $\Omega_{1,ri}$ will be used to denote the sets of $\underset{\sim}{\tilde{\beta}}$ as described above with the further restriction that $\sigma^2_{ri} \geq \Delta$. Formally, we write

$$\Omega_{0,ri} = \{\underset{\sim}{\tilde{\beta}} | \sigma^2_{ri} = 1\},$$

and

$$\Omega_{1,ri} = \{\underset{\sim}{\tilde{\beta}} | \sigma^2_{ri} \geq \Delta\},$$

where $i = 1,\ldots,k_r$; $r = 2,\ldots,p-1$, and let

$$\Omega_1 = \bigcup_{r=2}^{p-1} \bigcup_{i=1}^{k_r} \Omega_{1,ri}, \text{ and}$$

$$\Omega_0 = \bigcap_{r=2}^{p-1} \bigcap_{i=1}^{k_r} \Omega_{0,ri}.$$

Let $g_{\sigma^2_{ri}}(s_{ri})$ denote the probability density of S_{ri} depending on the parameter σ^2_{ri}, where $S_{ri} = \frac{SS_{ri}}{\nu_r}$, $i = 1,\ldots,k_r$; $r = 2,\ldots,p-1$.

Consider a family of hypotheses testing problems as follows:

(3) $$H_{0,ri}:\ \tilde{\underline{\beta}} \in \Omega_0 \quad \text{vs} \quad K_{ri}:\ \tilde{\underline{\beta}} \in \Omega_{1,ri};$$

$i = 1,\ldots,p-1$, $r = 2,\ldots,p-1$. A test of the hypotheses (3) will be defined to be a vector $(\varphi_1(\underline{y}),\ldots,\varphi_k(\underline{y}))$, where the elements of the vector are ordinary test functions; when $\underline{y}$ is observed we reject $H_{0,t}$ with probability $\varphi_t(\underline{y})$, $1 \le t \le k$. The power function of a test $(\varphi_1,\ldots,\varphi_k)$ is defined to be the vector $(p_1(\tilde{\underline{\beta}}),\ldots,p_k(\tilde{\underline{\beta}}))$ where

$$p_t(\tilde{\underline{\beta}}) = E_{\tilde{\underline{\beta}}}\ \varphi_t(\underline{Y}),$$

$1 \le t \le k$. Let $S(\gamma)$ be the set of all tests $(\varphi_1,\ldots,\varphi_k)$ such that

(4) $$E_{\tilde{\underline{\beta}}}\ \varphi_t(\underline{Y}) \le \gamma, \quad \tilde{\underline{\beta}} \in \Omega_0.$$

We define $\varphi^0 = (\varphi^0_1,\ldots,\varphi^0_k)$ as

$$\varphi^0_{ri}(\underline{y}) = \begin{cases} 1, & \text{if } g_\Delta(s_{ri}) \ge c\ g_1(s_{ri}), \\ 0, & \text{if } g_\Delta(s_{ri}) < c\ g_1(s_{ri}), \end{cases}$$

such that $E_{\tilde{\underline{\beta}}}\ \varphi^0_{ri}(\underline{Y}) = \gamma$, $\tilde{\underline{\beta}} \in \Omega_0$, where s_{ri} is the observed value of S_{ri}. It can be shown that φ^0 maximizes

$$\min_{1 \le t \le k} \inf_{\tilde{\beta} \in \Omega_{1,t}} E_{\tilde{\beta}} \varphi_t(\underline{Y})$$

among all tests $\varphi = (\varphi_1, \ldots, \varphi_k) \in S(\gamma)$ (cf. Gupta and Huang (1977)).

To determine the constant c, we proceed follows: for a given $\eta > 0$, there exists a smallest positive integer k_0 such that

$$\frac{a_{k_0}}{\eta} < 1 \quad \text{and} \quad \frac{a_{k_0+1}}{a_{k_0}} + \frac{a_{k_0}}{\eta} \le 1,$$

where

$$a_\ell(s_{ri}) = \frac{e^{-\lambda_r} \lambda_r^\ell}{\ell!} \left[\frac{\nu_r s_{ri}}{2}\right]^\ell \frac{\Gamma(\frac{1}{2}\nu_r)}{\Gamma(\frac{1}{2}\nu_r + \ell)},$$

$\ell = 0,1,2,\ldots;$ $\lambda_r = \frac{(\Delta-1)\nu_r}{2}$. For this k_0, it can be shown that

$$0 < \frac{g_\Delta(s_{ri})}{g_1(s_{ri})} - \sum_{\ell=0}^{k_0-1} a_\ell(s_{ri}) = \sum_{k=0}^{\infty} a_{k_0+k} \le \eta,$$

where

$$\frac{g_\Delta(s_{ri})}{g_1(s_{ri})} = \sum_{\ell=0}^{\infty} a_\ell .$$

Thus, approximately,

$$\frac{g_\Delta(s_{ri})}{g_1(s_{ri})} \approx \sum_{\ell=0}^{k-1} a_\ell(s_{ri})$$

with error less than η. For $\tilde{\beta} \in \Omega_0$,

$$E_{\tilde{\beta}}\, \varphi^0(\underline{Y}) = P_{\tilde{\beta}}\{g_\Delta(S_{ri}) \geq c\, g_1(S_{ri})\}$$

$$= P_{\tilde{\beta}}\{\sum_{\ell=0}^{k_0-1} a_\ell(S_{ri}) \geq c\}$$

$$= \int_0^\infty I_{[\sum_{\ell=0}^{k_o-1} a_\ell(s_{ri}) \geq c]}(s_{ri}) g_1(s_{ri}) ds_{ri} = \gamma,$$

where $g_1(s_{ri})$ is the central χ^2 with ν_r degrees of freedom and $I_A(x) = 0$ for $x \notin A$, $I_A(x) = 1$ for $x \in A$. The constant c can be determined.

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Recently, a number of methods have been developed for selecting the "best" or at least a "good" subset of variables in regression analysis. For various reasons, we may be interested in including only a subset, say, of size $r < p$, the number of independent variables. Various authors have considered this problem and a variety of techniques are presently being used to construct such subsets. In this paper, we are interested in deriving an optimal decision procedure based on residual mean squares to select a subset excluding all "inferior" independent variables. This kind of optimality criterion is related to the approach of Gupta and Huang (1977).

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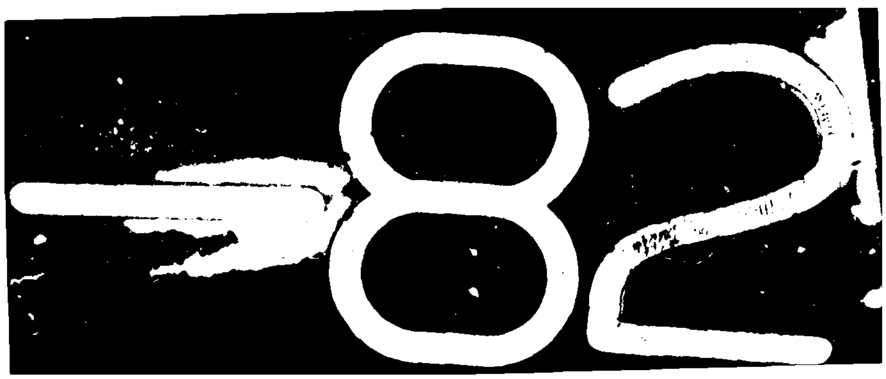
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