

# CS-E4895 Gaussian Processes

## Lecture 9: Deep GPs

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# Roadmap for today

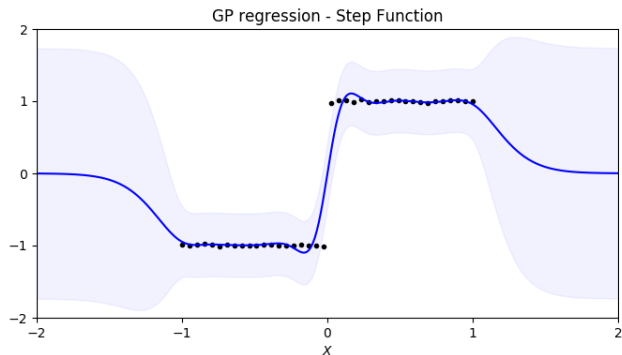
- 1 Introductions to Deep GPs
  - Limitations of standard GPs
  - Function Composition and Deep Learning
- 2 The Deep GP Model
  - Combining Layers of GPs
  - Deep GP Covariance
  - The Deep GP Posterior
- 3 Inference in Deep GPs
  - Stochastic Variational Inference
  - Alternative Approaches
  - Performance and Issues

# Section 1

## Introductions to Deep GPs

# Limitations of Standard GPs

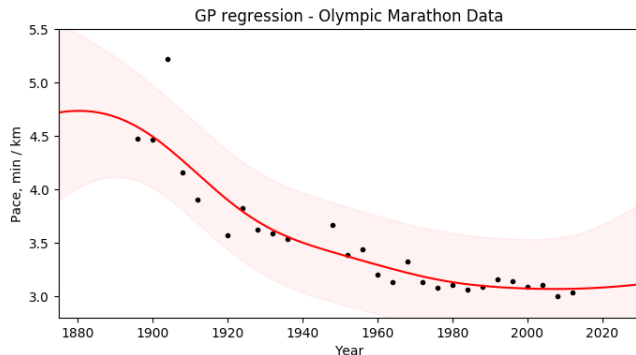
- **Discontinuities / jumps**



- A stationary GP fails to capture the sharp jump, and the variance is too large everywhere.

# Limitations of Standard GPs

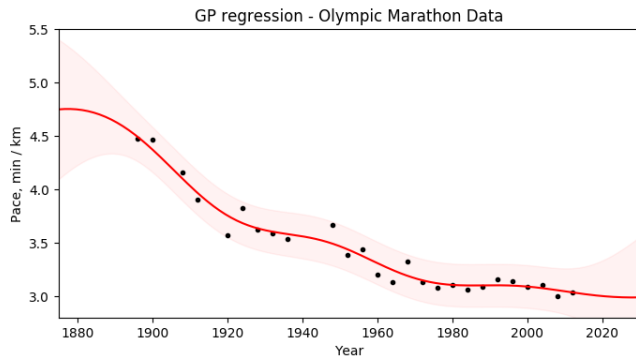
- Discontinuities / jumps
- Outliers



- The outlier has a *very* low probability under the model.
- To account for this, the model learns a likelihood variance that is too high for all other data points.

# Limitations of Standard GPs

- Discontinuities / jumps
- Outliers

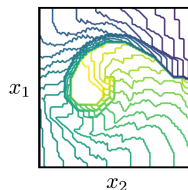


- Removing the outlier vastly improves the result. But we'd rather avoid such a manual intervention.

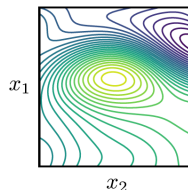
# Limitations of Standard GPs

- Discontinuities / jumps
- Outliers
- Non-stationarity

The previous two problems can be seen as issues arising due to a *stationary* model being applied to non-stationary data.



original data

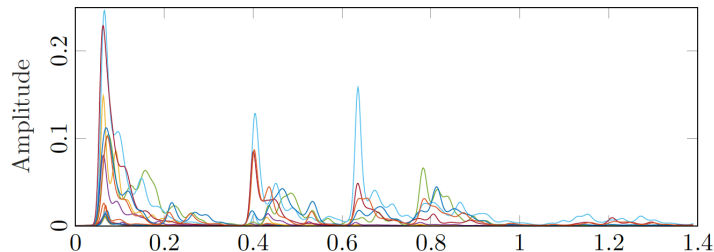


GP prediction  
MSE = 138

- Many real-world data sets do not have constant smoothness across the entire input space.

# Limitations of Standard GPs

- Discontinuities / jumps
- Outliers
- Non-stationarity
- Misalignment



- Multiple misaligned data streams cannot be modelling with a standard (multi-output) GP.
- The data must be aligned via a pre-processing step.
- Ideally this step should be incorporated into the probabilistic model, so that its uncertainty can be incorporated.



# Function Composition

- Function composition is at the heart of modern-day machine learning. Deep neural networks are made up of compositions of neural networks.
- Deep Gaussian processes work in an analogous way, whilst incorporating uncertainty and prior knowledge.

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Single GPs can model simple, stationary functions. The composition of multiple GPs,

$$f_3(f_2(f_1(\cdot))) = (f_3 \circ f_2 \circ f_1)(\cdot)$$

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can model more complex, nonstationary functions.

- We can view each “layer” as a warping of the inputs before feeding to the next layer.
- Function composition can be used to incorporate multiple layers of prior knowledge.

## Section 2

# The Deep GP Model

# Deep GP Intuition

Before writing down the model, let's gain some intuition about hierarchies of Gaussian processes.

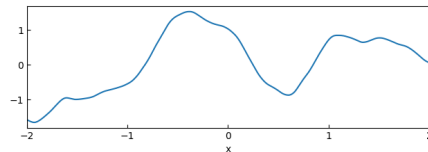
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Take inputs  $\mathbf{x}$ , and evaluate a GP,  $f_1(\cdot) \sim \mathcal{GP}(\mu_1(\cdot), \kappa_1(\cdot, \cdot))$ :

$$f_1(\mathbf{x}) \sim \mathcal{N}(\mu_1(\mathbf{x}), \kappa_1(\mathbf{x}, \mathbf{x}))$$

Draw a sample,  $\tilde{\mathbf{y}}_1$ , from this multivariate Gaussian:



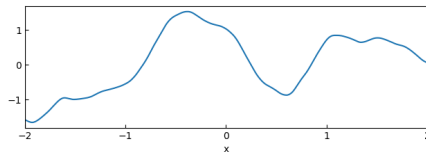
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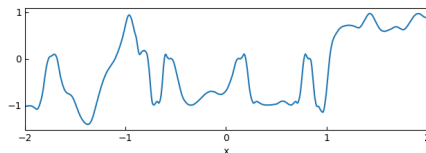
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Treat this sample as the input to *another* GP,  $f_2(\cdot) \sim \mathcal{GP}(\mu_2(\cdot), \kappa_2(\cdot, \cdot))$ :

$$f_2(\tilde{\mathbf{y}}_1) \sim \mathcal{N}(\mu_2(\tilde{\mathbf{y}}_1), \kappa_2(\tilde{\mathbf{y}}_1, \tilde{\mathbf{y}}_1))$$

and draw a sample,  $\tilde{\mathbf{y}}_2$ .



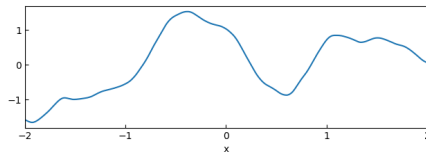
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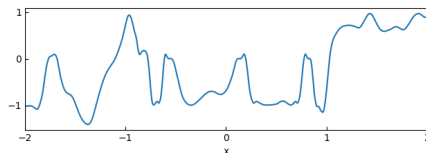
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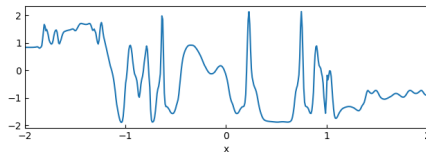
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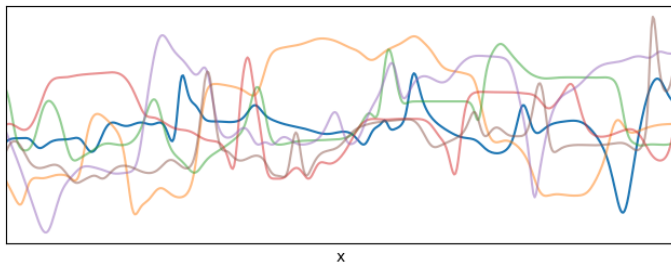
Repeat a third time for  $f_3(\cdot) \sim \mathcal{GP}(\mu_3(\cdot), \kappa_3(\cdot, \cdot))$ :

$$f_3(\tilde{\mathbf{y}}_2) \sim \mathcal{N}(\mu_3(\tilde{\mathbf{y}}_2), \kappa_3(\tilde{\mathbf{y}}_2, \tilde{\mathbf{y}}_2))$$



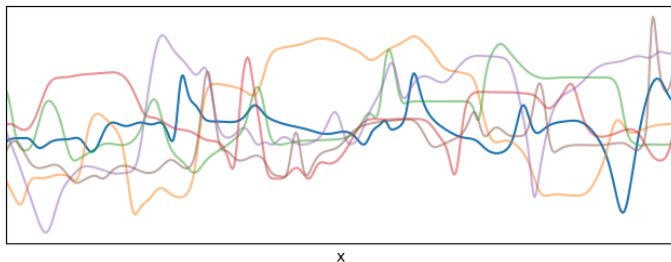


# Deep GP Intuition



These are samples from a 3-layer deep GP.

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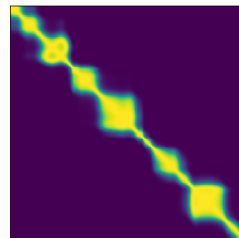
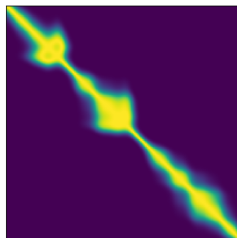
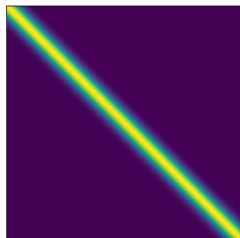
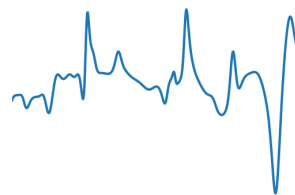


These are samples from a 3-layer deep GP.

- sharp jumps / discontinuities.
- highly nonstationary smoothness.
- rich space of function, high capacity
- how to avoid overfitting?

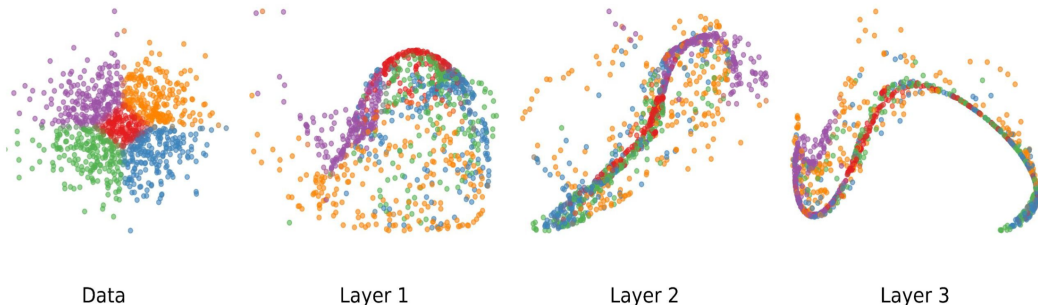
# Deep GP Covariance

As well as sampling, we can also plot the covariance matrix in each layer.  
Initially we have uniform input, later more 'clumping'



# Signal propagates through layers

(a) Draws from the DGP prior



- The color is only a visual aide
- In this example each layer maps  $\mathbb{R}^2$  to  $\mathbb{R}^2$
- The plot shows one sample path from the DGP
- Each sample path conserves neighborhoods to a degree
- No output layer shown

# The Deep GP Model

Now let's write down the deep GP model and look at its properties. Inference will come later.

$$\begin{aligned} f_\ell(\cdot) &\sim \mathcal{GP}(\mu_\ell(\cdot), \kappa_\ell(\cdot, \cdot)) , & \ell = 1, \dots, L \\ p(\tilde{\mathbf{y}}_\ell \mid f_\ell, \tilde{\mathbf{y}}_{\ell-1}) &= \prod_n \mathcal{N}(\tilde{y}_{\ell,n} \mid f_\ell(\tilde{\mathbf{y}}_{\ell-1,n}), \sigma_\ell^2) , & \tilde{\mathbf{y}}_1 = \mathbf{x} \\ p(\mathbf{y} \mid f_L, \tilde{\mathbf{y}}_{L-1}) &= \prod_n p(y_n \mid f_L(\tilde{\mathbf{y}}_{L-1,n})) \end{aligned}$$

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- We denote layer by  $\ell$  (**not** lengthscale!)
- $L$  layers of Gaussian process priors.
- $\tilde{\mathbf{y}}_\ell$  are latent variables - treated as input to layer  $\ell + 1$ .
- Typically include Gaussian noise between layers.

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where  $f_0 = \mathbf{x}$  and  $f_{L,n} = f_L(f_{L-1}(\dots(x_n)))$ .

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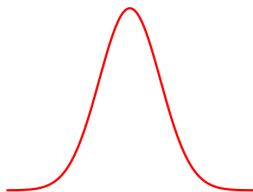
Using notation  $\mathbf{f}_\ell = f(\mathbf{f}_{\ell-1})$ , the full process has joint density

$$p(\mathbf{y}, \{\mathbf{f}_\ell\}_{\ell=1}^L) = \underbrace{\prod_{n=1}^N p(y_n \mid \mathbf{f}_{L,n})}_{\text{Likelihood}} \underbrace{\prod_{\ell=1}^L p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1})}_{\text{Deep GP Prior}}$$
$$p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1}) = \mathcal{N}(\mathbf{f}_\ell \mid \mu_{\ell-1}(\mathbf{f}_\ell), \mathbf{K}_\ell(\mathbf{f}_{\ell-1}, \mathbf{f}_{\ell-1}))$$

The  $\mathbf{f}_\ell$  is of size  $(N, M_\ell)$  where  $M_0 = D$  and  $M_L = \text{size}(\mathbf{y})$

# The Deep GP Posterior

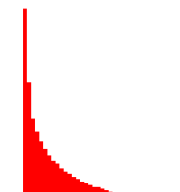
A Gaussian propagated through a nonlinearity is no longer Gaussian:



$$x \sim \mathcal{N}(x \mid \cdot, \cdot)$$



$$f(\cdot)$$



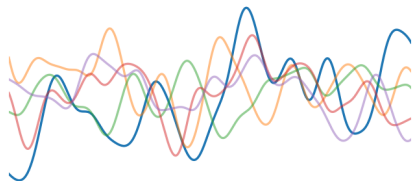
$$f(x) \sim ???$$

# The Deep GP Posterior

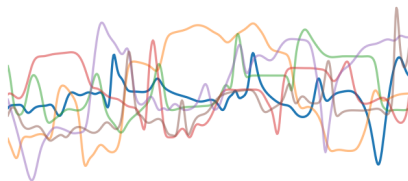
Similarly, a Gaussian process propagated through a nonlinearity (e.g., another GP) is no longer a Gaussian process (in the original inputs  $\mathbf{x}$ ).

$$f_1(\cdot) \sim \mathcal{GP}(\mu_1(\cdot), \kappa_1(\cdot, \cdot))$$

$$f_2(\cdot) \sim \mathcal{GP}(\mu_2(\cdot), \kappa_2(\cdot, \cdot))$$

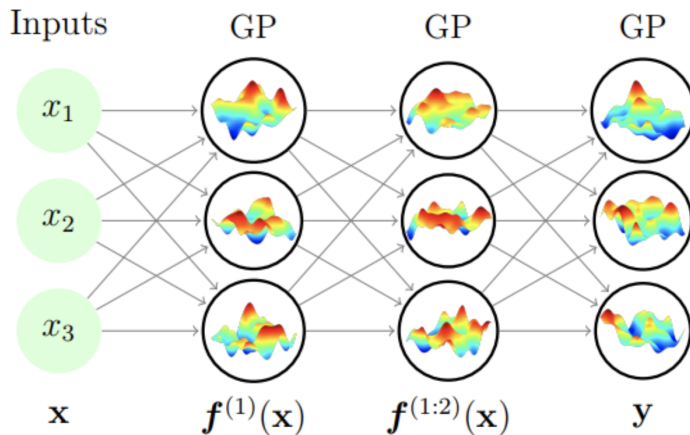


$$f_1(\mathbf{x}) \sim \mathcal{GP}(\cdot, \cdot)$$



$$f_2 \circ f_1(\mathbf{x}) = f_2(f_1(\mathbf{x})) \sim ???$$

# Deep GP illustration



- The size of each layer space can vary
- If observation  $y$  is a scalar, the final layer needs to map to scalar space

## Section 3

### Inference in Deep GPs

# Inference in Deep GPs

Since the posterior is not Gaussian, it is clear that we must resort to approximate inference.

- Various schemes have been proposed: Variational Inference, Expectation Propagation, Hamiltonian Monte Carlo.
- We will focus on **sparse, stochastic variational inference**.

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- Recall our joint probability:

$$p(\mathbf{y}, \{\mathbf{f}_\ell\}_{\ell=1}^L) = \prod_{n=1}^N \underbrace{p(y_n | \mathbf{f}_{L,n})}_{\text{likelihood}} \prod_{\ell=1}^L \underbrace{p(\mathbf{f}_\ell | \mathbf{f}_{\ell-1})}_{\text{DGP priors}}$$

where  $\mathbf{f}_0 = \mathbf{x}$ .

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where  $\mathbf{f}_0 = \mathbf{x}$ .

- We introduce inducing inputs  $\mathbf{z}_\ell$  and outputs  $\mathbf{u}_\ell = \mathbf{f}_\ell(\mathbf{z}_\ell)$  in each layer:

$$p(\mathbf{y}, \{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) = \prod_{n=1}^N p(y_n | \mathbf{f}_{L,n}) \prod_{\ell=1}^L p(\mathbf{f}_\ell | \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) p(\mathbf{u}_\ell)$$

where

$$p(\mathbf{f}_\ell | \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) = \mathcal{N}(\mathbf{f}_\ell | \mathbf{K}(\mathbf{f}_{\ell-1}, \mathbf{z}_\ell) \mathbf{K}(\mathbf{z}_\ell, \mathbf{z}_\ell)^{-1} \mathbf{u}_\ell, \mathbf{K}(\mathbf{f}_{\ell-1}, \mathbf{f}_{\ell-1}) - \mathbf{K}(\mathbf{f}_{\ell-1}, \mathbf{z}_\ell) \mathbf{K}(\mathbf{z}_\ell, \mathbf{z}_\ell)^{-1} \mathbf{K}(\mathbf{z}_\ell, \mathbf{f}_{\ell-1}))$$



# Stochastic VI for Sparse Deep GPs

True joint distribution

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- To construct a variational lower bound for the deep GP, we must first define an **approximate posterior**:

$$q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) = \prod_{\ell=1}^L \underbrace{p(\mathbf{f}_\ell | \mathbf{f}_{\ell-1}, \mathbf{u}_\ell)}_{\text{Gaussian conditional}} q(\mathbf{u}_\ell)$$

where  $q(\mathbf{u}_\ell) = \mathcal{N}(\mathbf{u}_\ell | \mathbf{m}_\ell, \mathbf{S}_\ell)$  are free-form Gaussians whose parameters are to be optimised. We also optimise inducing inputs  $\mathbf{z}_\ell$  as variational parameters.

# Stochastic VI for Sparse Deep GPs

- Recall the sparse variational bound for a single GP derived in previous lectures:

$$\begin{aligned}\ln p(\mathbf{y}) &\geq \sum_{n=1}^N \int q(f_n) \ln p(y_n | f_n) \mathrm{d}f_n - \mathbb{D}[q(\mathbf{u}) || p(\mathbf{u})] \\ &= \mathbb{E}_{q(\mathbf{f}, \mathbf{u})} [\ln p(\mathbf{y} | \mathbf{f})] + \mathbb{E}_{q(\mathbf{f}, \mathbf{u})} [\ln p(\mathbf{f}, \mathbf{u})] - \mathbb{E}_{q(\mathbf{f}, \mathbf{u})} [\ln q(\mathbf{f}, \mathbf{u})] \\ &= \mathbb{E}_{q(\mathbf{f}, \mathbf{u})} \left[ \ln \frac{p(\mathbf{y}, \mathbf{f}, \mathbf{u})}{q(\mathbf{f}, \mathbf{u})} \right]\end{aligned}$$

where

$$\begin{aligned}q(\mathbf{f}) &= \int p(\mathbf{f} | \mathbf{u}) q(\mathbf{u}) \mathrm{d}\mathbf{u} \\ &= \mathcal{N}(\mathbf{f} | \mathbf{A}\mathbf{m}, \mathbf{K}_{\mathbf{xx}} + \mathbf{A}(\mathbf{S} - \mathbf{K}_{\mathbf{zz}})^{-1}\mathbf{A}^T) \\ \mathbf{A} &= \mathbf{K}_{\mathbf{xz}}\mathbf{K}_{\mathbf{zz}}^{-1}\end{aligned}$$

- We will now derive a similar bound for the deep GP.

# Stochastic VI for Sparse Deep GPs

**joint:**  $p(\mathbf{y}, \{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) = \prod_{n=1}^N p(y_n \mid \mathbf{f}_{L,n}) \prod_{\ell=1}^L p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) p(\mathbf{u}_\ell)$

**approx. posterior:**  $q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) = \prod_{\ell=1}^L p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) q(\mathbf{u}_\ell)$

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- The variational bound is

$$\ln p(\mathbf{y}) \geq \mathcal{L}_{DGP} = \mathbb{E}_{q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)} \left[ \ln \frac{p(\mathbf{y}, \{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)}{q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)} \right]$$

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$$\text{approx. posterior: } q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) = \prod_{\ell=1}^L p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) q(\mathbf{u}_\ell)$$

- The variational bound is

$$\begin{aligned} \ln p(\mathbf{y}) \geq \mathcal{L}_{DGP} &= \mathbb{E}_{q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)} \left[ \ln \frac{p(\mathbf{y}, \{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)}{q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)} \right] \\ &= \int \int q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) \ln \left( \frac{p(\mathbf{y}, \{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)}{q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L)} \right) d\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L \\ &= \int \int q(\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L) \ln \left( \frac{\prod_{n=1}^N p(y_n \mid \mathbf{f}_{L,n}) \prod_{\ell=1}^L p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) p(\mathbf{u}_\ell)}{\prod_{\ell=1}^L p(\mathbf{f}_\ell \mid \mathbf{f}_{\ell-1}, \mathbf{u}_\ell) q(\mathbf{u}_\ell)} \right) d\{\mathbf{f}_\ell, \mathbf{u}_\ell\}_{\ell=1}^L \end{aligned}$$

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- This means that sampling from  $q(\mathbf{f}_{\ell,n})$  is cheap, and does not involve sampling from the full GP at each layer (in fact, it only requires sampling from univariate Gaussians).

# Optimising the Deep GP Bound

$$\ln p(\mathbf{y}) \geq \mathcal{L}_{DGP} = \sum_{n=1}^N \int q(f_{L,n}) \ln p(y_n | f_{L,n}) \mathrm{d}f_{L,n} - \sum_{\ell=1}^L \mathbb{D}[q(\mathbf{u}_\ell) || p(\mathbf{u}_\ell)]$$

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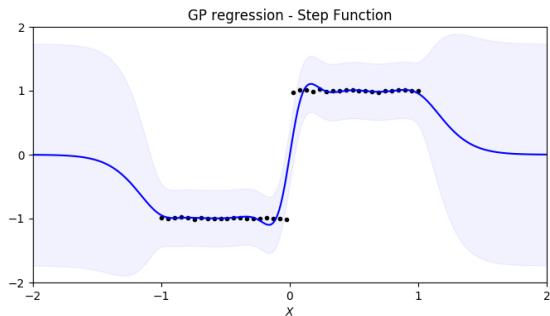
- For the second term, compute the KL divergence between  $q(\mathbf{u}_\ell)$  and  $p(\mathbf{u}_\ell)$  in each layer separately (this is available in closed form since both terms are Gaussian).
- This inference technique is called **doubly stochastic VI**, due to the two sources of stochasticity.

# Alternative Approaches

Other approaches to deep GP inference exist, but we won't go over them here:

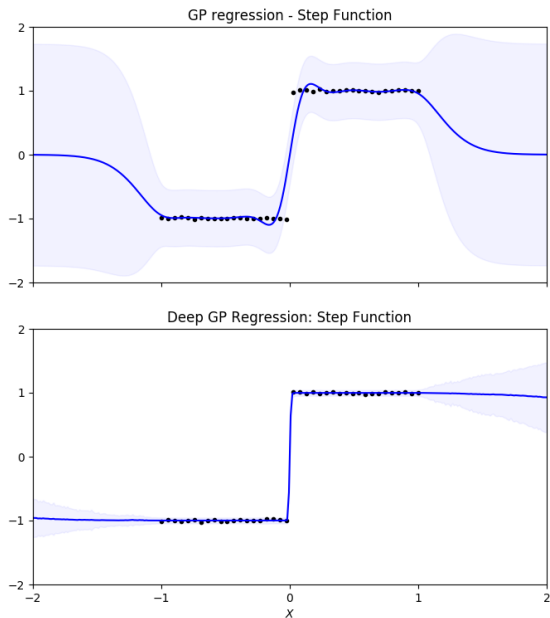
- **Deep GP Expectation Propagation** - similar to the above, but using EP for inference, and replacing the sampling procedure with Gaussian projections to approximate the marginals.
- **Importance-weighted VI with latent variables** - introduces additional latent variables which allow the model to represent non-Gaussian posteriors.
- **Hamiltonian Monte Carlo** - uses a sophisticated sampling approach to represent non-Gaussian posteriors.

- Discontinuities / jumps:

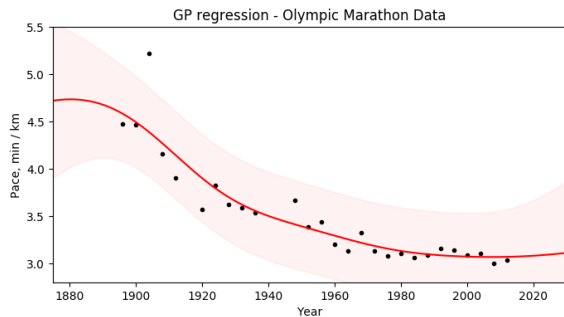


# Deep GP Performance

- **Discontinuities / jumps:**
- The deep GP captures the jump, whilst the variance elsewhere remains low.
- However, we would prefer that the variances increases in the region of the discontinuity.
- In the exercises, you will examine what happens in each layer.

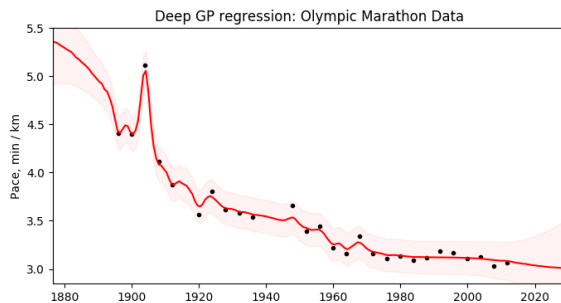
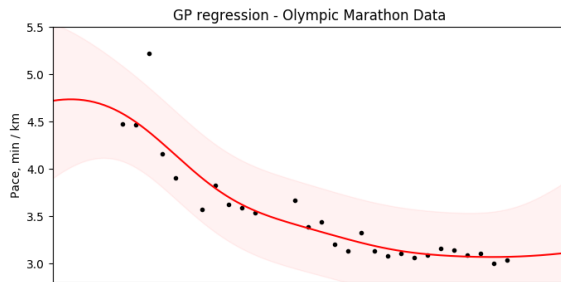


- **Outliers:**



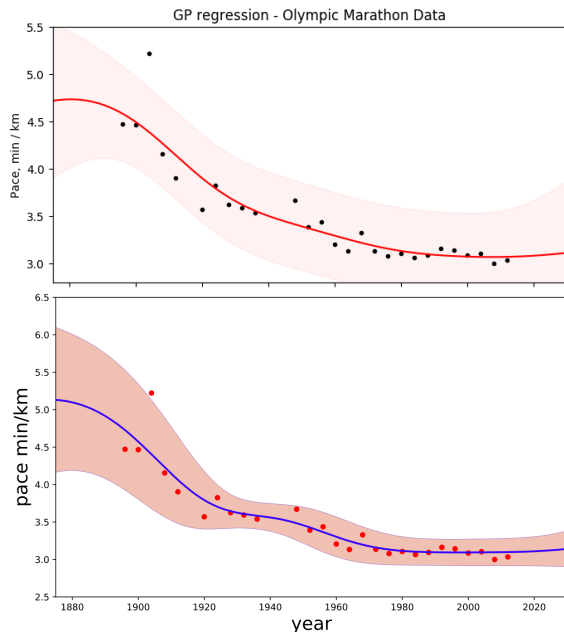
# Deep GP Performance

- **Outliers:**
- The deep GP seems to overfit the outlier.

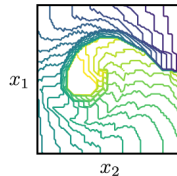
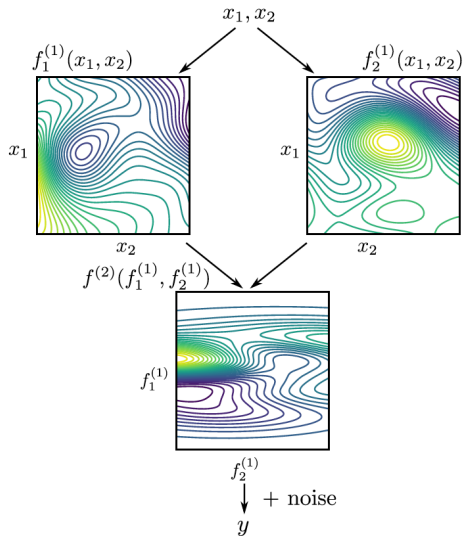


# Deep GP Performance

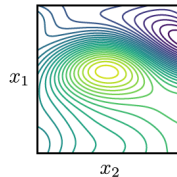
- **Outliers:**
- The deep GP seems to overfit the outlier.
- Whereas the originally proposed deep GP methods claim to solve these tasks well.
- But doubly stochastic VI reports superior performance on many machine learning tasks, potentially because it scales to large data.



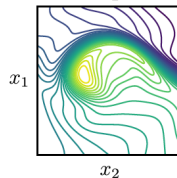
# Deep GP Performance



original data



GP prediction  
MSE = 138



DGP prediction  
MSE = 62.4



# Issues with Deep GPs

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- **Training can be slow**: we trade off the number of samples with accuracy.
- **Training is more prone to getting stuck in local minima** since there are many more parameters to optimise.
- **Current approaches to VI tend to “turn off” layers**, or reduce their variance to near-zero (such that they behave like deterministic mappings).

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- Performance matches *e.g.*, deep CNNs, but improves uncertainty quantification in predictions, *i.e.*, **the model is more aware when it is wrong.**
- So deep GPs have great potential. But, as we have seen, there is still much work to be done.



- Salimbeni and Deisenroth. Doubly Stochastic Variational Inference for Deep Gaussian Processes, NIPS 2017
  - Today's lecture followed the ds-DGP
- Duvenaud, Rippel, Adams, Ghahramani. Avoiding pathologies in very deep networks. AISTATS 2014
  - Interesting discussion of DGP risks of *rank collapse*

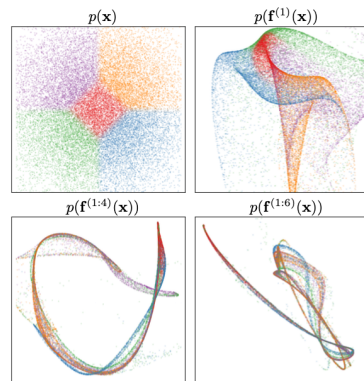


Figure 5: Visualization of draws from a deep GP. A 2-dimensional Gaussian distribution (top left) is warped by successive functions drawn from a GP prior. As the number of layers increases, the density concentrates along one-dimensional filaments.

# End of Today's Lecture

- Next time: I will give a lecture on latent models with GPs.