

# Bayesian Optimization

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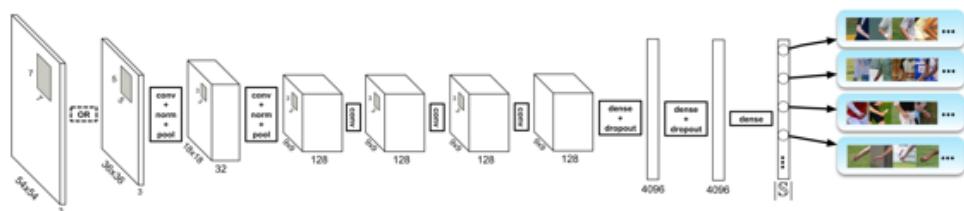
- Brochu et al.: *A Tutorial on Bayesian Optimization of Expensive Cost Functions, with Application to Active User Modeling and Hierarchical Reinforcement Learning*, arXiv:1012.2599, 2012
- Shahriari et al.: *Taking the Human Out of the Loop: A Review of Bayesian Optimization*, Proceedings of the IEEE, 2016

- Machine learning models are getting more and more complicated
  - ▶ Usually more parameters (e.g., deep neural networks)
- Non-convex and stochastic optimization methods have meta-parameters that are difficult to tune (learning rates, momentum parameters, ...)
  - ▶ Generally hard to apply modern techniques or reproduce results

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Goal: Automate the selection of critical meta-parameters  
(see also: [Automated Machine Learning \(AutoML\)](#))

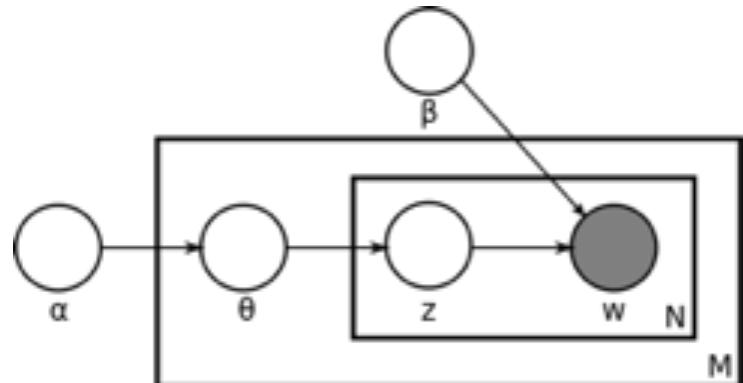
# Example: Deep Neural Networks



Huge interest in large neural networks

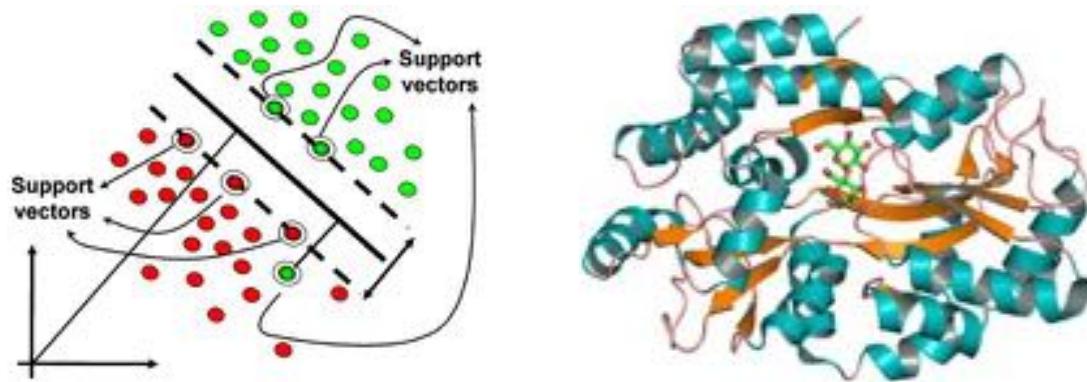
- When well-tuned, very successful for visual object identification, speech recognition, computational biology, ...
- Huge investments by Google, Facebook, Microsoft, etc.
- **Many choices:** number of layers, weight regularization, layer size, which nonlinearity, batch size, learning rate schedule, stopping conditions

# Example: Online Latent Dirichlet Allocation



- Hoffman et al. (2010): Approximate inference for [large-scale text analysis \(topic modeling\) with Latent Dirichlet Allocation](#)
- Good empirical results when well tuned
- **Hyper-parameters** tricky to set: Dirichlet parameters, number of topics, learning rate schedule, batch size, vocabulary size, ...

# Example: Classification of DNA Sequences



- Objective: Predict which DNA sequences will bind with which proteins
- Miller et al. (2012): [Latent Structural Support Vector Machine](#)
- **Hyper-parameters:** margin/slack parameter, entropy parameter, convergence criterion

# Search for Good Hyper-parameters

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- Standard search procedures:
  - Manual tuning
  - Grid search
  - Random search (very simple, works surprisingly well)
  - Black magic

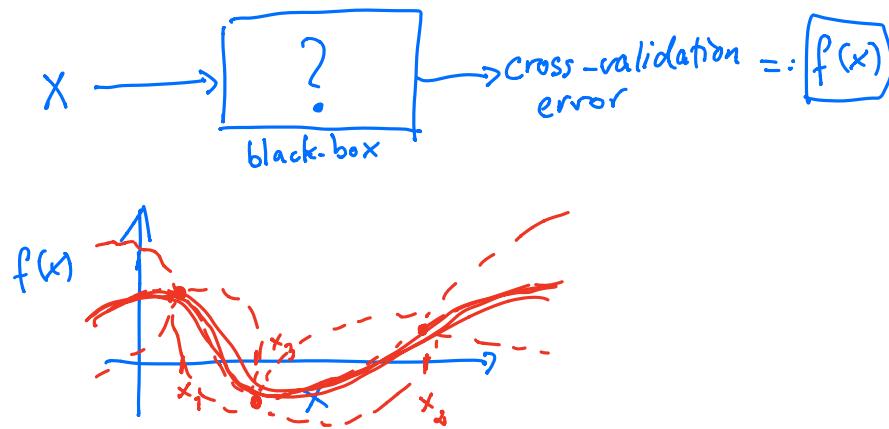
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- Standard search procedures:
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  - Random search (very simple, works surprisingly well)
  - Black magic
- Painful:
  - Evaluating the quality of the objective may be very expensive (e.g., time or money)
    - ▶ Imagine we would need to run a GPU/TPU cluster for 2 weeks
  - Many training cycles
  - Possibly noisy

## Setting

Globally optimize a black-box objective that is expensive to evaluate  
(e.g., cross-validation error for a massive neural network)

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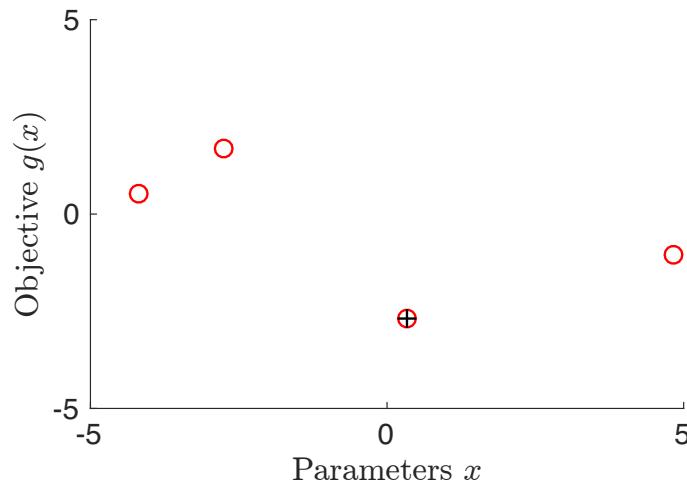
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- **Optimize cheap proxy** function to determine where to evaluate the true objective next
- Standard proxy: **Gaussian process**

- Objective: Find global minimum of objective function  $g$ :

$$\boldsymbol{x}_* = \arg \min_{\boldsymbol{x}} g(\boldsymbol{x})$$

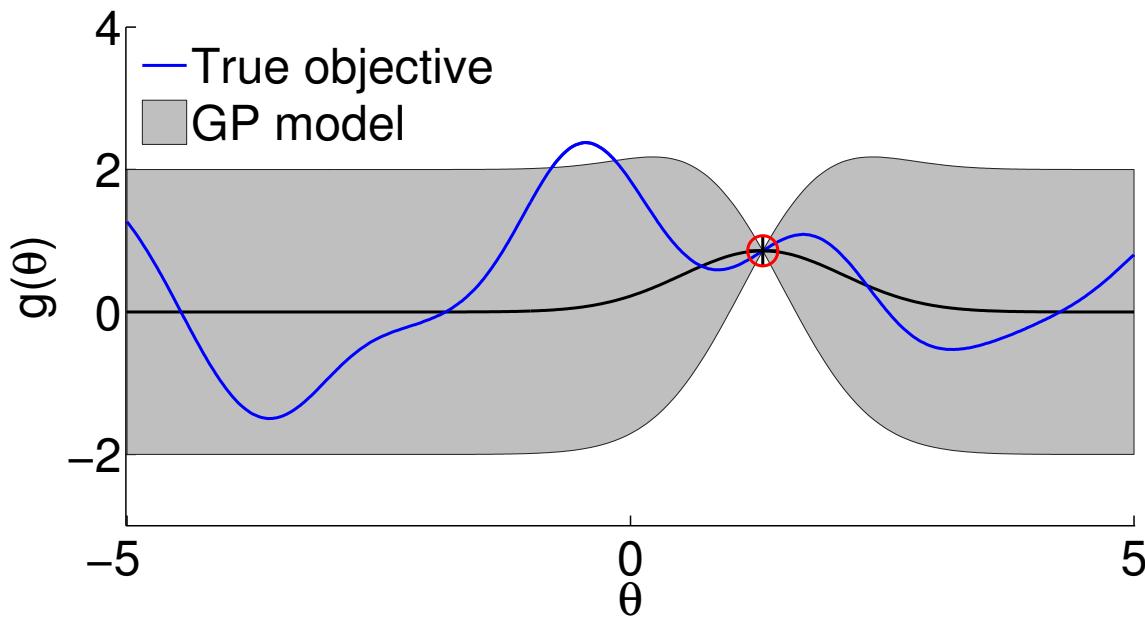
- We can evaluate the objective  $g$  pointwise, but do not have an easy functional form or gradients; observations may be noisy
- **Evaluating  $g$  is costly** (e.g., train a massive deep network)



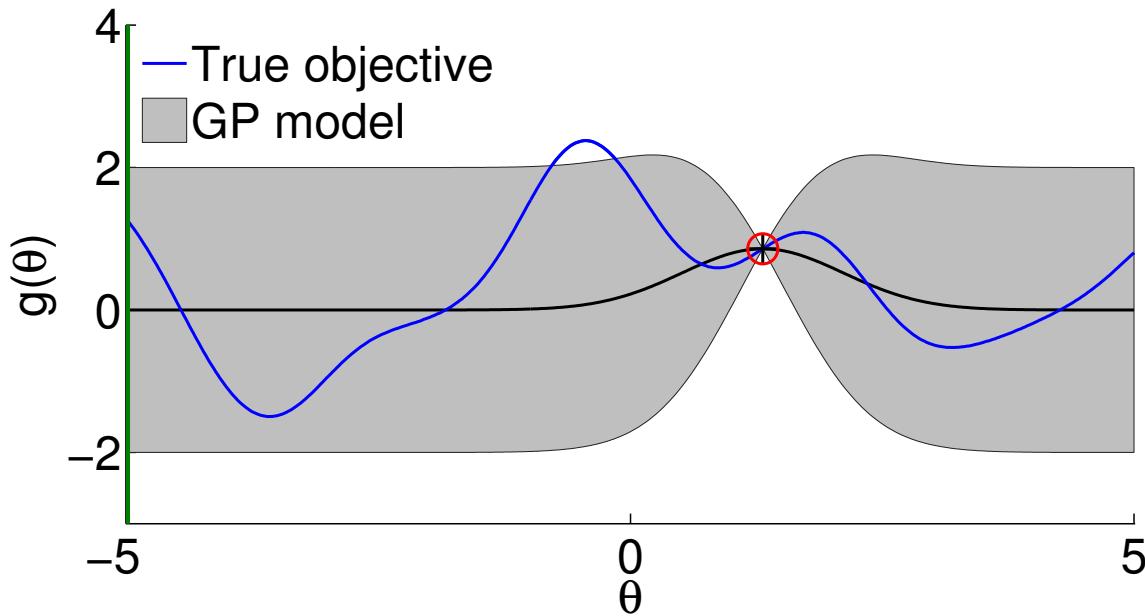
- To avoid evaluating  $g$  an excessive number of times, approximate it using a **proxy function**  $\tilde{g}$  (which is cheap to evaluate)
- Find a **global optimum**  $\tilde{g}(\mathbf{x}_*)$  of proxy function  $\tilde{g}$
- Evaluate true objective  $g$  at  $\mathbf{x}_*$
- Overall: Evaluate  $g$  only once

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- Overall: Evaluate  $g$  only once
- Works well if  $\tilde{g} \approx g$ .
- Usually not the case ➔ Repeat this cycle and keep updating  $\tilde{g}$

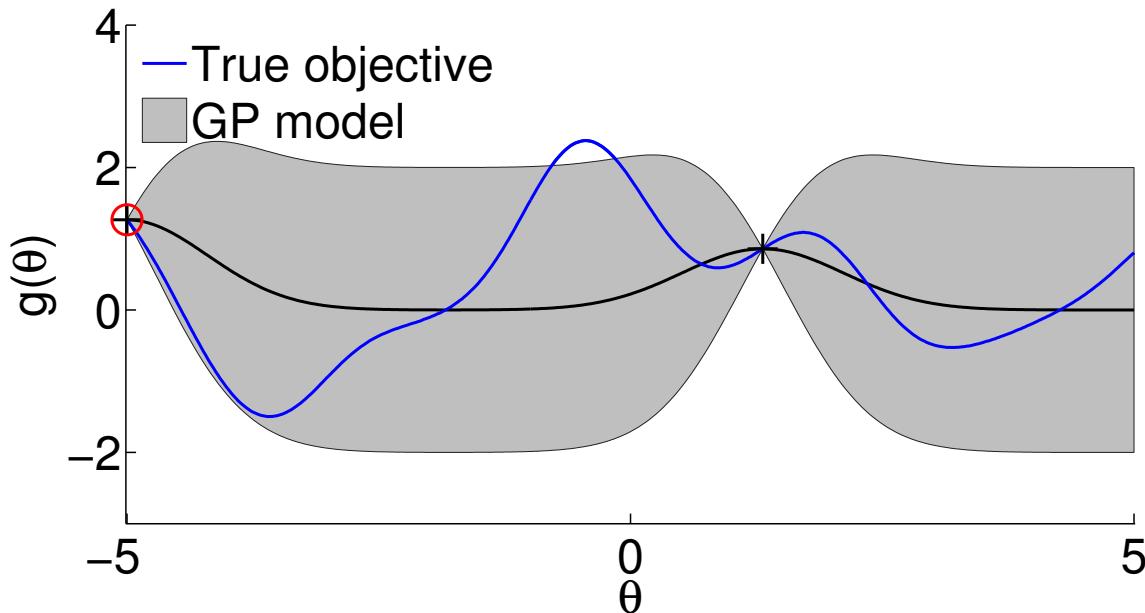
# Bayesian Optimization: Illustration



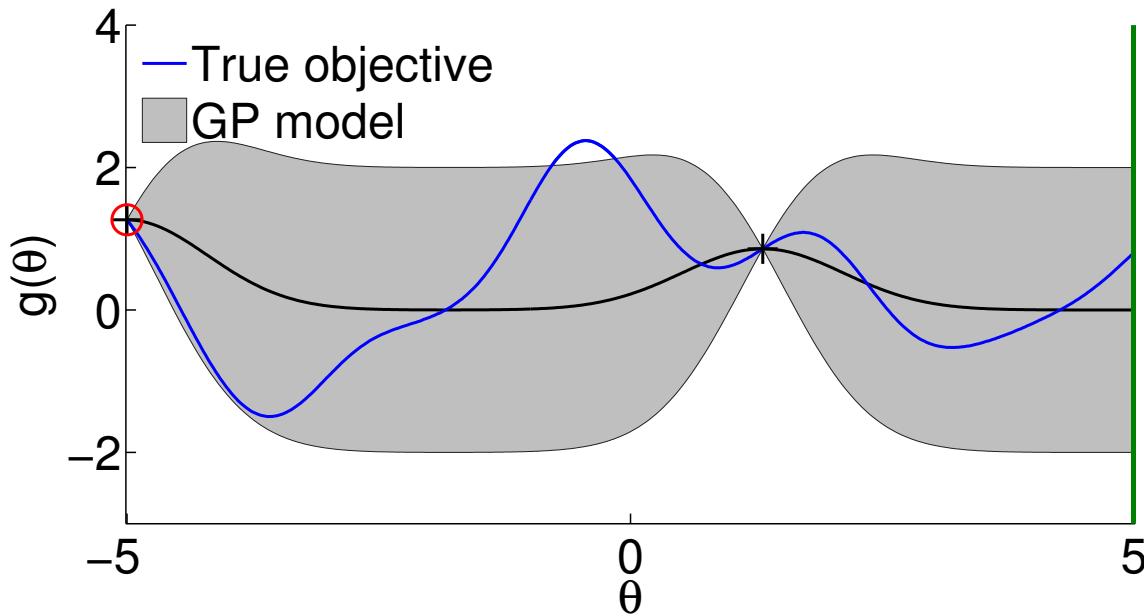
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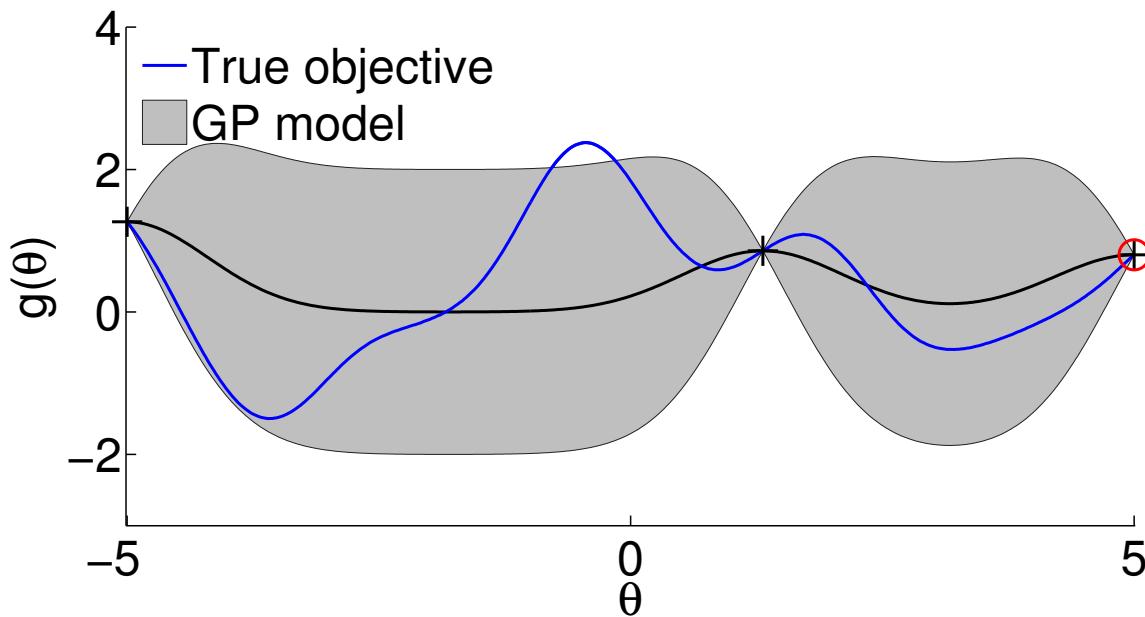
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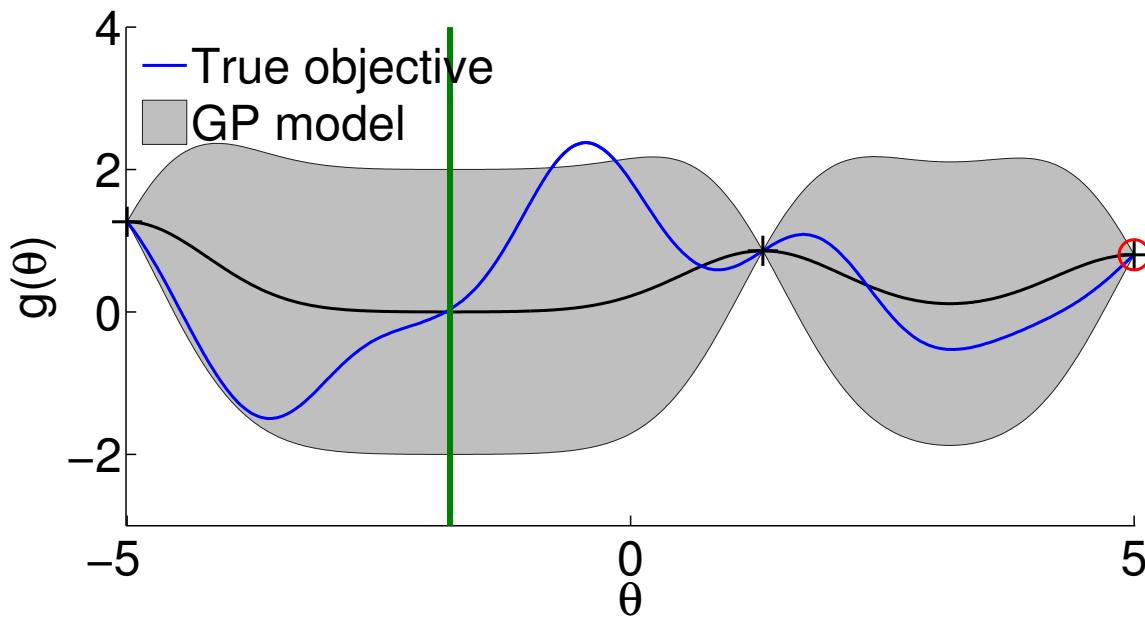
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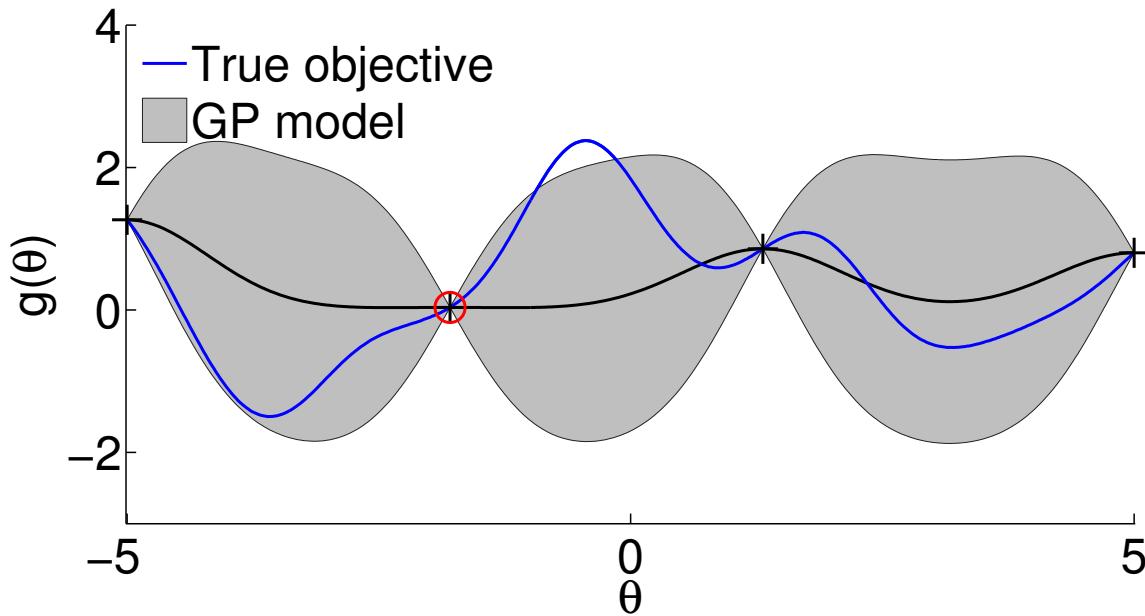
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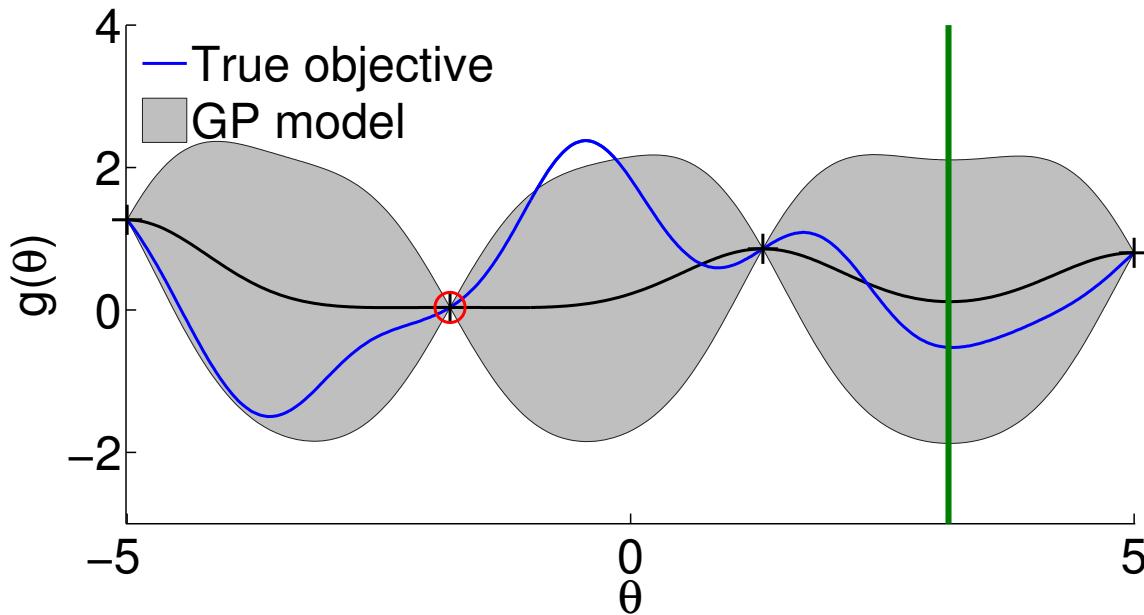
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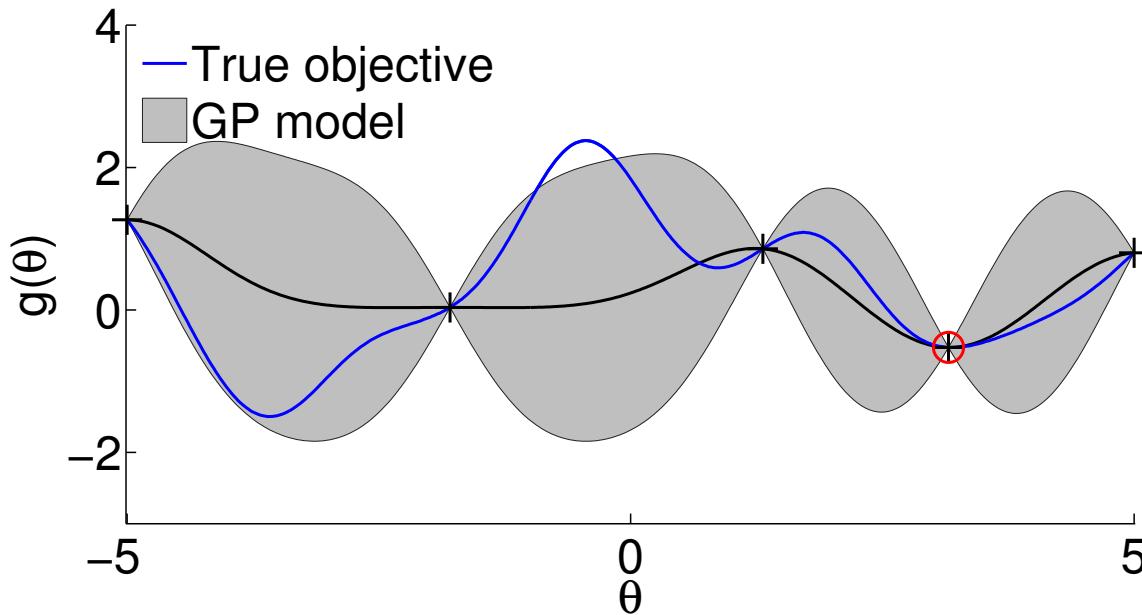
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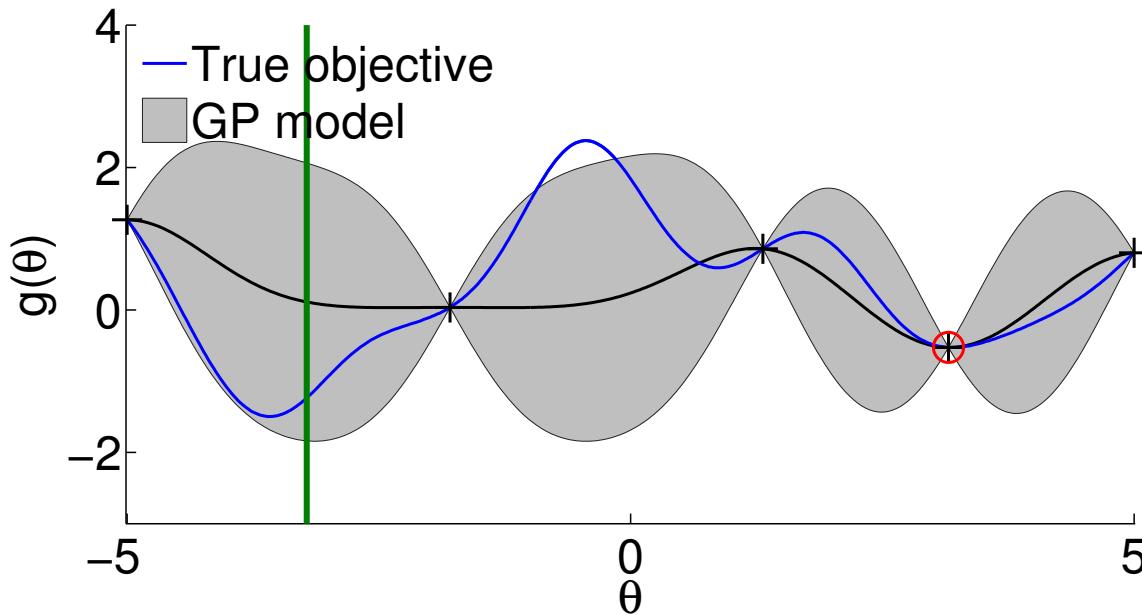
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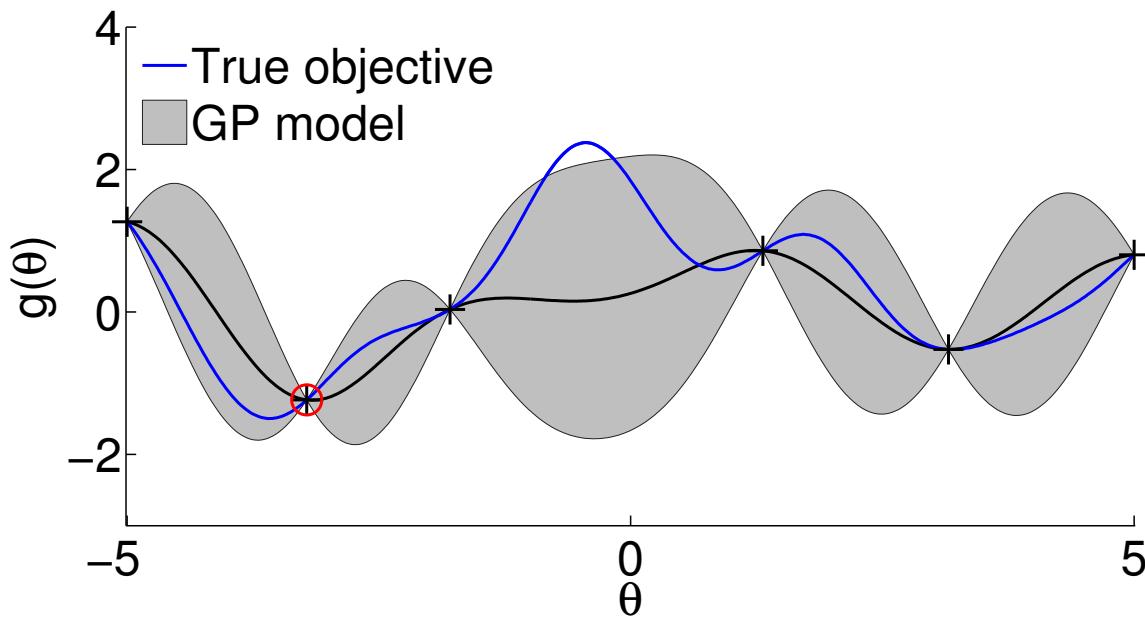
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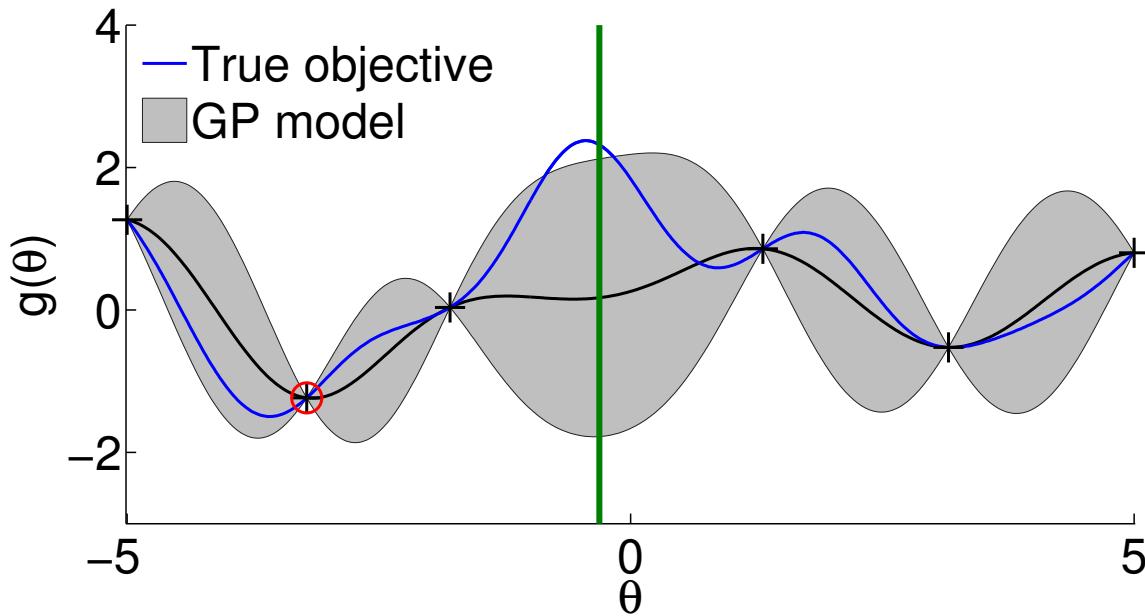
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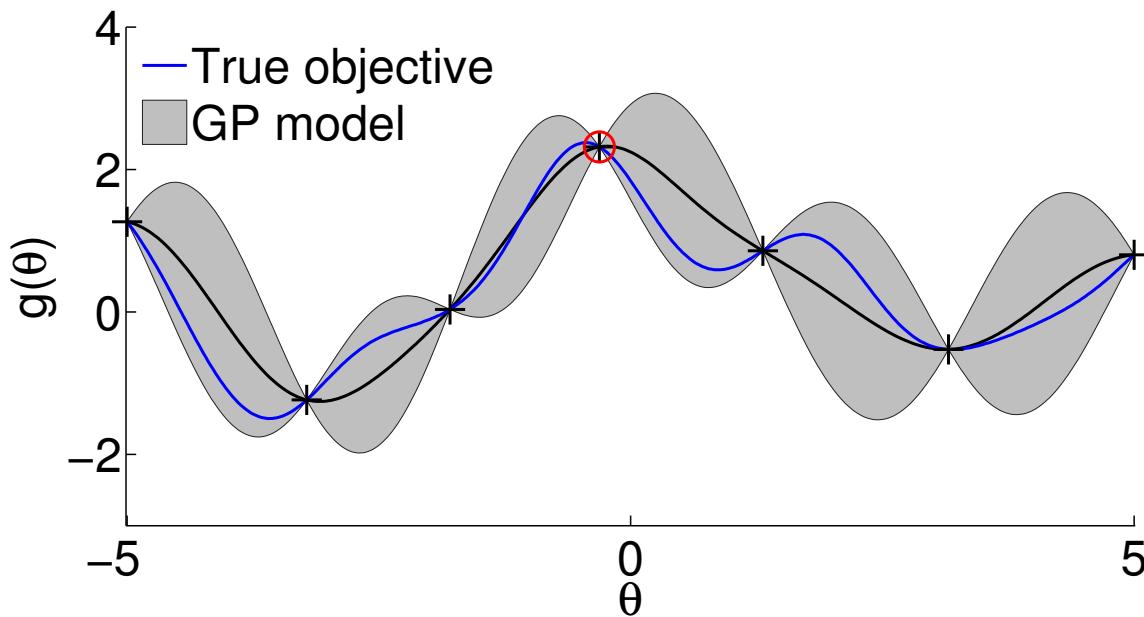
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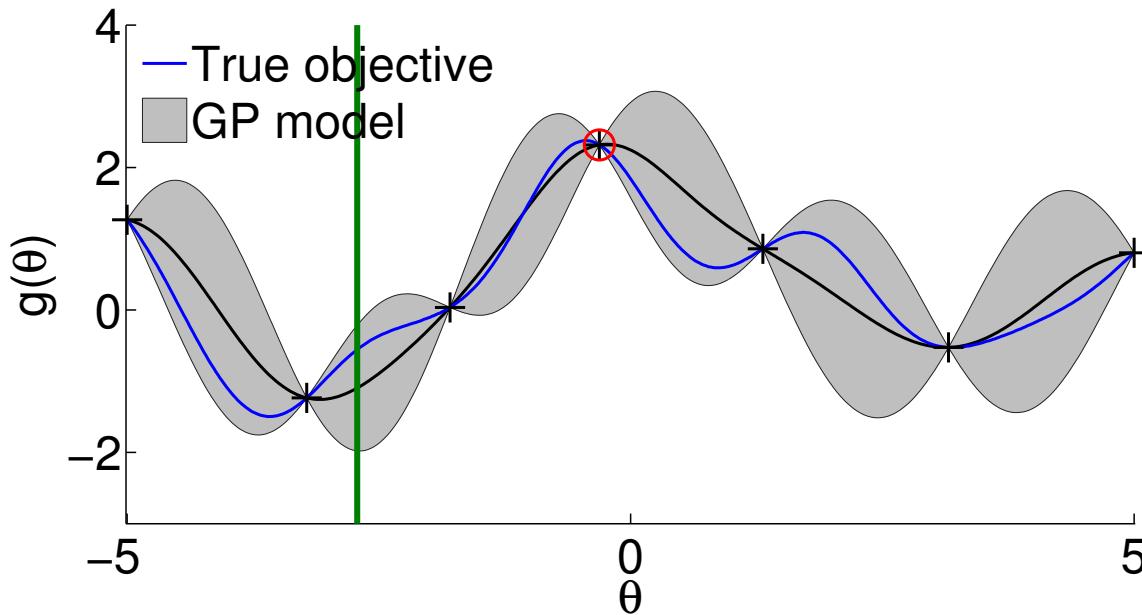
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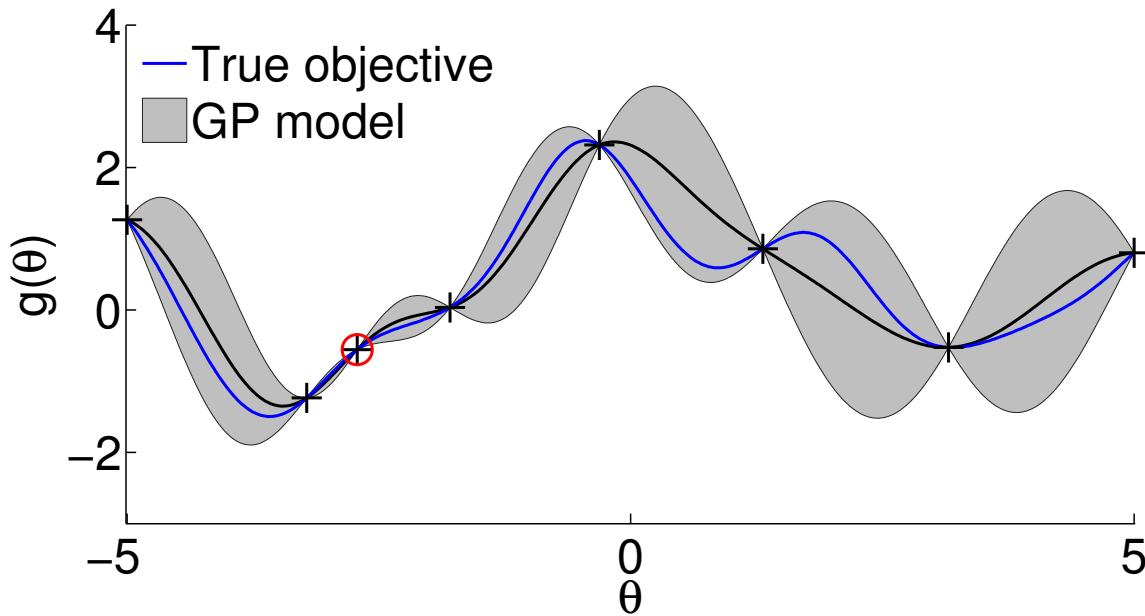
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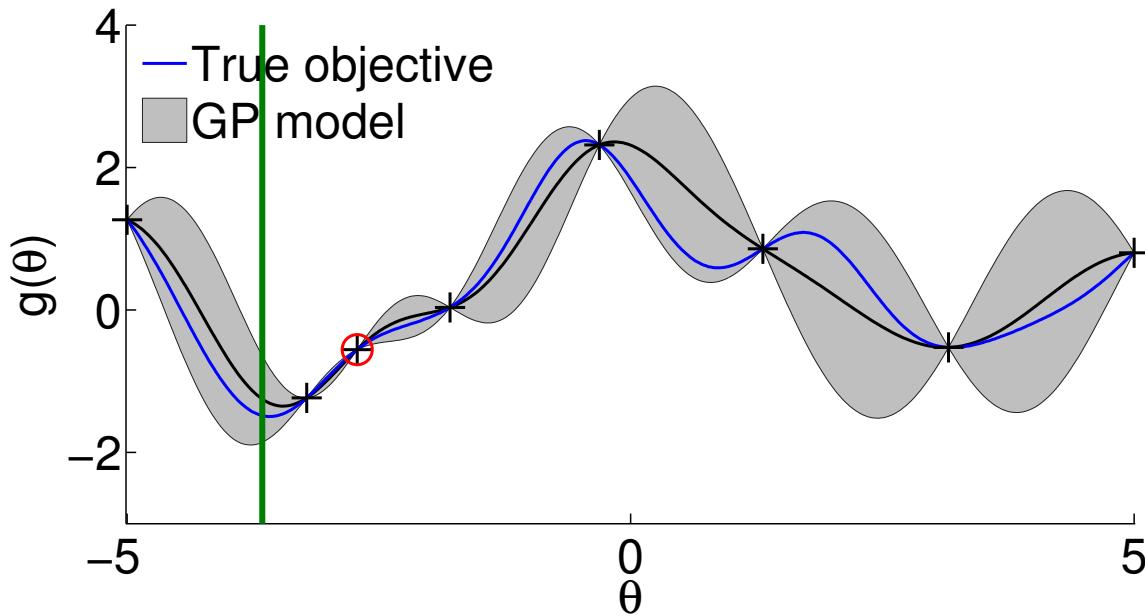
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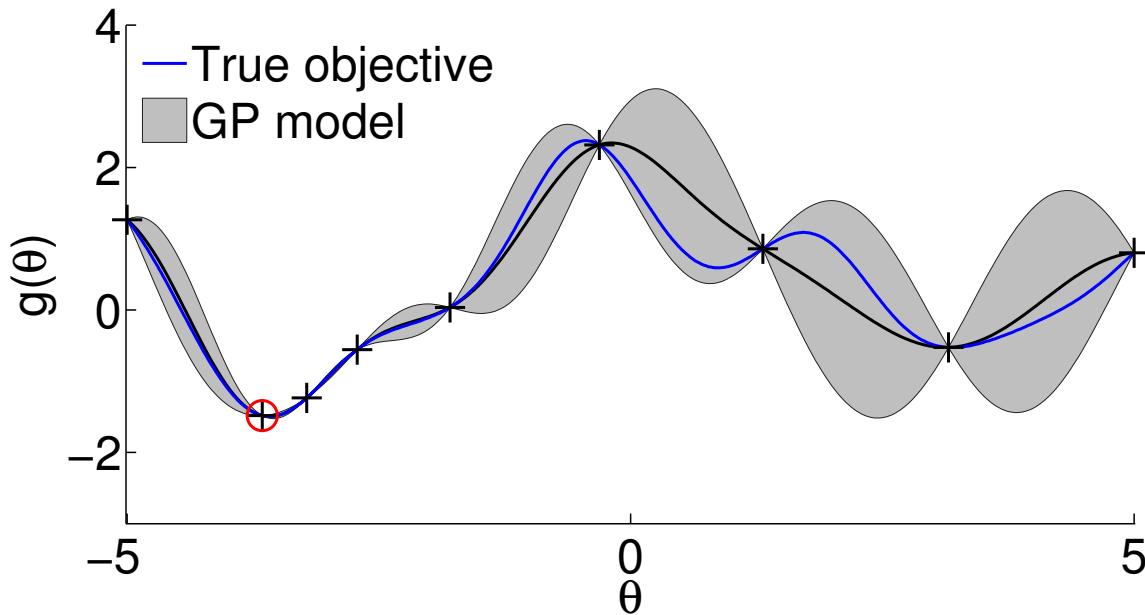
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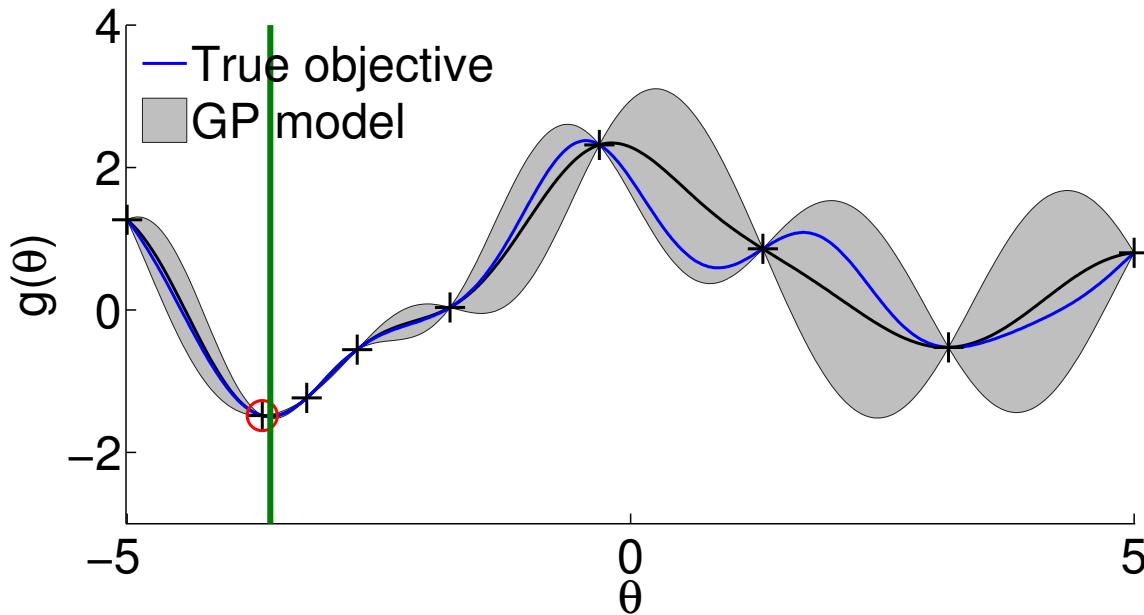
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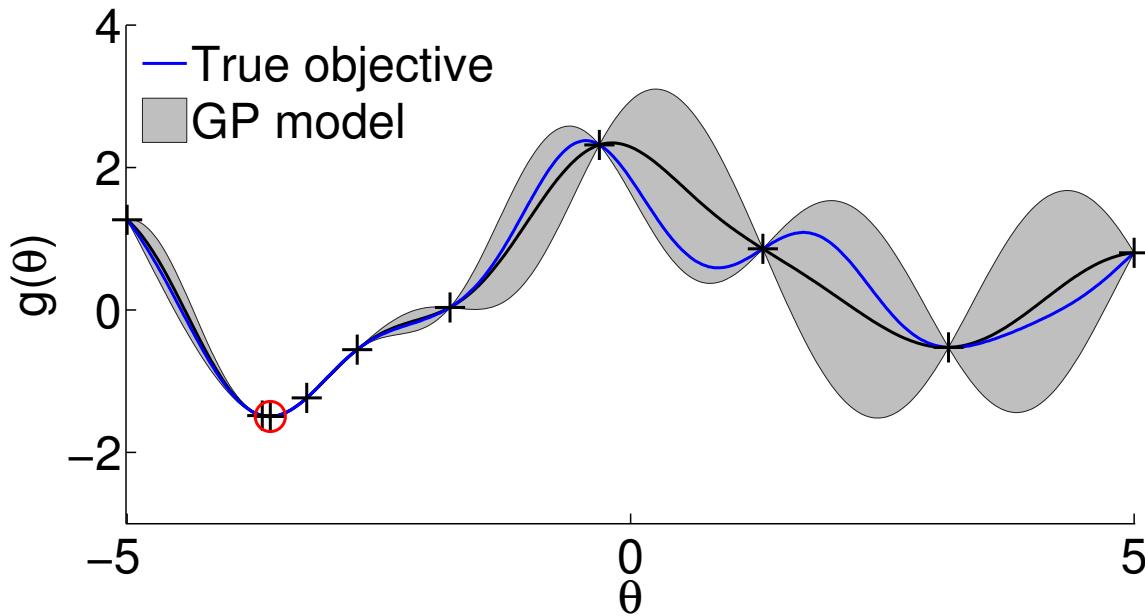
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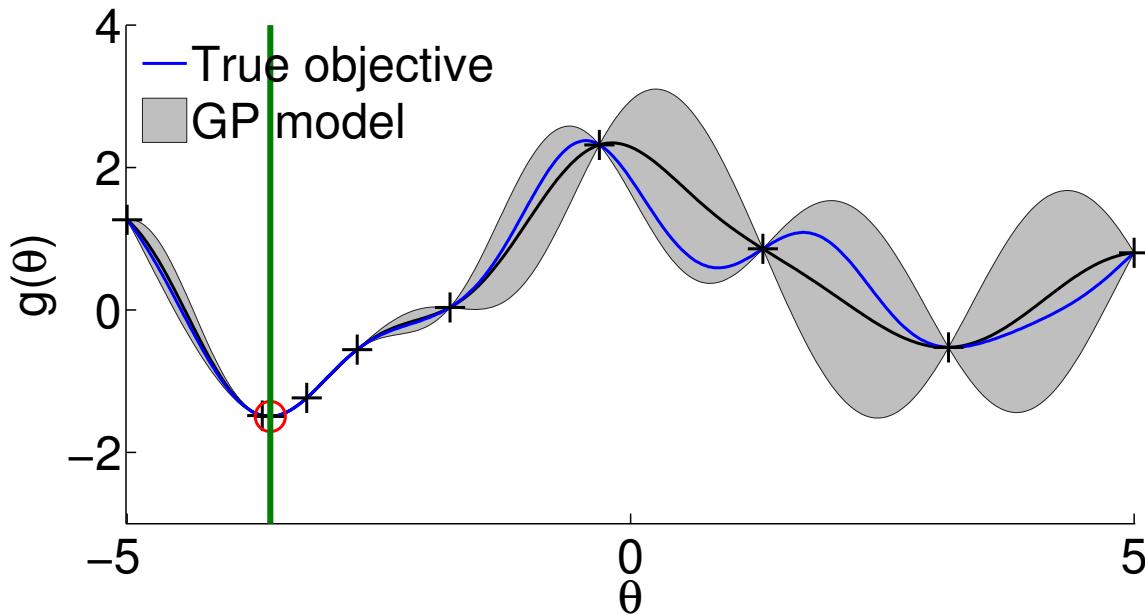
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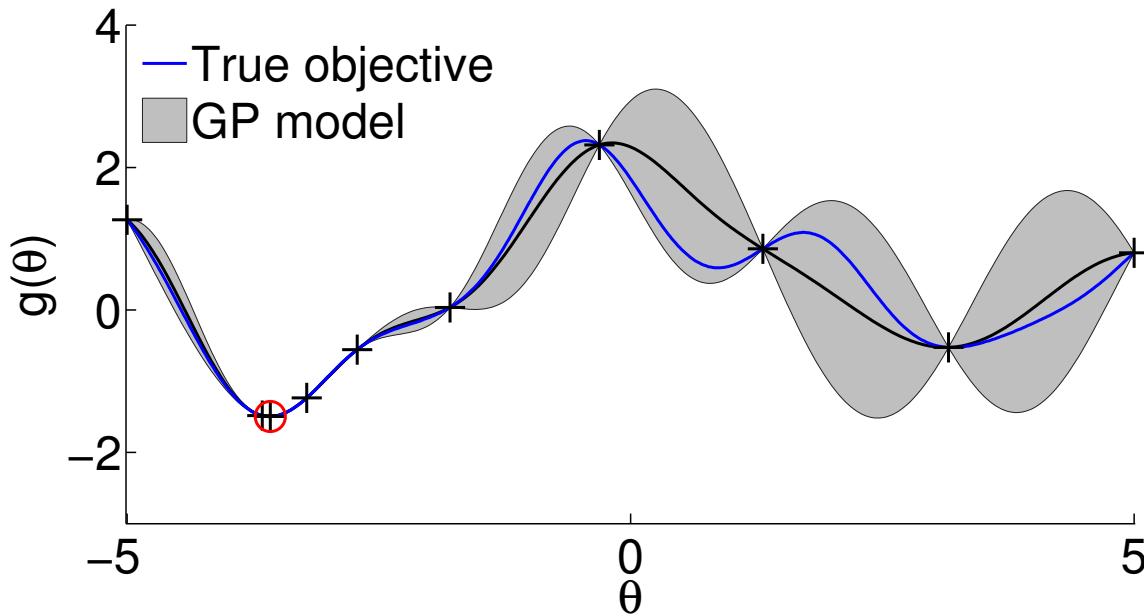
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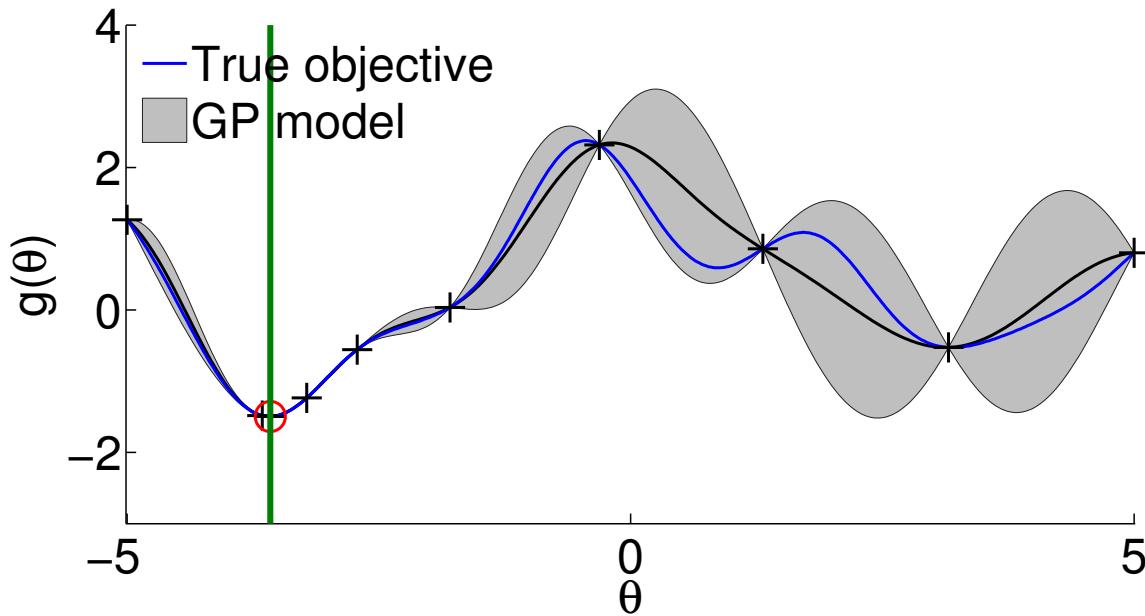
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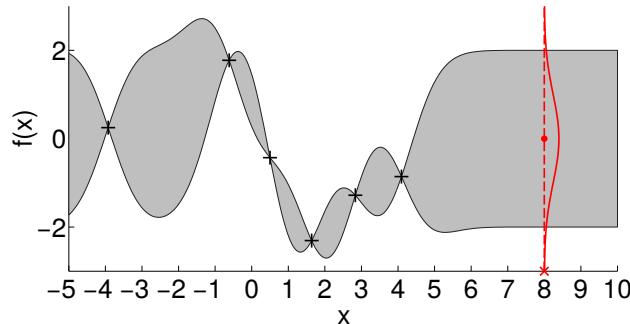


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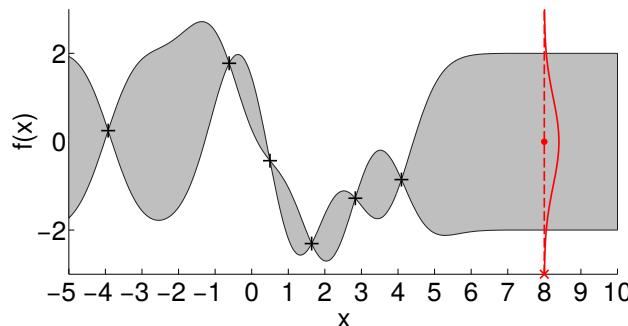
## Choosing the Next Point to Evaluate the True Objective: Acquisition Functions

# Using Uncertainty in Global Optimization



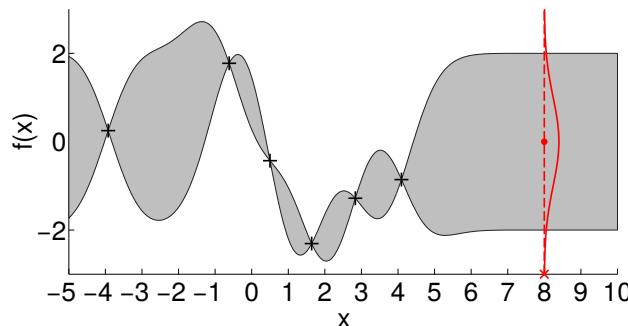
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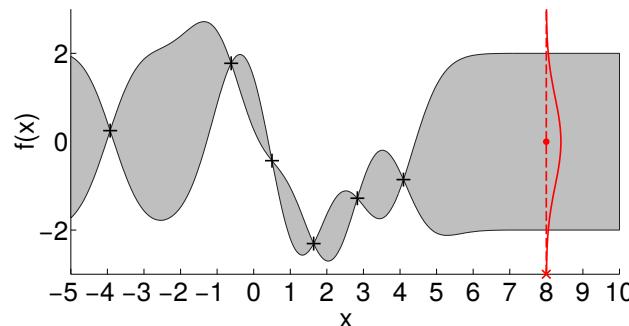
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# Using Uncertainty in Global Optimization



- Find a good (global) optimum
  - ▶ Need to get out of local optima
- Extrapolate from collected knowledge
- GP gives us closed-form means and variances
  - ▶ Trade off exploration and exploitation
    - **Exploration:** Seek places with high variance/uncertainty
    - **Exploitation:** Seek places with low mean

# Using Uncertainty in Global Optimization

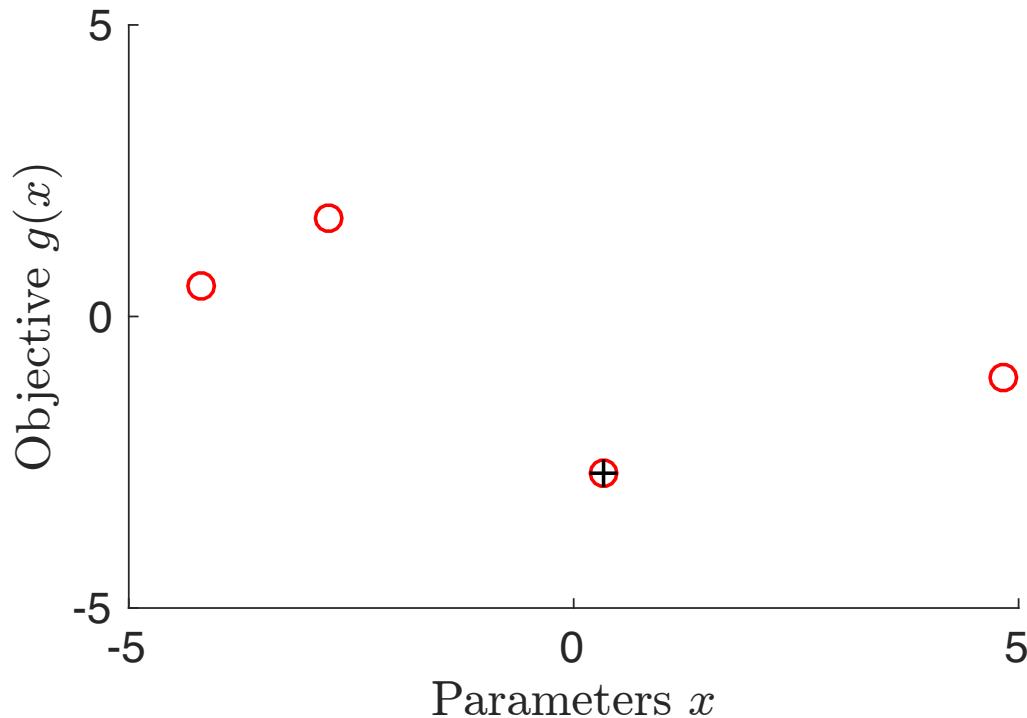


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    - **Exploration:** Seek places with high variance/uncertainty
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- **Acquisition function  $\alpha$**  trades off exploration and exploitation for our proxy optimization

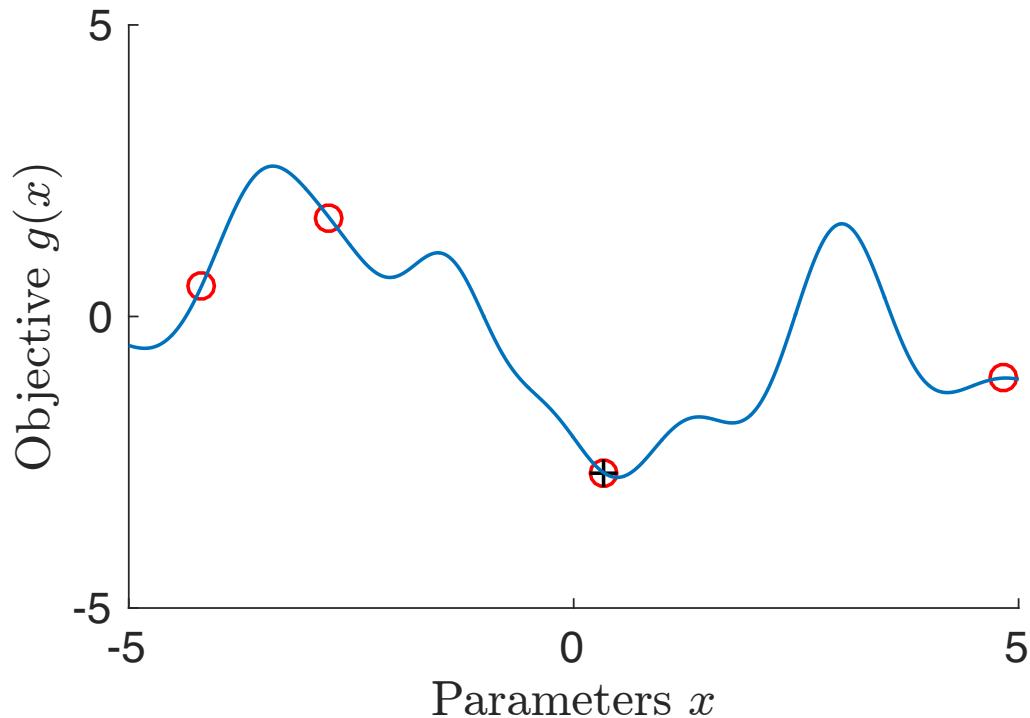
# Key Steps (Pseudo-Code)

```
1: Init: Data set  $\mathcal{D}_0 = \{\mathbf{X}_0, \mathbf{y}_0\}$ 
2: for iterations  $t = 1, 2, \dots$  do
3:   Update GP using data  $\mathcal{D}_{t-1}$ 
4:   Select  $\mathbf{x}_t = \arg \max_{\mathbf{x}} \alpha(\mathbf{x})$  by optimizing acquisition function
5:   Query true objective  $g$  at  $\mathbf{x}_t$ 
6:   Augment data set  $\mathcal{D}_t = \mathcal{D}_{t-1} \cup \{(\mathbf{x}_t, y_t)\}$ 
7: end for
8: Return best input in data set:  $\mathbf{x}^* = \arg \min_{\mathbf{x}} y(\mathbf{x})$ 
```

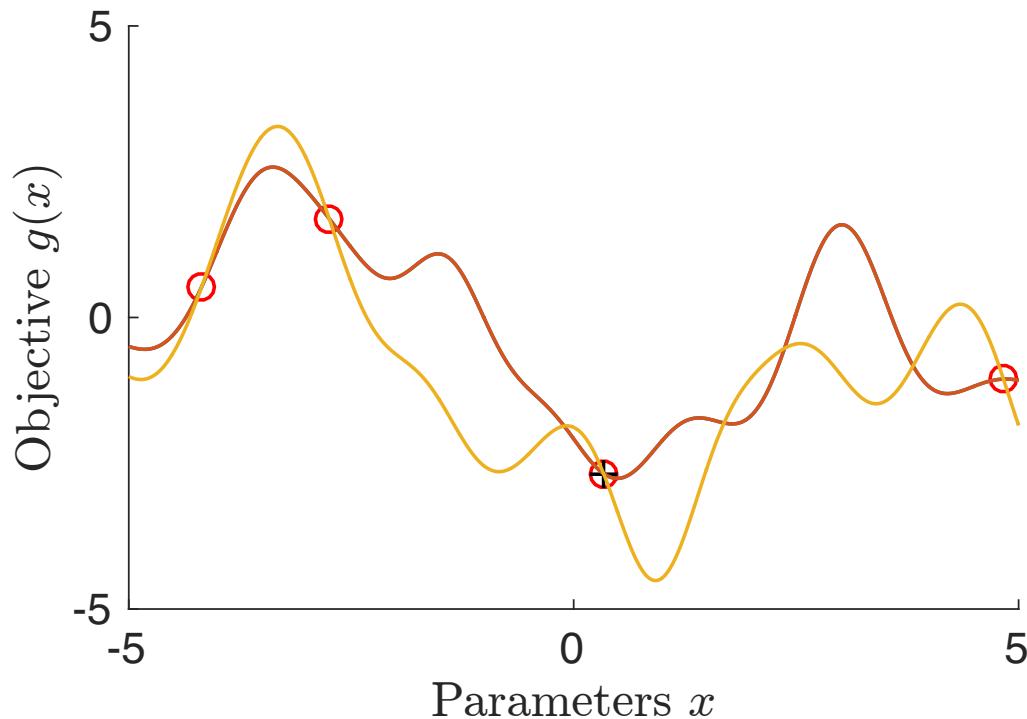
# Where to Evaluate Next?



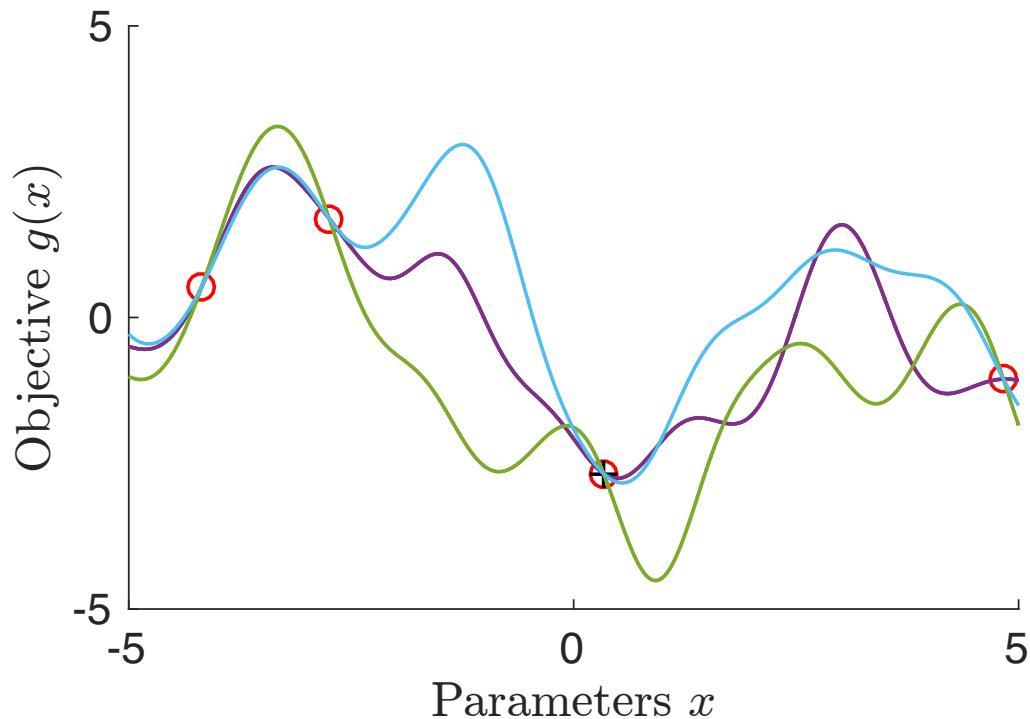
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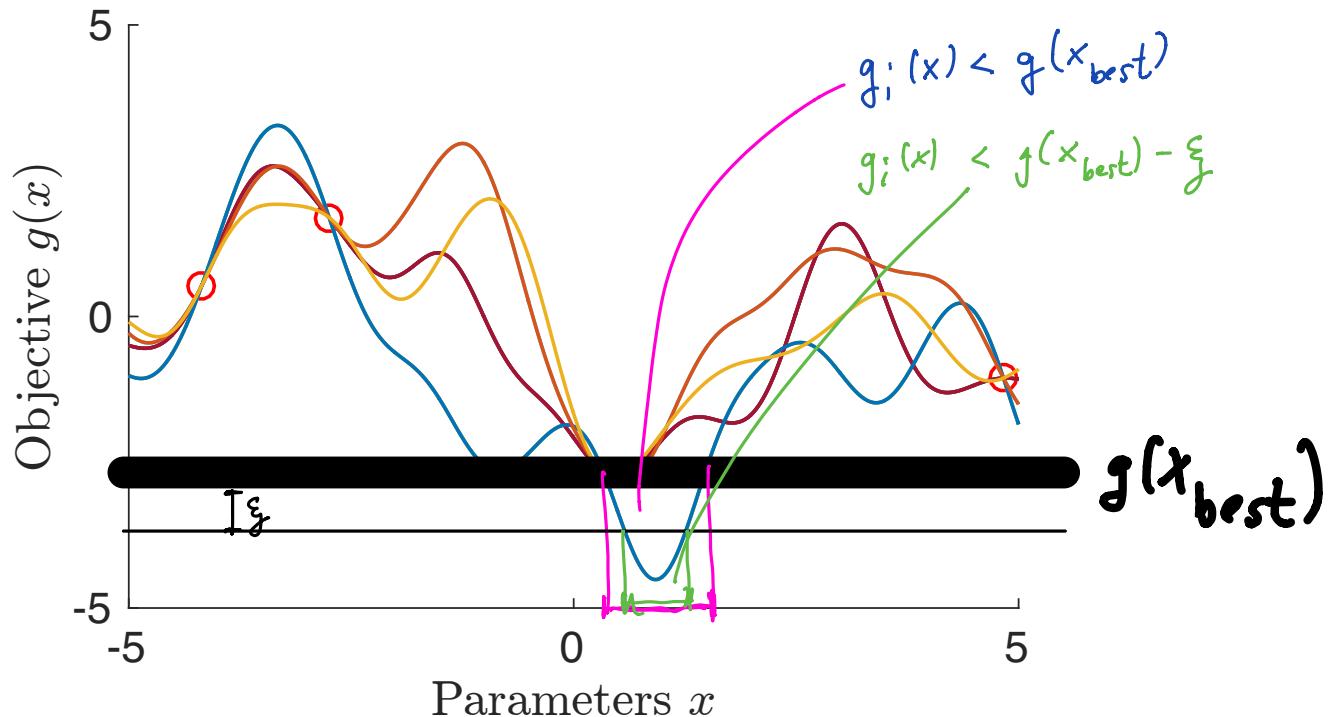
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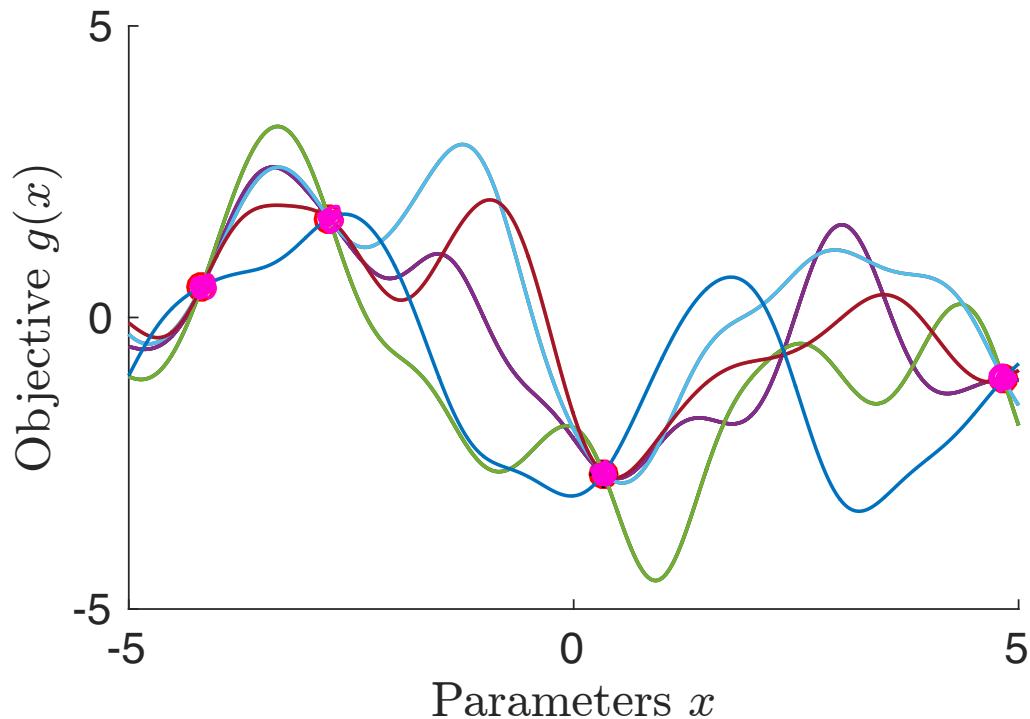
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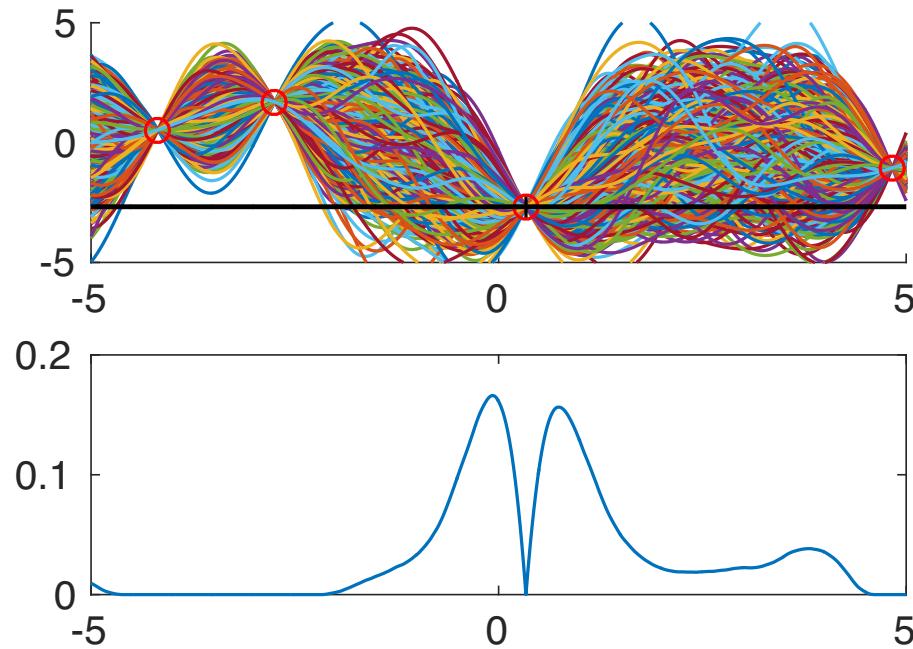
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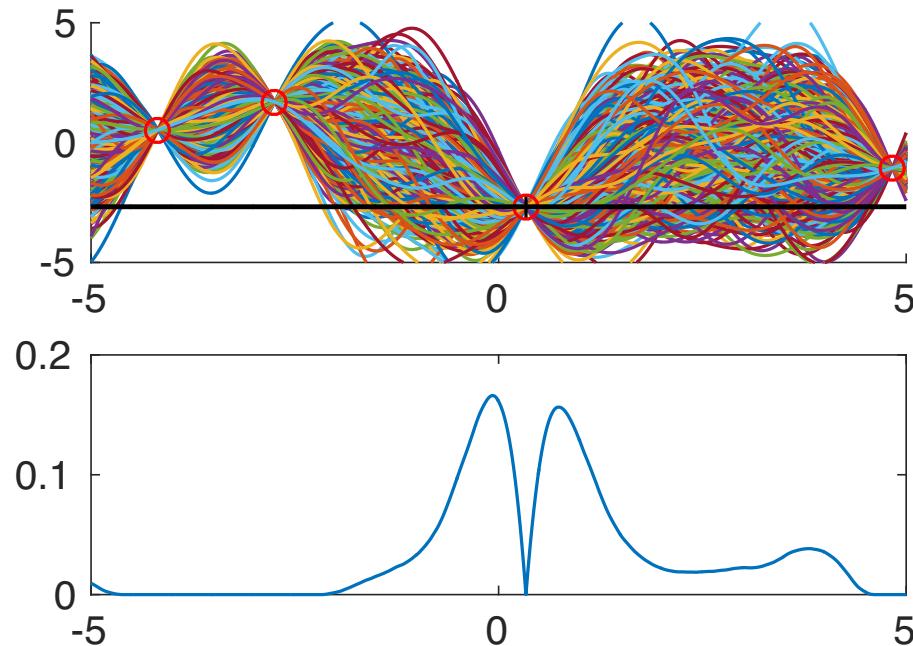


# Where to Evaluate Next to Improve Most?



- Upper panel: Samples from a probabilistic proxy  $\tilde{g}$

# Where to Evaluate Next to Improve Most?



- Upper panel: Samples from a probabilistic proxy  $\tilde{g}$
- Lower panel: Corresponding **expected improvement** over the best solution so far (black cross)
- ▶ Evaluate  $g$  at the maximum of the expected improvement

# Closed-Form Acquisition Functions

- For all  $x \in \mathbb{R}^D$  the GP posterior gives a predictive mean  $\mu(x)$  variance  $\sigma^2(x)$  of  $g(x)$
- Define

$$\gamma(x) = \frac{g(x_{\text{best}}) - \mu(x)}{\sigma(x)}$$

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$$\gamma(\mathbf{x}) = \frac{g(\mathbf{x}_{\text{best}}) - \mu(\mathbf{x})}{\sigma(\mathbf{x})}$$

- **Probability of Improvement (Kushner 1964):**

$$\alpha_{\text{PI}}(\mathbf{x}) = \Phi(\gamma(\mathbf{x}))$$

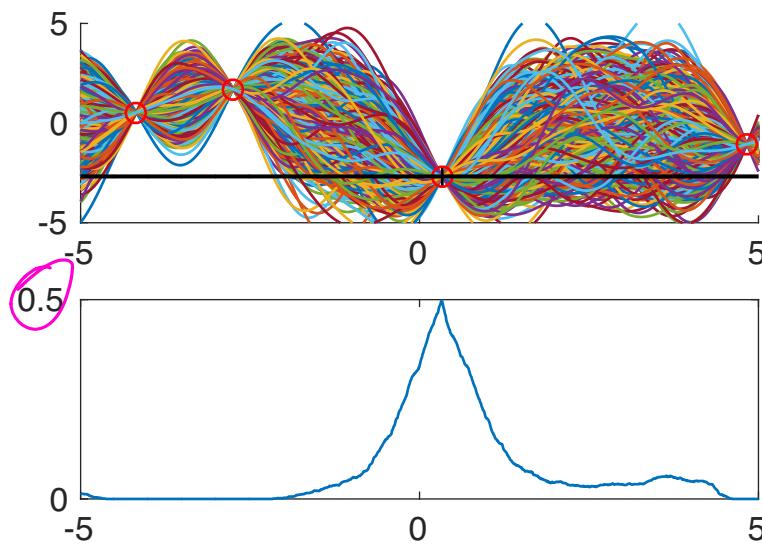
- **Expected Improvement (Mockus 1978):**

$$\alpha_{\text{EI}}(\mathbf{x}) = \sigma(\mathbf{x})(\gamma(\mathbf{x})\Phi(\gamma(\mathbf{x})) + \mathcal{N}(\gamma(\mathbf{x}) | 0, 1))$$

- **GP Lower Confidence Bound (Srinivas et al., 2010):**

$$\alpha_{\text{LCB}}(\mathbf{x}) = -(\mu(\mathbf{x}) - \kappa\sigma(\mathbf{x})), \quad \kappa > 0$$

# Probability of Improvement (1)

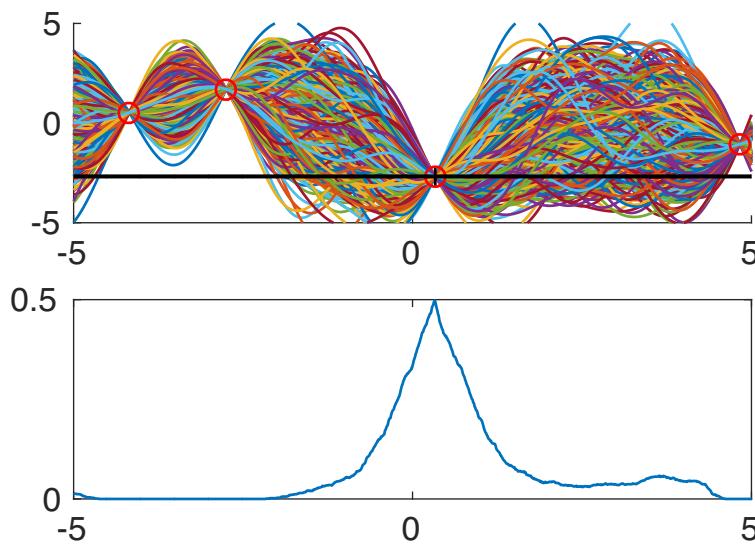


- **Idea:** Determine the probability that  $x_*$  leads to a better function value than the currently best one  $g(x_{\text{best}})$
- Sampling-based setting:  
Sample  $N$  functions  $g_i$ ; at every input  $x$  compute a Monte-Carlo estimate

$$\alpha_{\text{PI}}(x) = p(g(x) < g(x_{\text{best}})) \approx \frac{1}{N} \sum_{i=1}^N \delta(g_i(x) < g(x_{\text{best}}))$$

► Can lead to continued exploitation in an  $\epsilon$ -region around  $x_{\text{best}}$ .

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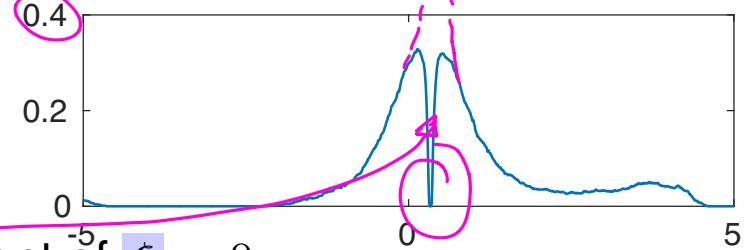
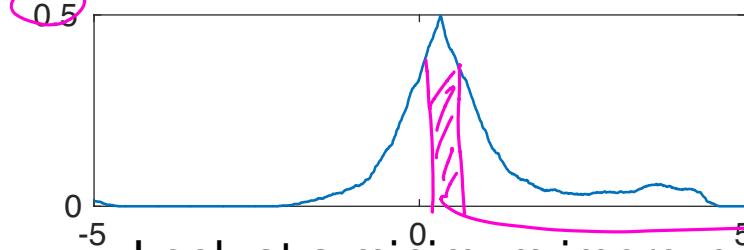
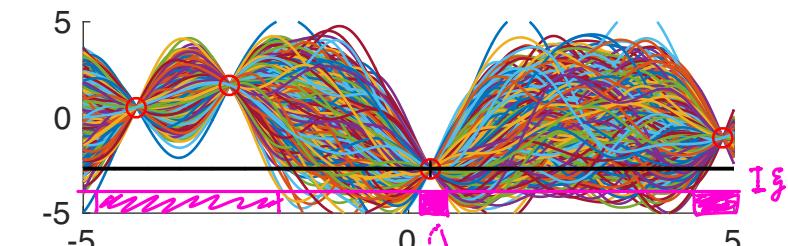
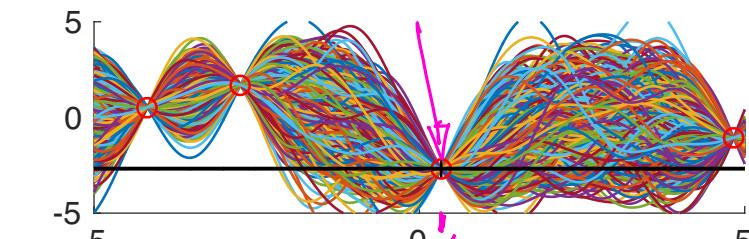


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- ▶ Can lead to continued exploitation in an  $\epsilon$ -region around  $x_{\text{best}}$ .
- ▶ Introduce a “slack variable”  $\xi$  for more aggressive exploration

# Probability of Improvement (2)



- Look at a minimum improvement of  $\xi > 0$ :

$$\alpha_{\text{PI}}(\mathbf{x}) = p(g(\mathbf{x}) < g(\mathbf{x}_{\text{best}}) - \xi) \approx \frac{1}{N} \sum_{i=1}^N \delta(g_i(\mathbf{x}) < g(\mathbf{x}_{\text{best}}) - \xi)$$

- If  $f \sim GP$  and  $p(g(\mathbf{x})) = \mathcal{N}(\mu(\mathbf{x}), \sigma(\mathbf{x}))$ :

$$\alpha_{\text{PI}}(\mathbf{x}) = \Phi(\gamma(\mathbf{x}, \xi)), \quad \gamma(\mathbf{x}, \xi) = \frac{g(\mathbf{x}_{\text{best}}) - \xi - \mu(\mathbf{x})}{\sigma(\mathbf{x})}$$

# Expected Improvement

- Idea: Quantify the amount of improvement

- Sampling-based scenario, where  $g_i \sim p(f)$ :

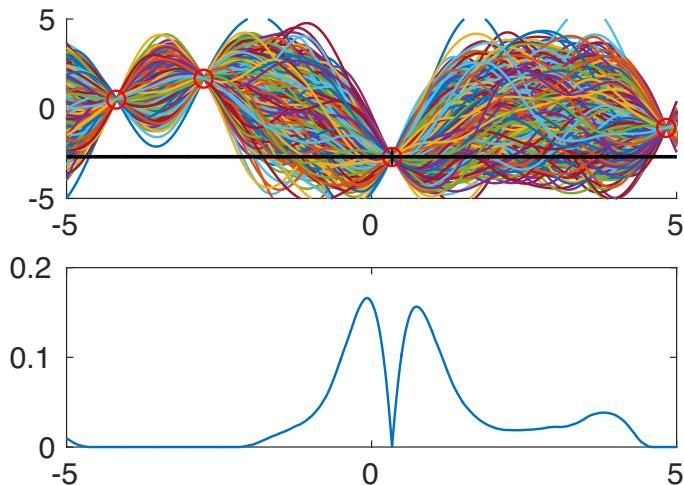
$$\alpha_{EI}(\mathbf{x}) = \mathbb{E}[\max\{0, g(\mathbf{x}_{\text{best}}) - g(\mathbf{x})\}]$$

$$\approx \frac{1}{N} \sum_{i=1}^N \max\{0, g(\mathbf{x}_{\text{best}}) - g_i(\mathbf{x})\}$$

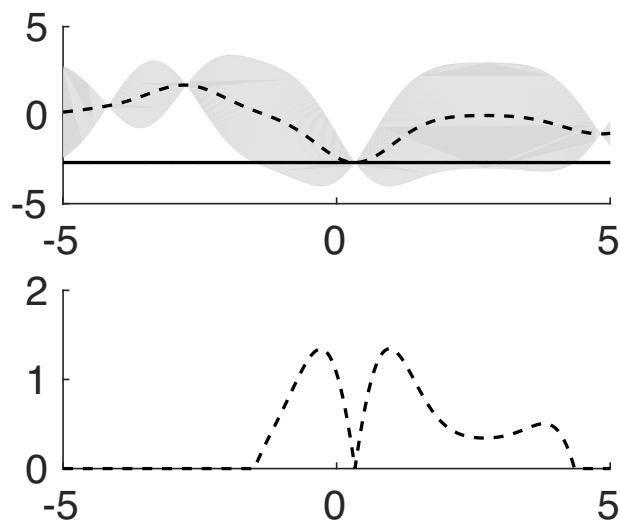
- If  $f \sim GP$ , we have a closed-form expression:

$$\alpha_{EI}(\mathbf{x}) = \sigma(\mathbf{x})(\gamma(\mathbf{x})\Phi(\gamma(\mathbf{x})) + \mathcal{N}(\gamma(\mathbf{x}) | 0, 1))$$

- Slack-variable approach also possible (similar to PI)



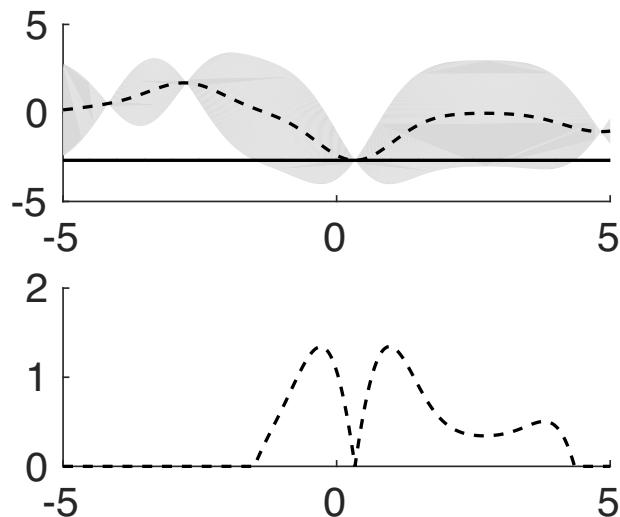
# GP-Lower Confidence Bound (1)



- Use the predictive mean  $\mu(x)$  and variance  $\sigma^2(x)$  of the GP prediction directly for targeted exploration by means of the acquisition function

$$\alpha_{LCB}(\mathbf{x}_t) = -(\mu(\mathbf{x}_t) - \sqrt{\kappa}\sigma(\mathbf{x}_t))$$

# GP-Lower Confidence Bound (2)



- More generally, we can get regret bounds for iteration-dependent  $\kappa$  (Srinivas et al., 2010)

$$\alpha_{\text{LCB}}(\mathbf{x}_t) = -(\mu(\mathbf{x}_t) - \sqrt{\kappa_t} \sigma(\mathbf{x}_t))$$

where  $\kappa_t \in \mathcal{O}(\log t)$  grows with the iteration  $t$

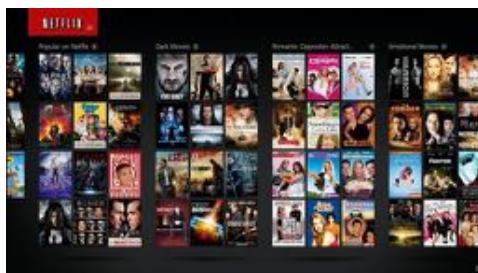
► Continue exploration

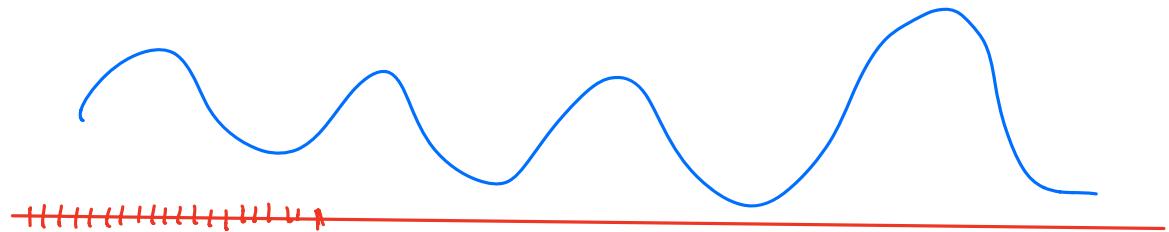
# Optimizing the Acquisition Function

- Optimizing the acquisition function **requires us to run a global optimizer inside Bayesian optimization**
- What have we gained?

# Optimizing the Acquisition Function

- Optimizing the acquisition function **requires us to run a global optimizer inside Bayesian optimization**
- What have we gained?
- Evaluating the acquisition function is cheap compared to evaluating the true objective
  - ▶ We can afford evaluating it many times



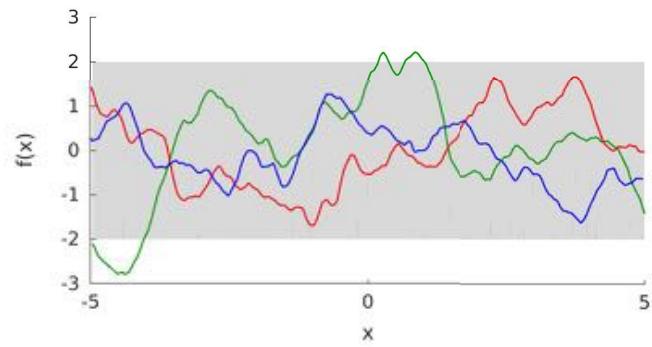
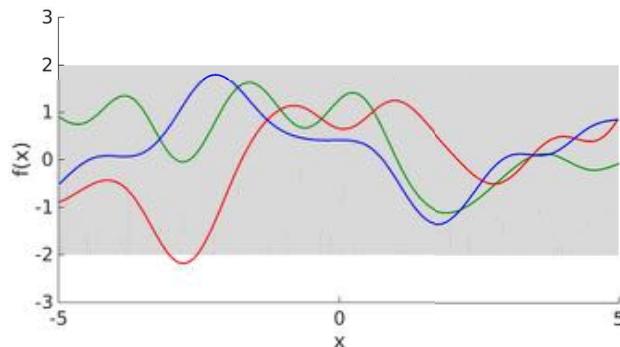


- Getting the function model (e.g., covariance function) wrong can be catastrophic
- Limited scalability in the number of dimensions and/or evaluations of the true objective function

Why?

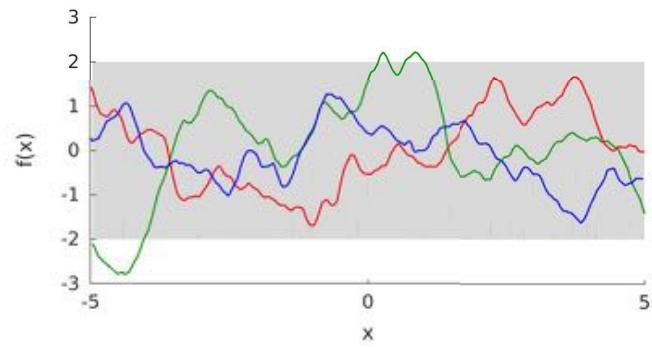
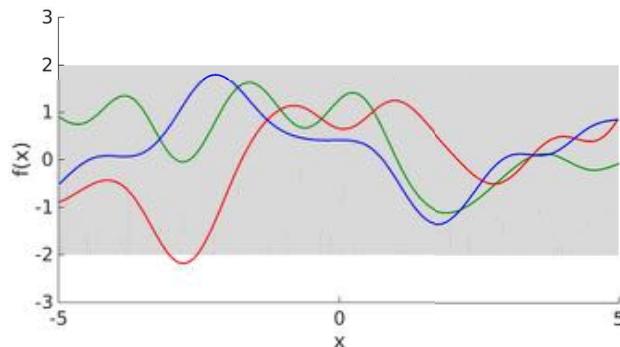
1 experiment = 1 training data point (GP)  
⇒ < 10 000 experiments  $O(N^3)$  training  
< 1000 experiments  
Run optimizer for acquisition function  
⇒ Query GP at many inputs  
→ Make predictions at many inputs  $O(N^2)$

# Poor Model Choice



- Covariance function selection is crucial for good performance
  - ▶ Choose a sufficiently flexible and adaptive kernel, e.g., Matérn (but not the squared exponential (Gaussian))

# Poor Model Choice



- Covariance function selection is crucial for good performance
  - ▶ Choose a sufficiently flexible and adaptive kernel, e.g., Matérn (but not the squared exponential (Gaussian))
- Nice side-effect of Matérn: Exploration is more encouraged than with the Gaussian kernel

# Choosing Covariance Functions

- Structured SVM for Protein Motif Finding (Miller et al., 2012)
- Optimize hyper-parameters of SSVM using BO (Snoek et al., 2012)

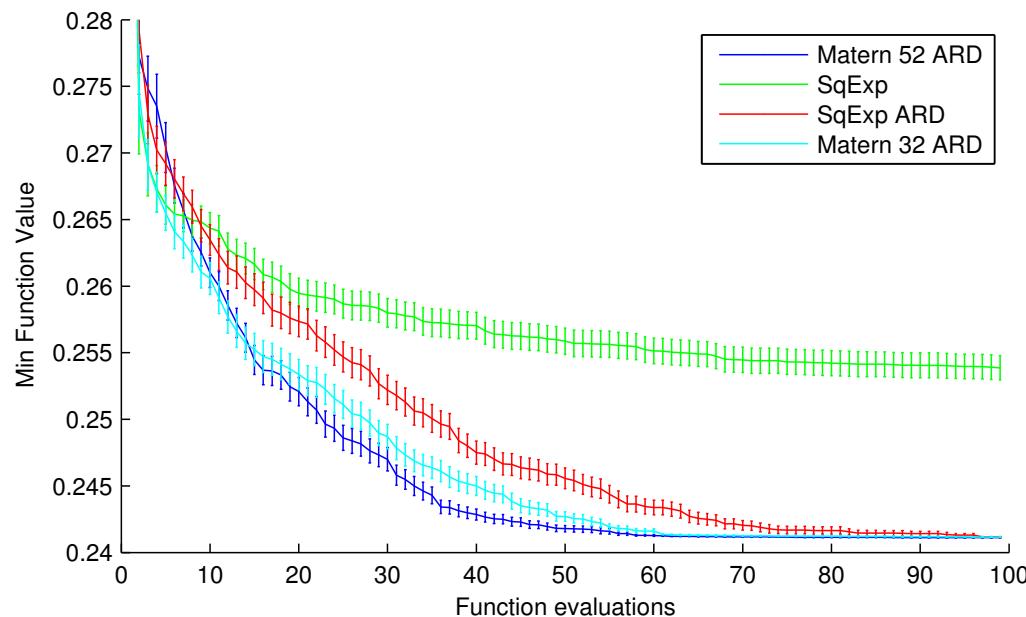


Figure: Figure from Snoek et al. (2012)

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# Gaussian Process Hyper-Parameters

- Empirical Bayes (maximize the marginal likelihood) can fail horribly, especially in the early stages of Bayesian optimization when we have only a few data points
- Solution: Integrate out the GP hyper-parameters  $\theta$  by [Markov Chain Monte Carlo \(MCMC\)](#) sampling (e.g., slice sampling)
- Look at [integrated acquisition function](#)

$$\begin{aligned}\alpha(\mathbf{x}) &= \mathbb{E}_{\boldsymbol{\theta}}[\alpha(\mathbf{x}, \boldsymbol{\theta})] = \int \alpha(\mathbf{x}, \boldsymbol{\theta}) p(\boldsymbol{\theta}) d\boldsymbol{\theta} \\ &\approx \frac{1}{K} \sum_{k=1}^K \alpha(\mathbf{x}, \boldsymbol{\theta}^{(k)}), \quad \boldsymbol{\theta}^{(k)} \sim \underbrace{p(\boldsymbol{\theta} | \mathbf{X}_n, \mathbf{y}_n)}_{\text{hyper-parameter posterior}}\end{aligned}$$

# Integrating out GP Hyper-parameters

- Online LDA (Hoffman et al., 2010) for topic modeling
- Two critical hyper-parameters that control the learning rate learned by BO (Snoek et al., 2012)

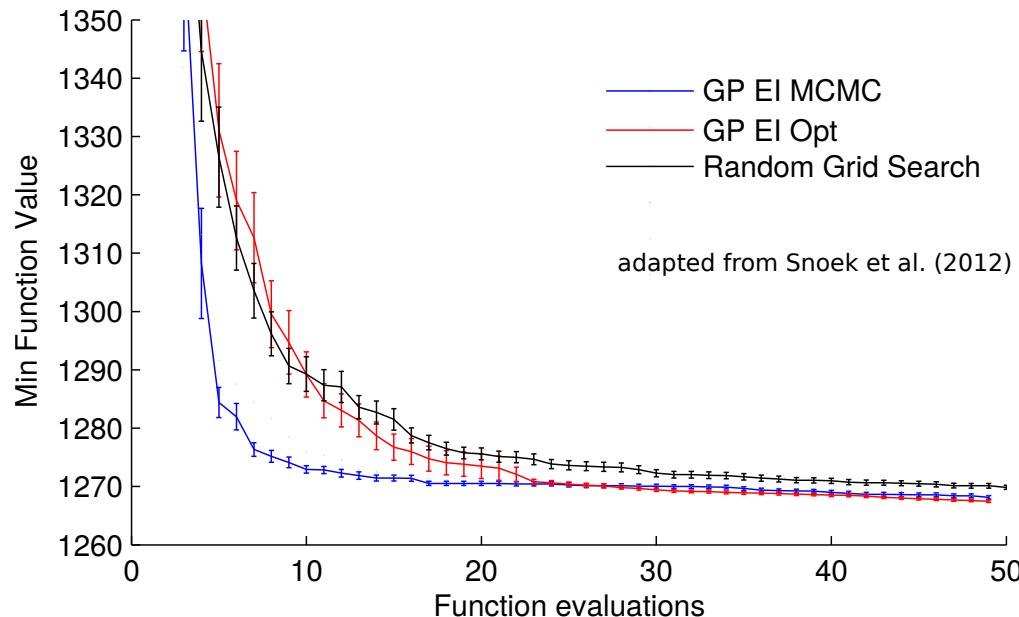
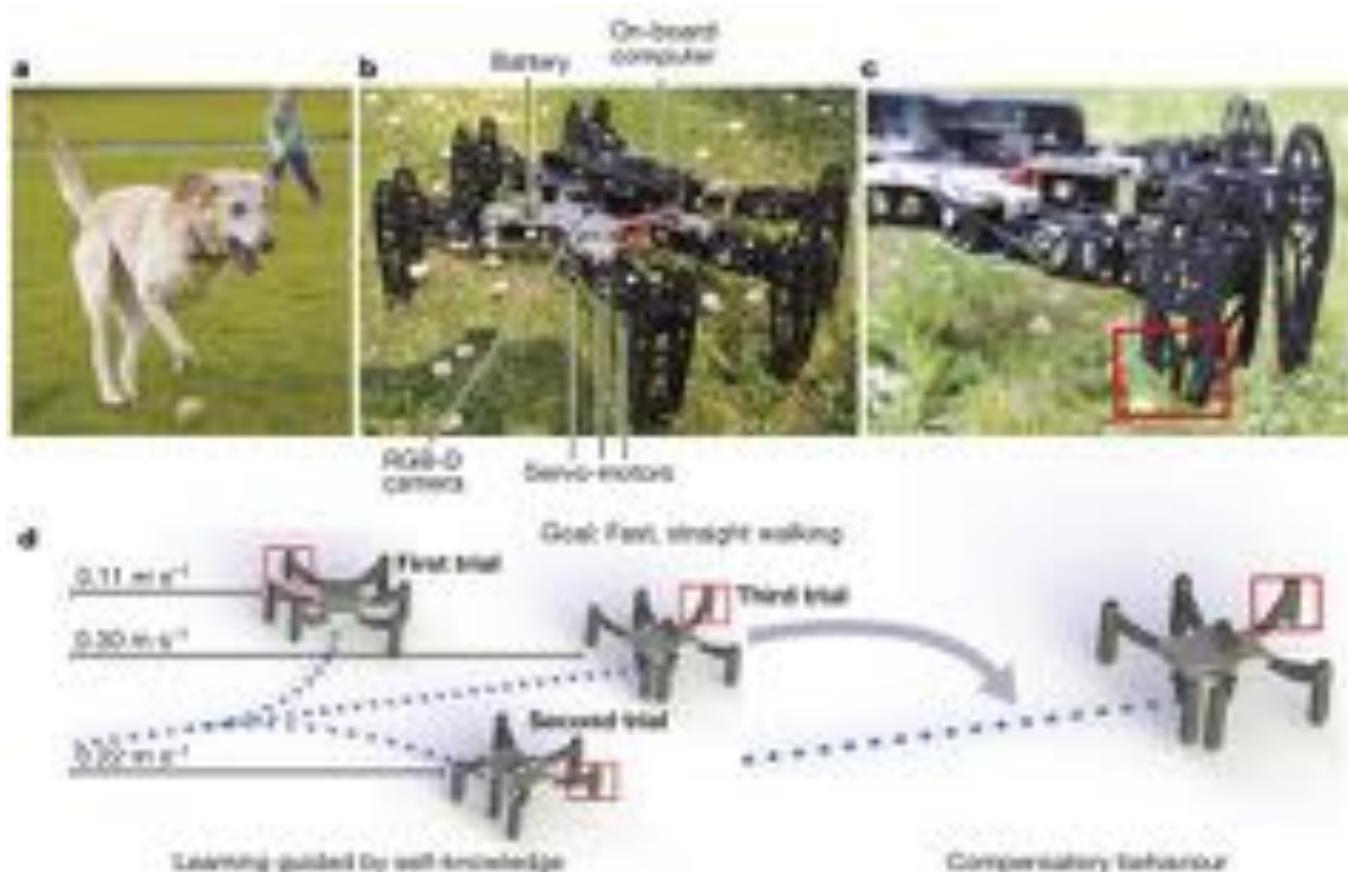


Figure: Figure from Snoek et al. (2012)

# Robots That Learn to Recover from Damage



Cully et al. (2015)

# Application Example: Controller Learning in

- Fragile bipedal robot
  - ▶ Only few experiments feasible
- Maximize robustness and walking speed
- 4 motors:
  - 2 actuated hips + 2 actuated knees
- Controller implemented as a finite-state-machine (8 parameters)



Calandra et al. (2015)

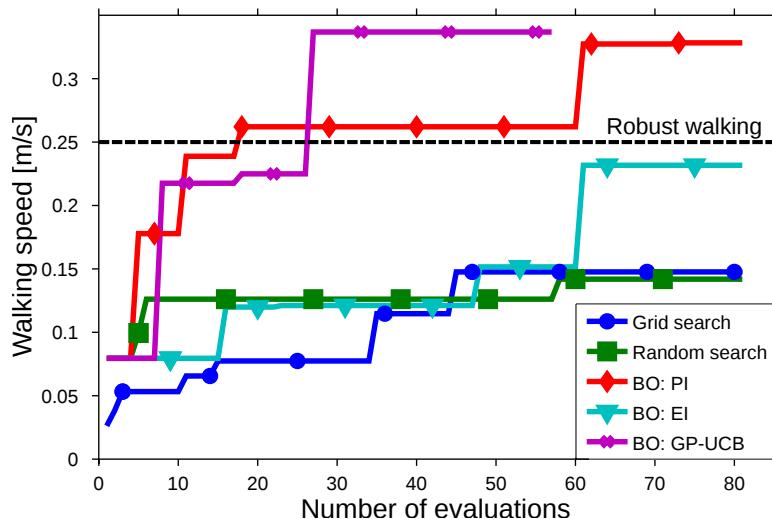
# Application Example: Controller Learning in

- Fragile bipedal robot
  - ▶ Only few experiments feasible
- Maximize robustness and walking speed
- 4 motors:
  - 2 actuated hips + 2 actuated knees
- Controller implemented as a finite-state-machine (8 parameters)
- Good parameters found after 80–100 experiments
- Substantial speed-up compared to manual parameter search



Calandra et al. (2015)

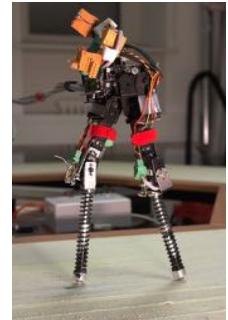
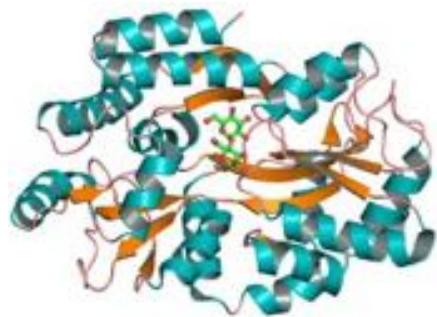
# Comparison



- Squared exponential covariance function
- Learned GP hyper-parameters (no MCMC for integrating them out)

- **Entropy-based acquisition functions:** Directly describe the distribution over the best input location (Hennig & Schuler, 2012; Hernández-Lobato et al., 2014)
- **Non-myopic** Bayesian optimization (e.g., Osborne et al., 2009)
- **High-dimensional** optimization (e.g., Wang et al., 2016)
- **Large-scale** Bayesian optimization (Hutter et al., 2014)
- **Efficient optimization of acquisition functions** (Wilson et al., 2018)
- **Non-GP** Bayesian optimization (Hutter et al., 2014; Snoek et al., 2015)
- **Constraints** (e.g., Gelbart et al., 2014)
- **Automated machine learning** (e.g., Feurer et al., 2015)
- **Multi-tasking, parallelizing, resource allocation, ...** (e.g., Swersky et al., 2014; Snoek et al., 2012; Wilson et al., 2018)

- **BoTorch** <https://github.com/pytorch/botorch>  
(Balandat et al., 2019)
- **BayesOpt**  
<https://bitbucket.org/rmcantin/bayesopt/>  
(Martinez-Cantin, 2014)
- **Spearmint** <https://github.com/HIPS/Spearmint>
- **Pybo** <https://github.com/mwhoffman/pybo> (Hoffman & Shariari)
- **GPyOpt** <https://github.com/SheffieldML/GPyOpt>  
(Gonzalez et al.)
- Matlab toolbox (bayesopt)



- Global optimization of black-box functions, which are expensive to evaluate ➤ Meta-challenges in machine learning, Auto-ML
- Use a probabilistic proxy model that is cheap to evaluate and use this to suggest next experiments
- Acquisition function trades off exploration and exploitation

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