## A Bayesian Approach to Model Selection

- In exploratory regression problems, we often must select which subset of our potential predictor variables produces the "best model."
- A Bayesian may consider the possible models and compare them based on their posterior probabilities.
- Note that if the value of coefficient β<sub>j</sub> is 0, then variable X<sub>j</sub> is not needed in the model.
- Let  $\beta_j = z_j b_j$  for each j, where  $z_j = 0$  or 1 and  $b_j \in (-\infty, \infty)$ .
- Then our model is

$$Y_i = z_0 b_0 + z_1 b_1 X_{i1} + z_2 b_2 X_{i2} + \dots + z_{k-1} b_{k-1} X_{i,k-1} + \epsilon_i, \ i = 1, \dots, n$$

where any  $z_j = 0$  indicates that this predictor variable does not belong in the model.

**Example**: Oxygen uptake example:

 $X_1 = \text{group}, X_2 = \text{age}, X_3 = \text{group} \times \text{age}$ :

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- For each possible value of the vector z, we calculate the posterior probability for that model:
- For any particular z\*, say:

$$\pi(\mathbf{z}^*|\mathbf{y}, \mathbf{X}) = \frac{p(\mathbf{z}^*)p(\mathbf{y}|\mathbf{X}, \mathbf{z}^*)}{\sum_{\mathbf{z}} p(\mathbf{z})p(\mathbf{y}|\mathbf{X}, \mathbf{z})}$$

- ► This involves a prior p(·) on each possible model a noninformative approach would be to let all these prior probabilities be equal.
- If there are a large number of potential predictors, we would use a method called **Gibbs sampling**) (more on this later) to search over the many models.

## Example of Bayesian Model Selection

- Example in R with Oxygen Data Set
- We can consider all possible subsets of set of predictor variables:

We can consider only certain subsets (here, we only consider including the interaction term when both first-order terms appear):

- Suppose we have built our Bayesian regression model using response data y and explanatory data matrix X.
- Suppose we consider future observations whose explanatory variable values are in the matrix X\*.
- What is the marginal distribution of the corresponding future response values y\*?
- This is the posterior predictive distribution

 $\pi(\mathbf{y}^*|\mathbf{y}, \mathbf{X}^*, \mathbf{X}).$ 

We will use this later as a tool for checking the fit of our regression model.

## The Posterior Predictive Distribution of the Data

► In our analysis with the noninformative priors, note that  $\pi(\mathbf{y}^*, \boldsymbol{\beta}, \sigma^2 | \mathbf{y}, \mathbf{X}^*, \mathbf{X}) = \pi(\mathbf{y}^* | \boldsymbol{\beta}, \sigma^2, \mathbf{X}^*) \pi(\boldsymbol{\beta}, \sigma^2 | \mathbf{X}, \mathbf{y})$ 

Then integrating out β and σ<sup>2</sup>, it can be shown that the posterior predictive distribution of y\* is multivariate-t with (n - k) degrees of freedom so that

$$E(\mathbf{y}^*) = \mathbf{X}^* \hat{\mathbf{b}} \text{ and}$$
  
covariance matrix  $= \frac{(n-k)\hat{\sigma}^2}{n-k-2} [\mathbf{I} + \mathbf{X}^* (\mathbf{X}' \mathbf{X})^{-1} \mathbf{X}^{*'}]$ 

- Intuition: Our original data are multivariate normal, given the model.
- Our future predictions are multivariate-t (reflects added uncertainty about the model).