

Bayesian Optimization : A New Sample Efficient Workflow for Reservoir Optimization under Uncertainty

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Abstract

Field development in E&P requires several level of decision making, where among the available options, the one with highest objective is decided. The decisions to make could be well configuration, well types , and well controls. The search to find the optimum decision, which called here optimization, a computationally demanding nonlinear problem. More challenging, in the case of optimization under uncertainty, such problems requires large number of flow simulation given different geological realizations. To alleviate this problem and to provide the feasible framework dealing with computational complexity in reservoir optimization, a new workflow is presented. The “Bayesian Optimization” in this work efficiently sample from the flow simulation based on the probabilistic modeling of the objective function, assisting to perform optimization process with as few as possible flow simulation, without losing the optimum solution. The paper begins with introduction to gaussian process and building a probabilistic view of the objective function as “Surrogate” and then provides the foundation for acquisition function, which decide to the next query point from the expensive function, here flow simulation. Further, a 1-D problem will be optimized through the Bayesian optimization which helps to better understanding of the workflow. Then, the workflow has been applied to 3D, “Egg model” to perform optimization in the realistic field scenario, considering geological uncertainty. Here, the comparison of the workflow with other two well-known algorithm in reservoir optimization literature, namely Genetic Algorithm (GA) and Particle Swarm Intelligence (PSO) has been performed. The results of the comparison show that the Bayesian Optimization workflow presented here could reach the same global optimum point achieved with GA and PSO, yet reduce computational complexity of the optimization 5X, which could be significant, in the case of the 3D flow simulation, potentially taking hours of running.

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1. Introduction:

While developing a field, prediction of reservoir response to change in the variable is an important factor to have an efficient reservoir management. The field of reservoir engineering is rich in the development and application of full-physics numerical simulations for forward modeling. However, the computational power needed to run such numerical simulators most of the time is huge. Especially, the framework of the Robust Optimization where uncertainty are considered through multiple of geological realization (thousand or multi-thousand), the practical applicability of such forward modeling is considerably limited. To address this challenge, the proxy-modeling for reservoir management has emerged to reduce the cost of forward simulation. The root of this field goes back to the work of the (Bruce 1943) where the analogy between flow of electricity in the electrical device and the response of the reservoir was constructed. (Albertoni and Lake 2003) Proposed the a Multivariate Linear Regression (MLR) to estimate the production from the well, where it claimed that total production at each is linear combination of injection rates with diffusivity filter. Building on the work of (Albertoni and Lake 2003), (Yousef 2006) proposed a new procedure to quantify communication between wells, where for each injector/producer pair, two coefficients, capacitance to quantify connectivity and time constant to quantify the degree of fluid storage were considered. (Sayarpour 2008) used superposition in time to analytically solve the numerical integration in CRM. (Zhao et al. 2015) articulated that CRM models are not applicable when production rates change significantly, mainly due to those models neglect to include interaction of production-production and injector-injector pair well. Formulating the model in the framework titled INSIM (interwell numerical simulation model), it was used as efficient replacement to reservoir simulator when history-matching phase rate data or water-cut data obtained during water flooding operations. (White and Royer, n.d.)

Separately, in sake of utilization of recent advancement in the world of Information technology, couple of research has been done on the development of “Surrogate Reservoir Models” - (Mohaghegh and Guruswamy 2006) proposed the workflow for SRM model where Fuzzy Pattern Recognition (FPR) technology was dimensionality reduction, in both static and dynamic parameters of the reservoir. Key Performance Indicator (KPI) was used to select the most important variable.- (Sampaio 2009) tried on use of feed-forward neural networks as nonlinear proxies of reservoir response. A few set of rock and fluid properties were considered as input (porosity, permeability in “x” direction, permeability in “y” direction, and rock compressibility) where in developing the neural network model, only one hidden layer was used. flow proxy modeling to replace full simulation in history matching, and built the proxy with a single hidden layer artificial neural network (ANN). To predict the oil production from the SAGD process, (Fedutenko et al. 2014) employed the RBF Neural Network to predict the cumulative oil production over the time horizon of production (10 years).

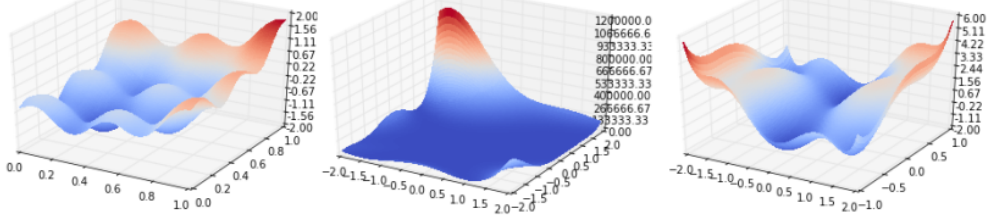
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2. Problem Statement

Consider a “well behaved” function $f : \chi \rightarrow \mathbb{R}$ where $\chi \subseteq \mathbb{R}^D$ is a bounded domain.

$$x_M = \operatorname{argmin} f(x) \quad (1)$$

reference equation 1



f is explicitly unknown and multimodal. Evaluations of f are expensive. Expensive functions, who doesn't have one?

In reservoir optimization, we can have different objectives: Recovery Factor, Net Present Value (NPV), ...

$$u_M = \operatorname{argmax}_{u \in \text{constraints}} \text{Objective Func}(u) \quad (2)$$

u is a decision variable to make: could be injection rate, well locations, order of drilling wells.

$$\text{Objective Func}(u) = J(u, G) = \sum_{k=1}^{n_T} \frac{q_o^k(u, G)P_o - q_w^k(u, G)P_{wp} - I^k(u, G)P_{wi}}{(1 + b)^{t_k/D}} \quad (3)$$

$$\text{Objective Func}(u) = \bar{J}(u) = \frac{\sum_{i=1}^{n_e} J_r(u, G_i)}{n_e} \quad (4)$$

Example case:.

- Lets' say you are optimizng the well control problem with dimension, $d = 10$
- Number of 3D geological realizations is $n_e = 100$ and running each possible u^* in commercial reservoir simulator, at each realization takes ~ 1 hour. Then time it takes to calculate $\bar{J}(u^*)$ is ~ 100 hours.
- With 6 month CPU running time budget, the total available budget is to run is ~ 50 .

- Availability of only 50 $\bar{J}(u^*)$ evaluation is a small budget, considering the 10 dimensional optimization problem.

Problem $\bar{J}(u)$ is expensive function, meaning the # times we can evaluate it severely limited. Solution Bayesian Optimization (BayesOpt) propose a new workflow to conduct optimization at the small $\bar{J}(u)$ budget, without affecting the optimum solution.

3. Bayesian Optimization Workflow

Overall View

Workflow to perform global optimization of multimodal black-box functions:

- Step 1. Choose some initial design points and build a probabilistic model over the space of possible objective f , this probabilistic model serves as prior.
- Step 2. Combine prior and the likelihood to get a posterior of probabilistic model over the objective given some observations.
- Step 3. Use the posterior to decide where to take the next evaluation \mathbf{x}^* according to some policy for decision making.
- Step 4. Evaluate the f at \mathbf{x}^* and augment it to the initial data, in step 1.

Iterate between 2 and 4 until the evaluation budget is over. ## 3.1 Gaussian Process

Step 1. Probabilistic Model as Prior

Gaussian Process (GP). reference to the book (**murphy2022?**)

Key Assumption in (GP) is that: the function values at a set of $M > 0$ inputs, $\mathbf{f} = [\mathbf{f}(\mathbf{x}_1), \dots, \mathbf{f}(\mathbf{x}_M)]$, is jointly Gaussian, with mean and Covariance

$$(\mu = m(x_1), \dots, m(x_M)) \sum_{i,j} = \kappa(x_i, x_j) \quad (5)$$

and κ is a positive definite (Mercer) kernel. Suppose we have a initial design points, set $\mathcal{D} = (x_n, y_n) : n = 1 : N$, where $y_n = f(x_n)$ is the noise-free observation of the function evaluated at x .

Now we consider the case of predicting the outputs for new inputs that may not be in \mathcal{D} .

$$\mathbf{f}_* = [\mathbf{f}(\mathbf{x}_1), \dots, \mathbf{f}(\mathbf{x}_{N_*})] \quad (6)$$

By definition of the GP, the joint distribution $p(\mathbf{f}_X, \mathbf{f} | \mathbf{X}, \mathbf{X}_*)$ has the following form:

$$\begin{bmatrix} \mathbf{f}_X \\ \mathbf{f}_* \end{bmatrix} \sim \mathcal{N} \left(\begin{bmatrix} \mu_X \\ \mu_* \end{bmatrix}, \begin{bmatrix} \mathbf{K}_{X,X} & \mathbf{K}_{X,*} \\ \mathbf{K}_{X,*}^\top & \mathbf{K}_{*,*} \end{bmatrix} \right) \quad (7)$$

$$\begin{aligned} \mu_X &= [m(x_1), \dots, m(x_N)] \\ \mu_* &= [m(x_1^*), \dots, m(x_N^*)] \end{aligned} \quad (8)$$

$$\begin{aligned} K_{X,X} &= \kappa(X, X; \theta), & size(N \times N) \\ K_{X,*} &= \kappa(X, X_*; \theta), & size(N \times N_*) \\ K_{*,*} &= \kappa(X_*, X_*; \theta), & size(N_* \times N_*) \end{aligned} \quad (9)$$

Covariance Kernels	assumeing $h = x - x' $
Gaussain	$\kappa(x, x') = \sigma_f^2 \exp(-\frac{h^2}{2\ell^2})$
Matern $\mu = \frac{5}{2}$	$\kappa(x, x') = \sigma_f^2 (1 + \frac{\sqrt{5} h }{\ell} \frac{5h^2}{3\ell^2}) \exp(-\frac{\sqrt{5} h }{\ell})$
Matern $\mu = \frac{3}{2}$	$\kappa(x, x') = \sigma_f^2 (1 + \frac{\sqrt{3} h }{\ell}) \exp(-\frac{\sqrt{3} h }{\ell})$
Exponetial	$\kappa(x, x') = \sigma_f^2 \exp(-\frac{ h }{\ell})$
Power-Exponetial	$\kappa(x, x') = \sigma_f^2 \exp(-(\frac{ h }{\ell})^p)$

$$\kappa(x, x'; ; \theta) = (1 + \frac{\sqrt{5}|h|}{\theta} + \frac{5h^2}{3\theta^2}) \exp(-\frac{\sqrt{5}|h|}{\theta}) \quad (10)$$

Covariance Kernel, Parameter estimation.

$$p(y|\mathbf{X}, \theta) = \int \mathbf{p}(y|\mathbf{f}, \mathbf{X}) \mathbf{p}(\mathbf{f}|\mathbf{X}, \theta) \quad (11)$$

$$\log p(y|\mathbf{X}, \theta) = \mathcal{L}(\zeta, \sigma_f^2) = -\frac{1}{2}(\mathbf{y} - \mu_{\mathbf{X}})^\top \mathbf{K}_{\mathbf{X}, \mathbf{X}}^{-1}(\mathbf{y} - \mu_{\mathbf{X}}) - \frac{1}{2} \log |\mathbf{K}_{\mathbf{X}, \mathbf{X}}| - \frac{n}{2} \log(2\pi) \quad (12)$$

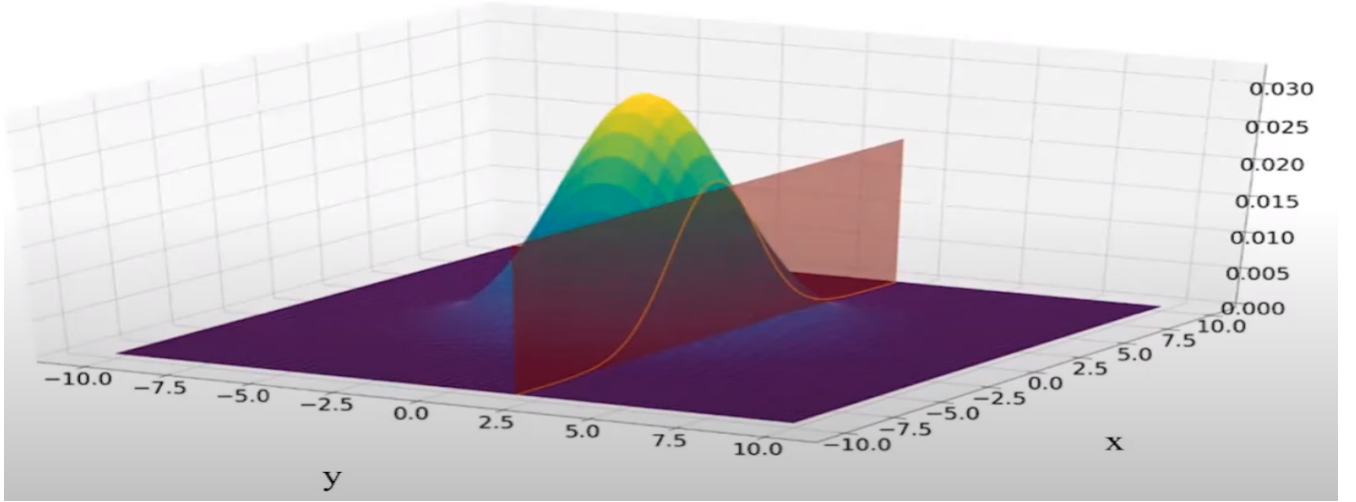
Where the dependence of the $\mathbf{K}_{\mathbf{X}, \mathbf{X}}$ on θ is implicit. The gradient-based optimizer is performed in order to:

$$[\zeta^*, \sigma_f^{2*}] = \operatorname{argmax} \mathcal{L}(\zeta, \sigma_f^2) \quad (13)$$

However, since the objective \mathcal{L} is not convex, local minima can be a problem, so we may need to use multiple restarts.

Step 2. Posterior of Probabilistic Model

Posterior of Gaussian Process, (conditioning on initial data).



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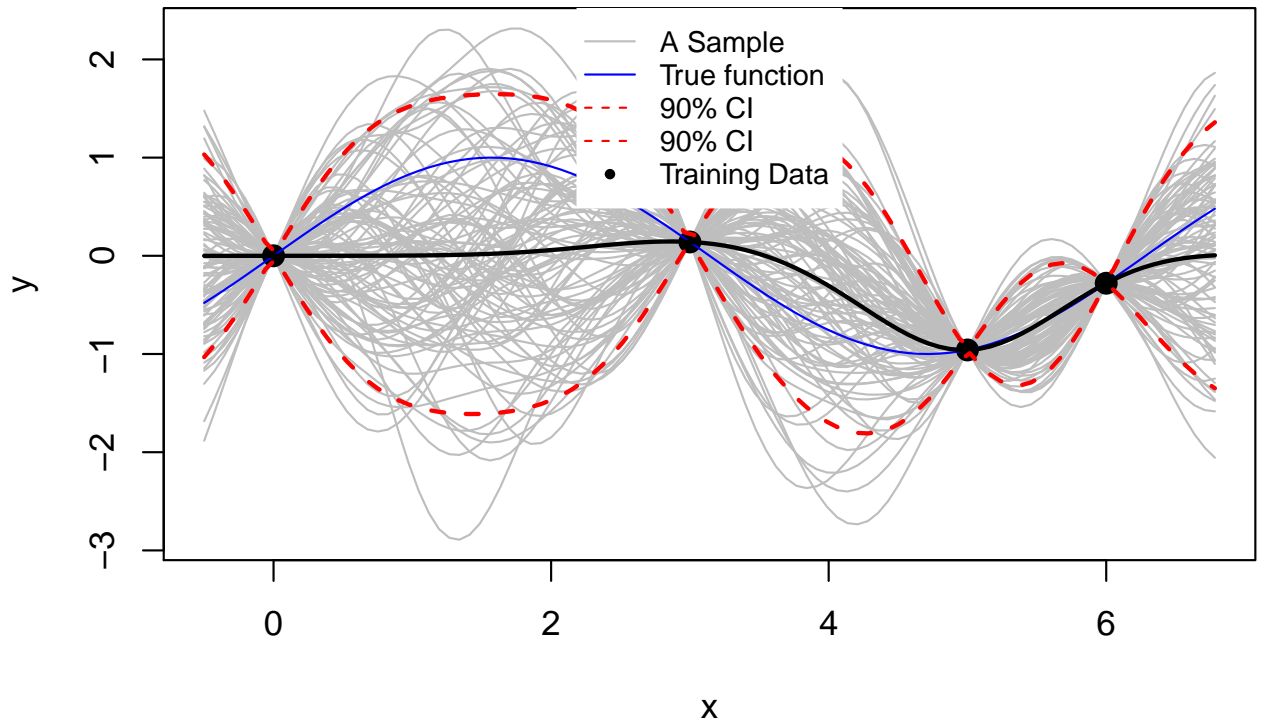
$$p(f_*|X_*, \mathcal{D}) = \mathcal{N}(f_*|\mu_*, \sum_*) \quad (14)$$

$$\begin{aligned}\mu_* &= m(\mathbf{X}_*) + \mathbf{K}_{\mathbf{X},*}^T \mathbf{K}_{\mathbf{X},\mathbf{X}}^{-1} (\mathbf{f}_{\mathbf{X}} - \mathbf{m}(\mathbf{X})) \\ \Sigma_* &= \mathbf{K}_{*,*} - \mathbf{K}_{\mathbf{X},*}^T \mathbf{K}_{\mathbf{X},\mathbf{X}}^{-1} \mathbf{K}_{\mathbf{X},*}\end{aligned}\tag{15}$$

Example of Step.1 and Step.2

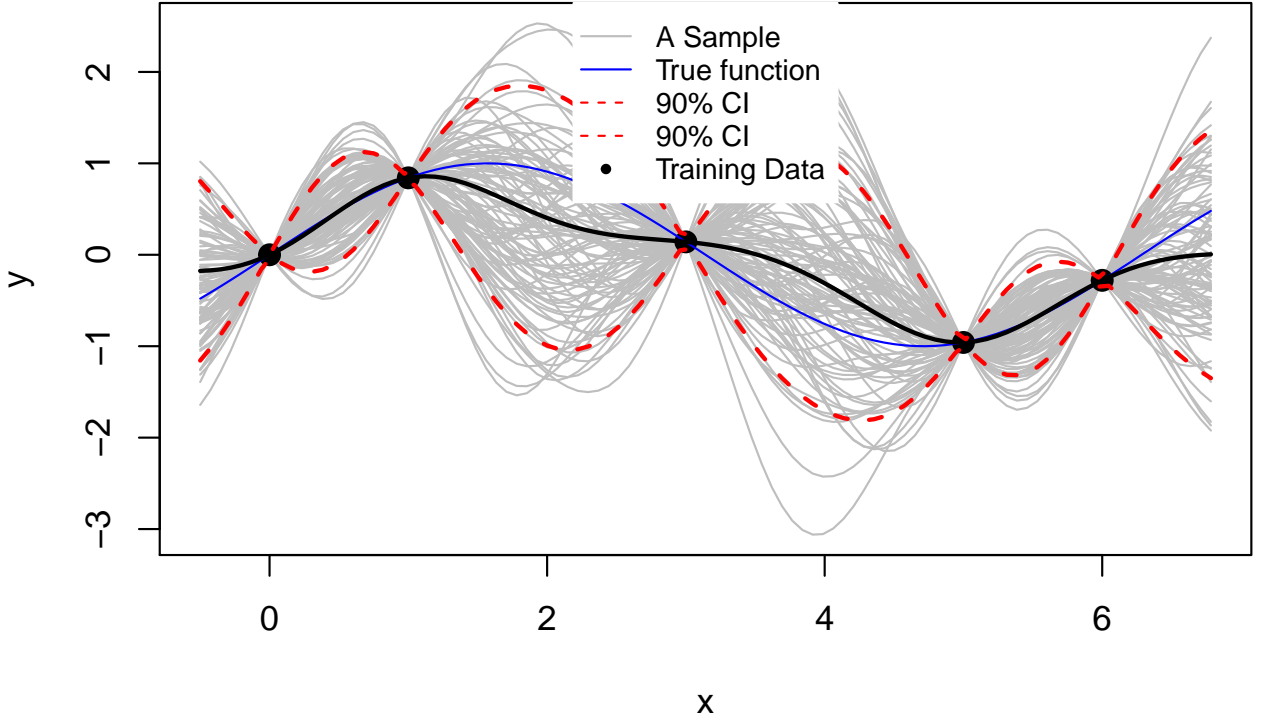
Assume $X = [0, 3, 5, 6]$ and $f_X = \sin(X)$, giving $\mathcal{D} = (X, f_X)$. What is $p(f_* | X_*, \mathcal{D})$

Gaussian Process Regression



Now we sample the point $X = 1$, and add to \mathcal{D}

Gaussian Process Regression



Step.3 Deciding on next \mathbf{x}^* based on Posterior

Posterior of the probabilistic model quantify the uncertainty over the space of the f . The question is what is the next \mathbf{x}^* to be sampled from the *expensive function*?

Define an utility function to collect new data points satisfying some optimality criterion: optimization as decision making.

There are a few of policies in the literature of Bayesopt, here the *Expected Improvement (EI)* policy will be used.

Expected Improvement as Policy for Decision Making. In Expected Improvement (EI) policy choose the next query point as the one which has the highest expected improvement over the space of the *expensive function*

$$utility(x; \theta, \mathcal{D}) = \alpha_{EI} = \int_y \max(0, y - f) p(y|x; \theta, \mathcal{D}) \quad (16)$$

$$utility(x; \theta, \mathcal{D}) = \alpha_{EI} = \int_y \max(0, y - f) p(y|x; \theta, \mathcal{D}) dy$$

However, we do not have access to the *expensive function*, f , therefore we replace the f with the best available solution found so far, y^+

$$utility(x; \theta, \mathcal{D}) = \alpha_{EI} = \int_y \max(0, y - y^\dagger) p(y|x; \theta, \mathcal{D}) dy \quad (17)$$

y^\dagger : The best solution found in the training dataset \mathcal{D}

The good news: The analytical form of the utility function is available for the gaussian process

$$\gamma(\mathbf{x}) = \frac{\mu(\mathbf{x}; \theta, \mathcal{D}) - y^\dagger}{\sigma(\mathbf{x}; \theta, \mathcal{D})} \quad (18)$$

$$utility(\mathbf{x}; \theta, \mathcal{D}) = \alpha_{EI}(x; \theta, \mathcal{D}) = (\mu(x; \theta, \mathcal{D}) - y^\dagger) \Phi(\gamma(x)) + \sigma(x; \theta, \mathcal{D}) \phi(\gamma(x)) \quad (19)$$

Where $\Phi(\cdot)$ and $\phi(\cdot)$ are CDF and PDF of standard Gaussian distribution.

It is too greedy in the context of the sequential decision making. Therefore, an explorative term is added as explorative" parameter ϵ .

$$\gamma(\mathbf{x}) = \frac{\mu(\mathbf{x}; \theta, \mathcal{D}) - y^\dagger - \epsilon}{\sigma(\mathbf{x}; \theta, \mathcal{D})} \quad (20)$$

$$\alpha_{EI}(x; \theta, \mathcal{D}) = (\mu(x; \theta, \mathcal{D}) - y^\dagger - \epsilon) \Phi(\gamma(x)) + \sigma(x; \theta, \mathcal{D}) \phi(\gamma(x)) \quad (21)$$

BO As a "mapping" between two problems. BO is an strategy to transform the problem

$$u_M = \underset{u \in \text{constraints}}{\operatorname{argmax}} \bar{J}(u) \quad (22)$$

unsolvabile!

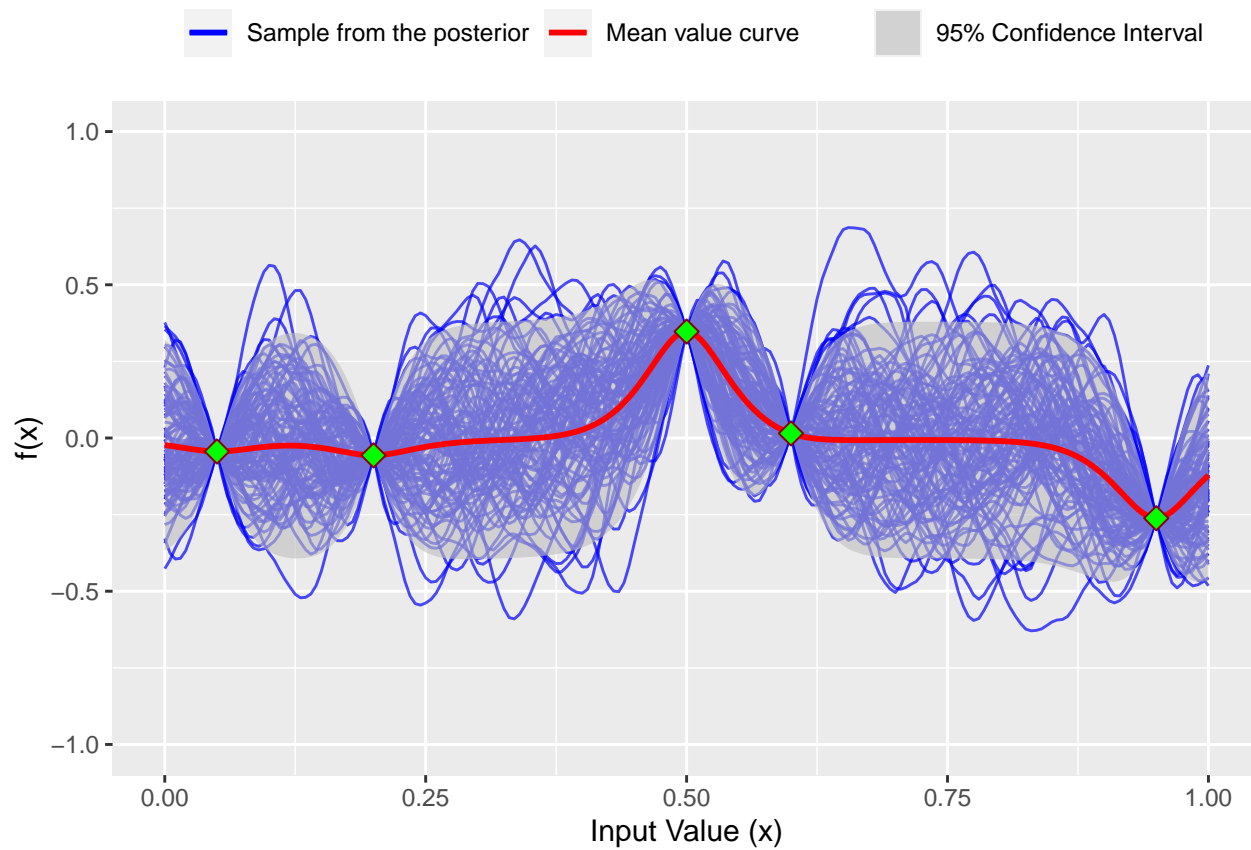
$$u^{next} = \underset{u \in \text{constraints}}{\operatorname{argmax}} \alpha_{EI}(u; \mathcal{D}_n, \theta^*) \quad (23)$$

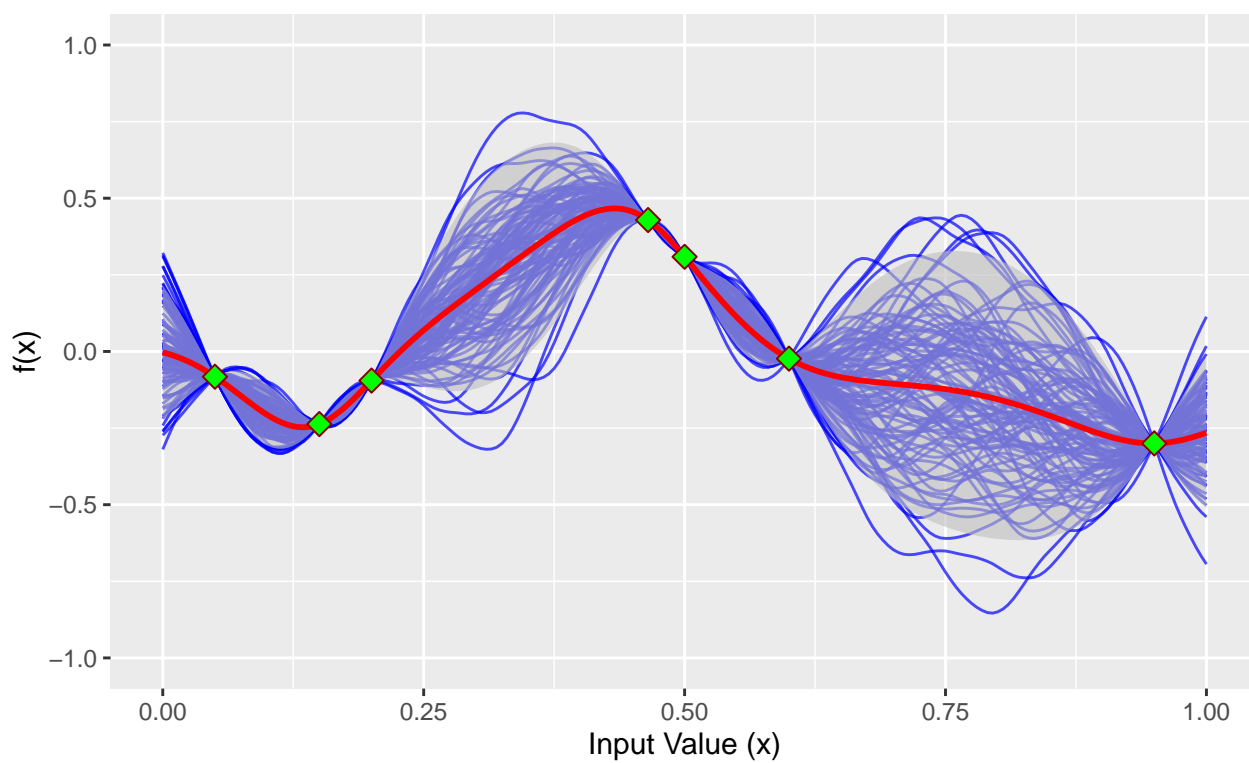
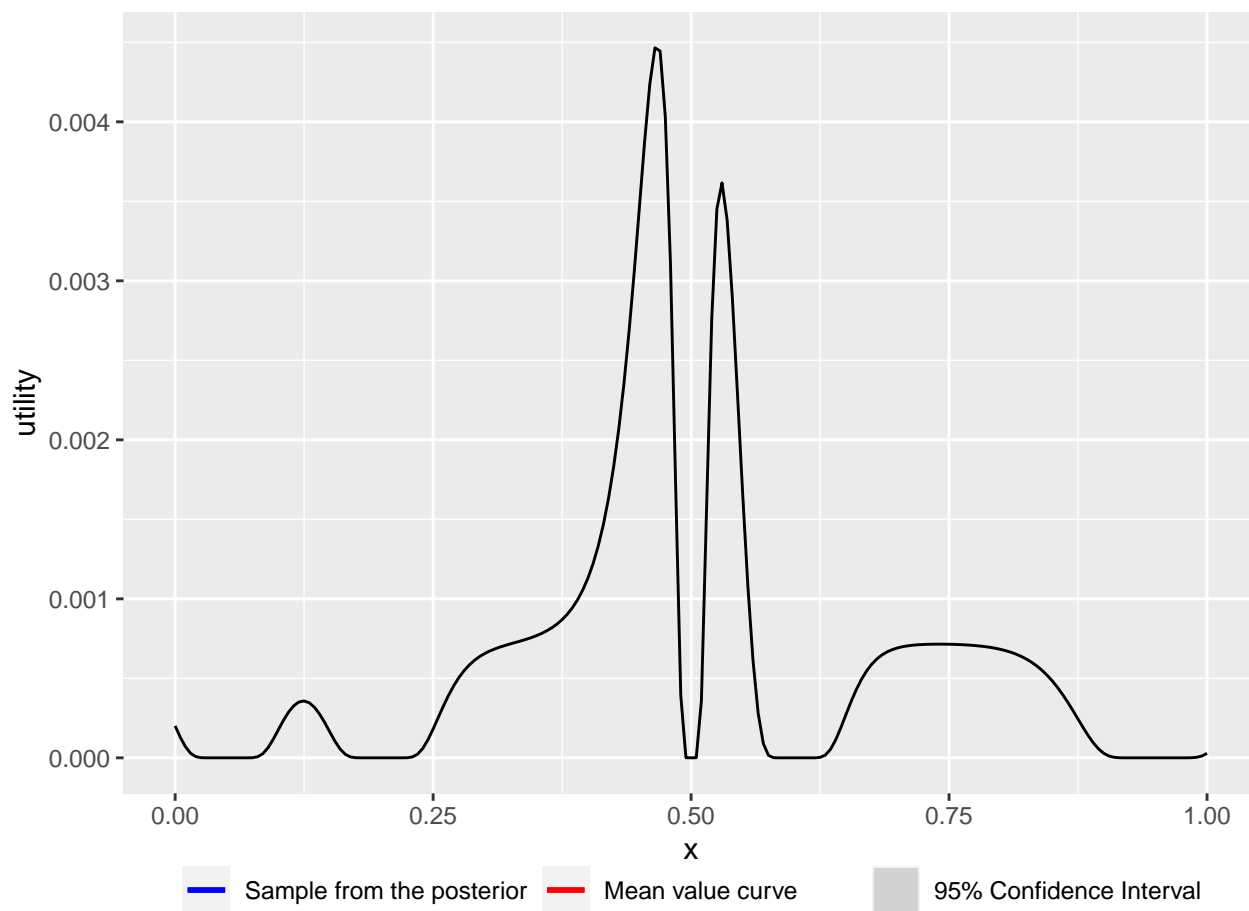
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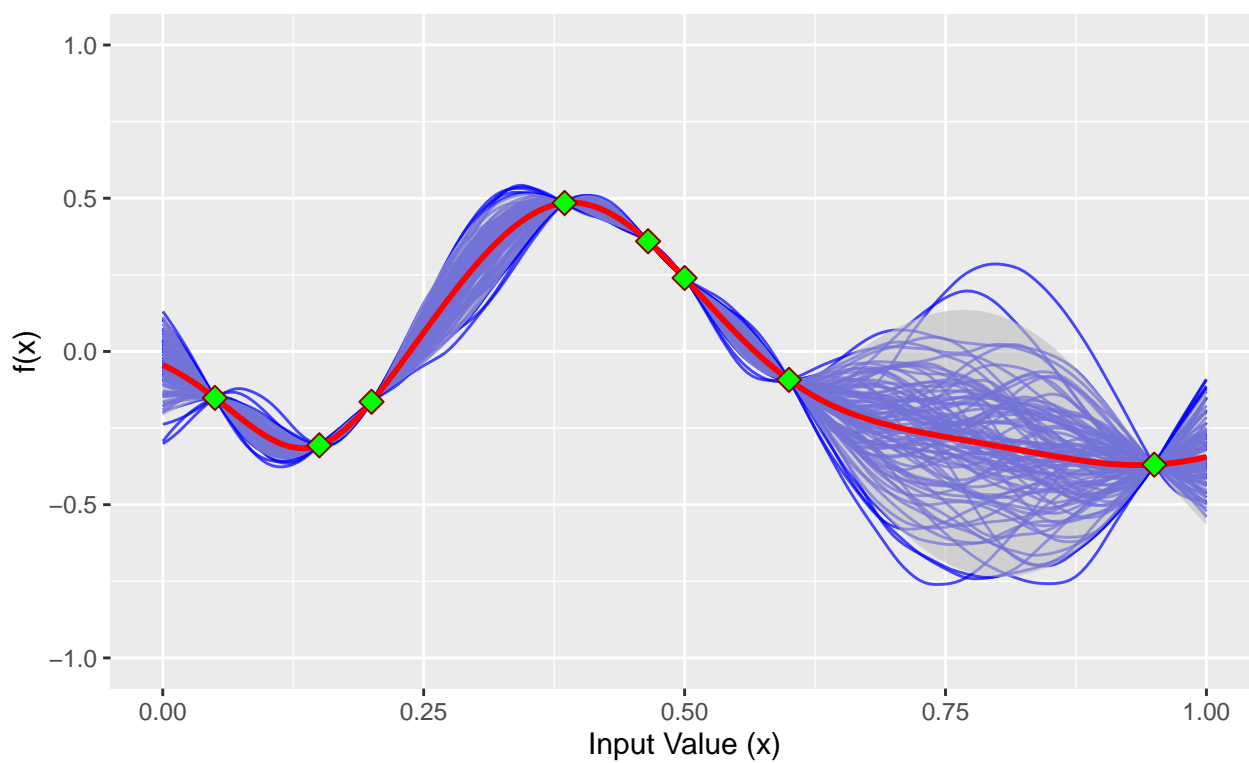
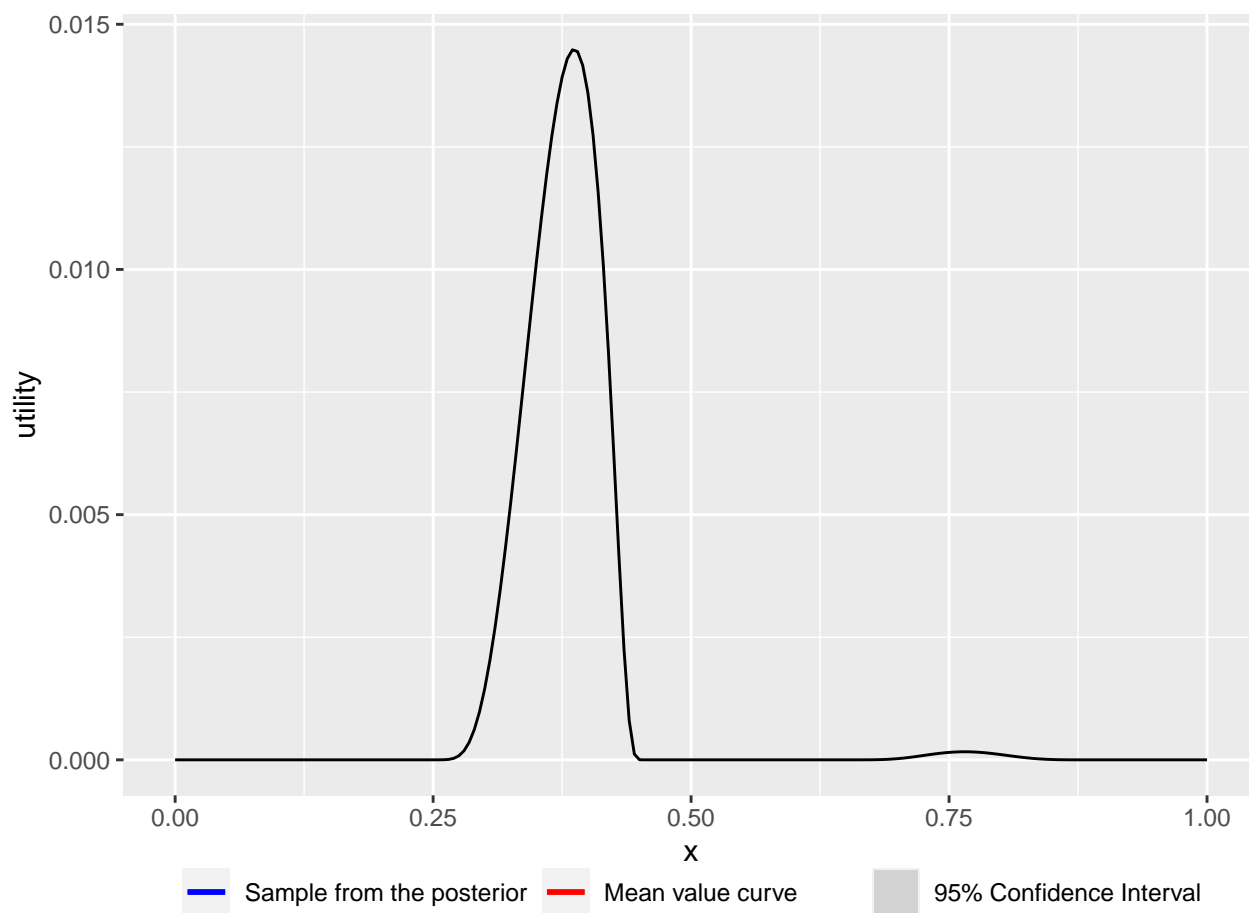
- $\alpha_{EI}(u)$ is inexpensive to evaluate.
- The analytical expression for gradient of $\alpha_{EI}(u)$ is available.
- Still need to find u^{next} , the multi-start BFGS is used for finding u^{next} .

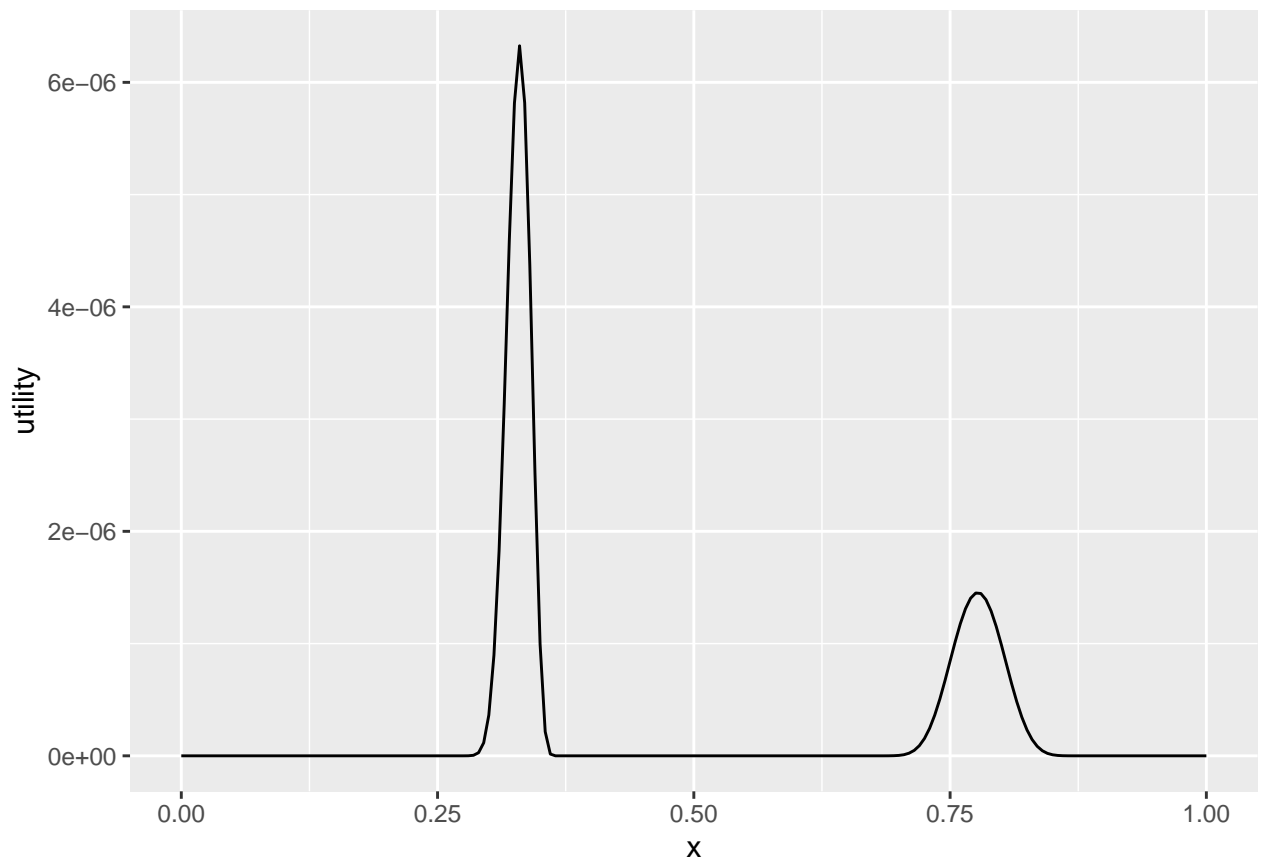
4. Example Cases

1-D toy Problem





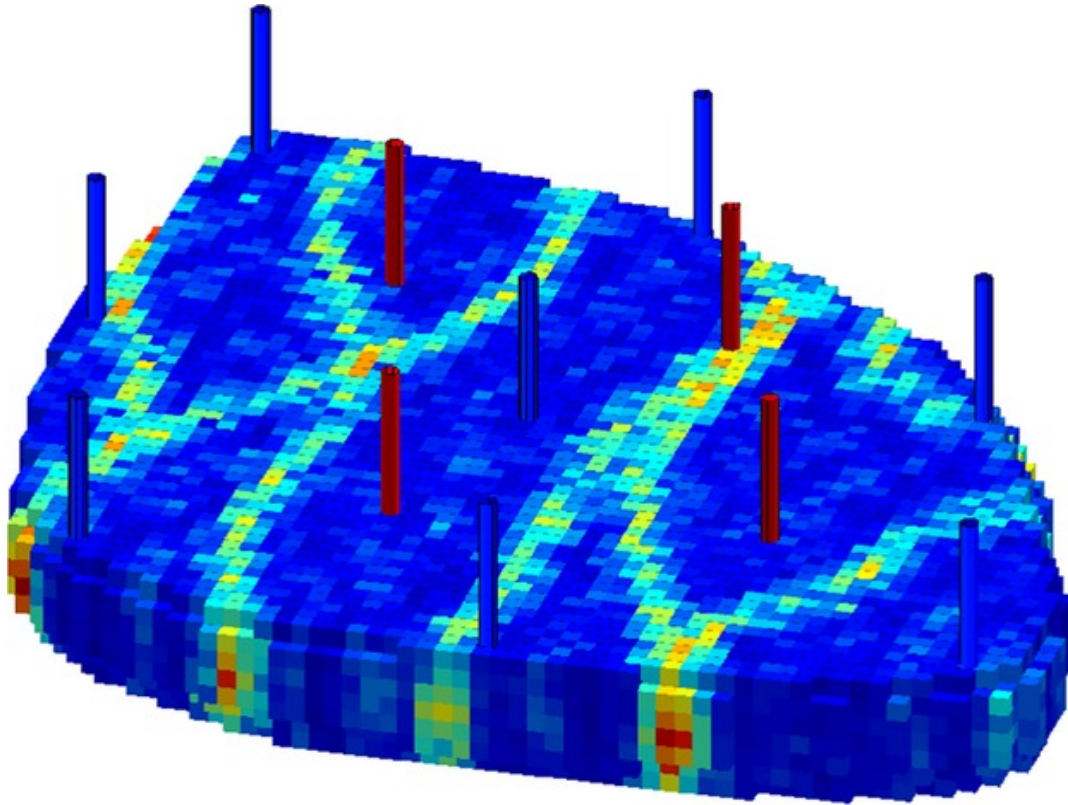




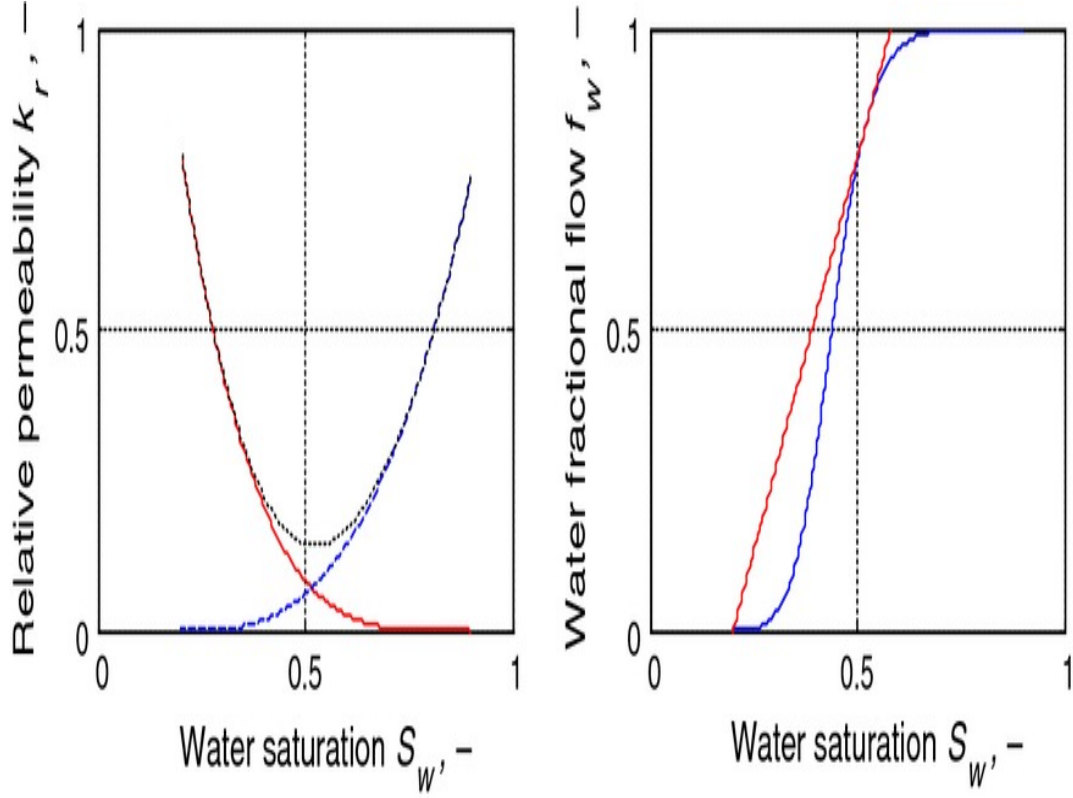
Field Scale

Egg Model (8 Injection wells, 4 Production wells)

- Egg Model



- Rel Perm Curves



BayesOpt Applied to Reservoir Case:

Optimization Objective

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$$J(u) = NPV(u) = \sum_{k=1}^{n_T} \frac{q_o^k(u)P_o - q_w^k(u)P_{wp} - I^k(u)P_{wi}}{(1+b)^{t_k/D}}$$

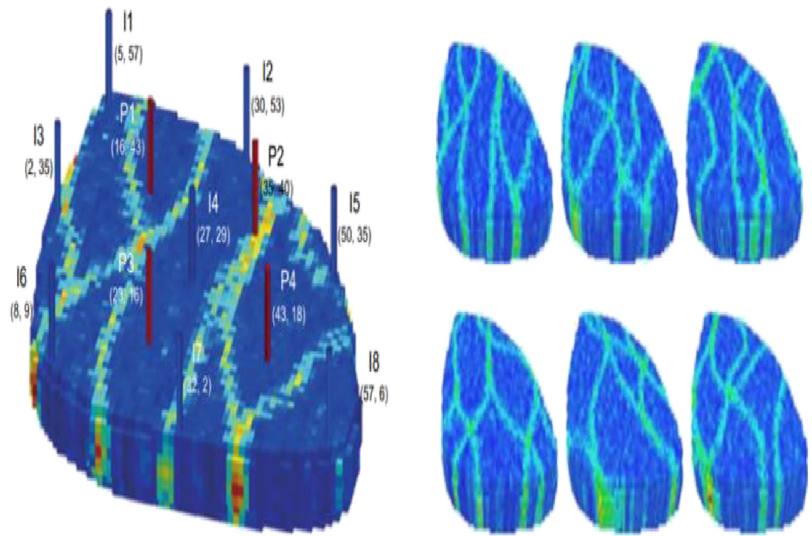
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$$\bar{J}(u) = \frac{\sum_{r=1}^{n_e} J_r(u)}{n_e}$$

• u is Injection rate for the each injection well,

• $u = [u_{inj1}, u_{inj2}, u_{inj3}, u_{inj4}, u_{inj5}, u_{inj6}, u_{inj7}, u_{inj8}]^T$

Item	Pric	Items	Value
P_o	315	b	8%
P_wp	47.5	D	365
P_wi	12.5	n_e	10



5. Coparsion with other optimization algorithm

6. Concluding Remarks

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Here are two sample references: (Dirac, 1953; Feynman and Vernon Jr., 1963).

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