

Probabilistic methods for Machine Learning

November 28
2023

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Topics For Today

- **Probability recap**
- Naïve Bayes
- Gaussian process
- Bayesian hyperparameter optimization

Probability: definitions

Sample Space Ω :

The set of all outcomes of a random experiment

Event space \mathcal{F} :

A set whose elements $A \in \mathcal{F}$ (called events) are subsets of Ω (i.e., $A \subseteq \Omega$ is a collection of possible outcomes of an experiment)

Probability measure:

A function $P : \mathcal{F} \rightarrow \mathbb{R}$ that satisfies the following properties:

- $P(A) \geq 0$, for all $A \in \mathcal{F}$
- $P(\Omega) = 1$
- If A_1, A_2, \dots are disjoint events (i.e., $A_i \cap A_j = \emptyset$ for $i \neq j$), then $P(\cup_i A_i) = \sum_i P(A_i)$

Probability: definitions

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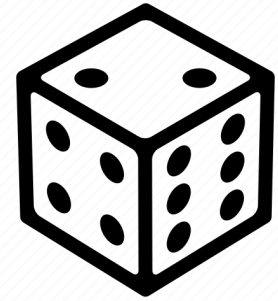
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Example: Tossing a six-sided die

$$\Omega = \{1, 2, 3, 4, 5, 6\}$$

$$\mathcal{F} = \{\emptyset, \Omega\} = \{\emptyset, 1, 2, 3, 4, 5, 6\}$$

$$P(\emptyset) = 0, P(\Omega) = 1$$

$$P(\{1\}) = \frac{1}{6}$$

$$P(\{1, 2, 3, 4\}) = \frac{4}{6}$$

$$P(\{2, 4, 6\}) = \frac{3}{6}$$

Probability: Conditional probability and independence

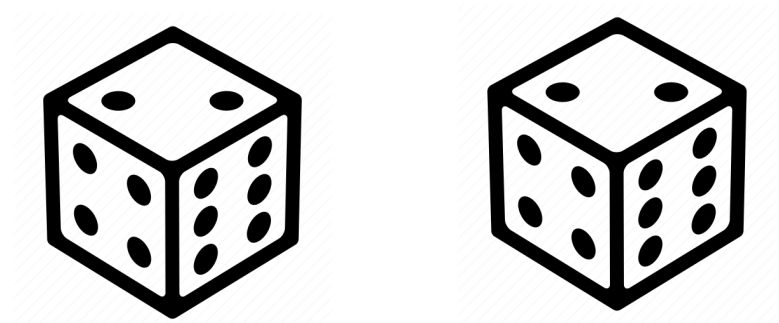
The conditional probability of any event A given B is

$$P(A|B) = \frac{P(A \cap B)}{P(B)}$$

Example: tossing two 6-sided die.

A = {We got same digits on first and second die}

B = {Sum of digits on first and second die more than 8}



Probability: Conditional probability and independence

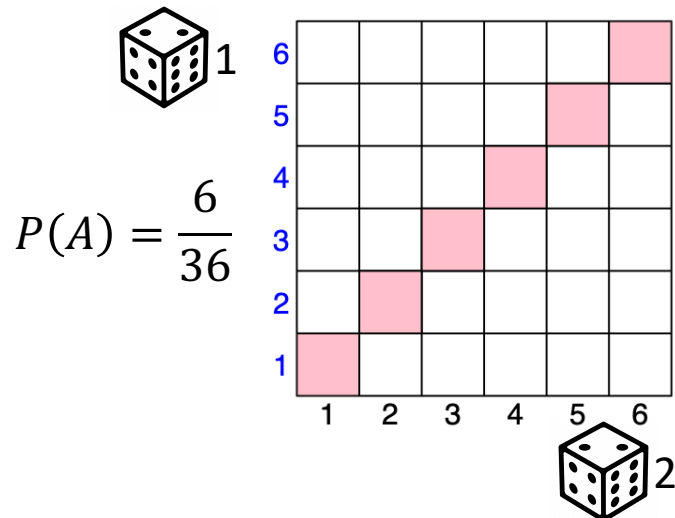
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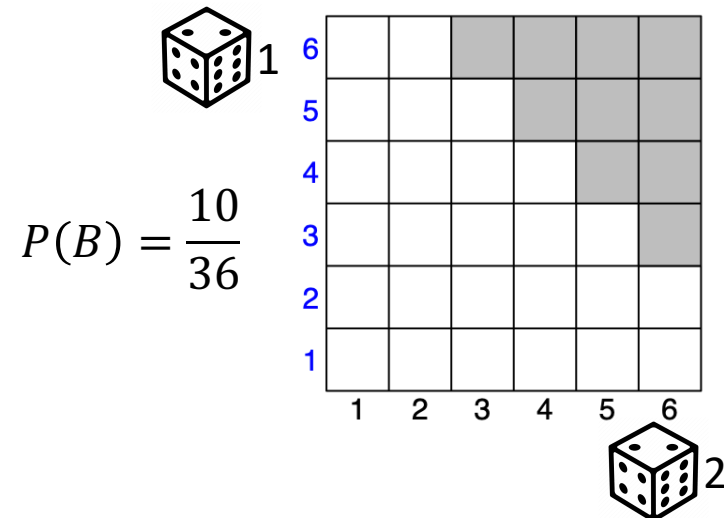
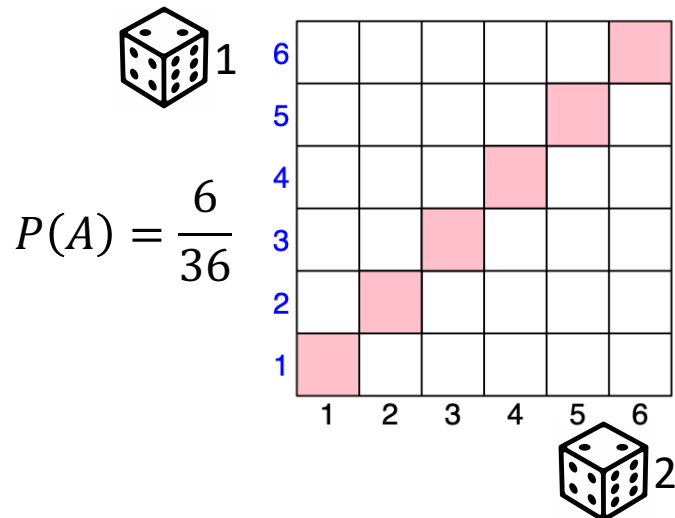
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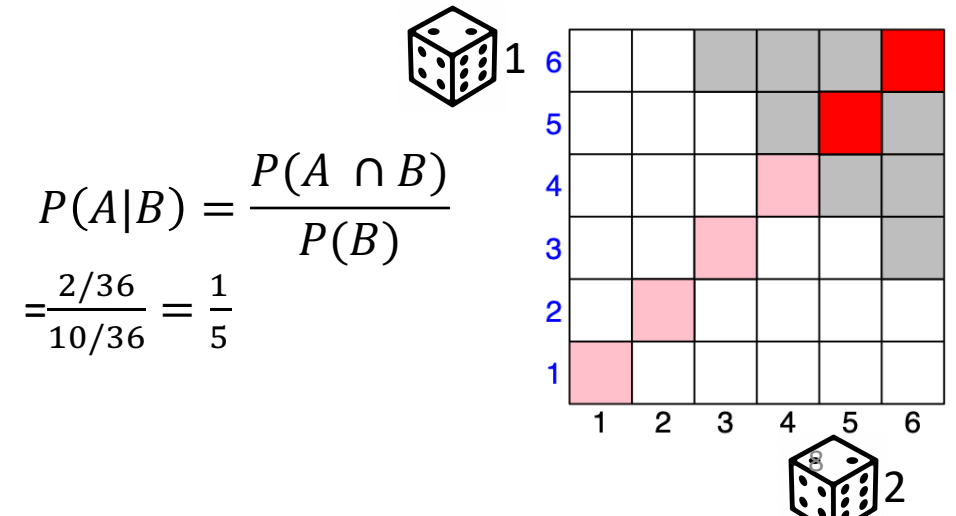
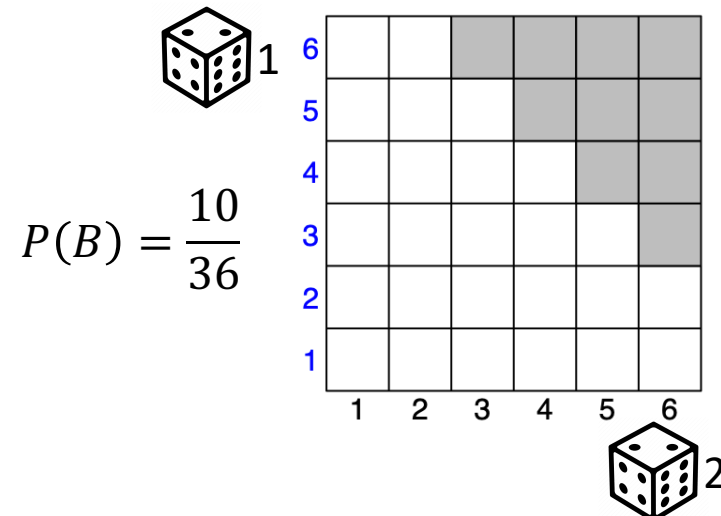
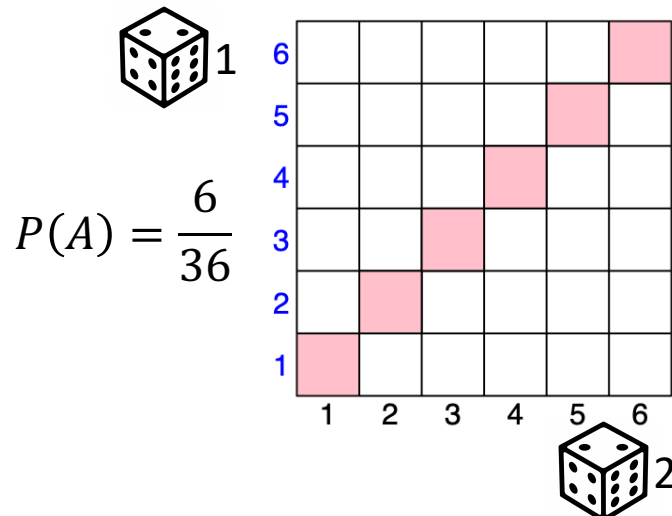
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Probability: Conditional probability and independence

The conditional probability of any event A given B is

$$P(A|B) = \frac{P(A \cap B)}{P(B)}$$

In other words, $P(A|B)$ is the probability measure of the event A after observing the occurrence of event B.

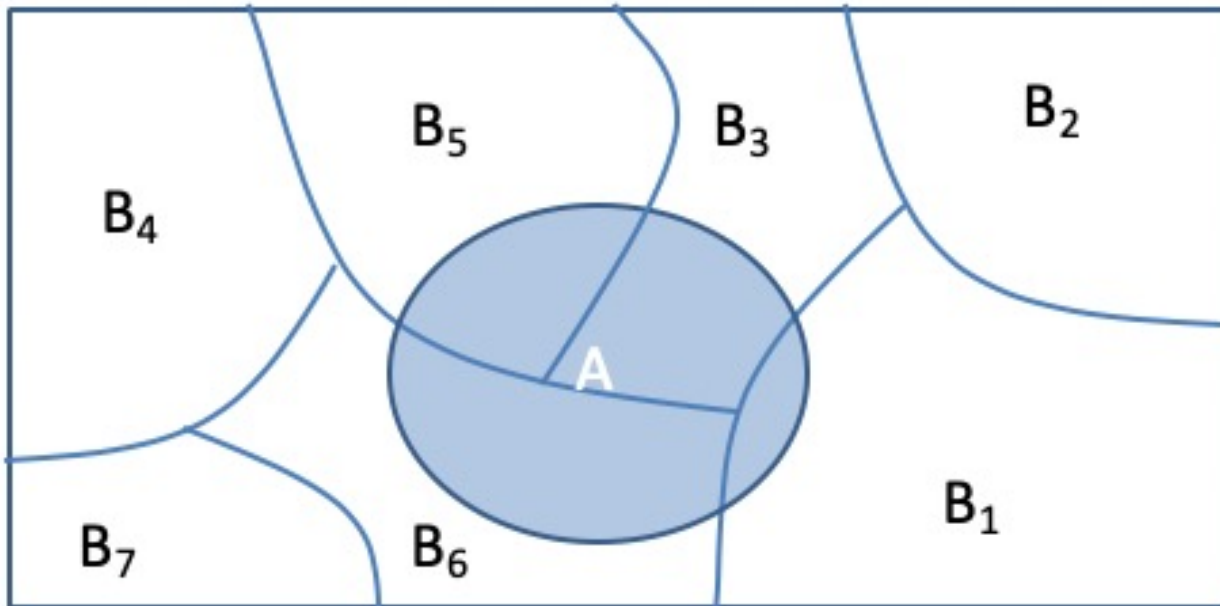
Two events are called **independent** if and only if

$$P(A \cap B) = P(A) \times P(B)$$

Independence is equivalent to saying that **observing B does not have any effect on the probability of A.**

Probability: Law of total probability

$$P(A) = \sum_{i=1}^n P(A \cap B_i) = \sum_{i=1}^n P(A|B_i)P(B_i)$$



Probability: Bayes' theorem

$$P(A|B) = \frac{P(B|A)P(A)}{P(B)}$$

Probability: Bayes' theorem

The diagram illustrates Bayes' theorem with the equation $P(A|B) = \frac{P(B|A)P(A)}{P(B)}$. The components are highlighted with colored boxes: $P(A|B)$ is in a yellow box, $P(B|A)$ is in an orange box, $P(A)$ is in a teal box, and $P(B)$ is in a pink box. Handwritten annotations in matching colors define each term: 'LIKELIHOOD' (orange) for $P(B|A)$, 'PRIOR' (teal) for $P(A)$, 'POSTERIOR' (yellow) for $P(A|B)$, and 'The probability of "B" being TRUE' (pink) for $P(B)$. Arrows point from each annotation to its corresponding term in the equation.

LIKELIHOOD
the probability of "B"
being TRUE given that "A" is TRUE

PRIOR
the probability of
"A" being TRUE

$$P(A|B) = \frac{P(B|A)P(A)}{P(B)}$$

POSTERIOR
the probability of "A"
being TRUE given that "B" is TRUE

The probability of "B" being TRUE

Topics

- Probability recap
- **Naïve Bayes**
- Gaussian process
- Bayesian hyperparameter optimization

Naïve Bayes: introduction

GAUSSIAN NAIVE BAYES CLASSIFIER

"Gaussian" because this is a normal distribution

This is our prior belief

$$P(\text{class} | \text{data}) = \frac{P(\text{data} | \text{class}) \times P(\text{class})}{P(\text{data})}$$

We don't calculate this in naive bayes classifiers

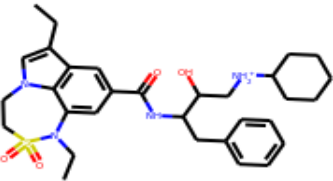
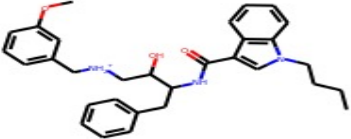
ChrisAlbon

Naïve Bayes: introduction

Problem statement:

- Given features X_1, X_2, \dots, X_n
- Given assumption about feature distribution
- Predict a label Y
- And give a level of [un]certainty about the prediction

Naïve Bayes: classification example [2 classes]

ROMol	MW	AMW	Sv	Se	Sp	H%	C%	N%		Class
	565.75	6.98	49.03	81.24	51.69	49.38	39.51	3.70	...	1
...										
	521.75	6.60	46.69	78.56	49.93	51.90	37.97	5.06	...	0

$X_1, \dots, X_n \in \mathbb{R}$ (features)

$Y \in \{1, 0\}$ (predict whether molecule active or inactive)

Naïve Bayes: Why “Bayes”?

Use Bayes Rule:

$$P(Y|X_1, \dots, X_n) = \frac{\overset{\text{Likelihood}}{P(X_1, \dots, X_n|Y)} \overset{\text{Prior}}{P(Y)}}{\underset{\text{Normalization Constant}}{P(X_1, \dots, X_n)}}$$

- Why did this help?

Well, we think that we might be able to specify how features are “generated” by the class label

Naïve Bayes: Why “Bayes”? [cont.]

For our example

$$P(Y = 0|X_1, \dots, X_n) = \frac{P(X_1, \dots, X_n|Y=0)P(Y=0)}{P(X_1, \dots, X_n)} = \frac{P(X_1, \dots, X_n|Y=0)P(Y=0)}{P(X_1, \dots, X_n|Y=1)P(Y=1) + P(X_1, \dots, X_n|Y=0)P(Y=0)}$$

$$P(Y = 1|X_1, \dots, X_n) = \frac{P(X_1, \dots, X_n|Y=1)P(Y=1)}{P(X_1, \dots, X_n)} = \frac{P(X_1, \dots, X_n|Y=1)P(Y=1)}{P(X_1, \dots, X_n|Y=1)P(Y=1) + P(X_1, \dots, X_n|Y=0)P(Y=0)}$$

To classify, we'll simply compute these two probabilities and predict based on which one is greater

Naïve Bayes: Why “Naïve”?

For the Bayes classifier, we need to “learn” two functions, the **likelihood** and the **prior**

But:

The diagram shows the Naïve Bayes formula with handwritten annotations in various colors:

- LIKELIHOOD** (orange): the probability of "B" being TRUE given that "A" is TRUE. An arrow points to the $P(B|A)$ term.
- PRIOR** (teal): the probability of "A" being TRUE. An arrow points to the $P(A)$ term.
- POSTERIOR** (green): the probability of "A" being TRUE given that "B" is TRUE. An arrow points to the $P(A|B)$ term.
- The probability of "B" being TRUE** (pink): An arrow points to the $P(B)$ term in the denominator.

$$P(A|B) = \frac{P(B|A) P(A)}{P(B)}$$

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Naïve Bayes: Why “Naïve”?

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But: How many parameters are required to specify the prior for our example?

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- POSTERIOR** (green): the probability of "A" being TRUE given that "B" is TRUE. An arrow points to the $P(A|B)$ term on the left side of the equation.
- The probability of "B" being TRUE** (pink): An arrow points to the $P(B)$ term in the denominator.

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Naïve Bayes: Why “Naïve”? [cont.]

For the Bayes classifier, we need to “learn” two functions, the **likelihood** and the **prior**

But: How many parameters are required to specify the prior for our example?

$$P(Y|X_1, \dots, X_n) = \frac{\overset{\text{Likelihood}}{P(X_1, \dots, X_n|Y)} \overset{\text{Prior}}{P(Y)}}{\underset{\text{Normalization Constant}}{P(X_1, \dots, X_n)}}$$

of parameters for modeling $P(X_1, \dots, X_n | Y)$: **$2(2^n - 1)$**

Naïve Bayes: Why “Naïve”? [cont. 2]

The ***Naïve Bayes Assumption***:

Assume that all features are *conditionally independent* **given the class label Y**

$$P(X_1, \dots, X_n | Y) = \prod_{i=1}^n P(X_i | Y)$$

$$P(Y | X_1, \dots, X_n) = \frac{\overset{\text{Likelihood}}{P(X_1, \dots, X_n | Y)} \overset{\text{Prior}}{P(Y)}}{\underset{\text{Normalization Constant}}{P(X_1, \dots, X_n)}}$$

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Naïve Bayes: Why “Naïve”? [And why 2n]

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of parameters for modeling $P(X_1 | Y), \dots, P(X_n | Y)$: **2n**

Gaussian naïve Bayes

assumption: features distributed normally :

$$p(x = v \mid C_k) = \frac{1}{\sqrt{2\pi\sigma_k^2}} e^{-\frac{(v-\mu_k)^2}{2\sigma_k^2}}$$

Naïve Bayes: Training

Training in Naïve Bayes is **easy** :

we have to select **feature distribution** (aka likelihood function)
[usually Gaussian distribution]

and **estimate parameters** of model of feature distribution

This corresponds to Maximum Likelihood estimation of model parameters.

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Is there other than Gaussian Naïve Bayes?

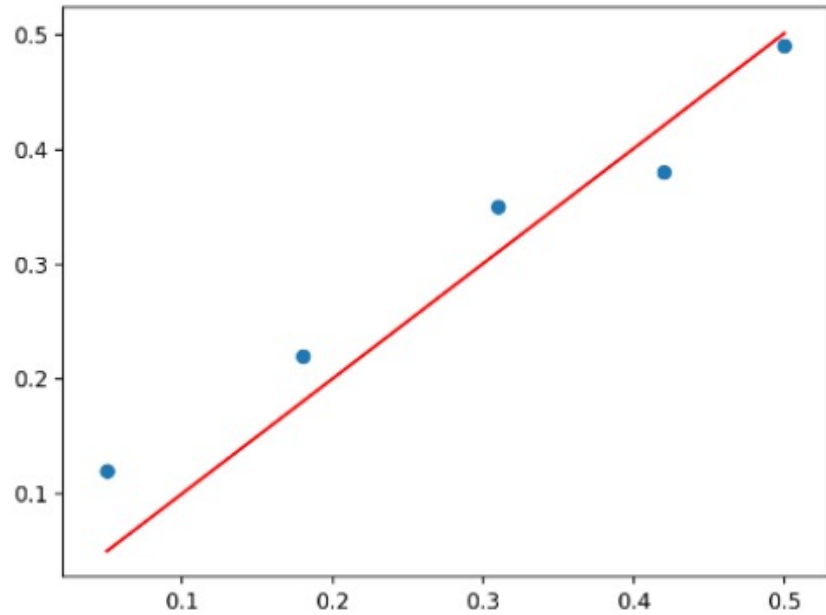
Sure, as a likelihood function, we could select different distributions.

In Scikit-learn there is implementation for Gaussian, Multinomial, Bernoulli etc.

Topics

- Probability recap
- Naïve Bayes
- **Gaussian process**
- Bayesian hyperparameter optimization

Gaussian Process: ML perspective



Gaussian Process: Definition

Continuous stochastic process — **random functions** — a set of random variables indexed by a continuous variable: $f(x)$

Set of 'inputs' $\mathbf{X} = \{x_1, x_2, \dots, x_N\}$; corresponding set of random function variables $\mathbf{f} = \{f_1, f_2, \dots, f_N\}$

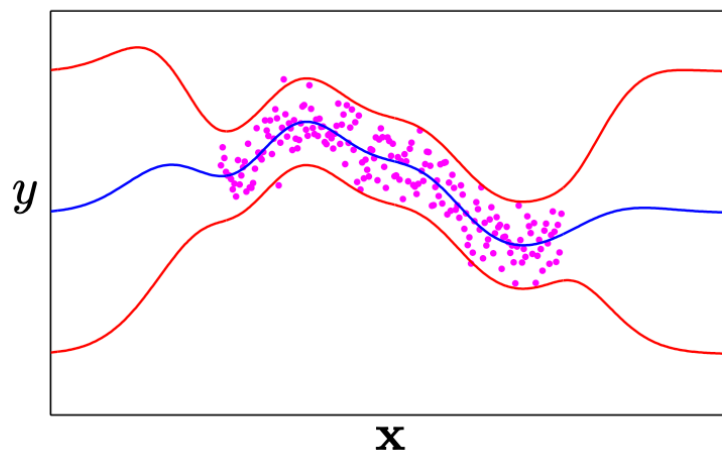
GP: Any set of function variables $\{f_n\}_{n=1}^N$ has joint (zero mean) Gaussian distribution:

$$p(\mathbf{f}|\mathbf{X}) = \mathcal{N}(\mathbf{0}, \mathbf{K})$$

Gaussian Process: ML perspective

Consider the problem of **nonlinear regression**:

You want to learn a **function** f with **error bars** from **data** $\mathcal{D} = \{\mathbf{X}, \mathbf{y}\}$



A **Gaussian process** is a prior over functions $p(f)$ which can be used for Bayesian regression:

$$p(f|\mathcal{D}) = \frac{p(f)p(\mathcal{D}|f)}{p(\mathcal{D})}$$

Gaussian Process: vs Gaussian distributions

Gaussian distributions

$$\mathcal{N}(\mu, \Sigma)$$

Distribution over vectors.

Fully specified by a mean and covariance.

The position of the random variable in the vector plays the role of the index.

Gaussian processes

$$\mathcal{GP}(m(\mathbf{x}), k(\mathbf{x}, \mathbf{x}'))$$

Distribution over functions.

Fully specified by a mean function and covariance function.

The argument of the random function plays the role of the index.

Gaussian Process: covariance matrix

Gaussian processes are merely based on the good old Gaussian

$$\mathcal{N}(\mathbf{x} \mid \boldsymbol{\mu}, \mathbf{K}) = \frac{1}{\sqrt{|2\pi \mathbf{K}|}} \exp \left[-\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu})^T \mathbf{K}^{-1} (\mathbf{x} - \boldsymbol{\mu}) \right]$$

Covariance matrix or kernel matrix

Covariance matrix constructed from *covariance function*:

$$\mathbf{K}_{ij} = K(x_i, x_j)$$

Covariance function characterizes correlations between different points in the process:

$$K(x, x') = \mathcal{E}[f(x)f(x')]$$

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Gaussian Process: kernel examples

Kernels -- are a crucial ingredient of GPs which determine the shape of prior and posterior of the GP.

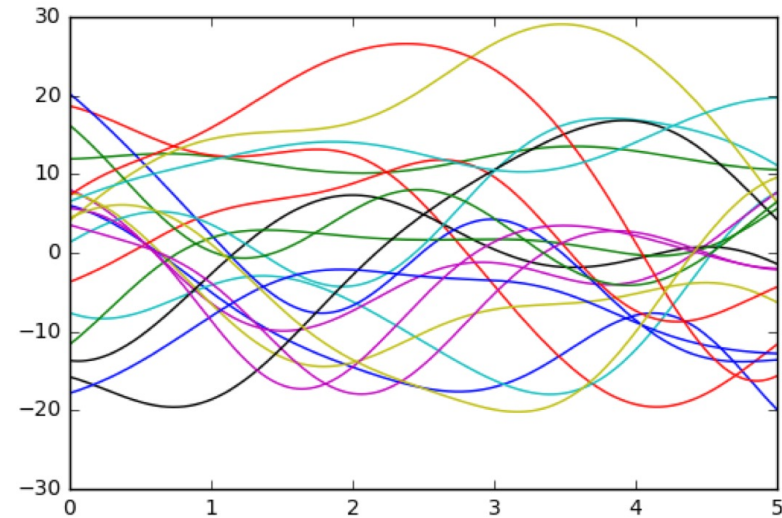
They **encode the assumptions** on the function being learned by defining the “similarity” of two datapoints combined with the assumption that similar datapoints should have similar target values.

Squared exponential (SE), RBF

$$K(x, x') = \sigma_0^2 \exp \left[-\frac{1}{2} \left(\frac{x - x'}{\lambda} \right)^2 \right]$$

Intuition: function variables close in input space are highly correlated, whilst those far away are uncorrelated

λ, σ_0 — hyperparameters. λ : lengthscale, σ_0 : amplitude



Gaussian Process: kernel examples

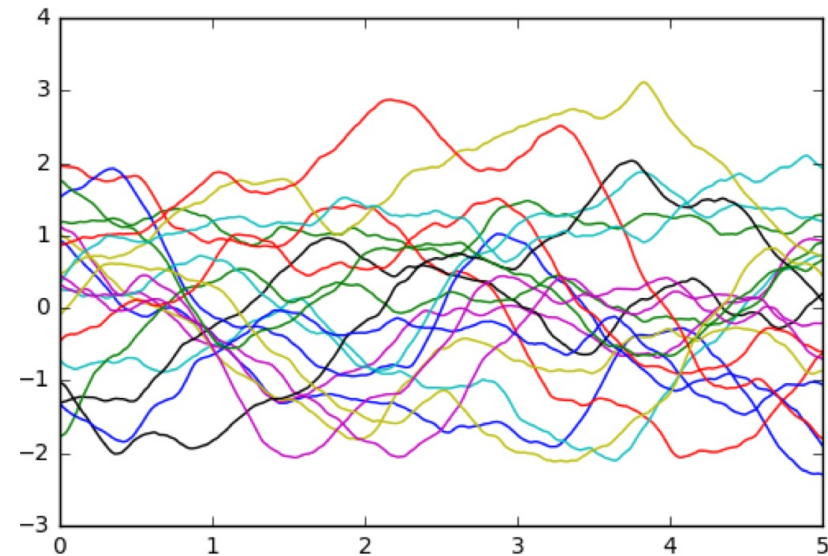
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Matérn class [usually $\nu = 1/2$ or $3/2$]

$$K(x, x') = \frac{2^{1-\nu}}{\Gamma(\nu)} \left(\frac{\sqrt{2\nu}|x - x'|}{\lambda} \right)^\nu K_\nu \left(\frac{\sqrt{2\nu}|x - x'|}{\lambda} \right)$$

where K_ν is a modified Bessel function.



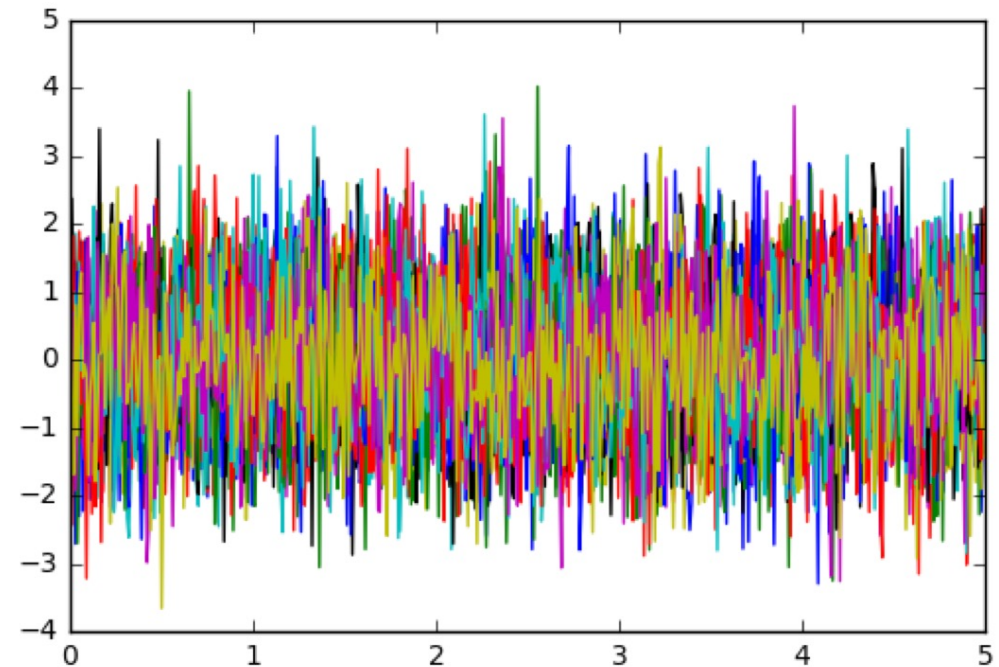
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White noise

$$k(x, x') = \begin{cases} 1 & \text{if } x = x' \\ 0 & \text{otherwise} \end{cases}$$



Gaussian Process: most standard kernel

Kernels -- are a crucial ingredient of GPs which determine the shape of prior and posterior of the GP.

They **encode the assumptions** on the function being learned by defining the “similarity” of two datapoints combined with the assumption that similar datapoints should have similar target values.

We can make new kernels by summing or multiplying kernels

Most default kernel $K = \text{RBF} + \text{White_noise}$

Gaussian Process Regression: training

$$\mathcal{GP}(m(\mathbf{x}), k(\mathbf{x}, \mathbf{x}'))$$

- I) Select mean function $m(\mathbf{x})$ [usually constant]
- II) Select kernel function $k(\mathbf{x}, \mathbf{x}')$ [usually RBF + White_noise]

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$$k(x, x') = \sigma_y^2 \exp\left(-\frac{(x - x')^2}{2\ell^2}\right) + \sigma_n^2 \delta_{ii'}$$

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- III) Find parameters [fit GP to data]

Gaussian Process: as a regression [in 2 words]

Covariance function defines the properties in the function space.

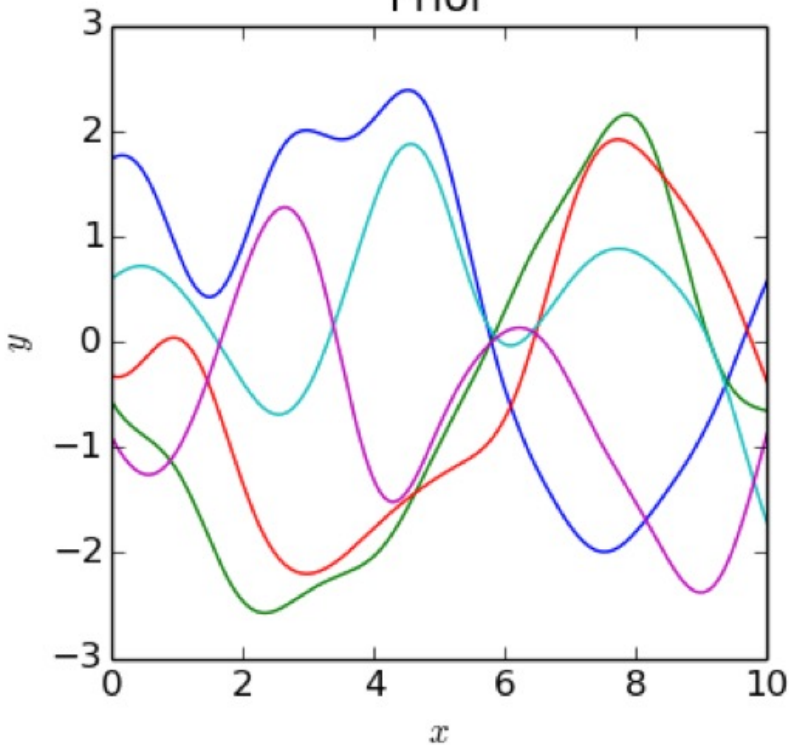
Data points "anchor" the function as specific locations.

Gaussian Process: as a regression

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Prior

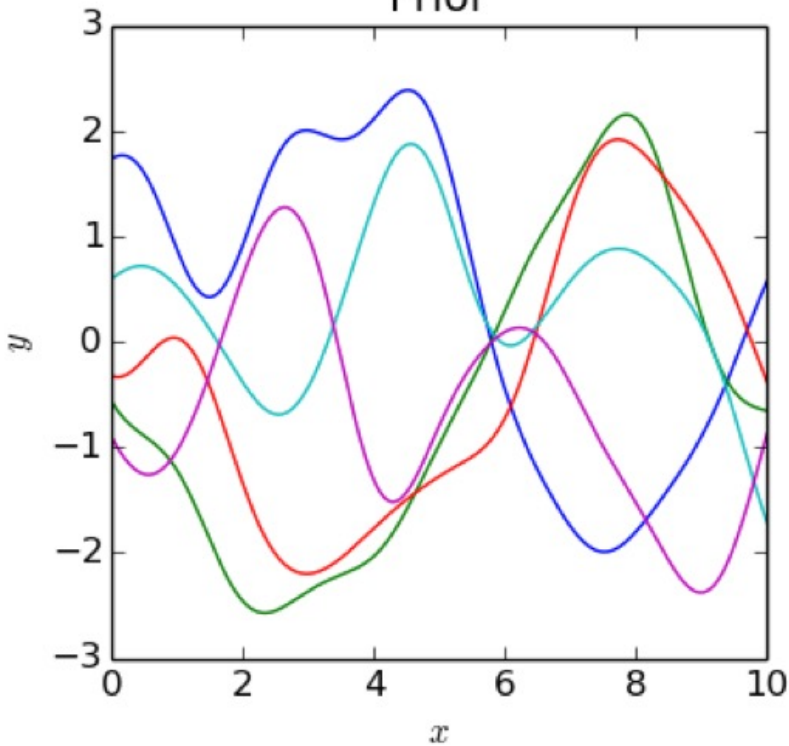


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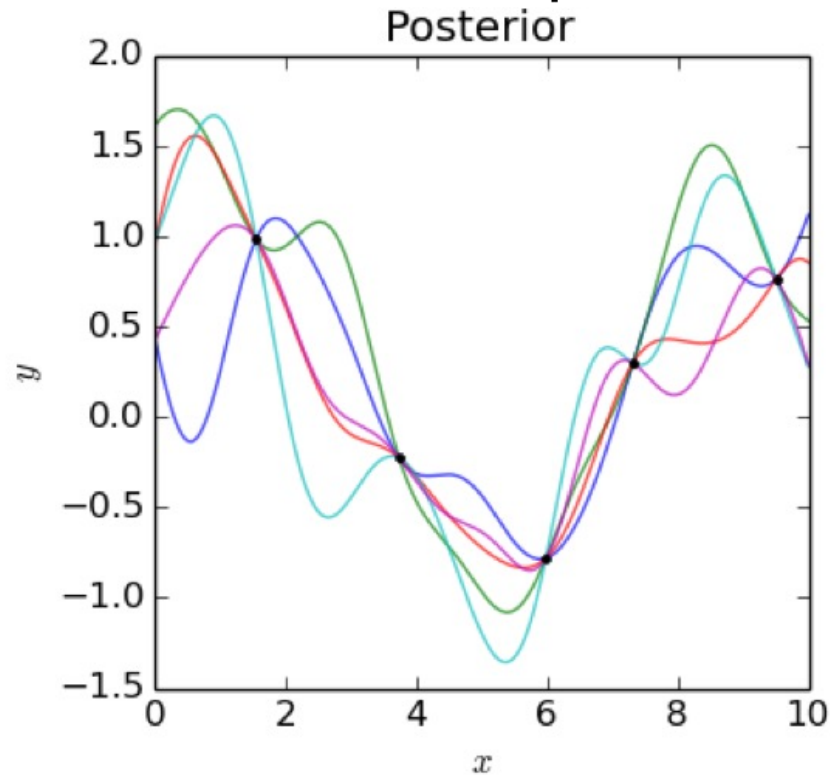
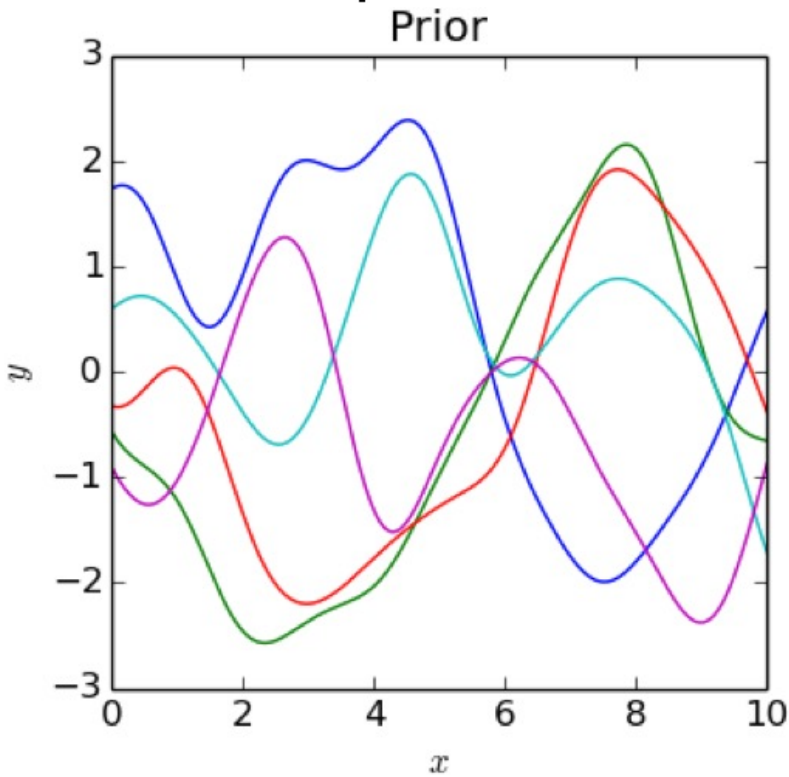


I) So we are fitting an **infinite number of functions** to the data

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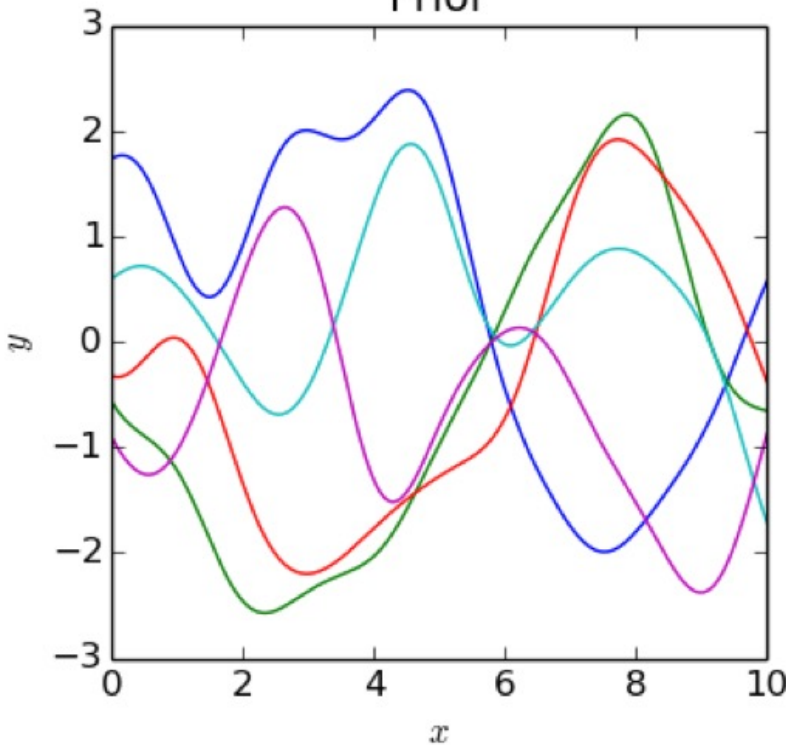
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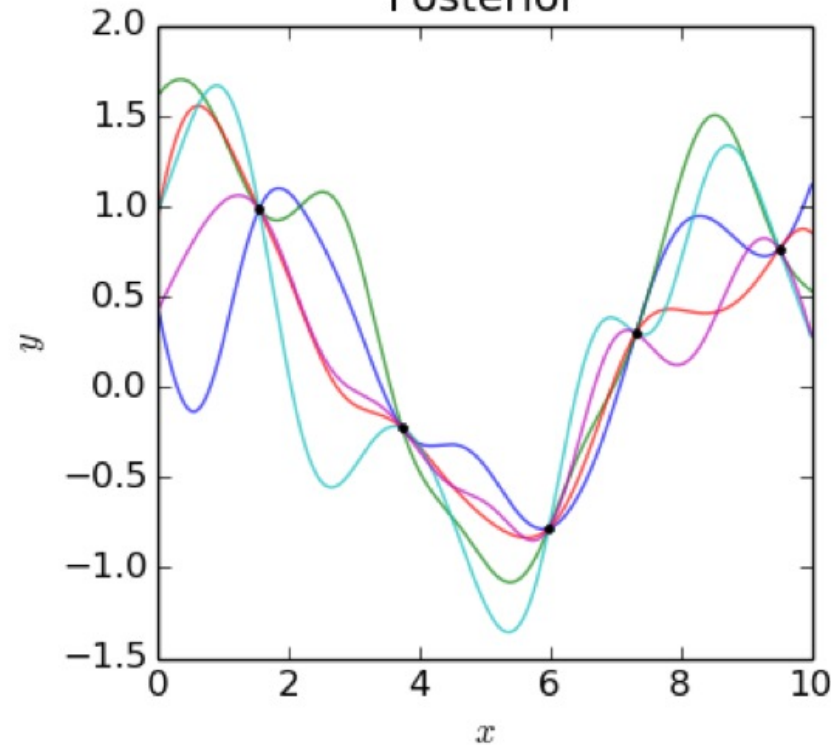
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Prior



Posterior



I) So we are fitting an **infinite number of functions** to the data

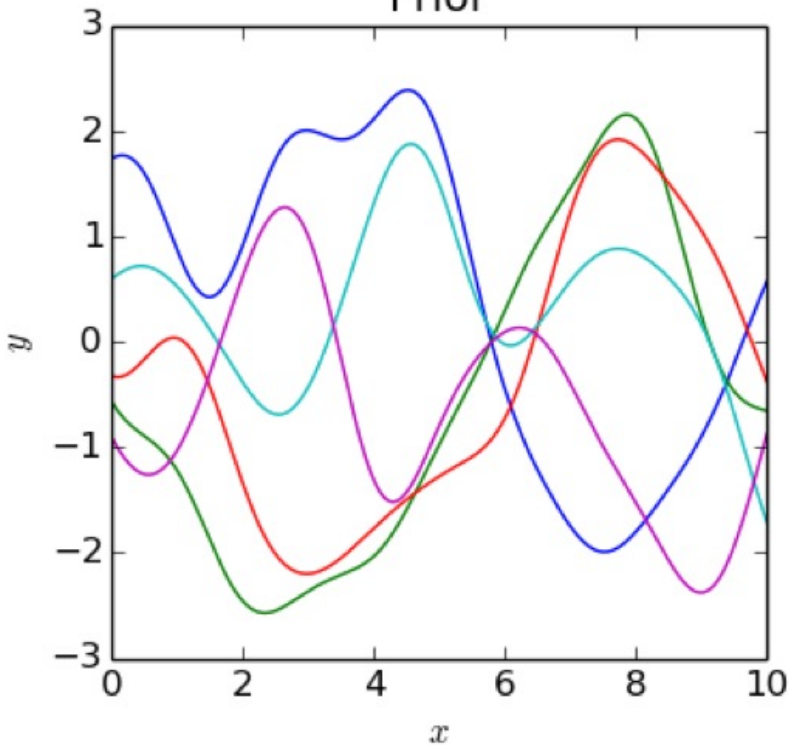
II) and then **sampling these functions from GP** and calculate values of those **sampled functions for particular input**

Gaussian Process: as a regression

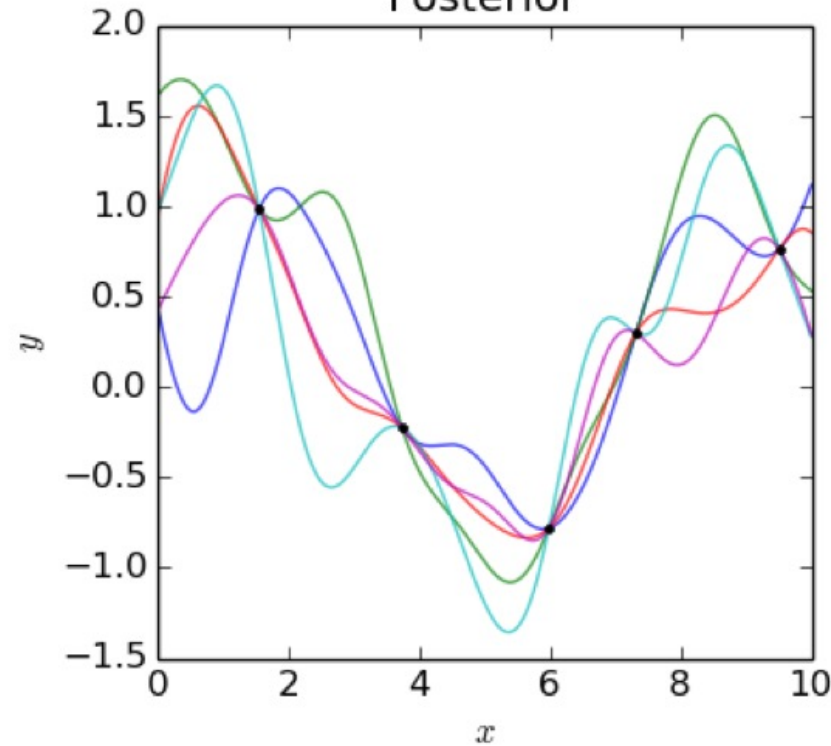
Covariance function defines the properties in the function space.

Data points "anchor" the function as specific locations.

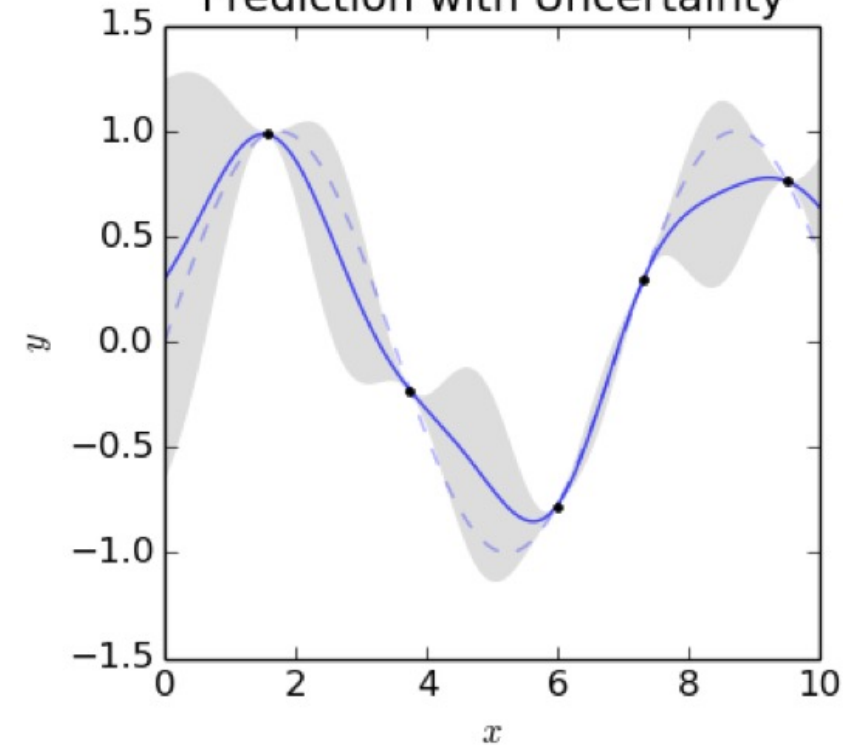
Prior



Posterior



Prediction with Uncertainty



I) So we are fitting an **infinite number of functions** to the data

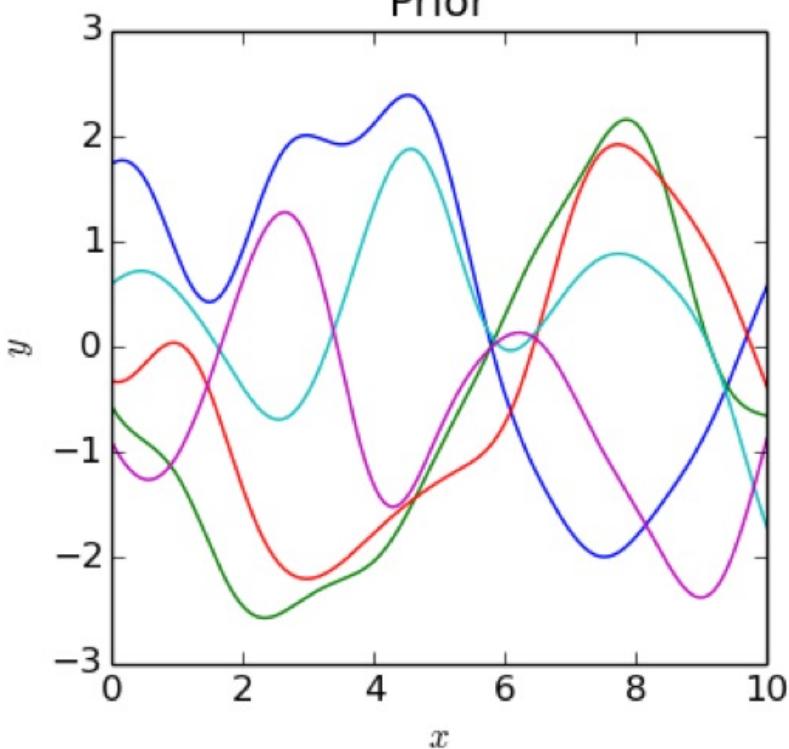
II) and then **sampling these functions from GP** and calculate values of those **sampled functions for particular input**

Gaussian Process: as a regression

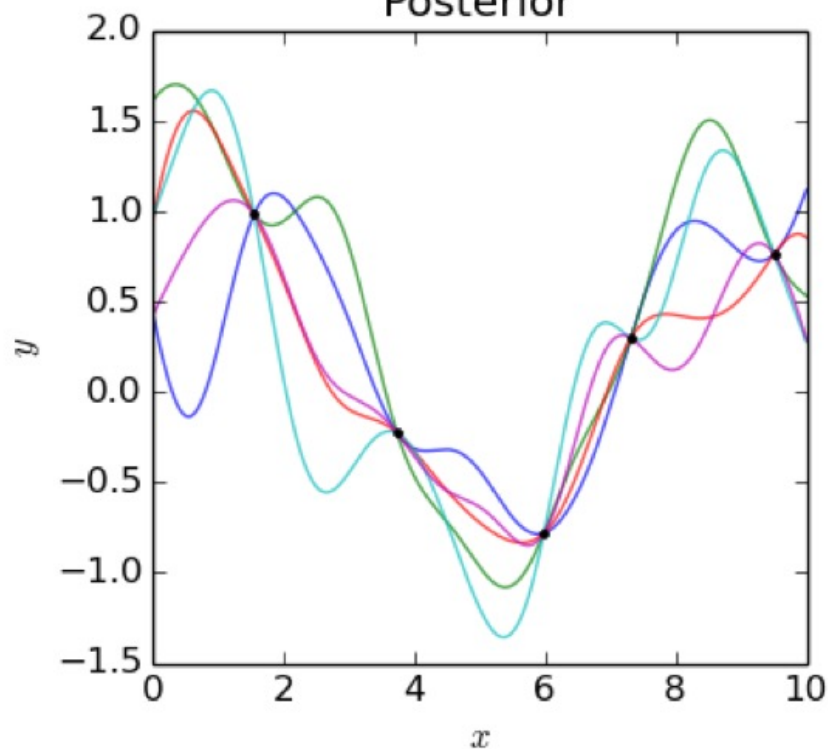
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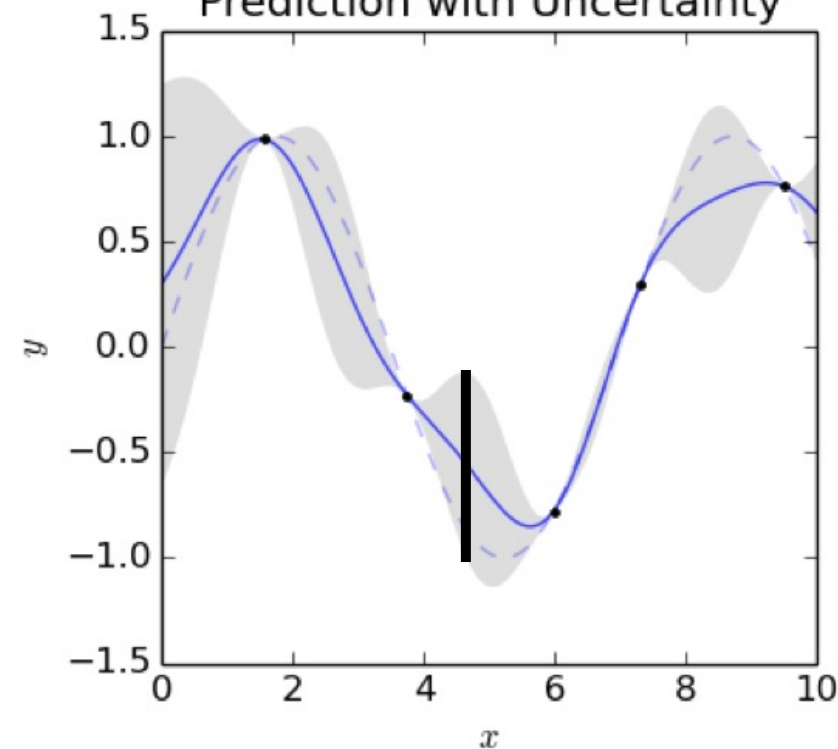
Prior



Posterior



Prediction with Uncertainty



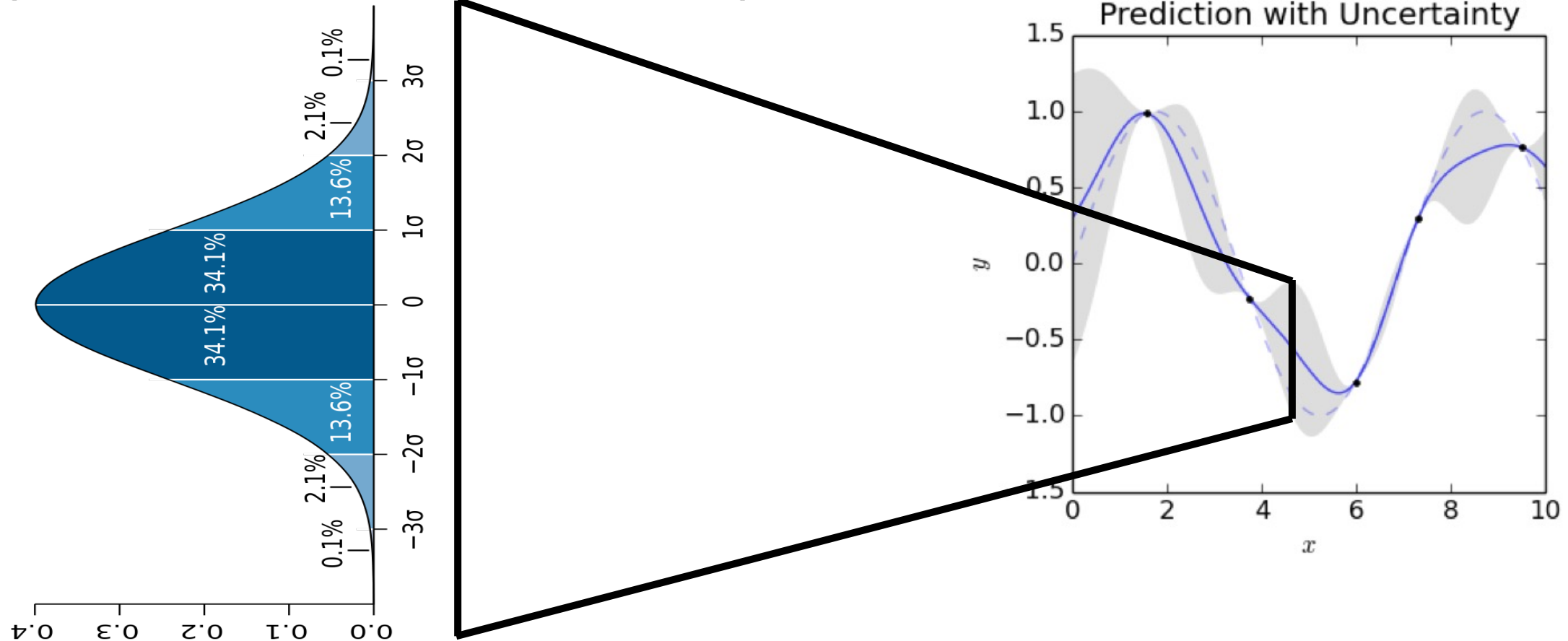
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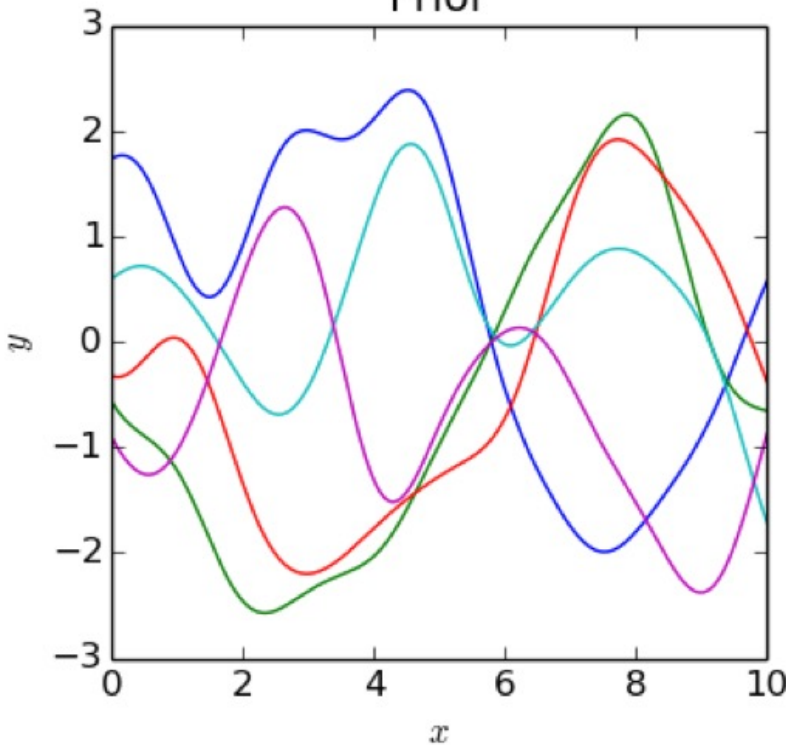


Gaussian Process: as a regression

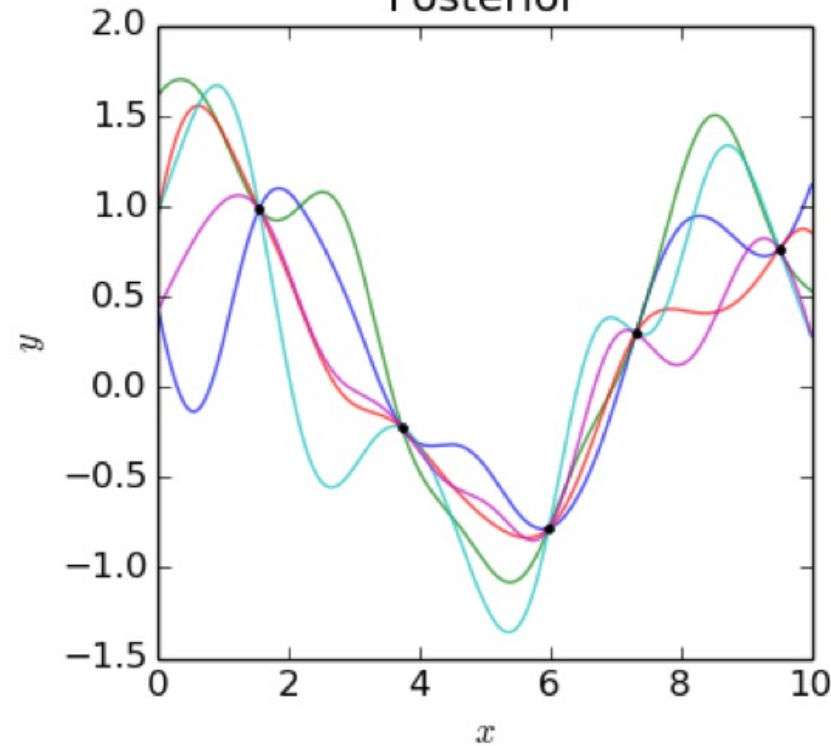
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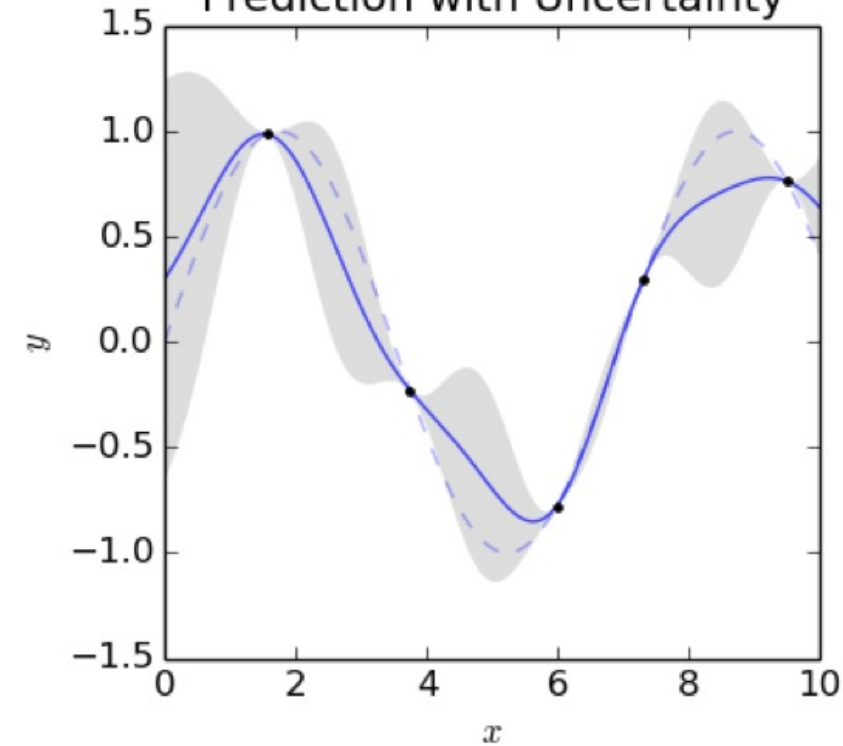
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Prediction with Uncertainty



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Topics

- Probability recap
- Naïve Bayes
- Gaussian process
- **Bayesian hyperparameter optimization**

Bayesian hyperparameter optimization

If you have an **experiment** with **set of hyperparameters** and some **objective function**, then you can optimize those hyperparameters with GP.

An experiment could be:

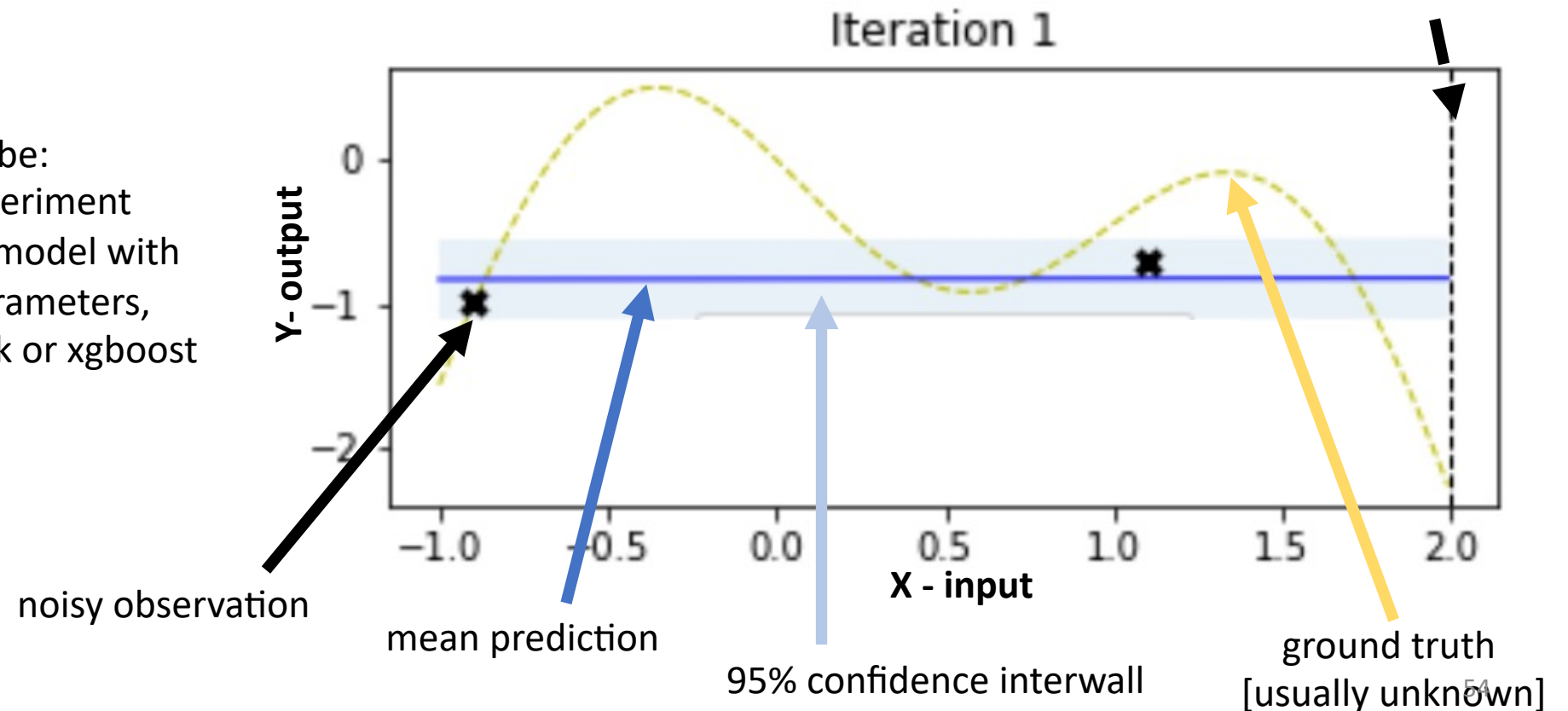
- Actual wet-lab experiment
- Machine learning model with big set of hyperparameters, like neural network or xgboost

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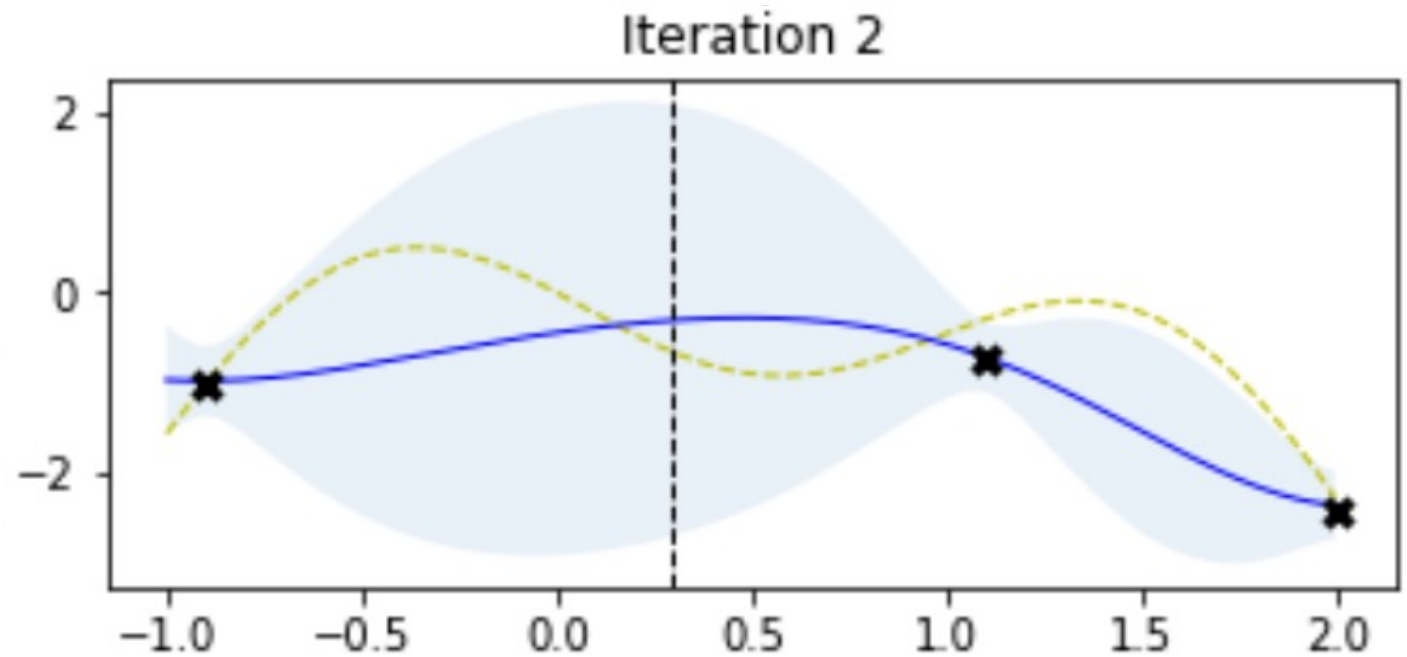


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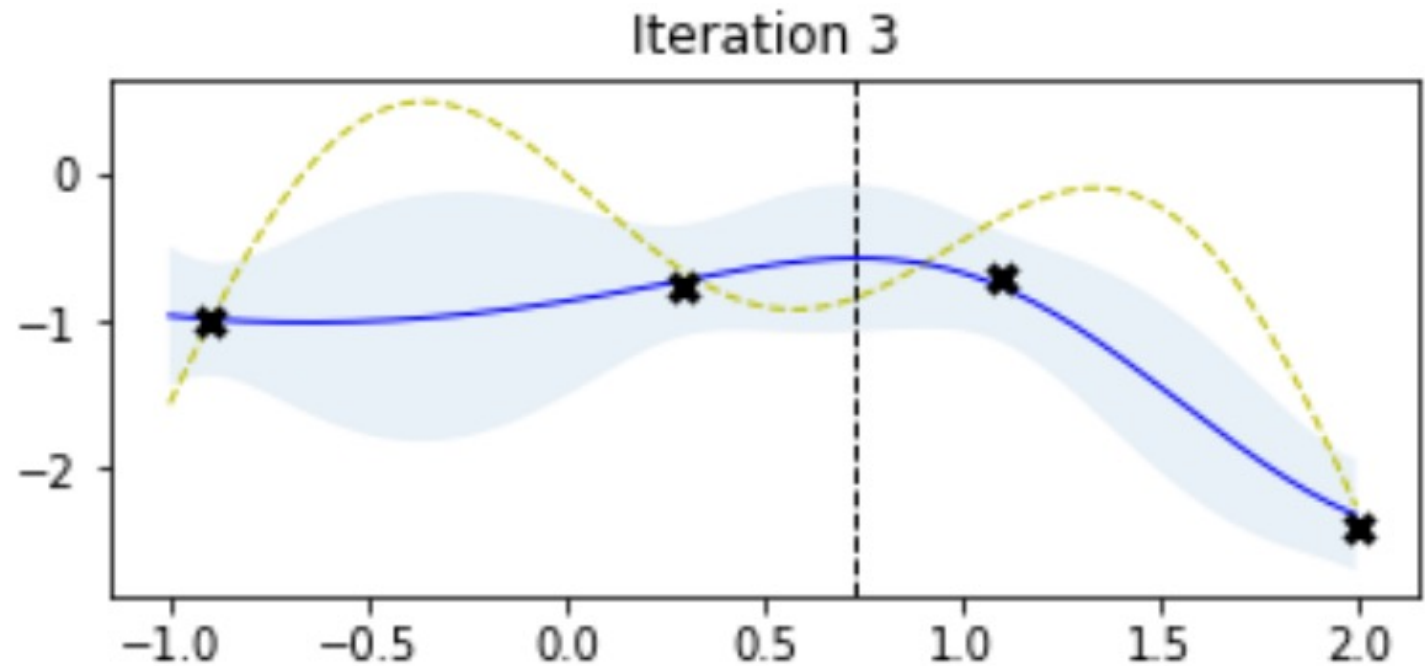


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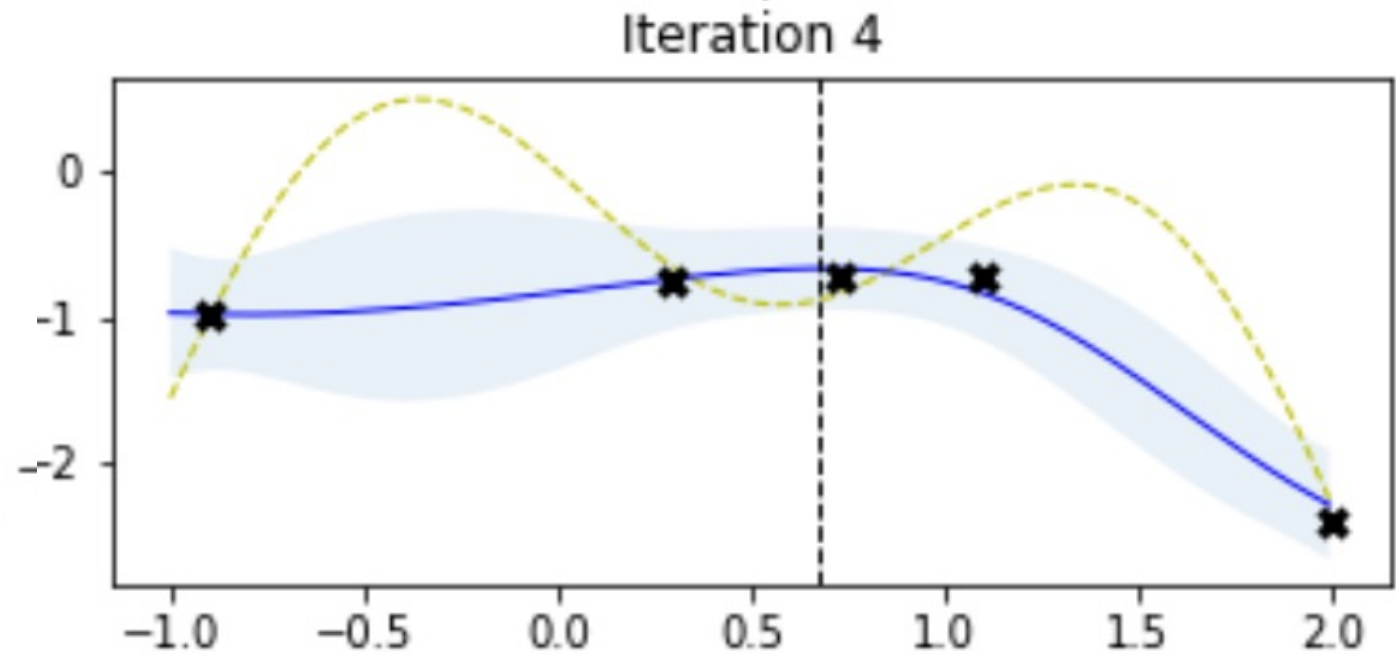


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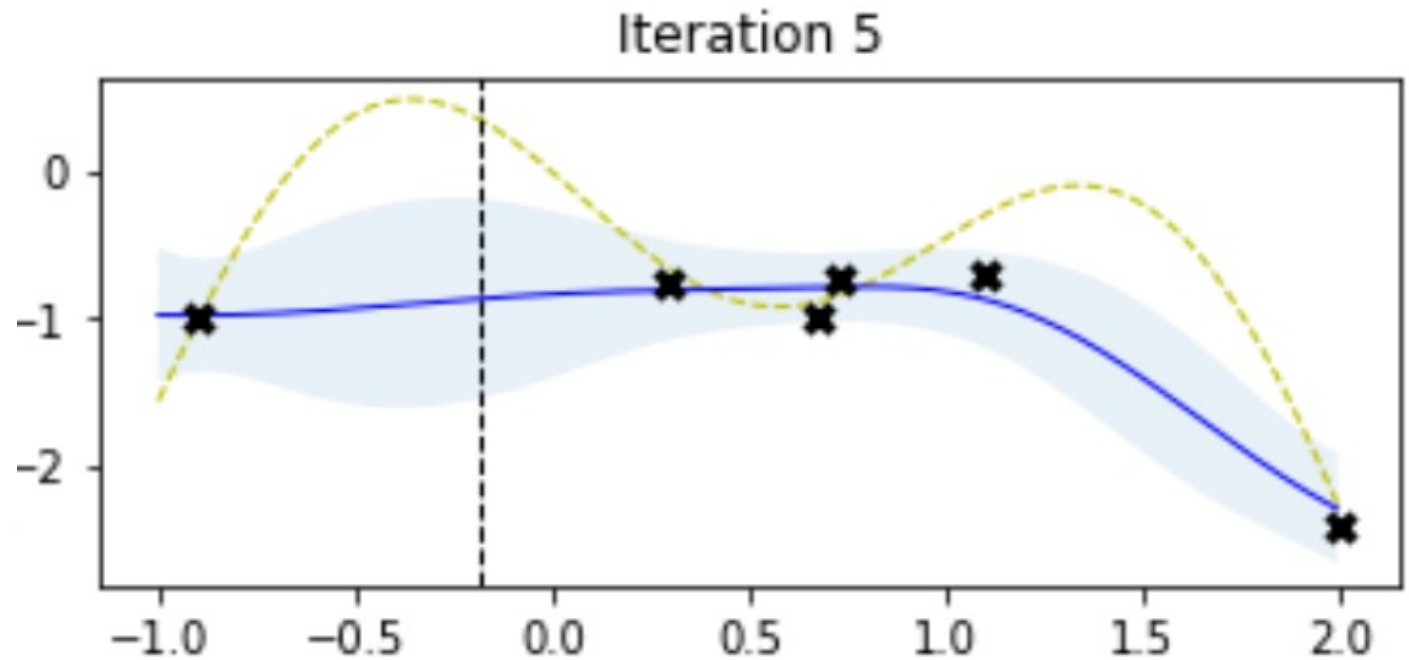


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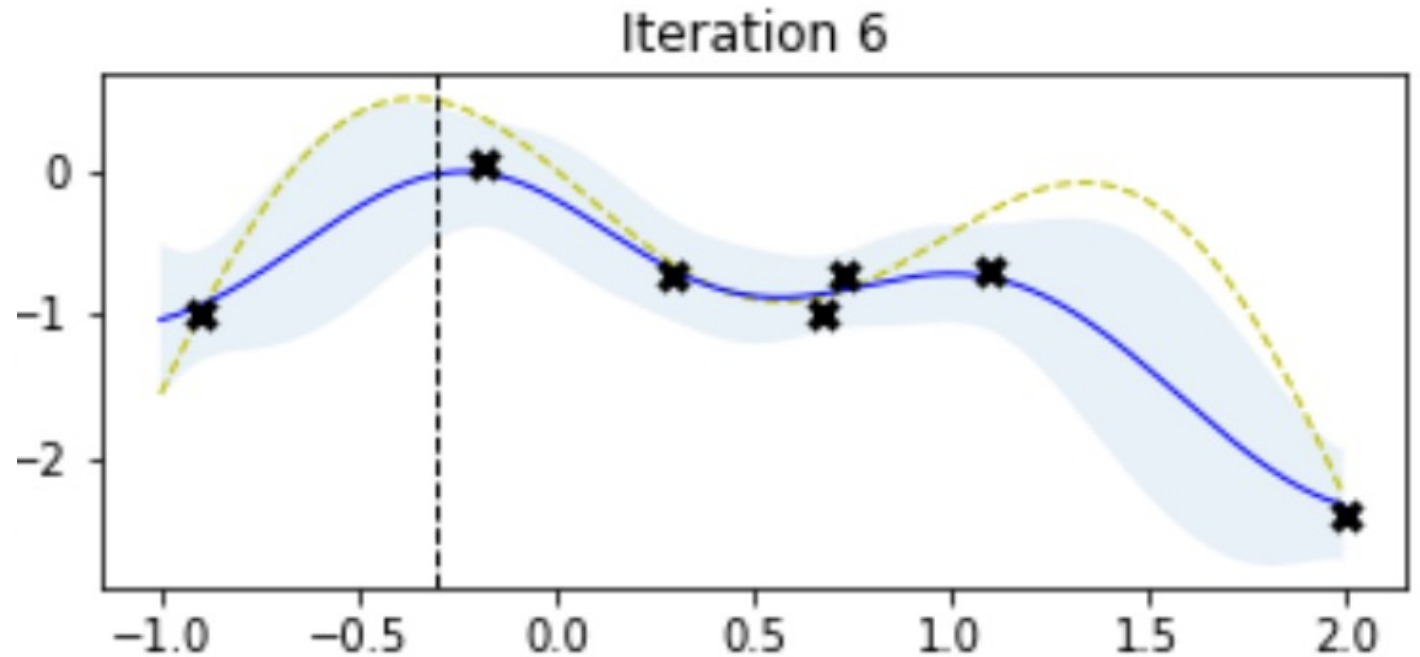


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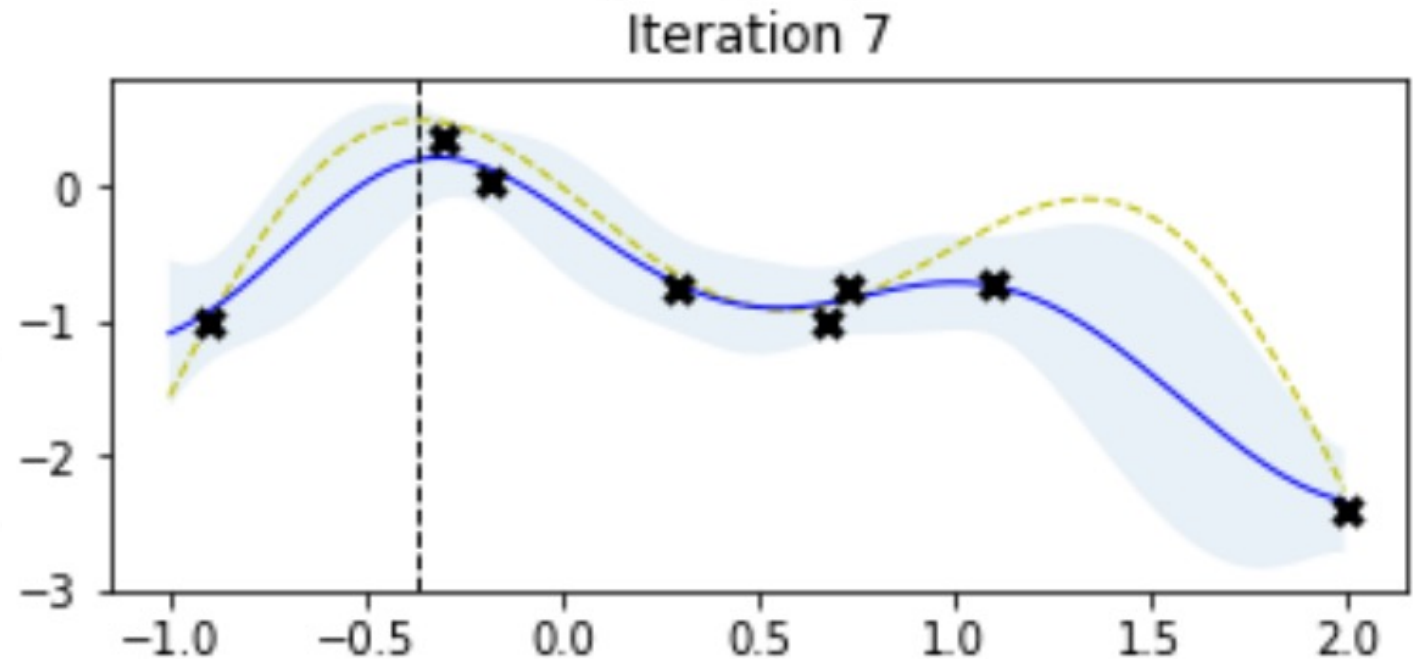


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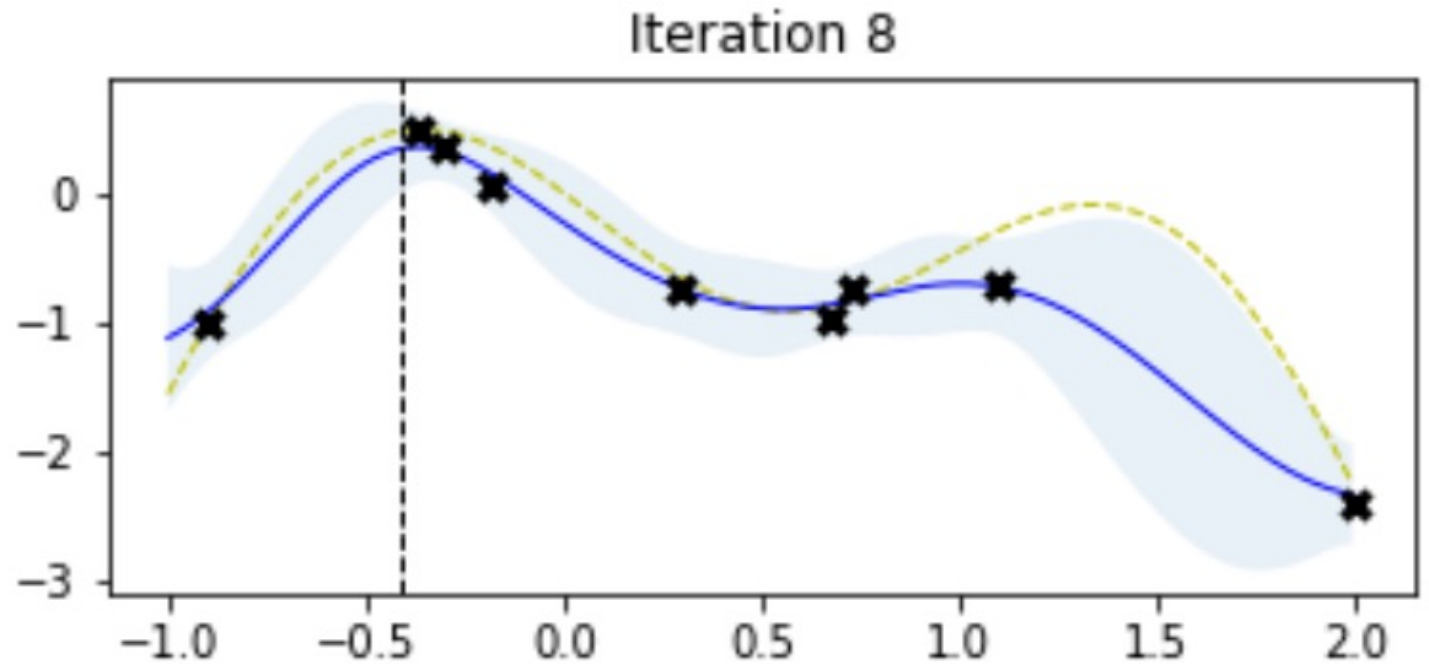


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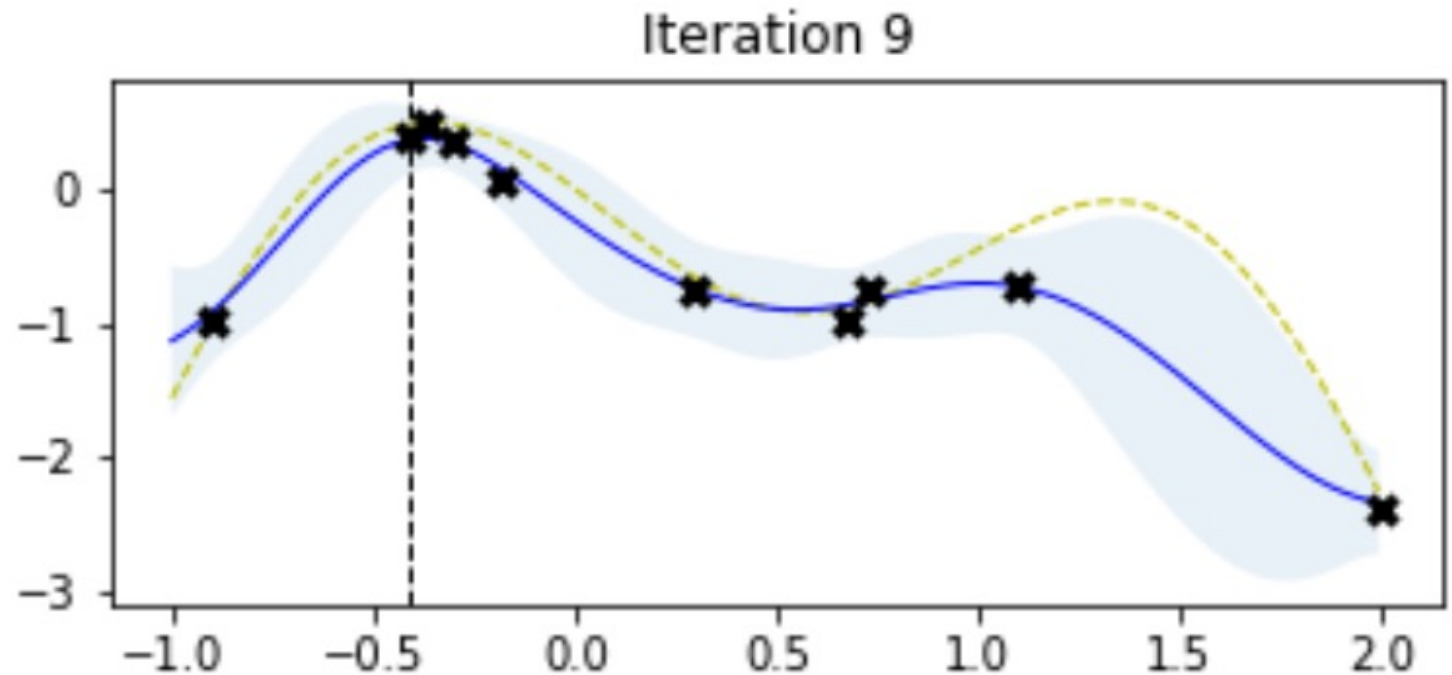


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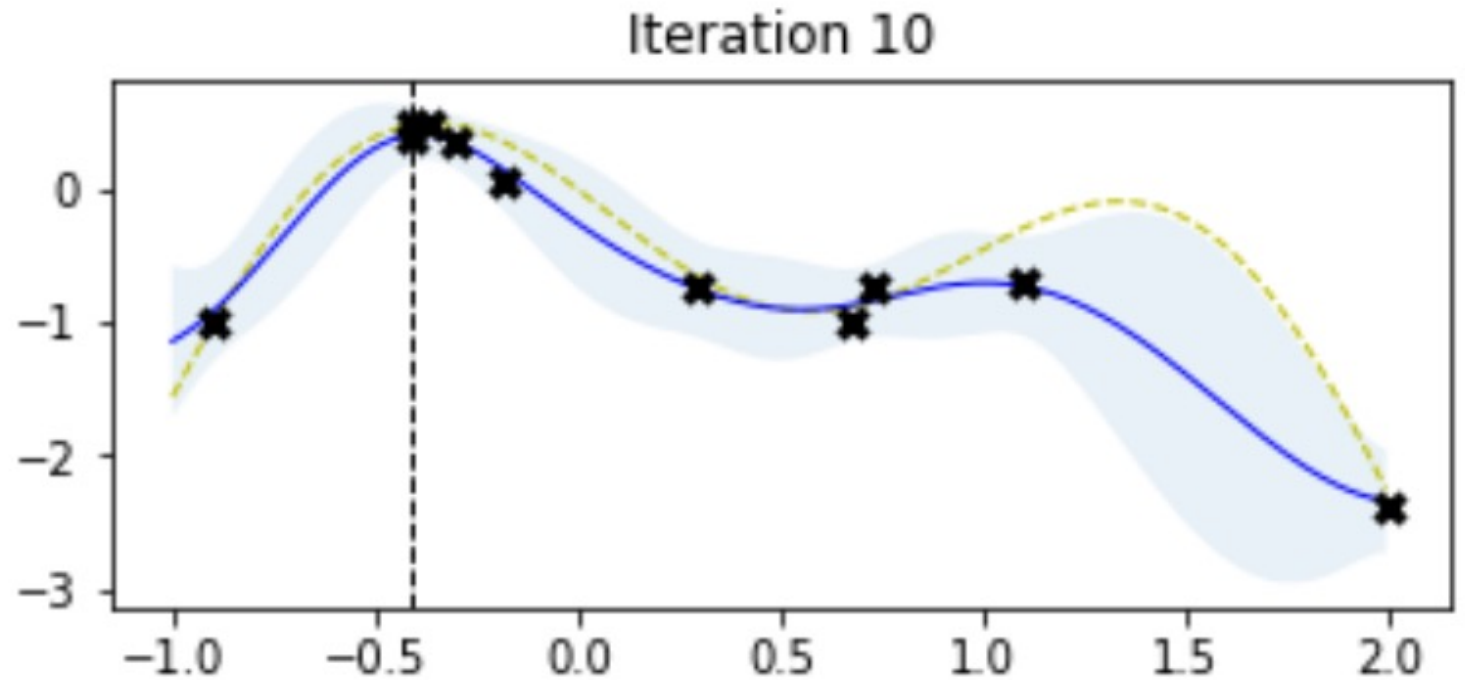


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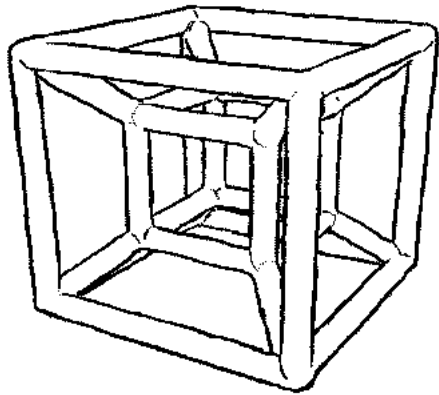
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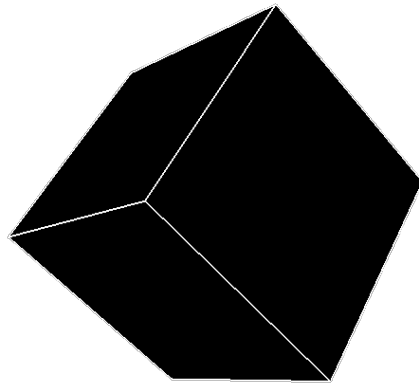
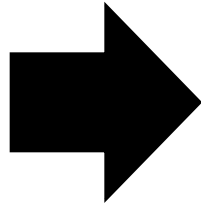


Bayesian hyperparameter optimization

Graduate student perspective



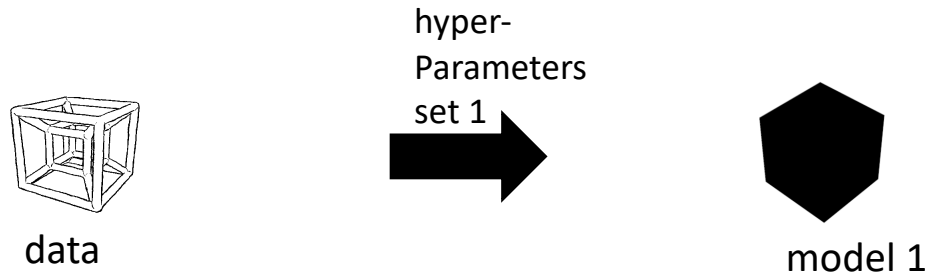
multidimensional
data



Machine learning
model

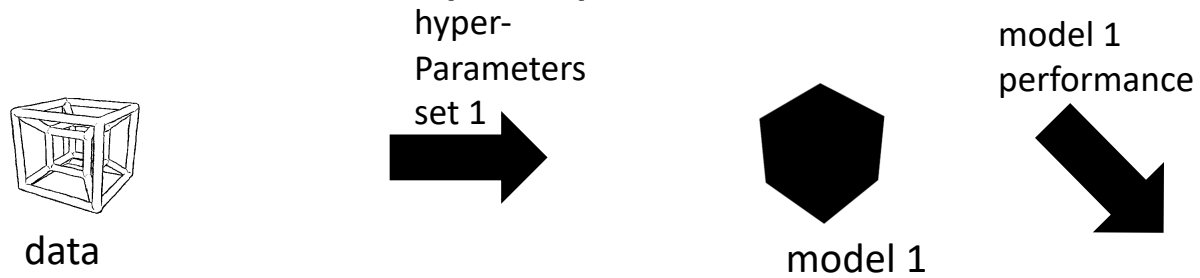
Bayesian hyperparameter optimization

Graduate student perspective



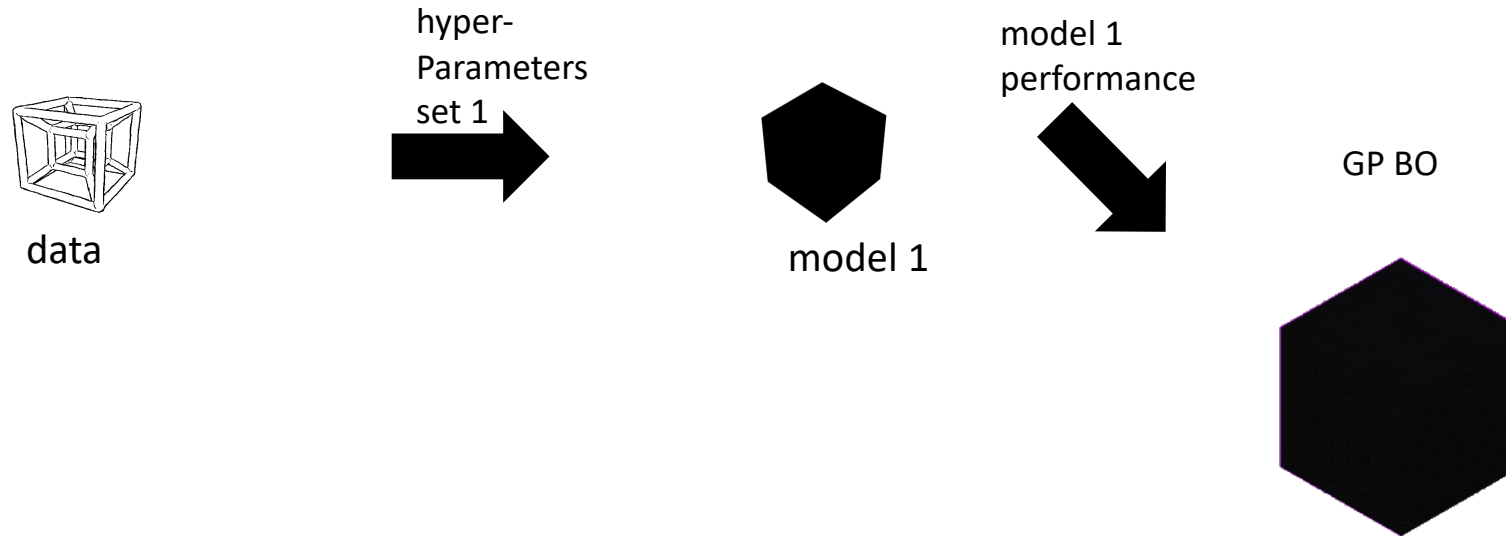
Bayesian hyperparameter optimization

Graduate student perspective



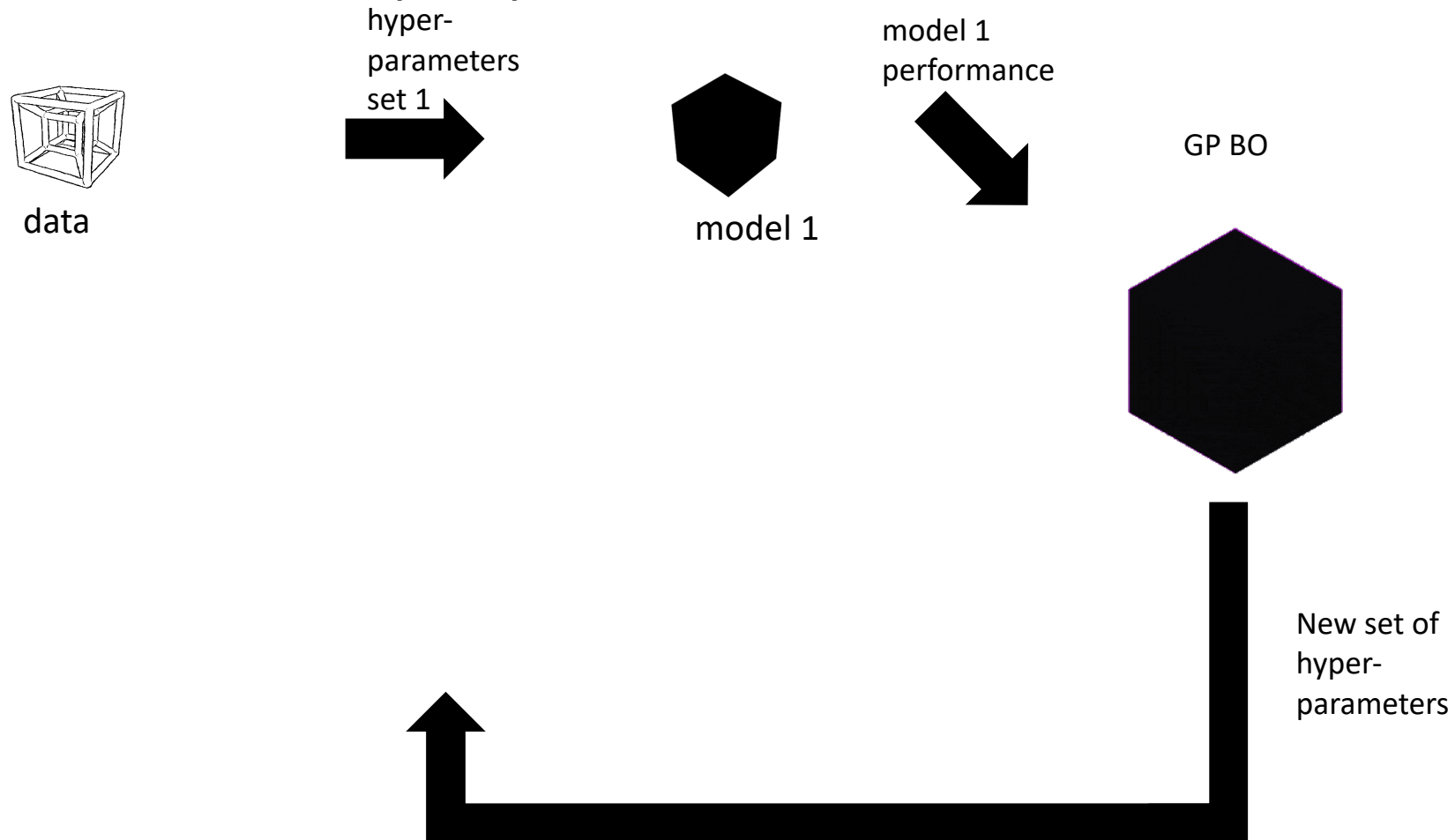
Bayesian hyperparameter optimization

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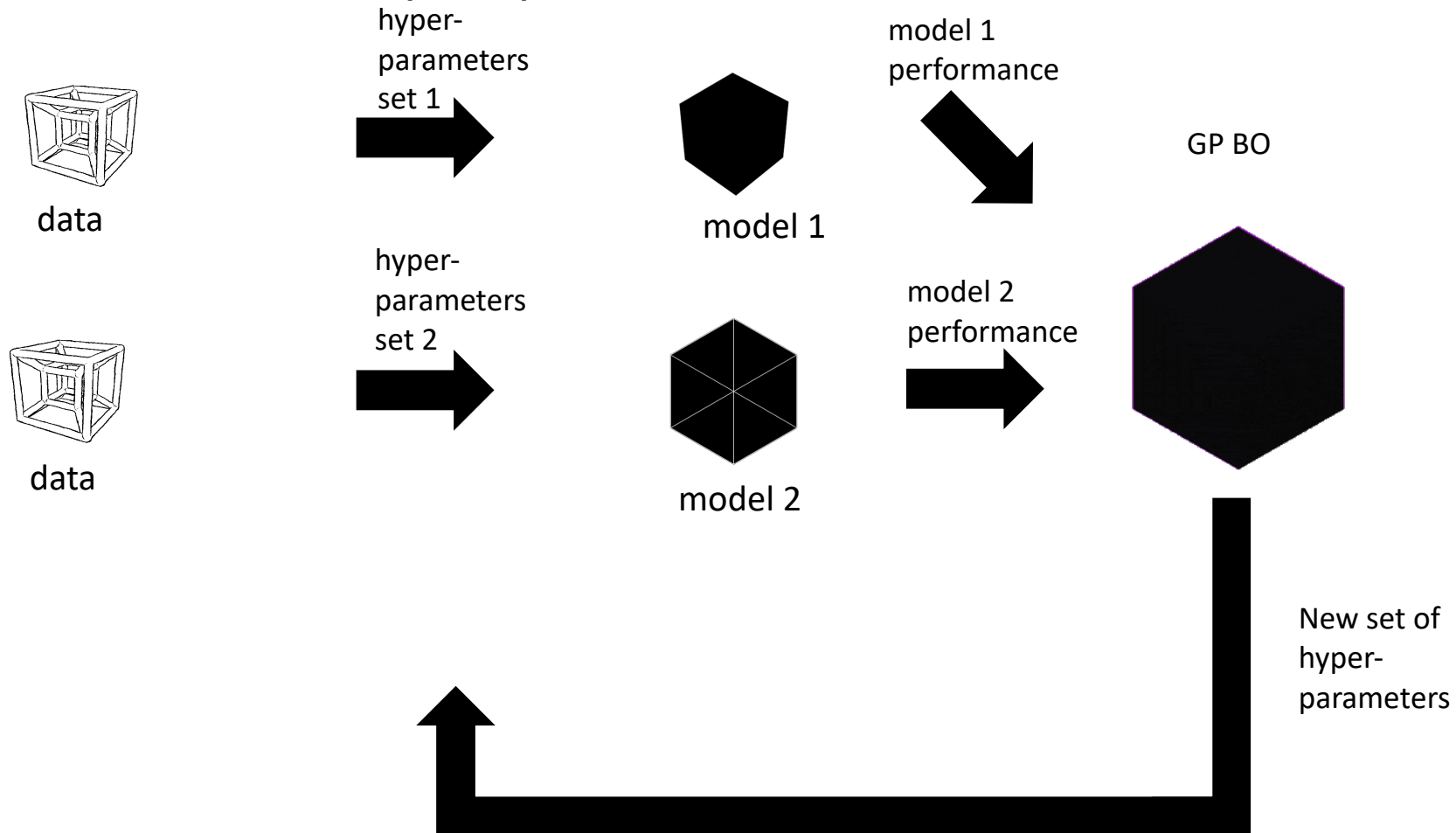
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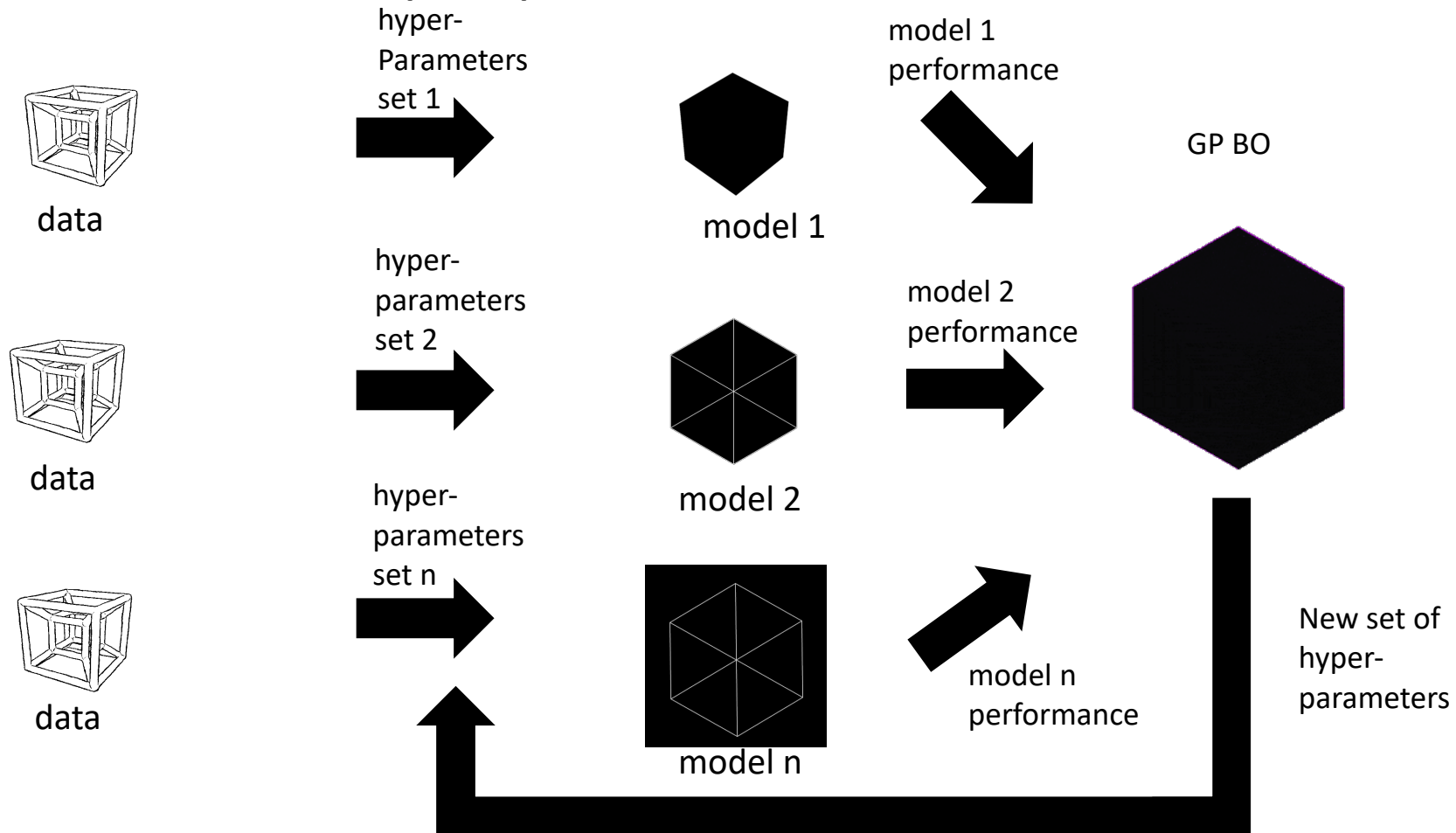
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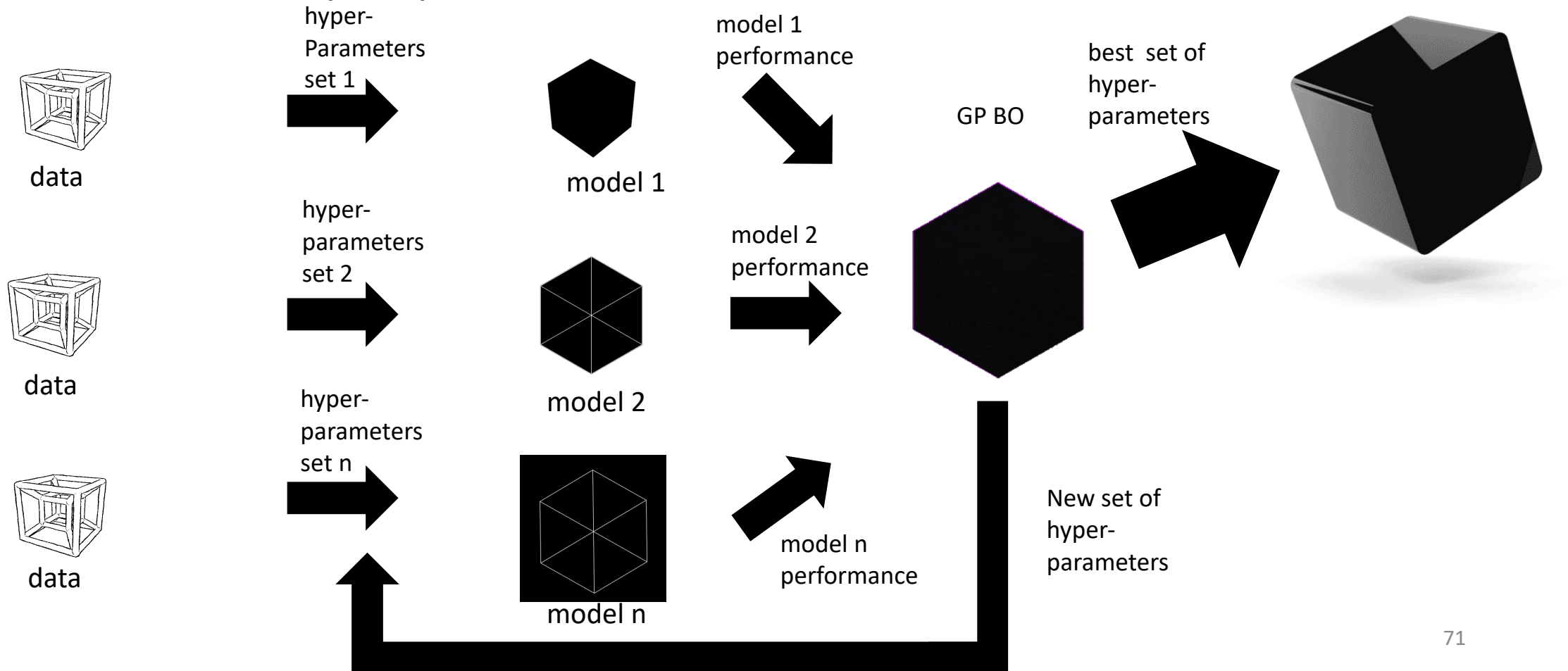
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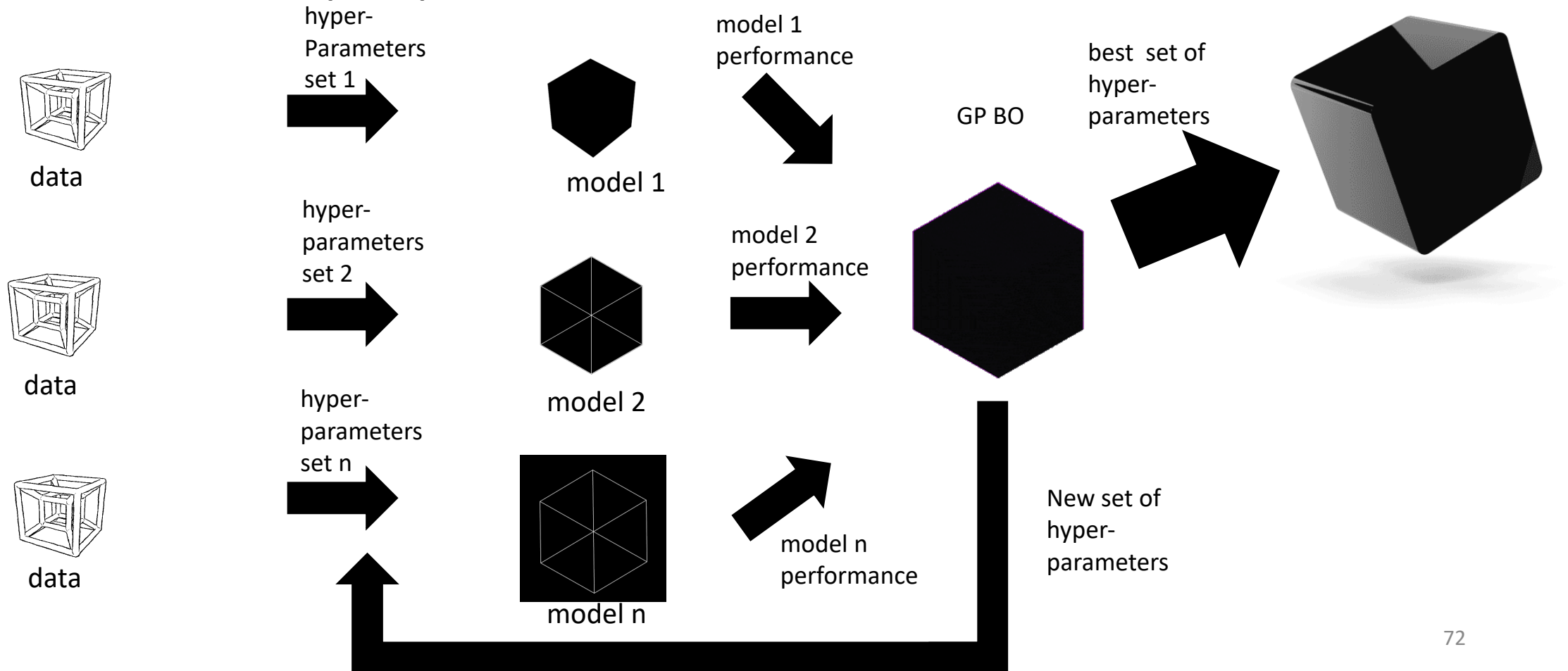
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That's all for today!

Thank you all for your attention!

