

Bayesian Optimisation

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Optimisation of expensive cost functions

In many situations, we want to optimise processes that we do not understand. They may be:

- ▶ Take long or costly to evaluate
(physical experiments, or long-running computer simulations)
- ▶ Black-box
(we have little a-priori knowledge about them)
- ▶ Impossible to evaluate gradients of
(can't back-propagate through reality)

Black-box cost functions

- ▶ Design of turbine blades
- ▶ Biological molecules
- ▶ Optimise chemical processes
- ▶ **Selecting Machine Learning hyperparameters**



Machine Learning Meta-Challenges

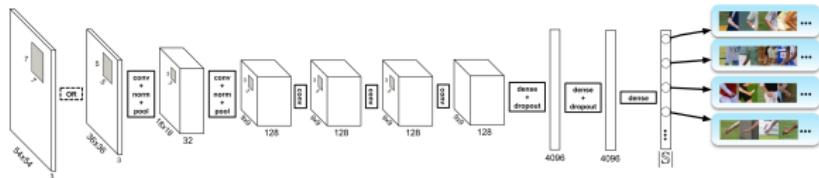
- ▶ Machine learning models are getting more and more complicated
 - ▶ Usually more parameters (e.g., deep neural networks)
- ▶ Non-convex and stochastic optimization methods have meta-parameters that are difficult to tune (learning rates, momentum parameters, ...)
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Goal: 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

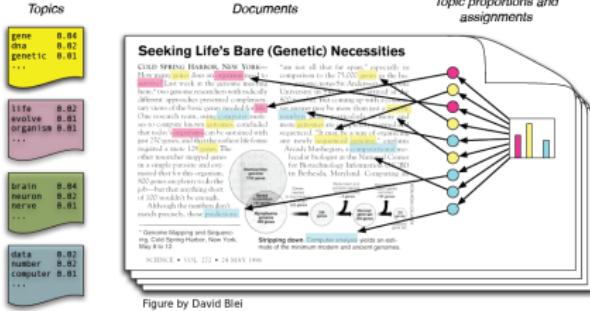
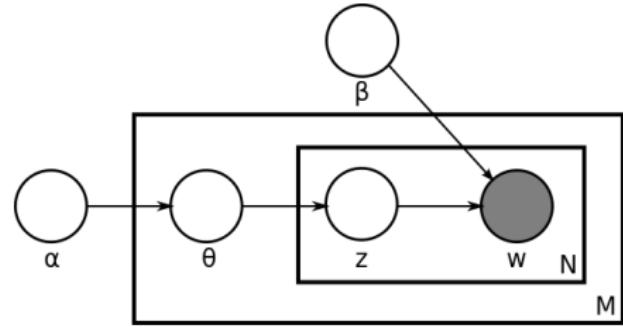
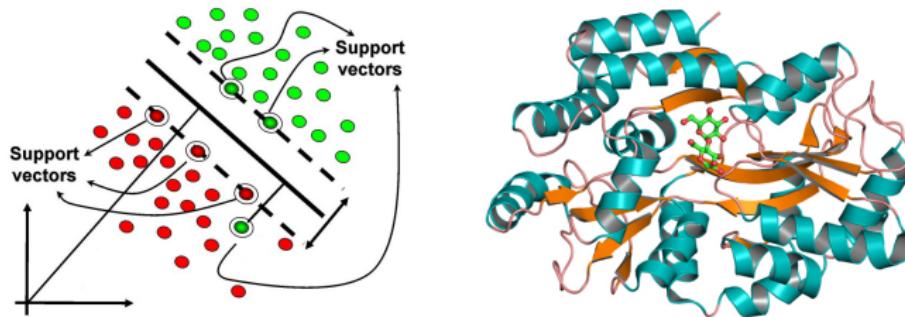


Figure by David Blei



- ▶ Hoffmann 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

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 - ▶ Grid search
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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

Alternative Approach: Bayesian Optimization

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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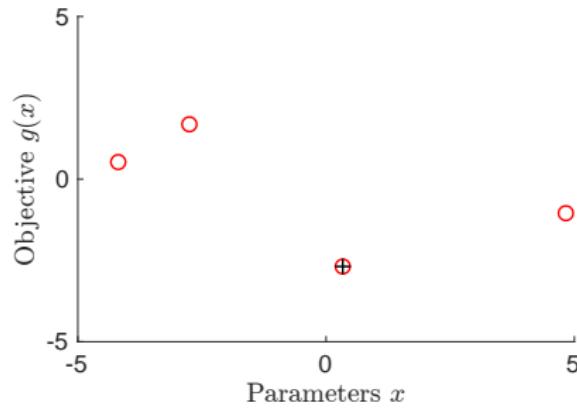
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- ▶ We can predict the outcome of the next experiment using the proxy model, which is much **cheaper to evaluate**
- ▶ **Optimize cheap proxy** function to determine where to evaluate the true objective next
- ▶ Standard proxy: **Gaussian process**

Setting (2)

- ▶ Objective: Find global minimum of objective function f :

$$\mathbf{x}_* = \arg \min_x f(\mathbf{x})$$

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



Bayesian optimisation

Let's phrase BayesOpt using decision theory:

$$L(f, \{\mathbf{x}_n\}_{n=1}^N) = Nc + \min_{1 \leq n \leq N} f(\mathbf{x}_n) \quad (1)$$

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- ▶ We decide how many evaluations we want to do N
- ▶ We decide where to evaluate $f(\cdot)$
- ▶ We incur a cost c for each evaluation
- ▶ We reduce loss by finding a lower $f(\mathbf{x})$

Bayesian optimisation – decision theory

For 3 observations:

$$\{\mathbf{x}_1^*, \mathbf{x}_2^*(\mathcal{D}_1), \mathbf{x}_3^*(\mathcal{D}_2)\} = \operatorname*{argmin}_{\{\mathbf{x}_n\}_{n=1}^3} \mathbb{E}_{p(f)} \left[Nc + \min_{1 \leq n \leq N} f(\mathbf{x}_n) \right] \quad (2)$$

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$$\mathbf{x}_1^* = -Nc + \operatorname{argmin}_{\mathbf{x}_1} \mathbb{E}_f \left[\min_{\mathbf{x}_2} \mathbb{E}_{f|\mathcal{D}_1} \left[\min_{\mathbf{x}_3} \mathbb{E}_{f|\mathcal{D}_2} \left[\min_{1 \leq n \leq 3} \{f(\mathbf{x}_n)\} \right] \right] \right] \quad (3)$$

Computational difficulties

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This is **very** difficult to evaluate.

- ▶ Difficult optimisations: closed-form optimum cannot be found
- ▶ Expectations of non-closed-form expressions
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Alternative approaches:

- ▶ Can try to approximate these quantities (fun to try for small-scale examples)
- ▶ Heuristic approaches

How to design a heuristic

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- ▶ Desired behaviour: Some sort of a trade-off between **exploration** and **exploitation**.

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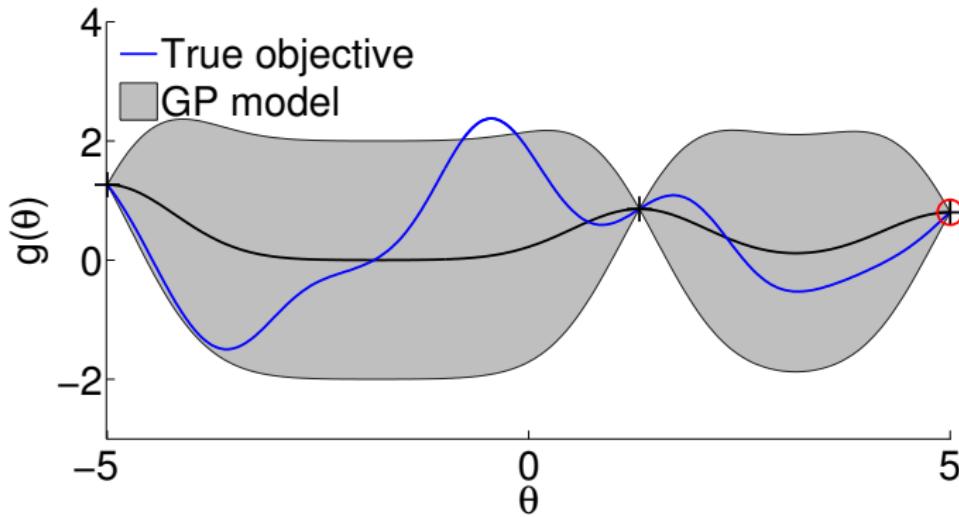
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- ▶ Computational constraint: Only use posterior of GP given **current** observations to choose current action
- ▶ Desired behaviour: Some sort of a trade-off between **exploration** and **exploitation**.
 - ▶ Want to try inputs that are different (exploration)
 - ▶ Want to focus on areas that are likely to have high values (exploitation)

Inadequate acquisition functions

Two inadequate acquisition functions:

- ▶ Minimise posterior mean (pure exploitation)
- ▶ Maximise posterior variance (pure exploration)



Common 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

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

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- Probability of Improvement (Kushner 1964):**

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

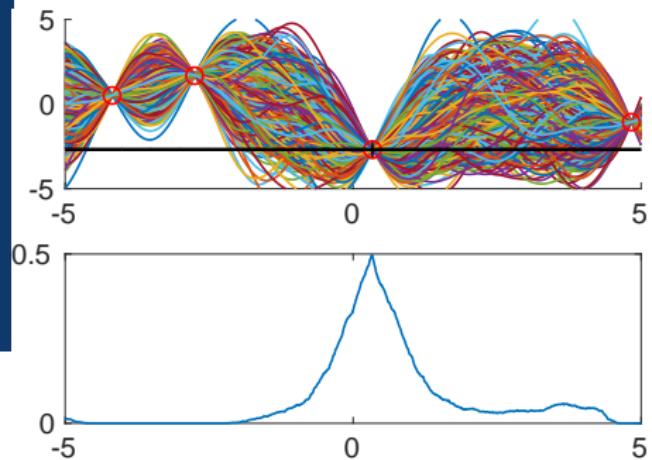
- Expected Improvement (Mockus 1978):**

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

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

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

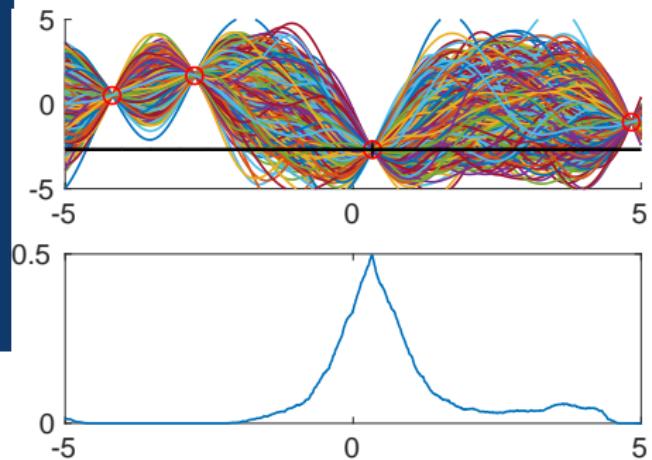
Probability of Improvement (1)



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

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

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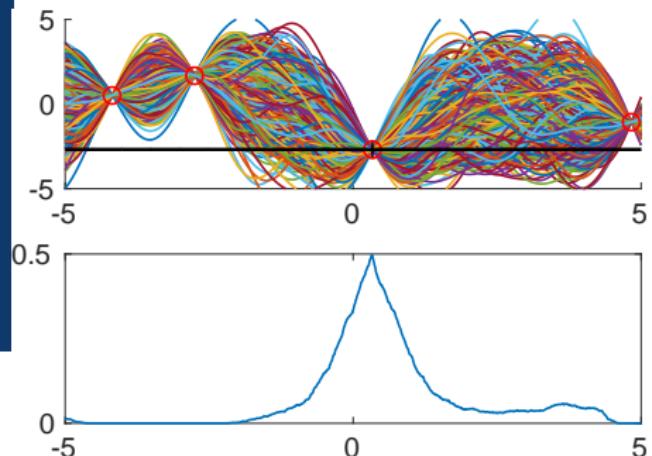


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- ▶ Can lead to continued exploitation in an ϵ -region around x_{best} .

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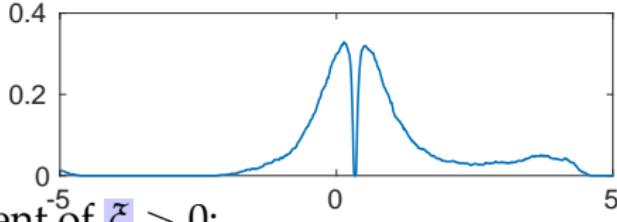
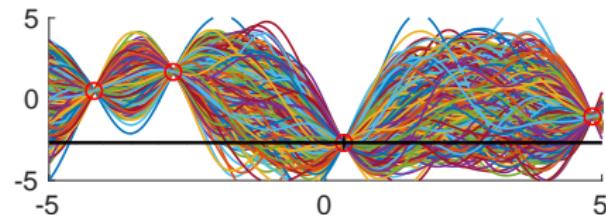
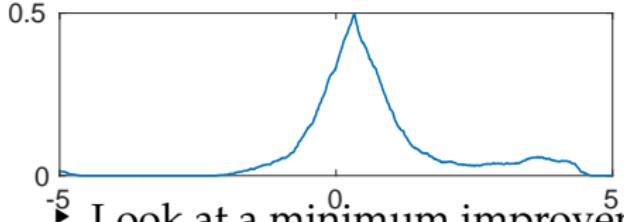
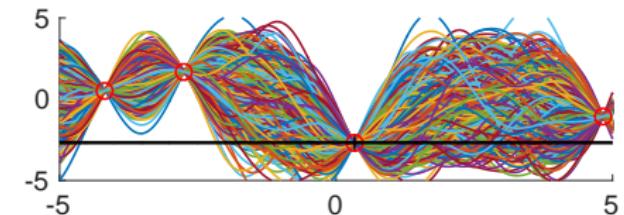


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- ▶ Can lead to continued exploitation in an ϵ -region around x_{best} .
- ▶ Introduce a “slack variable” ξ for more aggressive exploration

Probability of Improvement (2)



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

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

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

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

Expected Improvement

- ▶ Idea: Quantify the amount of improvement
- ▶ Sampling-based scenario, where $f_i \sim p(f)$:

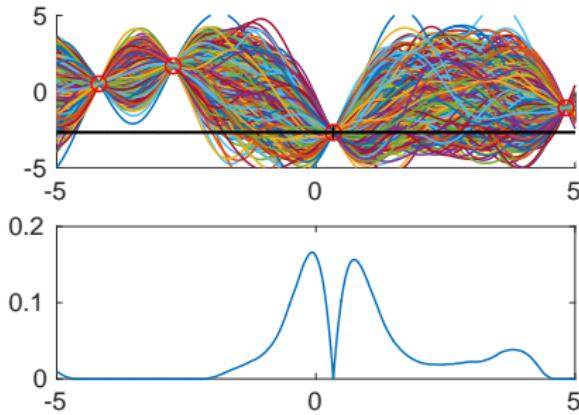
$$\alpha_{\text{EI}}(\mathbf{x}) = \mathbb{E}[\max\{0, f(\mathbf{x}_{\text{best}}) - f(\mathbf{x})\}]$$

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

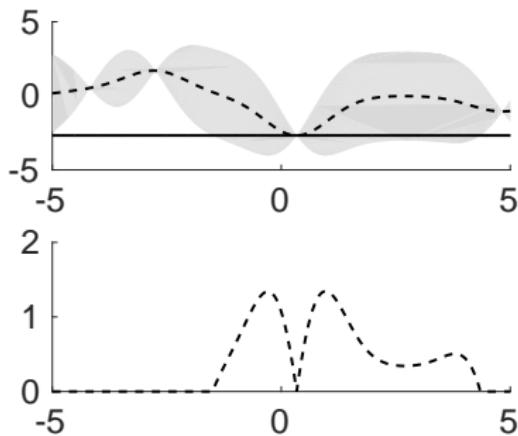
- ▶ If $f \sim GP$, we have a closed-form expression:

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

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



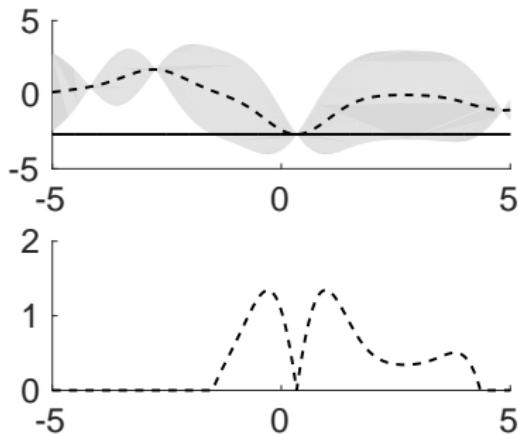
GP-Lower Confidence Bound (1)



- ▶ 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

$$\alpha_{LCB}(x_t) = -(\mu(x_t) - \sqrt{\kappa}\sigma(x_t))$$

GP-Lower Confidence Bound (2)



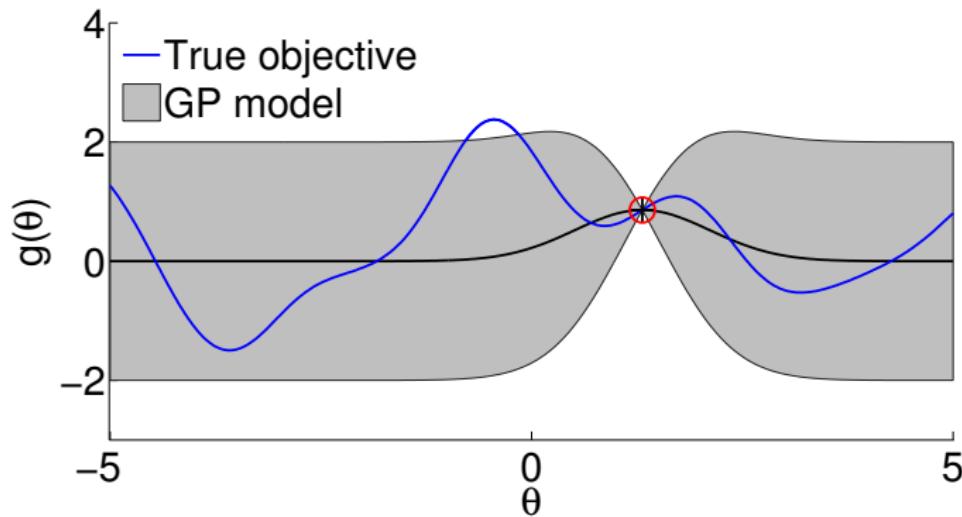
- More generally, we can get regret bounds for iteration-dependent κ (Srinivas et al., 2010)

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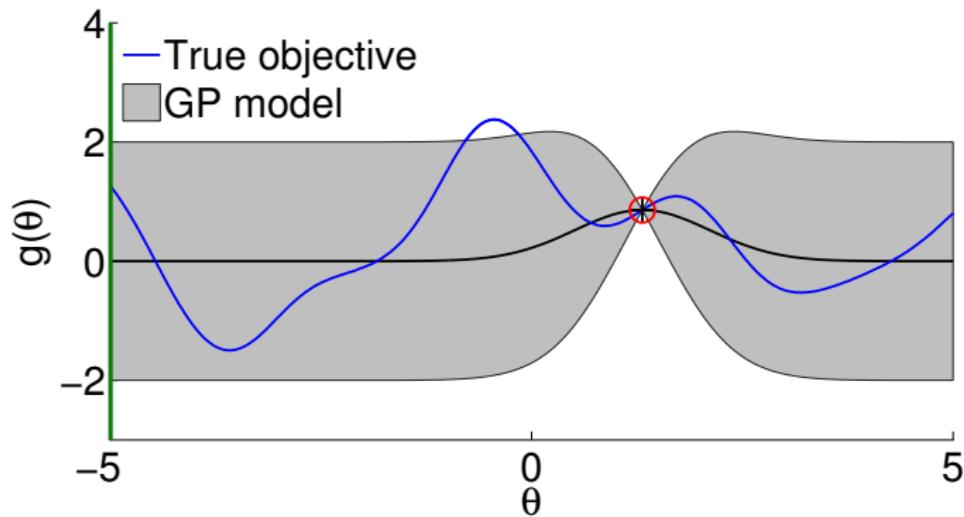
where $\kappa_t \in \mathcal{O}(\log t)$ grows with the iteration t

► Continue exploration

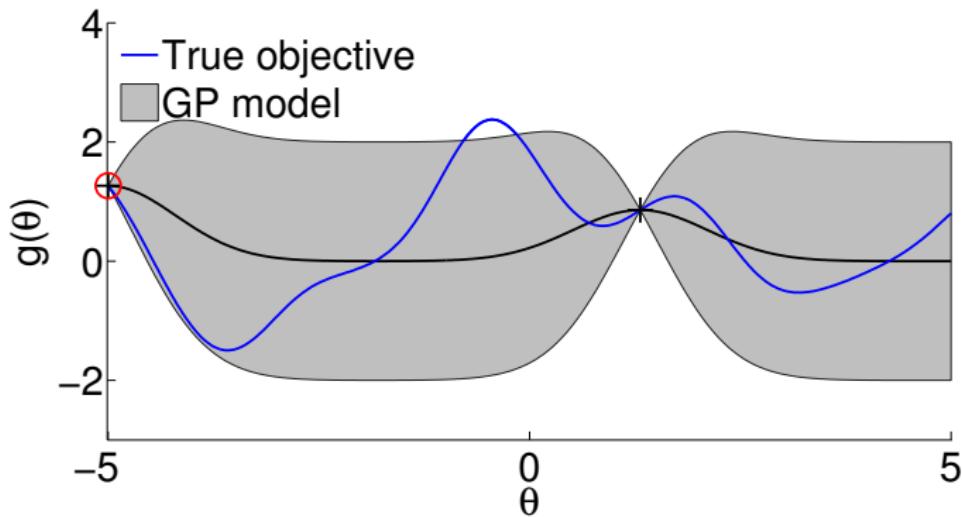
Bayesian Optimization: Illustration



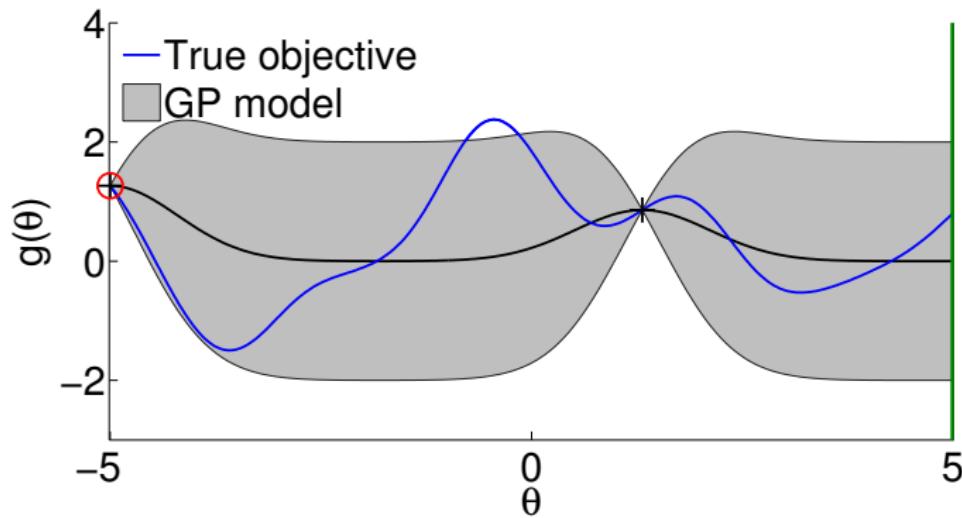
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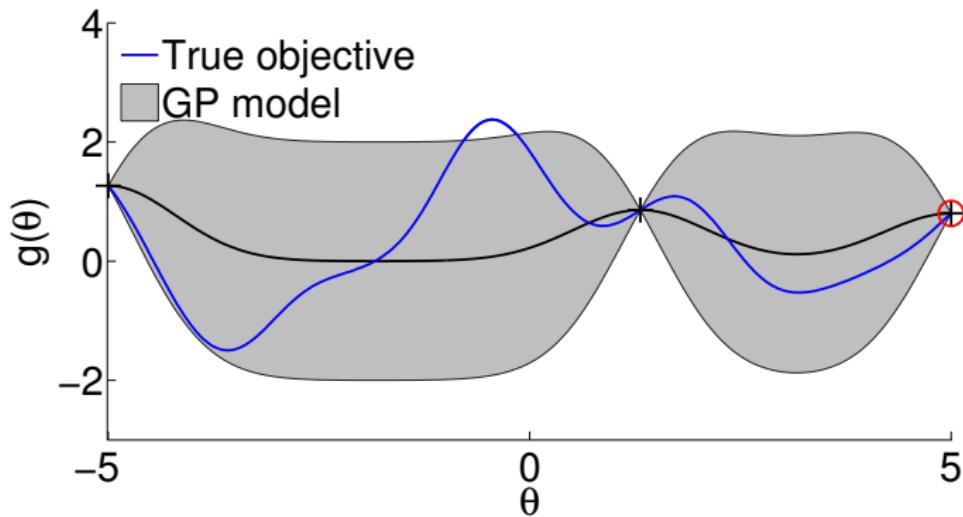
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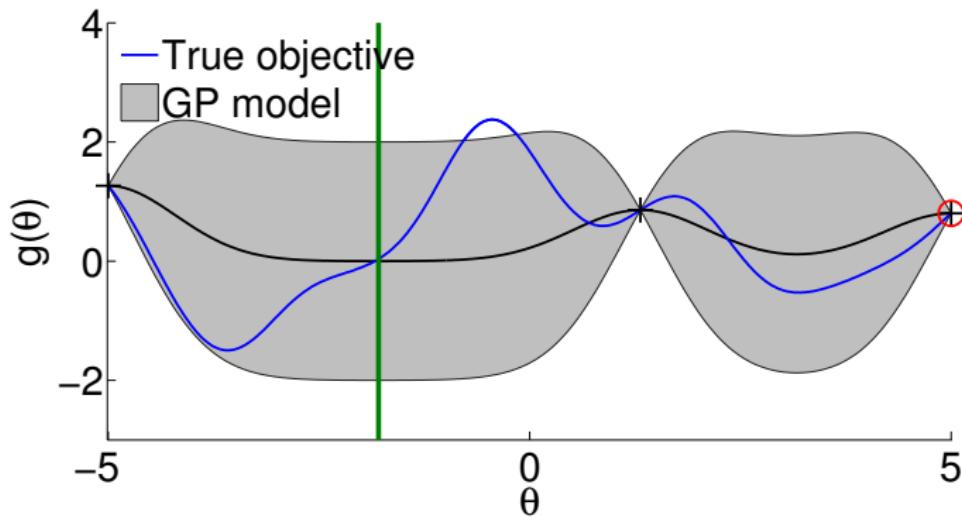
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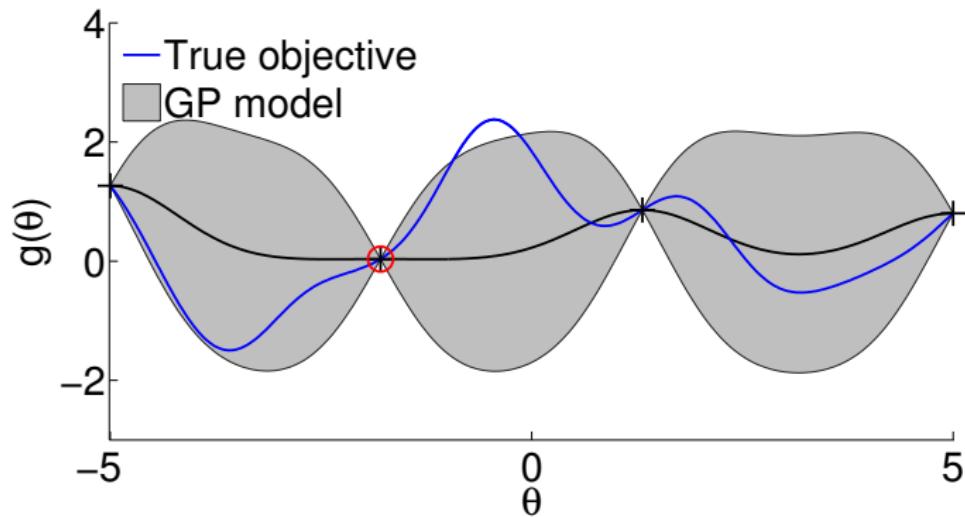
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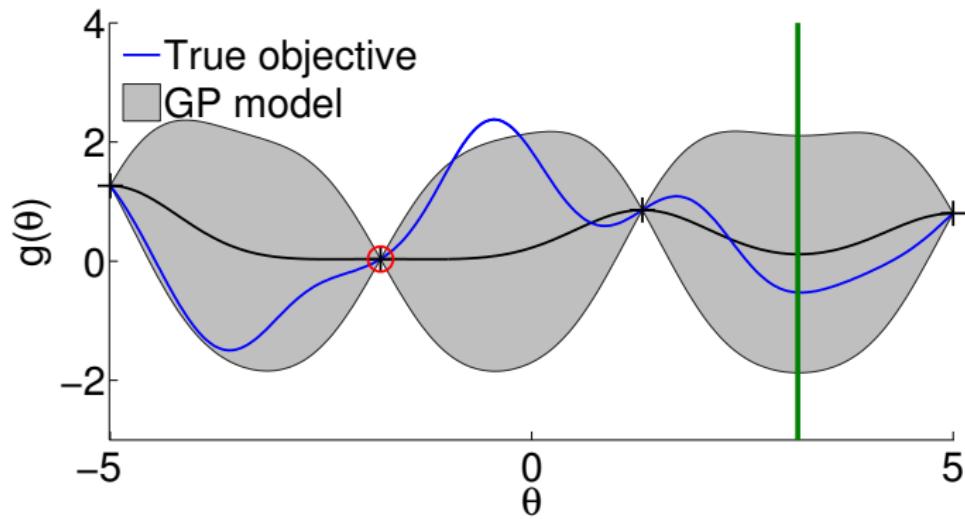
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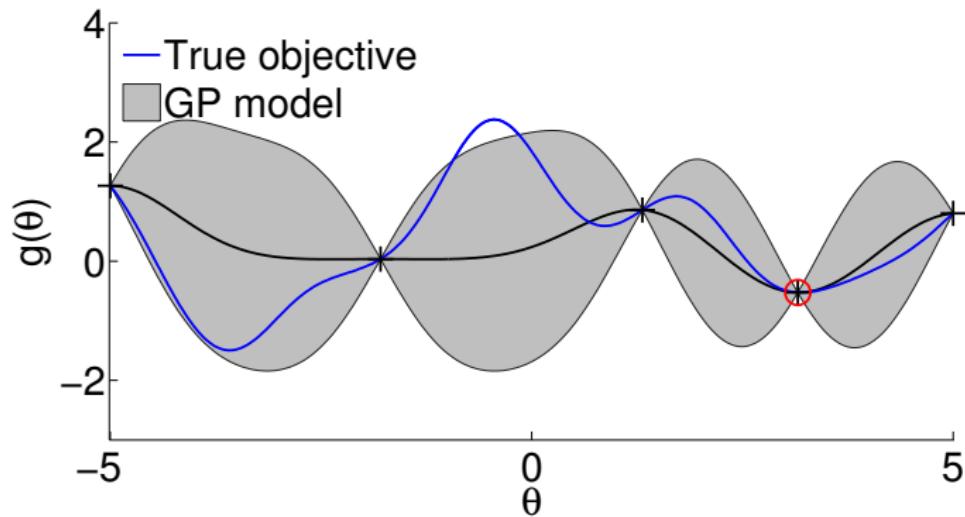
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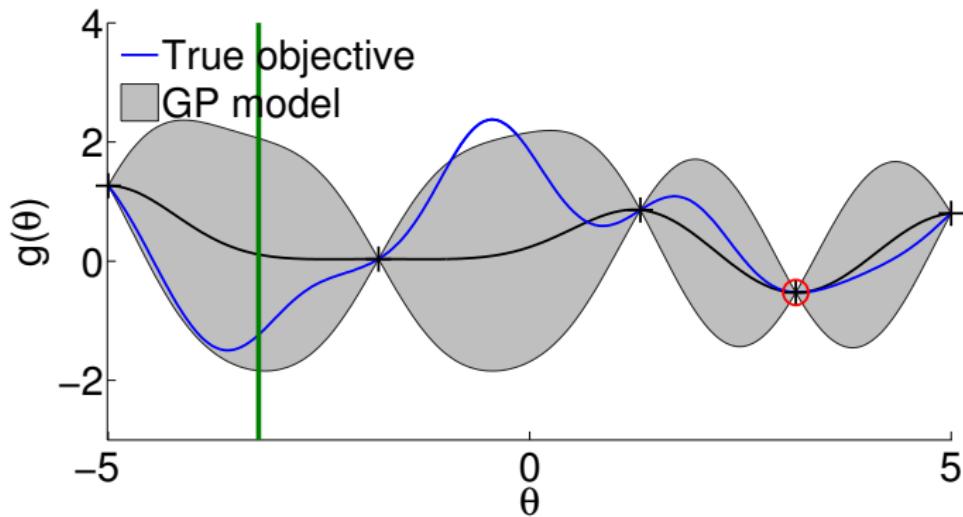
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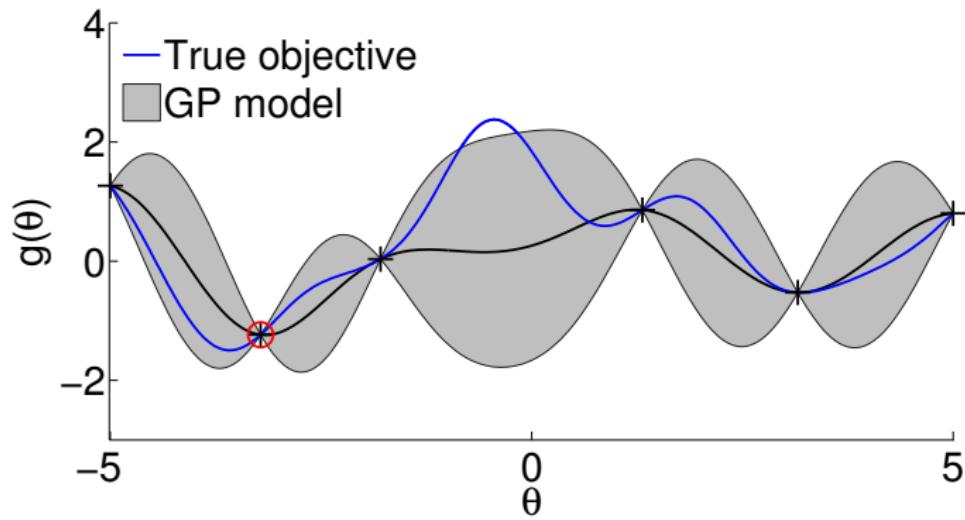
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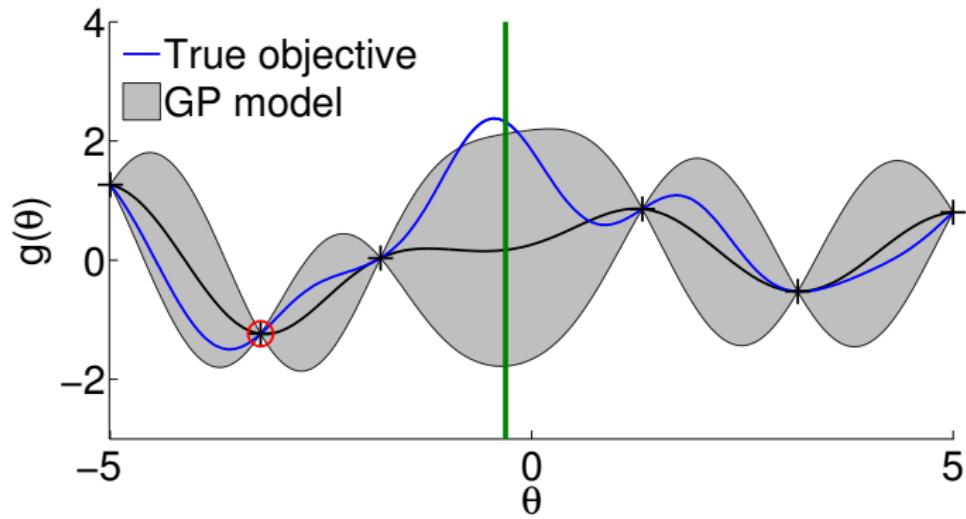
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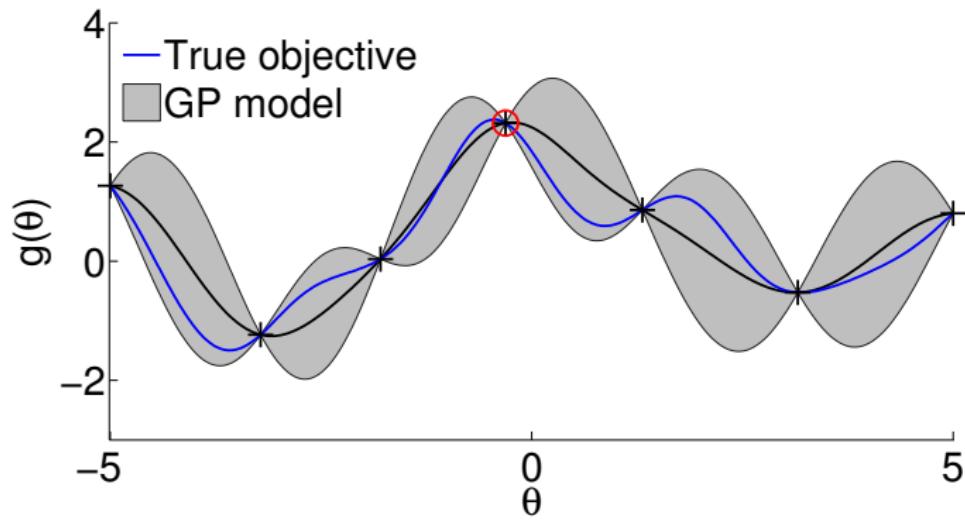
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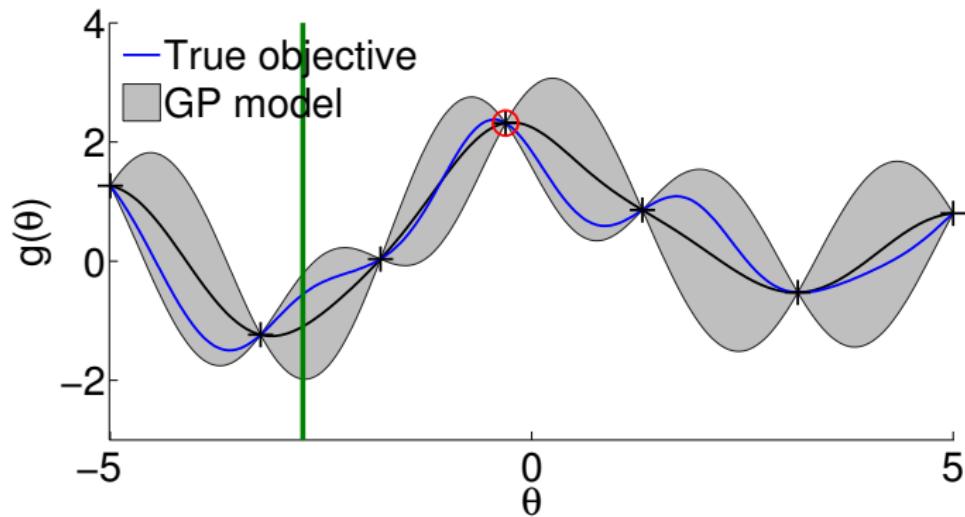
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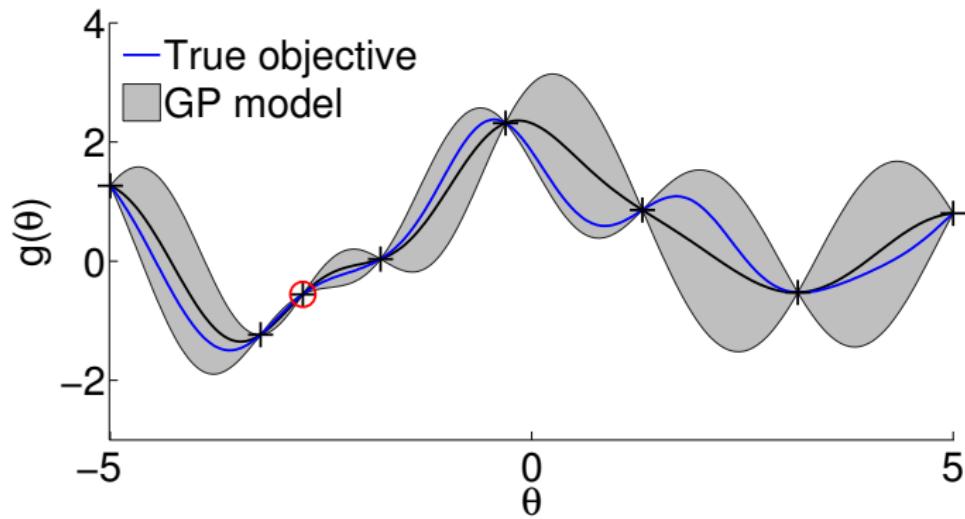
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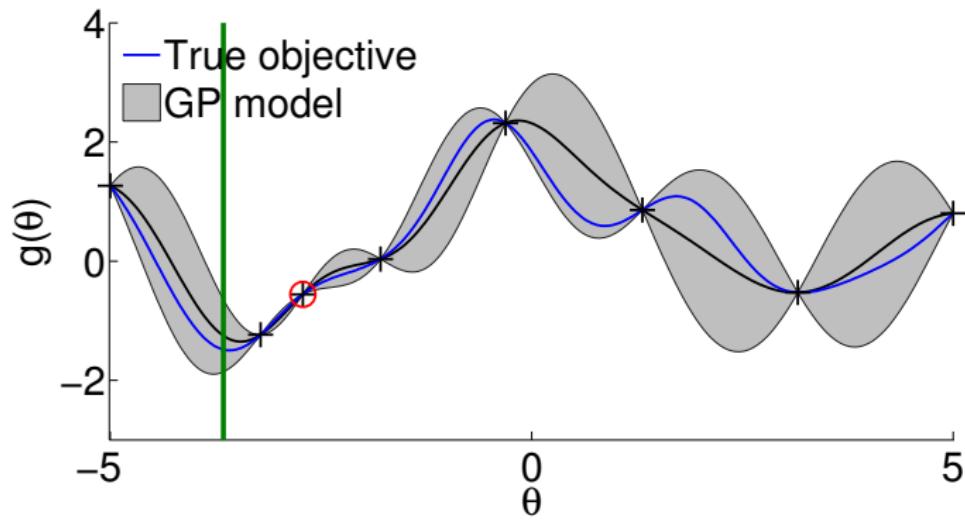
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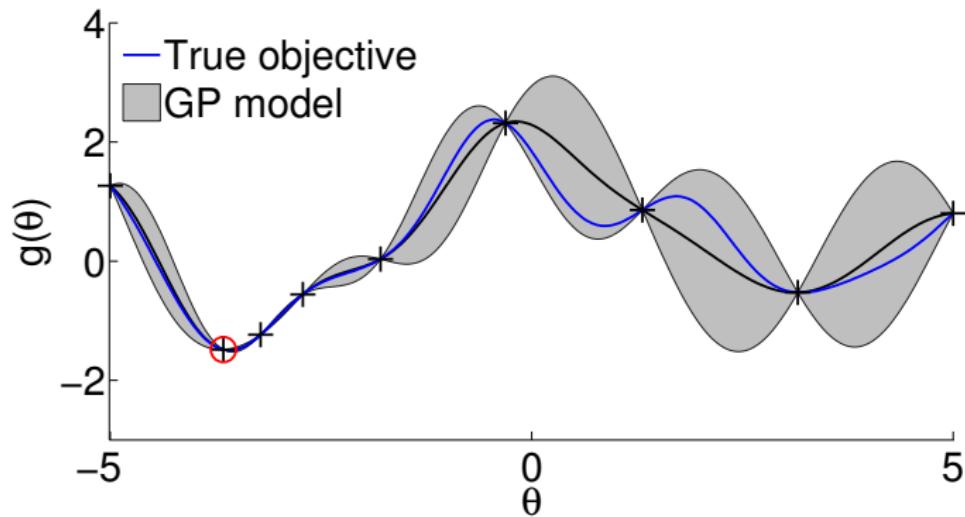
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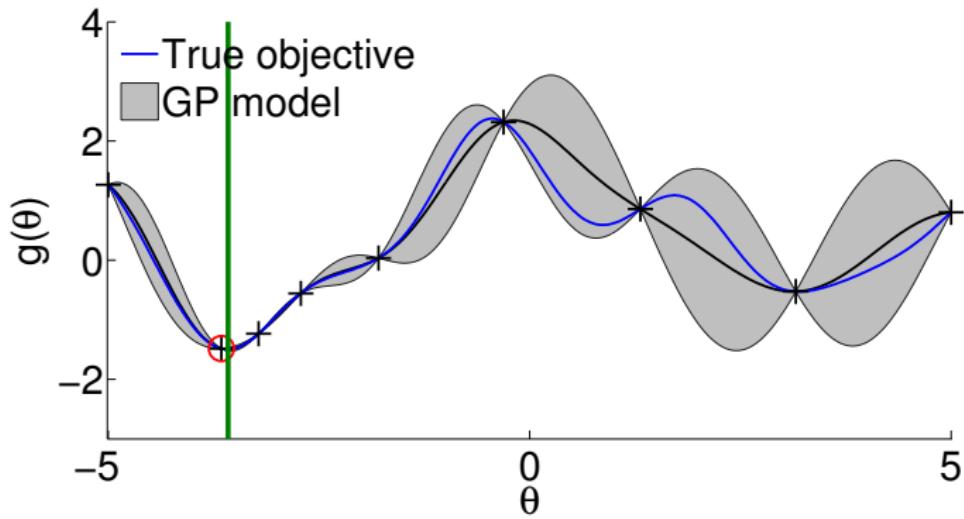
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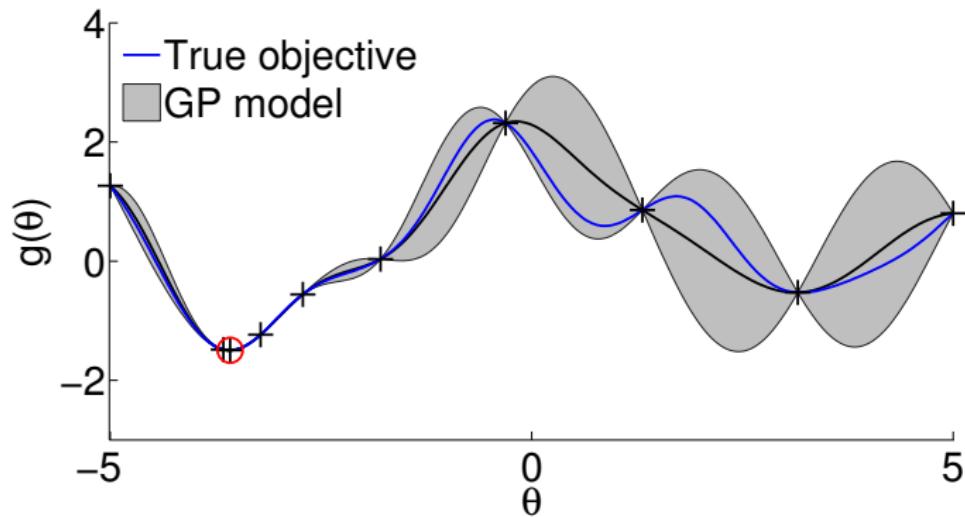
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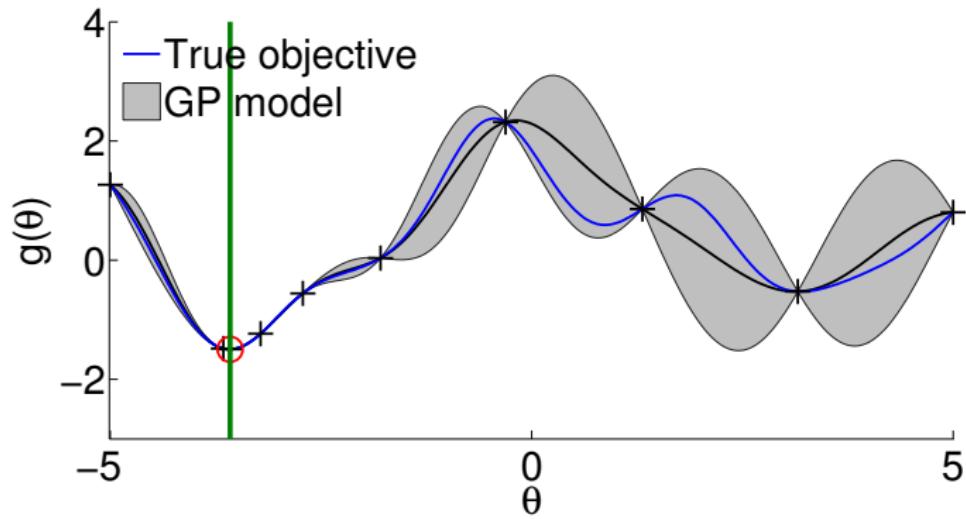
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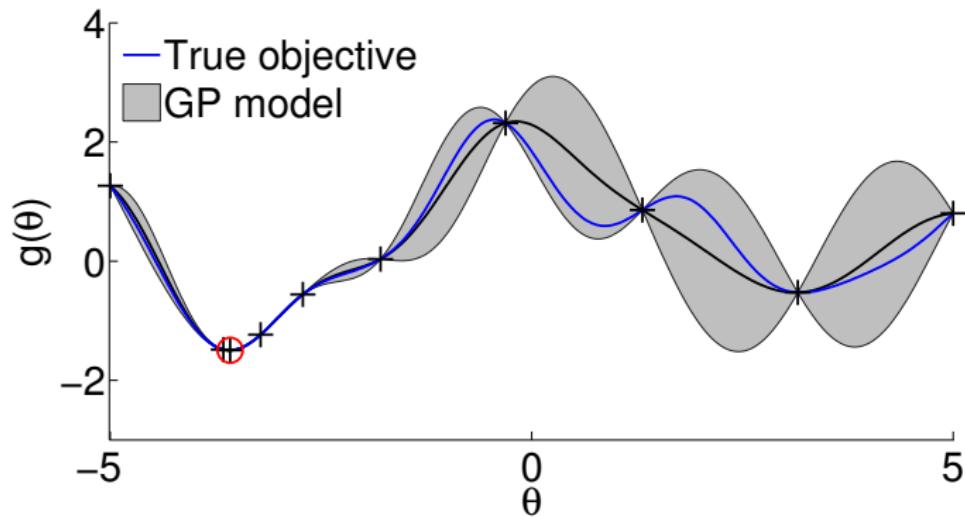
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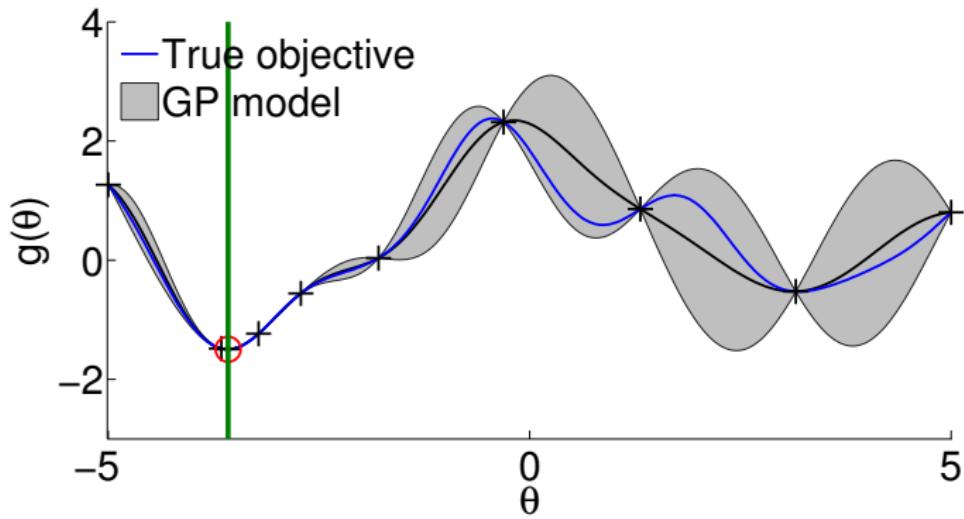
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Optimising the Acquisition Function

Notice: Acquisition function is multi-modal

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

Optimising the Acquisition Function

Notice: Acquisition function is multi-modal

- ▶ 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
 - ▶ Apply the usual tricks of many random restarts

Key Steps (Pseudo-Code)

- 1: **Init:** Data set $\mathcal{D}_0 = \{X_0, y_0\}$ (can be empty)
- 2: **for** iterations $t = 1, 2, \dots$ **do**
- 3: Update GP using data \mathcal{D}_{t-1}
- 4: Select $x_t = \arg \max_x \alpha(x)$ by optimising acquisition function
- 5: Query true objective g at x_t
- 6: Augment data set $\mathcal{D}_t = \mathcal{D}_{t-1} \cup \{(x_t, y_t)\}$
- 7: **end for**
- 8: **Return** best input in data set: $x^* = \arg \min_x y(x)$

Optimising acquisition function is itself an iterative process, done with a numerical gradient-based optimiser (e.g. BFGS).

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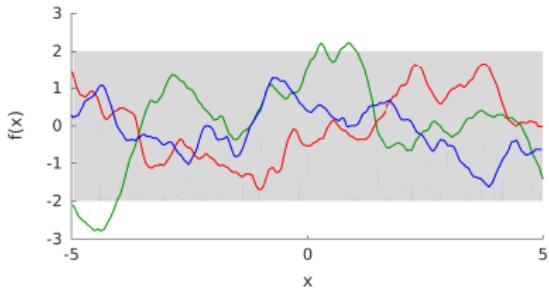
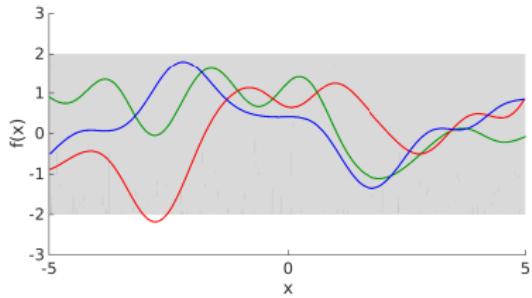
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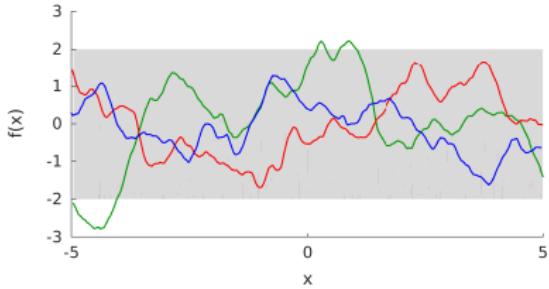
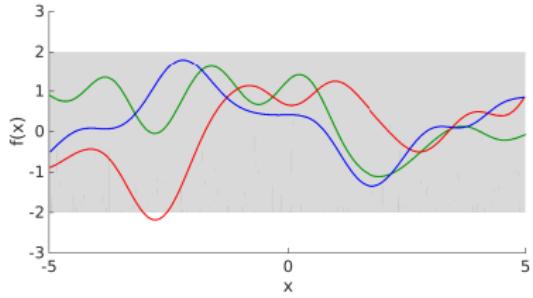
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- ▶ Covariance function selection is crucial for good performance
 - ▶ Choose a sufficiently flexible and adaptive kernel, e.g., Matérn (but not the squared exponential (Gaussian))
- ▶ Nice side-effect of Matérn: Exploration is more encouraged than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

Choosing Covariance Functions

Application:

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

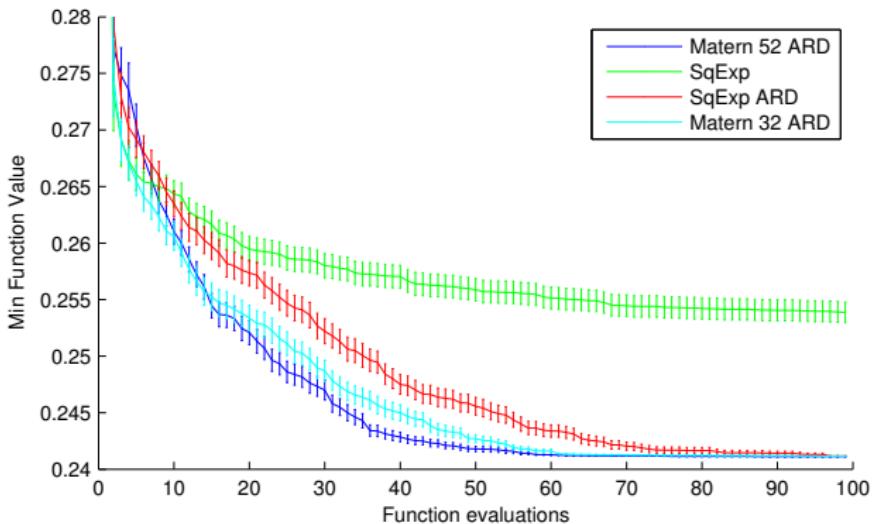


Figure: Figure from Snoek et al. (2012)

Gaussian Process Hyper-Parameters

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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 θ by Markov Chain Monte Carlo (MCMC) sampling (e.g., slice sampling)
- ▶ Look at integrated acquisition function

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

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)

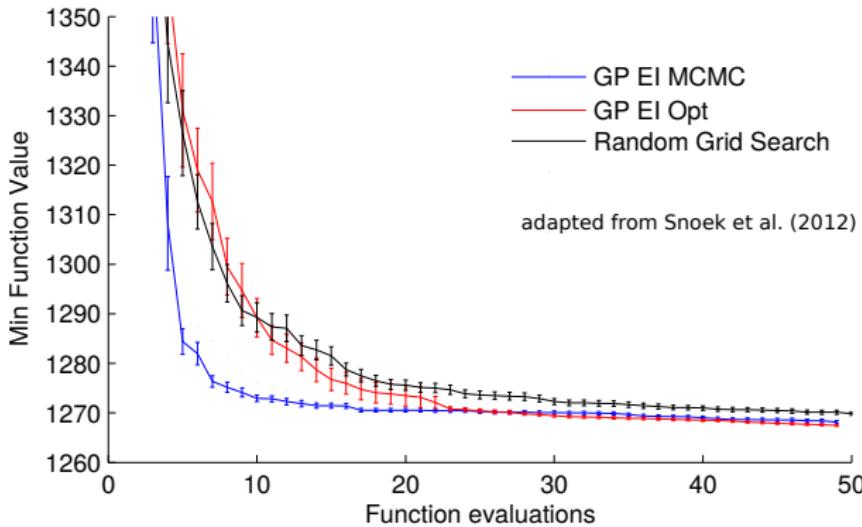
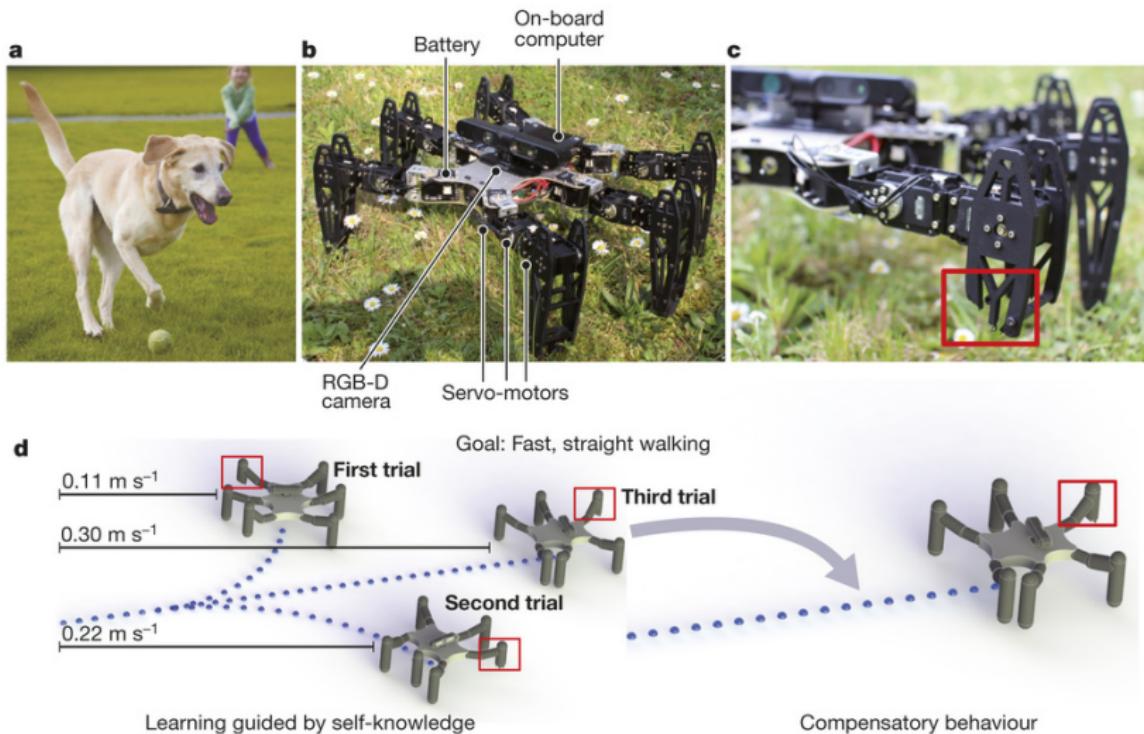


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Robots That Learn to Recover from Damage



Cully et al. (2015)

Application Example: Controller Learning in Robotics (Calandra et al., 2015)

- ▶ 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)

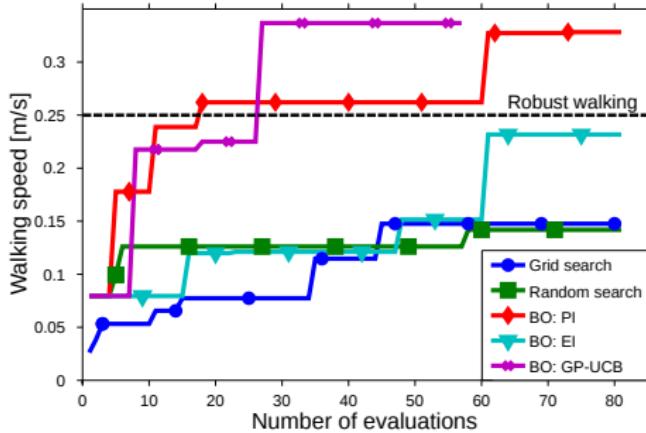
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 - 2 actuated hips + 2 actuated knees
- ▶ Controller implemented as a finite-state-machine (8 parameters)
- ▶ Good parameters found after 80–100 experiments
- ▶ Substantial speed-up compared to manual parameter search



Calandra et al. (2015)

Comparison



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

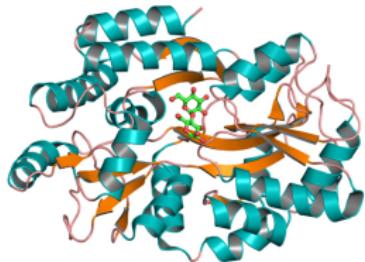
Further Topics in BO

- ▶ **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)

Software

- ▶ **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)

Summary



- ▶ 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

Further Reading

- ▶ 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

References I