Super-resolution using Gaussian Process Regression Final Year Project Interim Report

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Outline

- Introduction
- Gaussian Process Regression
 - Multivariate Normal Distribution
 - Gaussian Process
 - Regression
 - Training
- GPR for Super-resolution
 - Framework
 - Covariance Function

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The goal of **super-resolution (SR)** is to estimate a high-resolution (HR) image from one or a set of low-resolution (LR) images. It is widely applied in face recognition, medical imaging, HDTV etc.





Figure: Face recognition in video.

The goal of **super-resolution (SR)** is to estimate a high-resolution (HR) image from one or a set of low-resolution (LR) images. It is widely applied in face recognition, medical imaging, HDTV etc.

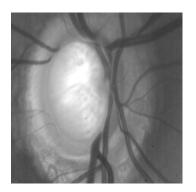




Figure: Super-resolution in medical imaging.

Super-resolution Methods

Interpolation-based methods

Fast but the HR image is usually blurred. E.g., bicubic interpolation, NEDI.

Learning-based methods

Hallucinate textures from the HR/LR image pair database.

Reconstruction-based methods

Formalize an optimization problem constrained by the LR image with various priors.

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Definition

A random vector $\mathbf{X} = (X_1, X_2, \dots, X_p)$ is said to be multivariate normally (MVN) distributed if every linear combination of its components $\mathbf{Y} = \mathbf{a}^T \mathbf{X}$ has a univariate normal distribution. Real-world random variables can often be approximated as following a multivariate normal distribution.

The probability density function of X is

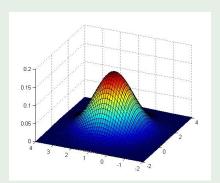
$$f(\mathbf{x}) = \frac{1}{(2\pi)^{(p/2)|\mathbf{\Sigma}|^{1/2}}} \exp\left\{\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \mathbf{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu})\right\}$$
(1)

where μ is the mean of **X** and Σ is the covariance matrix.

Example

Bivariate normal distribution

$$oldsymbol{\mu} = [1 \ 1]', \ oldsymbol{\Sigma} = \left[egin{array}{cc} 1 & 0 \ 0 & 1 \end{array}
ight].$$



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Property 1

The joint distribution of two MVN random variables is also an MVN distribution.

Given
$$\mathbf{X}_1 \sim \mathcal{N}(\boldsymbol{\mu}_1, \boldsymbol{\Sigma}_1)$$
, $\mathbf{X}_2 \sim \mathcal{N}(\boldsymbol{\mu}_2, \boldsymbol{\Sigma}_2)$ and $\mathbf{X} = \begin{bmatrix} \mathbf{X}_1 \\ \mathbf{X}_2 \end{bmatrix}$, we have $\mathbf{X} \sim \mathcal{N}_p(\boldsymbol{\mu}, \boldsymbol{\Sigma})$ with $\boldsymbol{\mu} = \begin{bmatrix} \boldsymbol{\mu}_1 \\ \boldsymbol{\mu}_2 \end{bmatrix}$, $\boldsymbol{\Sigma} = \begin{bmatrix} \boldsymbol{\Sigma}_{11} & \boldsymbol{\Sigma}_{12} \\ \boldsymbol{\Sigma}_{21} & \boldsymbol{\Sigma}_{11} \end{bmatrix}$.

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Property 2

The conditional distribution of the components of MVN are (multivariate) normal.

The distribution of X_1 , given that $X_2 = x_2$, is normal and has

Mean =
$$\mu_1 + \Sigma_{12} \Sigma_{22}^{-1} (\mathbf{x}_2 - \mu_2)$$
 (2)

Covariance =
$$\mathbf{\Sigma}_{11} - \mathbf{\Sigma}_{12} \mathbf{\Sigma}_{22}^{-1} \mathbf{\Sigma}_{21}$$
 (3)

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Gaussian Process

Definition

Gaussian Process (GP) defines a distribution over the function f, where f is a mapping from the input space \mathcal{X} to \mathfrak{R} , such that for any finite subset of \mathcal{X} , its marginal distribution $P(f(\mathbf{x}_1), f(\mathbf{x}_2), ... f(\mathbf{x}_n))$ is a multivariate normal distribution.

$$\mathbf{f}|\mathbf{X} \sim \mathcal{N}(m(\mathbf{x}), K(\mathbf{X}, \mathbf{X}))$$
 (4)

where

$$\mathbf{X} = \{\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_n\} \tag{5}$$

$$m(\mathbf{x}) = E[f(\mathbf{x})] \tag{6}$$

$$k(\mathbf{x}_i, \mathbf{x}_j) = E\left[(f(\mathbf{x}_i) - m(\mathbf{x}))(f(\mathbf{x}_i)^T - m(\mathbf{x}^T)) \right]$$
(7)

and $K(\mathbf{X}, \mathbf{X})$ denotes the covariance matrix such that $\mathbf{K}_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$.

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Gaussian Process

Formally, we write the Gaussian Process as

$$f(\mathbf{x}) \sim \mathcal{GP}(m(\mathbf{x}), k(\mathbf{x}_i, \mathbf{x}_j))$$
 (8)

Without loss of generality, the mean is usually taken to be zero.

- Parameterized by the *mean* function $m(\mathbf{x})$ and the *covariance* function $k(\mathbf{x}_i, \mathbf{x}_j)$
- Infer in the function space directly

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Gaussian Process Regression

Model:

$$f(\mathbf{x}) \sim \mathcal{GP}(m(\mathbf{x}), k(\mathbf{x}_i, \mathbf{x}_j))$$
 (9)

Given the inputs X_* , the output f_* is

$$\mathbf{f}_* \sim \mathcal{N}(\mathbf{0}, \mathcal{K}(\mathbf{X}_*, \mathbf{X}_*))$$
 (10)

According to the Gaussian prior, the joint distribution of the training outputs \mathbf{f} , and the test outputs \mathbf{f}_* is

$$\begin{bmatrix} \mathbf{f} \\ \mathbf{f}_* \end{bmatrix} \sim \mathcal{N} \left(\mathbf{0}, \begin{bmatrix} \mathcal{K}(\mathbf{X}, \mathbf{X}) & \mathcal{K}(\mathbf{X}, \mathbf{X}_*) \\ \mathcal{K}(\mathbf{X}_*, \mathbf{X}) & \mathcal{K}(\mathbf{X}_*, \mathbf{X}_*) \end{bmatrix} \right). \tag{11}$$

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Noisy Model

In reality, we do not have access to true function values but rather noisy observations. Assuming independent indentically distributed noise, we have the noisy model

$$y = f(\mathbf{x}) + \varepsilon, \ \varepsilon \sim \mathcal{N}(0, \sigma_n^2)$$
 (12)

$$f(\mathbf{x}) \sim \mathcal{GP}(m(\mathbf{x}), K(\mathbf{X}, \mathbf{X}))$$
 (13)

$$Var(y) = Var(f(\mathbf{x})) + Var(\varepsilon) = K(\mathbf{X}, \mathbf{X}) + \sigma_n^2 I$$
 (14)

Thus, the joint distribution for prediction is

$$\begin{bmatrix} \mathbf{y} \\ \mathbf{f}_* \end{bmatrix} \sim \mathcal{N} \left(\mathbf{0}, \begin{bmatrix} K(\mathbf{X}, \mathbf{X}) + \sigma_n^2 I & K(\mathbf{X}, \mathbf{X}_*) \\ K(\mathbf{X}_*, \mathbf{X}) & K(\mathbf{X}_*, \mathbf{X}_*) \end{bmatrix} \right)$$
(15)



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Prediction

Referring to the previous property of the conditional distribution, we can obtain

$$\mathbf{f}_* \sim \mathcal{N}(\mathbf{\bar{f}}, V(\mathbf{f}_*))$$
 (16)

$$\overline{\mathbf{f}}_* = K(X_*, X)[K(X, X) + \sigma_n^2 I]^{-1} \mathbf{y}, \tag{17}$$

$$V(\mathbf{f}_{*}) = K(X_{*}, X_{*}) - K(X_{*}, X)[K(X, X) + \sigma_{n}^{2}I]^{-1}K(X, X_{*}).$$
 (18)

y are the training outputs and \mathbf{f}_* are the test outputs, which are predicted as the mean $\overline{\mathbf{f}}$.

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Marginal Likelihood

GPR model:

$$\mathbf{y} = \mathbf{f} + \boldsymbol{\epsilon} \tag{19}$$

$$\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}), \mathbf{K})$$
 (20)

$$\epsilon \sim \mathcal{N}(\mathbf{0}, \sigma_n^2 \mathbf{I})$$
 (21)

y is an n-dimensional vector of observations. Without loss of generality, let $m(\mathbf{x}) = 0$. Thus $y | \mathbf{X}$ follows a normal distribution with

$$E(y|\mathbf{X}) = 0 (22)$$

$$Var(y|\mathbf{X}) = K(\mathbf{X}, \mathbf{X}) + \sigma_n^2 \mathbf{I}$$
 (23)

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Marginal Likelihood

Let $\mathbf{K}_y = Var(y|\mathbf{X})$,

$$p(\mathbf{y}|\mathbf{X}) = \frac{1}{(2\pi)^{n/2}|\mathbf{K}_y|^{1/2}} \exp\left\{-\frac{1}{2}\mathbf{y}^T \mathbf{K}_y^{-1} \mathbf{y}\right\}$$
(24)

The log marginal likelihood is

$$\mathcal{L} = \log p(\mathbf{y}|\mathbf{X}) = -\frac{n}{2}\log 2\pi - \frac{1}{2}\log |\mathbf{K}_y| - \frac{1}{2}\mathbf{f}^{\mathsf{T}}\mathbf{K}_y^{-1}\mathbf{f}$$
 (25)

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Maximum a posteriori

Matrix derivative:

$$\frac{\partial}{\partial x} \mathbf{Y} = -\mathbf{Y}^{-1} \frac{\partial \mathbf{Y}}{\partial \theta_i} \mathbf{Y}^{-1}$$
 (26)

$$\frac{\partial}{\partial x} \mathbf{Y} = -\mathbf{Y}^{-1} \frac{\partial \mathbf{Y}}{\partial \theta_i} \mathbf{Y}^{-1}$$

$$\frac{\partial}{\partial x} \log |\mathbf{Y}| = \operatorname{tr}(\mathbf{Y}^{-1} \frac{\partial \mathbf{Y}}{\partial \theta_i})$$
(26)

Gradient ascent:

$$\frac{\partial \mathcal{L}}{\partial \theta_i} = \frac{1}{2} \mathbf{y}^T K^{-1} \frac{\partial K}{\partial \theta_i} K^{-1} \mathbf{y} - \frac{1}{2} tr(K^{-1} \frac{\partial K}{\partial \theta_i})$$
 (28)

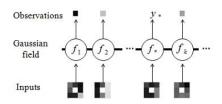
 $\frac{\partial K}{\partial \theta}$ is a matrix of derivatives of each element.

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Graphical Representation



- Model: $y = f(\mathbf{x}) + \varepsilon$
- Squares: observed pixels
- Circles: unknown Gaussian field
- Inputs (x): neighbors (predictors) of the target pixel
- Outputs (y): pixel at the center of each 3×3 patch
- Thick horizontal line: a set of fully connected nodes.

Stage 1: interpolation Input LR patch



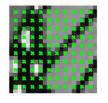
Stage 1: interpolation Sample training targets





Stage 1: interpolation SR based on Bicubic Interpolation



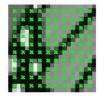




Stage 2: deblurring

Stage 1: interpolation





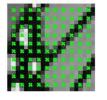


Stage 2: deblurringSample training targets



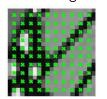
Stage 1: interpolation







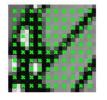
Stage 2: deblurringObtain neighbors from the downsampled patch





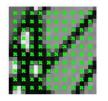
Stage 1: interpolation







Stage 2: deblurringSR based on the simulated blurring process







Covariance Equation

- defines the similarity between two points (vectors)
- indicate the underlying distribution of functions in GP
- Squared Exponential covariance function

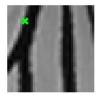
$$k(\mathbf{x}_i, \mathbf{x}_j) = \sigma_f^2 \exp\left(-\frac{1}{2} \frac{(\mathbf{x}_i - \mathbf{x}_j)'(\mathbf{x}_i - \mathbf{x}_j)}{\ell^2}\right)$$
(29)

 σ_f^2 represents the signal variance and ℓ defines the *characteristic* length scale.

Given an image \mathbf{I} , the covariance between two pixels $\mathbf{I}_{i,j}$ and $\mathbf{I}_{m,n}$ is calculated as $k(\mathbf{I}_{(i,j),N},\mathbf{I}_{(m,n),N})$, where N means to take the 8 nearest pixels around the pixel. Therefore, the similarity is based on the Euclidean distance between the pixels' neighbors.

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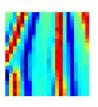
Covariance Equation







(b) Training patch



(c) Covariance matrix

- **Local similarity**: high responses (red regions) from the training patch are concentrated on edges
- Global similarity: high-responsive regions also include other similar edges within the patch
- Conclusion: pixels embedded in a similar structure to that of the target pixel in terms of the neighborhood tend to have higher weights during prediction

Hyperparameters:

- σ_f^2 : signal variance
- σ_n^2 : noise variance
- ℓ : characteristic length scale











- (a) Test
- - $\sigma_n = .01$ $\sigma_n = .001$ $\sigma_n = .14$
- (b) Training (c) $\ell = .50$, (d) $\ell = .05$, (e) $\ell = 1.65$,
- (c): MAP estimation
- (d): Quickly varying field with low noise
- (e): Slowly varyin field with high noise

Hyperparameters:

- σ_f^2 : signal variance
- σ_n^2 : noise variance
- ℓ: characteristic length scale











- (a) Test
- (b) Training (c) $\ell = .50$, (d) $\ell = .05$, (e) $\ell = 1.65$,
- - $\sigma_n = .01$ $\sigma_n = .001$ $\sigma_n = .14$

- (c): MAP estimation
- (d): Quickly varying field with low noise (high-frequncy artifacts)
- (e): Slowly varyin field with high noise

Hyperparameters:

- σ_f^2 : signal variance
- σ_n^2 : noise variance
- ℓ: characteristic length scale











- (a) Test
- (b) Training (c) $\ell = .50$, (d) $\ell = .05$, (e) $\ell = 1.65$,

 - $\sigma_n = .01$ $\sigma_n = .001$ $\sigma_n = .14$

- (c): MAP estimation
- (d): Quickly varying field with low noise (high-frequncy artifacts)
- (e): Slowly varyin field with high noise (too smooth)

Log marginal likelihood:

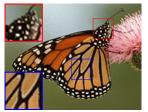
$$\log p(\mathbf{y}|X,\boldsymbol{\theta}) = -\frac{1}{2}\mathbf{y}^T K_y^{-1} \mathbf{y} - \frac{1}{2}\log|K_y| - \frac{n}{2}\log 2\pi$$
 (30)

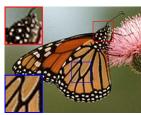
Maximize a posteriori (gradient descent):

$$\frac{\partial \mathcal{L}}{\partial \theta_i} = \frac{1}{2} \mathbf{y}^T K^{-1} \frac{\partial K}{\partial \theta_i} K^{-1} \mathbf{y} - \frac{1}{2} tr(K^{-1} \frac{\partial K}{\partial \theta_i})$$
(31)

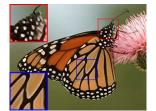
heta denotes the parameter set.

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(a) Bicubic (MSSIM=0.84) (b) GPP (MSSIM=0.84)



result

(c) Our (MSSIM=0.86)



(d) Ground truth





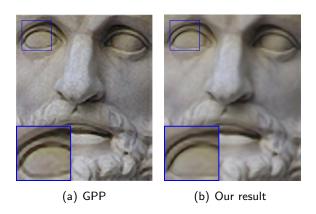
(a) Input (b) $3\times$ direct magnification



(c) $10 \times$ our result



(d) $10\times$ detail synthesis







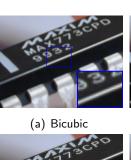


(b) Edge statistics



(c) Patch redundancy









(b) Edge statistics



(c) Patch redundancy

(d) Ours







(b) Edge statistics







(b) Edge statistics



(a) Bicubic

(b) Edge statistics