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Bayesian and evolutionary optimization: searching from both sides

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11 July 2023

talk at the Hasso-Plattner Institut, Potsdam

Context: global optimization

$$\min_{x \in \mathcal{S}} f(x)$$

\mathcal{S} : search space, continuous, discrete, mixed, others (graphs?).
Default $\mathcal{S} \in \mathbb{R}^d$ (hyper-rectangle). d is the dimension.

Both evolutionary and Bayesian optimization algorithms (EO & BO) apply.
EO & BO have different scientific communities (computer science vs. applied math, but it is changing).

Goal of the talk : introduce BO through a comparison with EO and draw common perspectives.

Context: BO for costly optimization

$$\min_{x \in \mathcal{S}} f(x)$$

Costly: one call to f takes more CPU than the rest of the optimization algorithm. Examples: nonlinear partial differential equations (finite elements), training of a neural network, real experiment . . .

To save calls to f , build a model of it based on previous evaluations and rely on it whenever possible \rightarrow metamodel / surrogate based optimization. **Gaussian process as metamodel : Bayesian Optimization (BO)**.

Beginnings and references

EO & BO both date from the 60s or 70s with the work of

- [Fogel et al., 1965], [Holland, 1973], [Rechenberg, 1973] and [Schwefel, 1977] for EO,
- and [Kushner, 1962], [Močkus, 1972] and [Saltenis, 1971] for BO.

Reference texts for BO : [Frazier, 2018],[Garnett, 2023].

Algorithms skeletons I

BO

- 1 make an initial design of experiments (DoE) \mathbb{X}^t and calculate $\mathbb{F}^t = f(\mathbb{X}^t)$, $t = \text{length}(\mathbb{F})$
- 2 build a Gaussian Proc. from $(\mathbb{X}^t, \mathbb{F}^t) \rightarrow GP^t$
- 3 $x^{t+1} = \arg \max_{x \in \mathcal{S}} \text{AcquisCrit}(x; GP^t)$
- 4 $(\mathbb{X}^{t+1}, \mathbb{F}^{t+1}) = (\mathbb{X}^{t+1}, \mathbb{F}^{t+1}) \cup (x^{t+1}, f(x^{t+1}))$, increment t
- 5 stop ($t > t^{\max}, \dots$) or go to 2.

EO

- 1 make an initial DoE \mathbb{X} and calculate $\mathbb{F} = f(\mathbb{X})$, $t = \text{length}(\mathbb{F})$
- 2 Apply variation operators (selection, mutation, crossover) to get a new DoE:
 $\mathbb{X}' = \text{Variation}(\mathbb{X}, \mathbb{F})$
- 3 calculate $\mathbb{F}' = f(\mathbb{X}')$,
 $t = t + \text{length}(\mathbb{F}')$
- 4 Create a new (\mathbb{X}, \mathbb{F}) from old (\mathbb{X}, \mathbb{F}) and $(\mathbb{X}', \mathbb{F}')$
- 5 stop ($t > t^{\max}, \dots$) or go to 2.

Algorithms skeletons II

BO

- 1 make an initial design of experiments (DoE) \mathbb{X}^t and calculate $\mathbb{F}^t = f(\mathbb{X}^t)$, $t = \text{length}(\mathbb{F})$
- 2 build a Gaussian Proc. from $(\mathbb{X}^t, \mathbb{F}^t) \rightarrow GP^t$
- 3 $x^{t+1} = \arg \max_{x \in \mathcal{S}} \text{AcquisCrit}(x; GP^t)$
- 4 $(\mathbb{X}^{t+1}, \mathbb{F}^{t+1}) = (\mathbb{X}^t, \mathbb{F}^t) \cup (x^{t+1}, f(x^{t+1}))$, increment t
- 5 stop ($t > t^{\max}, \dots$) or go to 2.

EO

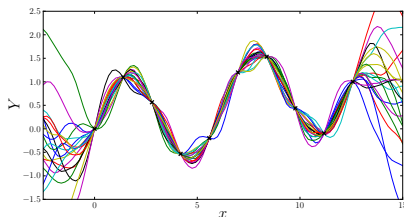
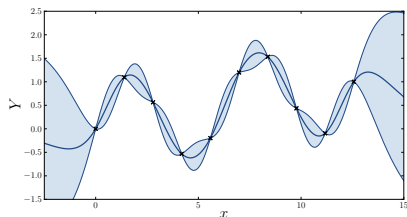
- 1 make an initial DoE \mathbb{X} and calculate $\mathbb{F} = f(\mathbb{X})$, $t = \text{length}(\mathbb{F})$
- 2 Apply variation operators (selection, mutation, crossover) to get a new DoE: $\mathbb{X}' = \text{Variation}(\mathbb{X}, \mathbb{F})$
- 3 calculate $\mathbb{F}' = f(\mathbb{X}')$, $t = t + \text{length}(\mathbb{F}')$
- 4 Create a new (\mathbb{X}, \mathbb{F}) from old (\mathbb{X}, \mathbb{F}) and $(\mathbb{X}', \mathbb{F}')$
- 5 stop ($t > t^{\max}, \dots$) or go to 2.

Differences:

- Maximization of an **acquisition criterion** that depends on a Gaussian Process instead of the variation operators;
- EO is population oriented while all points are kept in BO. BO would be memory consuming for large t but not its typical range of use. The state of EO is implicitly stored in its population, efficient but loss prone.

A SHORT INTRODUCTION TO BAYESIAN OPTIMIZATION

Gaussian Process Regression (kriging) – I



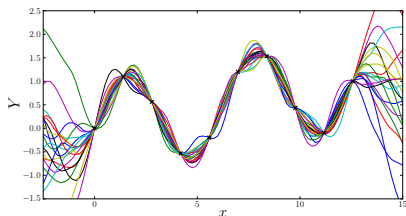
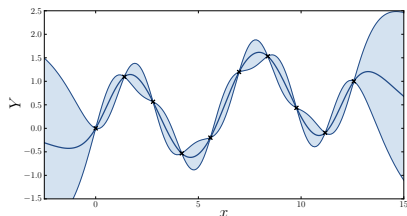
$Y(x)$ is $\mathcal{N}(\mu(x), k(x, x'))$

$Y(x) | Y(\mathbb{X}) = \mathbb{F} \sim \mathcal{N}(m(x), c(x, x'))$ is also Gaussian, interpolating and depends on the **kernel** $k(., .)$ and $\mu(.)$ through parameters θ .

Example: $k(x, x') = \sigma^2 \exp(-\sum_{i=1}^d \frac{(x_i - x'_i)^2}{2\theta_i^2})$.

Note: kernels are defined over all kinds of spaces. Example of mixed continuous-discrete BO in [Cuesta-Ramirez et al., 2022].

Gaussian Process Regression (kriging) – II



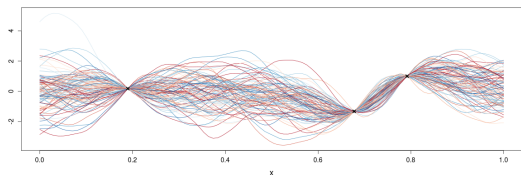
$Y(x) | Y(\mathbb{X}) = \mathbb{F} \sim \mathcal{N}(m(x), c(x, x'))$, interpolating and depends on the kernel $k(., .)$ and $\mu(.)$ through parameters θ .

$$m(x) = \mathbb{E}[Y(x) | Y(\mathbb{X}) = \mathbb{F}] = \mu(x) + k(x, \mathbb{X})k(\mathbb{X}, \mathbb{X})^{-1}(\mathbb{F} - \mu(\mathbb{X})\mathbb{1})$$
$$c(x, x') = \text{Cov}[Y(x), Y(x') | Y(\mathbb{X}) = \mathbb{F}] = k(x, x') - k(x, \mathbb{X})k(\mathbb{X}, \mathbb{X})^{-1}k(\mathbb{X}, x')$$

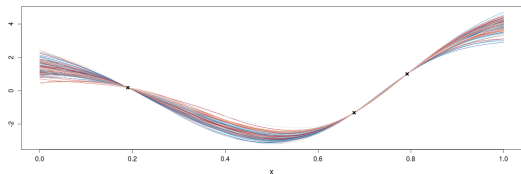
Learn the GP typically by max. (log) likelihood or leave-one-out crossvalidation, $\theta^* = \arg \max_{\theta} LL(\theta; \mathbb{F})$.

Gaussian Process Regression (kriging)

θ 's as length scales, $k(x, x') = \sigma^2 \prod_{i=1}^d \text{correlation}_i \left(\frac{|x_i - x'_i|}{\theta_i} \right)$



$\theta = 0.1$



$\theta = 0.5$

(Matérn kernel, $\sigma = 1$)

An example of acquisition criterion: the $\mathbb{E}I$

Measure of progress: the improvement,

$$I(x) = \max(0, (\min(\mathbb{F}) - Y(x) \mid Y(\mathbb{X})=\mathbb{F})).$$

Acquisition criterion: $\mathbb{E}I$ [Saltens, 1971, Schonlau, 1997], to maximize at each iteration.

The $\mathbb{E}I$ is analytically known and does not involve calls to f ,

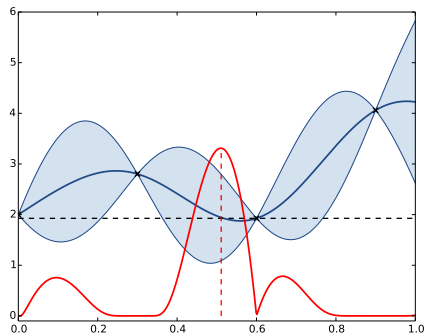
$$\mathbb{E}I(x) = \int_{-\infty}^{+\infty} I(x) dy(x) = \dots =$$

$$\sqrt{c(x, x)} [w(x)\text{cdf}_{\mathcal{N}}(w(x)) + \text{pdf}_{\mathcal{N}}(w(x))] \\ \text{with } w(x) = \frac{\min(F) - m(x)}{\sqrt{(c(x, x))}}.$$

A parameter-less intensification (1st term) – exploration (2nd term) compromise.

Expected Improvement

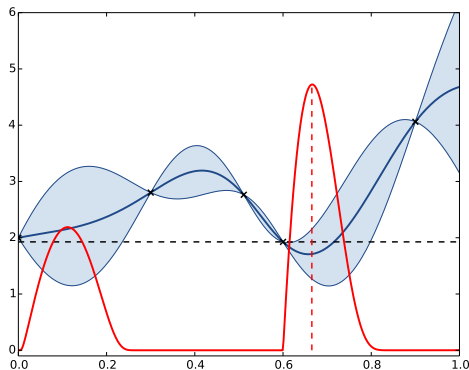
Let's see how it works... iteration 4



Expected Improvement

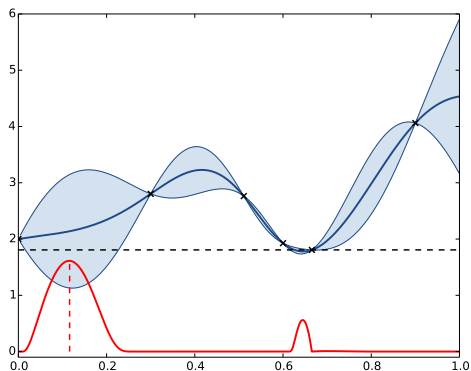
$$x^{t+1} = \arg \max_{x \in \mathcal{S}} \mathbb{E}I(x)$$

Let's see how it works... iteration 4+1



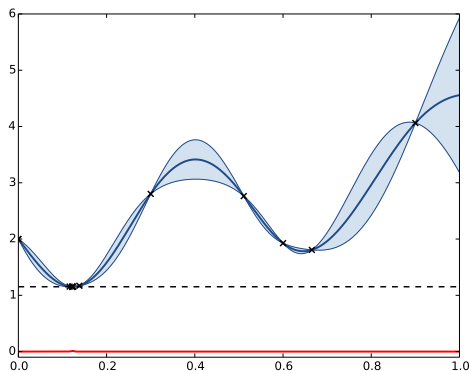
Expected Improvement

$$x^{t+1} = \arg \max_{x \in S} \mathbb{E}I(x) \dots \text{iteration } 4+2$$



Expected Improvement

$$x^{t+1} = \arg \max_{x \in S} \mathbb{E}I(x) \dots \text{iteration } 4+5$$



Expected Improvement: comments

There are other acquisition criteria:

- Upper Confidence Bound [Kushner, 1962, Srinivas et al., 2010],
 $= m(x) - \alpha c(x, x)$. Need to choose α .
- Knowledge gradient [Moćkus, 1972, Frazier and Powell, 2007],
 $\min_{x \in \mathcal{S}} m^t(x) - \mathbb{E}[\min_{x \in \mathcal{S}} m^{t+1}(x)]$. No analytical expression.
- Criteria based on the reduction in entropy of the optimum [Villemonteix et al., 2009, Hernández-Lobato et al., 2014]. No analytical expression, complex.
- ...
- $\mathbb{E}I$ is a myopic criterion compared to the above criteria. It is still a good entry criterion thanks to its simplicity and ease of interpretation.

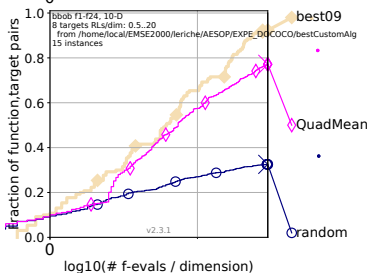
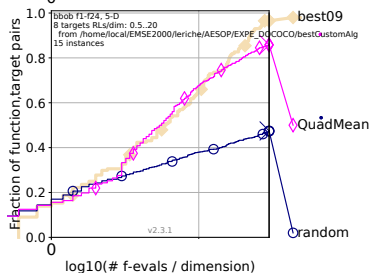
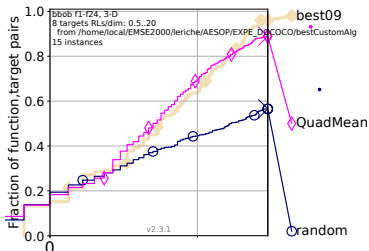
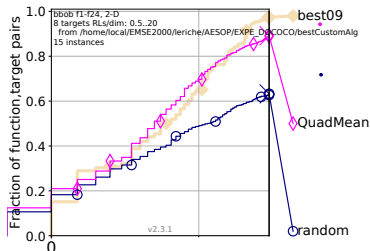
Bayesian optimization and dimension

Bayesian optimizers are competitive at low number of function evaluations but they lose this advantage with dimension.

- ⇐ Loss of GP accuracy?
- ⇐ EI sample too often at boundary?

Next slide: COCO tests [Hansen et al., 2016] from [Le Riche and Picheny, 2021]

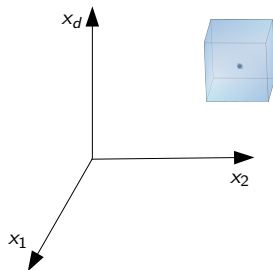
Effect of dimension on Bayesian Optimization



QuadMean = BO with quadratic mean, runs length = 30 d.

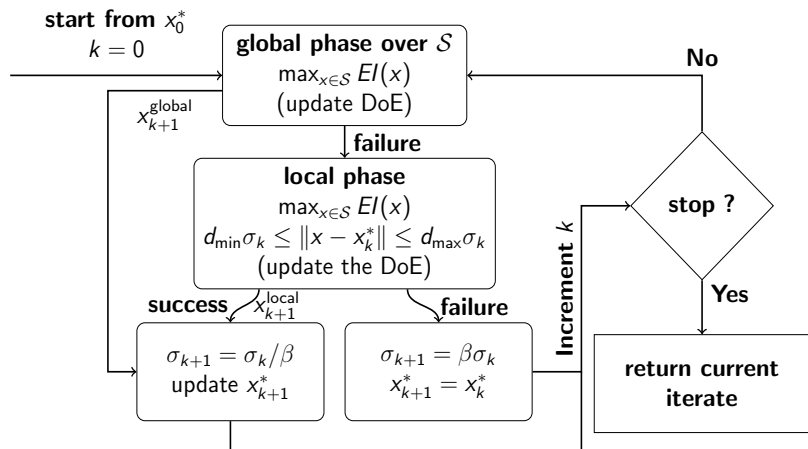
BO and trust regions

Principle: counteract the effect of increasing dimension (volume) by restricting the search to a smaller (controlled) trust region.



- TRIKE, Trust-Region Implementation in Kriging-based optimization with Expected Improvement, [Regis, 2016].
- TURBO, a TrUst-Region BO solver, [Eriksson et al., 2019].
- TREGO, a Trust-Region framework for EGO, [Diouane et al., 2023] : mix searches inside (local) and outside (global) the trust region.

TREGO algorithm



Parameters : $\sigma_0, \beta < 1$

Sufficient decrease condition for success of the local phase,

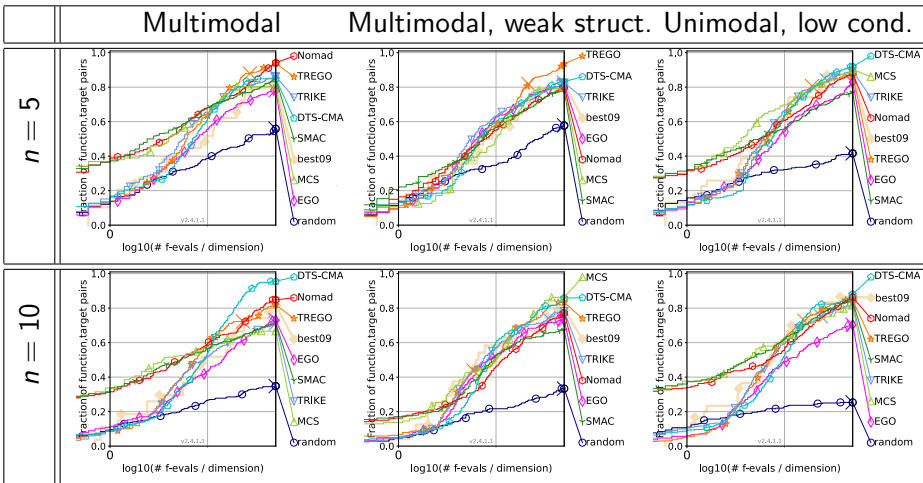
$$f(x_{k+1}^{\text{local}}) \leq f(x_k^*) - 10^{-4} \sigma_k^2$$

TREGO properties

From [Diouane et al., 2023],

- TREGO iterates converge to a local minimum : by assuming f is bounded below, Lipschitz continuous near the point of convergence, and by considering a subsequence of the local iterates. No assumption on GP or x_0^* .
- Empirical COCO tests:
 - more local than global steps (4 to 1) is beneficial
 - TREGO is robust to the values of σ_0 and β
 - A local GP was thought an asset for non stationary functions. But it is a drawback on badly conditioned functions. Not kept.

TREGO performance



Trust regions solve BO's oversampling of the boundaries in high-dim. while helping on unimodal functions (not the natural target for BO).

RELATIONS BETWEEN BAYESIAN AND EVOLUTIONARY OPTIMIZATION

The probabilistic representations of f

BO and EO represent f differently

BO

Global model of the function
(in \mathbb{R})

$$p(f | x) = \mathcal{N}(m(x), c(x, x))$$

EO

Model the distribution of good
points (in \mathcal{S})

$$p(x | f(x) \leq \text{threshold})$$

$$\approx \frac{1}{\text{card}(\mathbb{X})} \sum_{x_i \in \mathbb{X}} \delta_{x_i}(x)$$

Both views are tied by Bayes rule (see later).

Fundamental degrees of freedom

BO

Choosing and learning the kernel = choosing the feature space of the $\phi(x)$'s: by Mercer's theorem, $k(x, x') = \sum_{i=1}^D \lambda_i \phi_i(x) \phi_i(x')$ (N possibly infinite).
Quality metrics: regression metrics (Q2, LOO normality test).

Acquisition criterion: controls the exploration-intensification tradeoff.

Quality metrics: optimization metrics (convergence, performance profiles, ...).

EO

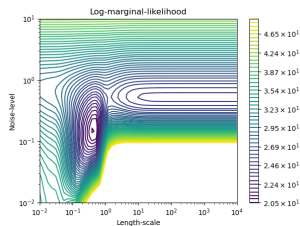
Choosing and learning the representation and variation operators.
Quality metrics: ergodicity (completeness), unbiasedness, locality (genotype-phenotype local correlation).

Both views will be linked later (perspectives).

EO for BO

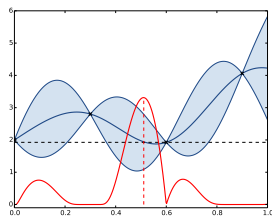
BO has 2 internal optimizations that are not costly and multimodal: they benefit from EO (CMA-ES is a typical choice for maximizing the $\mathbb{E}I$).

$$\max_{\theta} \text{likelihood}(\theta; \mathbb{X}, \mathbb{F})$$



(from `scikit-learn.org`)

$$\max_x \mathbb{E}I(x; GP)$$



BO for EO: adding a GP – I

Principle : EO algorithms need to order new x 's for selection.
Replace calls to f by calls to GPs.

Like surrogate-based EO, but with prediction uncertainties. Ranking from

- UCB [Büche et al., 2005], $m(x) - \alpha c(x, x)$,
- prob. of improvement [Ulmer et al., 2003],
 $\text{cdf}_{\mathcal{N}} \left((\min \mathbb{F} - m(x)) / \sqrt{c(x, x)} \right)$,
- partial (Pareto) ranks of (function prediction,- function uncertainty) [Volz et al., 2017], $(m(x), -\sqrt{c(x, x)})$,
- ... cf. [Bajer et al., 2019] for other criteria.

BO for EO: adding a GP – II

Algorithms are variants around the following sequence:

- $\mathbb{X}' = \text{Variation}(\mathbb{X})$
- Rank points in \mathbb{X}' using the GP and one of the ranking criteria (UCB, PI, EI, m , ...)
- Check the quality of the GP:
 - number of generations without GP update
 - OR compare ranking from true f and from GP on a small portion of \mathbb{X}'
- GP update from a reduced number of points chosen in $\mathbb{X}' \cup \mathbb{X}$ where true f is calculated

a sequence which is decomposed, repeated, varied ... inside EO implementations.

The aforementioned interfaces between BO and EO are external.

PERSPECTIVES:

A DEEPER LINK BETWEEN EO AND BO

through 2 examples.

Crossover as a barycenter in feature space I

Summary

- Feature learning (from data, through the kernels) can be seen as a search for an appropriate representation where linear reasoning works.
- Interpretate crossover as a barycenter in feature space.

A kernel can always be seen as an inner product in feature space[†], $k(x, x') = \langle \phi(x), \phi(x') \rangle$, where the feature is a map $\phi : \mathcal{S} \rightarrow \mathbb{R}^D$ and D can be infinite.

[†]For example it can be the eigendecomposition of Mercer's theorem seen earlier, $k(x, x') = \sum_{i=1}^{\infty} \sqrt{\lambda_i} \phi_i(x) \sqrt{\lambda_i} \phi_i(x')$ with the usual L2 scalar product, or the canonical decomposition in the RKHS, $k(x, x') = \langle k(x, \cdot), k(x', \cdot) \rangle_{\mathcal{H}}$.

Crossover as a barycenter in feature space II

Definition (Crossover as barycenter of features)

Let $x_C \in \mathcal{S}$ be the result of the crossover of $x_A \in \mathcal{S}$ with $x_B \in \mathcal{S}$. The features of x_C are the α -barycenter of the features of x_A and x_B , $0 \leq \alpha \leq 1$:

$$x_C := \arg \min_{x \in \mathcal{S}} \alpha \|\phi(x) - \phi(x_A)\|^2 + (1 - \alpha) \|\phi(x) - \phi(x_B)\|^2$$

Crossover as a barycenter in feature space III

Calculation:

$$\begin{aligned}\|\phi(x) - \phi(x_A)\|^2 &= \langle \phi(x) - \phi(x_A), \phi(x) - \phi(x_A) \rangle \\ &= \langle \phi(x), \phi(x) \rangle + \langle \phi(x_A), \phi(x_A) \rangle - 2\langle \phi(x), \phi(x_A) \rangle \\ &= k(x, x) + k(x_A, x_A) - 2k(x, x_A)\end{aligned}$$

and idem with $\|\phi(x) - \phi(x_B)\|^2$.

Assume stationarity, $k(x, x) = k(x_A, x_A) = k(x_B, x_B) = \sigma^2$, then

$$x_C = \arg \min_{x \in \mathcal{S}} 2\sigma^2 - 2\alpha k(x, x_A) - 2(1 - \alpha)k(x, x_B)$$

or equivalently

$$x_C = \arg \max_{x \in \mathcal{S}} \alpha k(x, x_A) + (1 - \alpha)k(x, x_B)$$

i.e., maximize weighted similarities with x_A and x_B as seen by the (learned) kernels.

For a crossover, choose $\alpha \sim \mathcal{U}[0, 1]$.

A Bayesian variation operator I

Link the probabilistic representations of BO and EO (cf. earlier slide) to define a variation (crossover and mutation) operator.

- T , a threshold (e.g., $T = \min \mathbb{F}$)
- $F(x) := Y(x) \mid f(\mathbb{X}) = \mathbb{F} \sim \mathcal{N}(m(x), c(x, x))$, the GP
- $F(x) < T$, a logical random variable. It can be written $F(X) < T \mid X = x$, X random variable in \mathcal{S}
- Bayes:
$$p(F(X) < T \mid X = x) \cdot p(X = x) = p(X = x \mid F(X) < T) \cdot p(F(X) < T)$$
- GP : $p(F(X) < T \mid X = x) = \text{cdf}_{\mathcal{N}}\left(\frac{(T - m(x))}{\sqrt{c(x, x)}}\right)$
- $p(X = x) = \mathcal{U}[\mathcal{S}]$. Other more intensifying choices possible, at the risk of premature convergence.

A Bayesian variation operator II

$p(X = x | F(X) < T)$ is the model for the variation operator of EO. It is sampled through

$$p(X = x | F(X) < T) \propto \text{cdf}_{\mathcal{N}}\left(\frac{(T - m(x))}{\sqrt{c(x, x)}}\right) p(X = x)$$

by a Metropolis-Hastings (or any appropriate MCMC) scheme.

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