Model Selection for Geostatistical Models

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Muddy Creek- tributary to Sun River in Central Montana

Muddy Creek: surveyed every 15.24 meters, total of 5456m; 358 measurements

Degree | AIC<sub>c</sub>
--- | ---
0 | 1455
1 | 294.3
2 | 251.3
3 | 47.1
4 | 34.0
5 | 35.5
Muddy Creek: residuals from poly(d=4) fit

Minimum AIC\textsubscript{c} ARMA model: ARMA(1,1)

\[ Y_t = 0.574 \, Y_{t-1} + \varepsilon_t - 0.311 \, \varepsilon_{t-1}, \{\varepsilon_t\} \sim \text{WN}(0,0.0564) \]

Some theory:

- ARMA(1,1) model:
  \[ Y_t = 1.743 \, Y_{t-1} + \varepsilon_t - 0.311 \, \varepsilon_{t-1} \]

LS estimates of trend parameters are **asymptotically efficient**.

- LS estimates are **asymptotically indep** of cov parameter estimates.
Summary of models fitted to Muddy Creek bed elevation:

<table>
<thead>
<tr>
<th>Degree</th>
<th>$\text{AIC}_c$</th>
<th>ARMA</th>
<th>$\text{AIC}_c$</th>
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</thead>
<tbody>
<tr>
<td>0</td>
<td>1455</td>
<td>(1,2)</td>
<td>59.67</td>
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<td>1</td>
<td>294.3</td>
<td>(2,1)</td>
<td>26.98</td>
</tr>
<tr>
<td>2</td>
<td>251.3</td>
<td>(2,1)</td>
<td>26.30</td>
</tr>
<tr>
<td>3</td>
<td>47.1</td>
<td>(1,1)</td>
<td>7.12</td>
</tr>
<tr>
<td>4</td>
<td>34.0</td>
<td>(1,1)</td>
<td>2.78</td>
</tr>
<tr>
<td>5</td>
<td>35.5</td>
<td>(1,1)</td>
<td>4.68</td>
</tr>
</tbody>
</table>
Muddy Creek (cont)

Simulated series: polynomial degree 4 + ARMA(1,1):
Recap of Traditional Model Fitting

Suppose \( \{Y_t\} \) follows the linear model with time series (or spatial) errors given by

\[
Y_t = \beta_0 + X_{t1} \beta_1 + \ldots + X_{tp} \beta_p + W_t,
\]

where \( \{W_t\} \) is a stationary (ARMA) time series.

- Estimate \((\beta_0, \ldots, \beta_p)\) by ordinary least squares (OLS).
- Identify and estimate ARMA parameters using the estimated residuals,

\[
Y_t - (\hat{\beta}_0 + X_{t1}\hat{\beta}_1 + \cdots + X_{tp}\hat{\beta}_p)
\]

- Re-estimate \((\beta_0, \ldots, \beta_p)\) and ARMA parameters using full MLE.

Limitations of this approach for model selection:

- Ignore potential confounding between explanatory variables and correlation in \( \{W_t\} \).
- Ignoring autocorrelation function can mask importance of explanatory variables.
PROGRAM

➢ Motivation

➢ Introduction

➢ Model Selection for Geostatistical Models
  • Basic setup

➢ AIC for Geostatistical Models
  • Basics

➢ Minimum Description Length (MDL) for Geostatistical Models

➢ Simulation Results

➢ Prediction

➢ Application to Orange-Throated Whiptail Lizard
Introduction

Model selection setup: Suppose \( \{Z(s), s \in D\} \) is a random field where \( D \) is some subset of \( \mathbb{R}^2 \).

Spatial data: observations \( Z(s_1), \ldots, Z(s_n) \) at locations \( s_1, \ldots, s_n \).

Explanatory variables (covariates): \( p \) explanatory variables \( X_1(s), \ldots, X_p(s) \) are available at each location \( s \).

Linear model for \( Z \):

\[
Z(s) = \beta_0 + X_1(s) \beta_1 + \ldots + X_p(s) \beta_p + \delta(s)
\]

where \( \delta(s) \) is a mean 0 stationary (isotropic) Gaussian random field.

Model selection issue.

- Which explanatory variables should be included in the model?
- What family of covariance functions should be used to model \( \delta(s) \)?
Model Selection

Problem: How does one choose the best set of covariates and family of covariance functions?

Some Objectives of Model Selection.

1. Choose the correct order model (*consistency*).
   - There exists a true model and the model selection procedure will choose the correct set of covariates and the right family of covariance functions as sample size increases.

2. Choose the model that performs best for prediction (*efficiency*).
   - Find the model that predicts (or interpolates) well at unobserved locations.

3. Choose the model that maximizes data compression.
   - Find a model that summarizes the data in the most compact fashion, yet retains the salient features present in the data.
The Geostatistical Model

Model:

\[ Z(s) = \beta_0 + X_1(s) \beta_1 + \ldots + X_p(s) \beta_p + \delta(s) \]

- \( X(s) = (1, X_1(s), \ldots, X_p(s))^T \) is a vector of explanatory variables observed at location \( s \).
- \( \beta = (\beta_0, \beta_1, \ldots, \beta_p)^T \) is a \((p + 1)\) dimensional parameter vector.
- \( \delta(s) \) is the unobserved regression error at location \( s \).

Assumptions on \( \delta(s) \).

- \( \delta(s) \) is a stationary, isotropic Gaussian process with mean 0 and covariance function

\[ \text{Cov}(\delta(s), \delta(t)) = \sigma^2 \rho(||s-t||, \theta). \]

- \( \sigma^2 \) is the variance of the process
- \( ||s-t|| \) is the euclidean distance between locations \( s \) and \( t \).
- \( \rho(\cdot, \theta) \) is an isotropic correlation function parameterized by a \( k\)-dim’l vector \( \theta \).
Some standard autocorrelation functions (these are a bit limiting).

1. Exponential

\[ \rho(d; \theta) = \exp(-d / \theta_1), \quad d > 0 \]

2. Gaussian

\[ \rho(d; \theta) = \exp(-d^2 / \theta_1^2), \quad d > 0 \]

3. Matern

\[ \rho(d; \theta) = \frac{1}{2^{\theta_2 - 1} \Gamma(\theta_2)} \left( \frac{2d \sqrt{\theta_2}}{\theta_1} \right)^{\theta_2} K_{\theta_2} \left( \frac{2d \sqrt{\theta_2}}{\theta_1} \right), \quad d > 0, \]

where \( K_\theta(.) \) is the modified Bessel function.

- \( \theta_1 \) is the range parameter controlling the rate of decay of correlations.
- \( \theta_2 \) is the smoothness parameter controlling the smoothness of the random field.
Autocorrelation functions (cont)

Exponential $\theta_1 = 1.0$

Gaussian $\theta_1 = 4.0$

Matern $\theta_1 = 1.5, \theta_2 = .75$

Matern $\theta_1 = 2.5, \theta_2 = 1.0$

Matern $\theta_1 = 3.5, \theta_2 = 2.0$
**Estimation**

Log-likelihood: For the data \( Z = (Z(s_1), \ldots, Z(s_n))^T \)

\[
\log L_Z(\beta, \theta, \sigma^2) = -0.5 \log |\sigma^2 \Omega| - 0.5 \sigma^{-2} (Z - X\beta)^T \Omega^{-1} (Z - X\beta)^T, 
\]

where \( \Omega = \rho(||s_i - s_j||; \theta) \) is the matrix of correlations for the data vector \( Z \).

MLE estimate: \((\hat{\beta}, \hat{\theta}, \hat{\sigma}^2)\) maximizes the log-likelihood.

Note.

- **MLE estimates** can be difficult to compute for large sample sizes.
- **Restricted maximum likelihood** (REML) estimates often have more desirable sampling properties, but performance for model selection is not clear.
AIC for Geostatistical Models

Background on AIC (Burnham and Anderson (1998) and McQuarrie and Tsai (1998).):

Suppose

- \( Z \sim f_T \)
- \( \{ f(\cdot;\psi), \psi \in \Psi \} \) is a family of candidate probability density functions.

The Kullback-Leibler information between \( f(\cdot;\psi) \) and \( f_T \) is

\[
I(\psi) = \int -2 \log \left( \frac{f(z|\psi)}{f_T(z)} \right) f_T(z) dz
\]

has the interpretation as a measure of the

- distance between \( f(\cdot;\psi) \) and \( f_T \)
- loss of information when \( f(\cdot;\psi) \) is used as the model instead of \( f_T \)
AIC for Geostatistical Models

By Jensen’s inequality,

\[ I(\psi) \geq -2 \log \int \left( \frac{f(z \mid \psi)}{f_T(z)} \right) f_T(z) dz = 0 \]

with equality holding if and only if \( f(z \mid \psi) = f_T(z) \), a.e.

Basic idea: minimize the Kullback-Leibler index

\[ \Delta(\psi) = \int -2 \log(f(z \mid \psi)) f_T(z) dz \]
\[ = E_T(-2 \log L_Z(\psi)), \]

where \( L_Z(\psi) \) is the likelihood based on the data \( Z \).

Problem: Cannot compute \( \Delta(\psi) \) or \( \Delta(\hat{\psi}) \), where \( \hat{\psi} \) is the MLE of \( \psi \).

Strategy: Search for an unbiased estimator of

\[ E_\psi(\Delta(\hat{\psi})) \]

and find the candidate model which minimizes this statistic as a function of the model.
AIC for Geostatistical Models

Note: The expectation above can be written as the double expectation,

\[ E_\psi (\Delta(\hat{\psi})) = E_\psi (E_T (-2 \log L_Y (\hat{\psi}) \mid Z)) = E_\psi (-2 \log L_Y (\hat{\psi})), \]

where \( Y \) is independent of \( Z \) with the same distribution. Thus \(-2 \log L_Y (\hat{\psi})\) is an unbiased estimate of \( \Delta(\hat{\psi}) \). But \( Y \) is unobserved and that is where the AIC correction factor comes into play.

Applied to the Geostatistical model: Parameter vector \( \psi = (\beta, \theta, \sigma^2)^T \) and

\[-2 \log L_Y (\hat{\beta}, \hat{\theta}, \hat{\sigma}^2) = -2 \log L_Z (\hat{\beta}, \hat{\theta}, \hat{\sigma}^2) + \hat{\sigma}^{-2} S_Y (\hat{\beta}, \hat{\theta}) - n, \]

where

\[ S_Y (\beta, \theta) = (Y - X\beta)^T \Omega^{-1} (\theta)(Y - X\beta) \]

and

\[ \hat{\sigma}^2 = (Y - X\hat{\beta})^T \Omega^{-1} (\hat{\theta})(Y - X\hat{\beta}) / n. \]

So just need to approximate

\[ E_\psi (\hat{\sigma}^{-2} S_Y (\hat{\beta}, \hat{\theta})). \]
AIC for Geostatistical Models

In order to compute $E_\psi (\hat{\sigma}^{-2} S_Y (\hat{\beta}, \hat{\theta}))$, we assume \textit{standard asymptotics} hold:

- $(\hat{\beta}, \hat{\theta})$ is $\text{AN}((\beta, \theta)^T, I_n^{-1})$, where $I_n$ is the Fisher information.

- For large $n$, $I_n$ can be approximated by

$$V(\beta, \theta) = 1/2\sigma^{-2} E_\psi \left( \frac{\partial^2 S_Y}{\partial(\beta, \theta)^T \partial(\beta, \theta)^T} (\beta, \theta) \right).$$

- For $n$ large, $n\hat{\sigma}^2 = S_Z (\hat{\beta}, \hat{\theta})$ is approximately distributed as $\sigma^2 \chi^2 (n-p-1-k)$.

- $\hat{\sigma}^2$ is approximately independent of $(\hat{\beta}, \hat{\theta})$.

Then

$$E_\psi (\hat{\sigma}^{-2} S_Y (\hat{\beta}, \hat{\theta})) - n \sim \sigma^2 (n + 1 + p + k) E_\psi \hat{\sigma}^{-2} - n$$

$$\sim \sigma^2 (n + 1 + p + k) \left(\sigma^2 \frac{n - p - 1 - k - 2}{n}\right)^{-1} - n$$

$$= \frac{2(p + 2 + k)n}{n - p - k - 3}.$$
AIC for Geostatistical Models

The quantity, referred to as the corrected AIC, is then

\[ AIC_c = -2 \log L_Z(\hat{\beta}, \hat{\theta}, \hat{\sigma}^2) + \frac{2(p + 2 + k)n}{n - p - k - 3}. \]

The standard AIC statistic is given by

\[ AIC = -2 \log L_Z(\hat{\beta}, \hat{\theta}, \hat{\sigma}^2) + 2(p + 2 + k). \]
Model Selection Using Minimum Description Length

Basics of MDL:
Choose the model which maximizes the compression of the data or, equivalently, select the model that minimizes the code length of the data (i.e., amount of memory required to encode the data).

\[ M = \text{class of operating models for } y = (y_1, \ldots, y_n) \]

\[ L_F(y) = \text{code length of } y \text{ relative to } F \in M \]

Typically, this term can be decomposed into two pieces (two-part code),

\[ L_F(y) = L(\hat{F}|y) + L(\hat{\epsilon} | \hat{F}) , \]

where

\[ L(\hat{F}|y) = \text{code length of the fitted model for } F \]

\[ L(\hat{\epsilon} | \hat{F}) = \text{code length of the residuals based on the fitted model} \]
Encoding the data: \((x_1, y_1), \ldots, (x_n, y_n)\)

1. "Naïve" case

\[
L("naive") = L(x_1, \ldots, x_n) + L(y_1, \ldots, y_n) = L(x_1) + \cdots + L(x_n) + L(y_1) + \cdots + L(y_n)
\]

2. Linear model; suppose \(y_i = a_0 + a_1 x_i, \ i = 1, \ldots, n\). Then

\[
L("p = 1") = L(x_1, \ldots, x_n) + L(a_0, a_1) = L(x_1) + \cdots + L(x_n) + L(a_0) + L(a_1)
\]

3. Linear model with noise; suppose \(y_i = a_0 + a_1 x_i + \varepsilon_i, \ i = 1, \ldots, n\), where \(\{\varepsilon_i\} \sim \text{IID } N(0, \sigma^2)\). Then

\[
L("p = 1") = L(x_1) + \cdots + L(x_n) + L(\hat{a}_0) + L(\hat{a}_1) + L(\hat{\sigma}^2) + L(\hat{\varepsilon}_1, \ldots, \hat{\varepsilon}_n | \hat{a}_0, \hat{a}_1, \hat{\sigma}^2)
\]

If \(A < L(y_1) + \ldots + L(y_n)\), then "p=1" encoding scheme dominates the "naïve" scheme.
Applied to the Geostatistical model:

\[ Z(s) = \beta_0 + X_1(s) \beta_1 + \ldots + X_p(s) \beta_p + \delta(s) \]

First term \( L(\hat{F}|y) \):

\[ L(\hat{F}|y) = L(\hat{\beta}_0, \ldots, \hat{\beta}_p) + L(\hat{\theta}_1, \ldots, \hat{\theta}_k) \]

Encoding:

- integer \( I \): \( \log_2 I \) bits (if \( I \) unbounded)
- \( \log_2 I_U \) bits (if \( I \) bounded by \( I_U \))

- MLE \( \hat{\theta} \): \( \frac{1}{2} \log_2 N \) bits (where \( N = \) number of observations used to compute \( \hat{\theta} \); Rissanen (1989))
Model Selection Using Minimum Description Length (cont)

So,

\[
L(\hat{F}|y) = \sum_{j=0}^{p} \frac{1}{2} \log_2 n + \sum_{j=1}^{k} \frac{1}{2} \log_2 n = \frac{1}{2} (p + 1 + k) \log_2 n
\]

Second term \( L(\hat{e} | \hat{F}) \): Using Shannon’s classical results on information theory, Rissanen demonstrates that the code length of \( \hat{e} \) can be approximated by the negative of the log-likelihood of the fitted model, i.e., by

\[
L(\hat{e} | \hat{F}) \approx -\log \hat{L}_Z(\hat{\beta}, \hat{\theta}, \hat{\sigma}^2)
\]

MDL for the model \( F \) is then

\[
MDL = \frac{1}{2} (p + 1 + k) \log_2 n - \log_2 \hat{L}_Z(\hat{\beta}, \hat{\theta}, \hat{\sigma}^2)
\]

The strategy is to find the model that minimizes \( MDL \).

Note:

\[
MDL = \frac{1}{2} \left( -2 \log_2 \hat{L}_Z(\hat{\beta}, \hat{\theta}, \hat{\sigma}^2) + (p + 1 + k) \log_2 n \right)
\]

Penalty coefficient \((\log_2 n)\) is larger than that \((2)\) for AIC.
Model Selection and Spatial Correlation

Traditional approach to model selection:

1. Select explanatory variable to model the large scale variation. Here one might use $\text{AIC}_c$ assuming independence in the noise term.

2. Estimate correlation function parameters using residuals from model in previous step.

3. Re-estimate regression parameters using GLS.

4. Iterate steps 2 and 3 until convergence.

Limitations of this approach:

- Ignore potential confounding between explanatory variables and correlation in spatial process.

- Ignoring autocorrelation function can mask importance of explanatory variables.
Goal of simulation study: compare model selection performance of AIC versus traditional method.

1. **Sampling design:** 100 sites randomly located on \([0,10] \times [0,10]\).

2. **Explanatory variables:** Five potential explanatory variables, 
   \(X_1(s_j), X_2(s_j), \ldots, X_5(s_j), \; j =1, \ldots, 100, \; \text{IID sqrt}(12/10)\cdot t_{12}\)

3. **Response variables:**
   \[Z = 2 + 0.75 \; X_1 + 0.50 \; X_2 + 0.25 \; X_3 + \delta,
   \]
   where \(\delta\) is Gaussian with mean 0, \(\sigma^2 = 50\), and Matern autocorrelation function with parameters \(\theta_1 = 4\) and \(\theta_2 = 1\).

4. **Replicates:** 500 replicates were simulated with a new Gaussian random field generated for each replicate.

5. **AIC:** Computed for \(2^5 = 32\) possible models per replicate.
Model Selection: Random Pattern Sampling Design
Model Selection: Simulation Results for the Random Pattern

- Independent AIC and Spatial AIC report the percentage of simulations that each model was selected.

- Of the 32 possible models, the results given here include only those with 10% or more support for one of the models.

<table>
<thead>
<tr>
<th>Variables in Model</th>
<th>Spatial AIC</th>
<th>Independent AIC</th>
<th>MDL</th>
</tr>
</thead>
<tbody>
<tr>
<td>$X_1, X_2, X_3$</td>
<td>56.0</td>
<td>2.4</td>
<td>40.4</td>
</tr>
<tr>
<td>$X_1, X_2, X_3, X_5$</td>
<td>14.4</td>
<td>0.2</td>
<td>4.2</td>
</tr>
<tr>
<td>$X_1, X_2, X_3, X_4$</td>
<td>10.8</td>
<td>0.2</td>
<td>0.8</td>
</tr>
<tr>
<td>$X_1, X_2$</td>
<td>10.2</td>
<td>8.4</td>
<td>46.4</td>
</tr>
<tr>
<td>Intercept only</td>
<td>0.0</td>
<td>26.8</td>
<td>0.0</td>
</tr>
<tr>
<td>$X_1$</td>
<td>0.4</td>
<td>14.2</td>
<td>1.2</td>
</tr>
<tr>
<td>$X_2$</td>
<td>0.0</td>
<td>13.8</td>
<td>0.2</td>
</tr>
</tbody>
</table>
Independent Model AIC values
Spatial Model AIC values
Sampling Patterns

- Highly Clustered
- Lightly Clustered
- Random Pattern
- Regular Pattern
- Grid Design
# Effect of Sampling Design

<table>
<thead>
<tr>
<th>Variables in Model</th>
<th>Highly Clustered</th>
<th>Lightly Clustered</th>
<th>Random</th>
<th>Regular Pattern</th>
<th>Grid Design</th>
</tr>
</thead>
<tbody>
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<td>73</td>
<td>65</td>
<td>46</td>
<td>43</td>
<td>16</td>
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<td>$X_1, X_2$</td>
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<td>2</td>
<td>18</td>
<td>21</td>
<td>35</td>
</tr>
<tr>
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<td>13</td>
<td>8</td>
<td>8</td>
<td>3</td>
</tr>
<tr>
<td>$X_1, X_2, X_3, X_4, X_5$</td>
<td>10</td>
<td>13</td>
<td>11</td>
<td>7</td>
<td>7</td>
</tr>
</tbody>
</table>

- Each column reports the percentage of simulations that each model was selected.
- Of the 32 possible models, the results given here include only those with 10% or more support for one of the models.
### Effect of Smoothness Parameter

<table>
<thead>
<tr>
<th>Variables in Model</th>
<th>$\theta_2 = 0.50$</th>
<th>$\theta_2 = 0.75$</th>
<th>$\theta_2 = 1.00$</th>
<th>$\theta_2 = 4.00$</th>
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<td></td>
<td>spat</td>
<td>ind</td>
<td>spat</td>
<td>ind</td>
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<td>3</td>
<td>35</td>
<td>1</td>
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<td>0</td>
<td>3</td>
<td>0</td>
</tr>
<tr>
<td>$X_1, X_2, X_3, X_4$</td>
<td>3</td>
<td>0</td>
<td>7</td>
<td>0</td>
</tr>
<tr>
<td>$X_1, X_2$</td>
<td>22</td>
<td>7</td>
<td>20</td>
<td>4</td>
</tr>
<tr>
<td>Intercept only</td>
<td>4</td>
<td>30</td>
<td>0</td>
<td>32</td>
</tr>
<tr>
<td>$X_1$</td>
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<td>17</td>
<td>12</td>
<td>20</td>
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<tr>
<td>$X_2$</td>
<td>4</td>
<td>12</td>
<td>0</td>
<td>11</td>
</tr>
</tbody>
</table>

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Effect of smoothness (cont)

Matern $\theta_1=4.0$, $\theta_2=.5$
Matern $\theta_1=4.0$, $\theta_2=.75$
Matern $\theta_1=4.0$, $\theta_2=1.0$
Matern $\theta_1=4.0$, $\theta_2=4.0$
Effect of Range

<table>
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<th>Variables in Model</th>
<th>$\theta_1 = 2$</th>
<th>$\theta_1 = 4$</th>
<th>$\theta_1 = 6$</th>
<th>$\theta_1 = 8$</th>
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</thead>
<tbody>
<tr>
<td></td>
<td>spat</td>
<td>ind</td>
<td>spat</td>
<td>ind</td>
</tr>
<tr>
<td>$X_1, X_2, X_3$</td>
<td>15</td>
<td>2</td>
<td>56</td>
<td>2</td>
</tr>
<tr>
<td>$X_1, X_2, X_3, X_5$</td>
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<td>0</td>
<td>14</td>
<td>0</td>
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<tr>
<td>$X_1, X_2, X_3, X_4$</td>
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<td>0</td>
<td>11</td>
<td>0</td>
</tr>
<tr>
<td>$X_1, X_2$</td>
<td>22</td>
<td>8</td>
<td>10</td>
<td>8</td>
</tr>
<tr>
<td>Intercept only</td>
<td>0</td>
<td>22</td>
<td>0</td>
<td>27</td>
</tr>
<tr>
<td>$X_1$</td>
<td>14</td>
<td>20</td>
<td>0</td>
<td>14</td>
</tr>
<tr>
<td>$X_2$</td>
<td>1</td>
<td>8</td>
<td>0</td>
<td>14</td>
</tr>
</tbody>
</table>

- Each column reports the percentage of simulations that each model was selected.
- Of the 32 possible models, the results given here include only those with 10% or more support for one of the models.
Effect of range (cont)

Matern $\theta_1=4.0, \theta_2=1.0$
Matern $\theta_1=4.0, \theta_2=1.0$
Matern $\theta_1=6.0, \theta_2=1.0$
Matern $\theta_1=8.0, \theta_2=1.0$
Prediction

Efficient prediction:

- **Time series** ([Shibata (1980), Brockwell and Davis (1991)]). AIC is an efficient order selection procedure for autoregressive models.
- **Regression** (see [McQuarrie and Tsai (1998)]).
- **Other notions of efficiency**, e.g., Kullback-Leibler efficiency and $L_2$ efficiency (see [McQuarrie and Tsai (1998)]).
Prediction Error

Simulations:

• Model selection and estimation using 100 observations.
• Used fitted model to predict at 100 additional locations.
• Evaluated performance using mean square prediction error

$$MSPE = \frac{1}{100} \sum_{j=1}^{100} (Z_j - \hat{Z}_j)^2,$$

where $\hat{Z}_j$ is the universal kriging predictor for the $j^{th}$ prediction location.
**MSPE**

- Spatial AICC is **16.9%** improvement—methods agreed only 11 times.
- Spatial AICC is **39.6%** better over independent AICC with indep noise.
- Predictive coverage was about the same (.92 for spatial AIC, .95 for indep error AIC)
Example: Lizard abundance

Abundance for the orange-throated whiptail lizard in southern California

Data:

- 147 locations
- $Z(s_i) = \log$ (average number of lizards caught per day at location $s_i$)
- Explanatory variables. 37 variables ($2^{37} = 1.374 \times 10^{11}$ models) reduced to the following 6.
  - ant abundance (three levels)
  - log (% sandy soils)
  - elevation
  - barerock indicator
  - % cover
  - log(% chaparral plants)
- 160 possible models
Locations (Southern California) of Whiptail Lizards
Lizard abundance (cont)

<table>
<thead>
<tr>
<th>Predictors</th>
<th>AIC</th>
<th>Spatial Rank</th>
<th>Ind Rank</th>
</tr>
</thead>
<tbody>
<tr>
<td>Ant₁, % sand</td>
<td>54.1</td>
<td>1</td>
<td>66</td>
</tr>
<tr>
<td>Ant₁, Ants₂ , % sand</td>
<td>54.8</td>
<td>2</td>
<td>56</td>
</tr>
<tr>
<td>Ant₁, % sand, % cover</td>
<td>55.7</td>
<td>3</td>
<td>59</td>
</tr>
<tr>
<td>Ant₁, Ant₂ , % sand, % cover, elevation, barerock, % chaparral</td>
<td>92.2</td>
<td>41</td>
<td>1</td>
</tr>
<tr>
<td>Ant₁, Ant₂ , % sand, elevation, barerock, % chaparral</td>
<td>95.3</td>
<td>33</td>
<td>2</td>
</tr>
<tr>
<td>Ant₁, % sand, % cover, elevation, barerock</td>
<td>95.7</td>
<td>38</td>
<td>3</td>
</tr>
</tbody>
</table>

Note: MDL chooses the same model as AICC.
Lizard abundance—simulation

Simulation of model: use covariates $\text{Ant}_1$, % sand, with Matern covariance function
Key: 1 = $\text{ant}_1$, 2 = $\text{ant}_2$, 3 = $\text{ant}_3$, 4 = % sand, 5 = elevation, 6 = barerock, 7 = cover, 8 = chaparral plants
Summary Remarks

1. *Spatial correlation* should not be ignored when selecting explanatory variables.

2. *Model choice* for prediction should involve *joint selection* of the explanatory variables and the form of the autocorrelation function.

3. *Sampling patterns* can severely impact *model selection*.

4. *MDL* offers a nice philosophical alternative to AIC for modeling complex phenomena.