Myopic Posterior Sampling for Adaptive Goal Oriented Design of Experiments

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Abstract

Bayesian methods for adaptive decision-making, such as Bayesian optimisation, active learning, and active search have seen great success in relevant applications. However, real world data collection tasks are more broad and complex, as we may need to achieve a combination of the above goals and/or application specific goals. In such scenarios, specialised methods have limited applicability. In this work, we design a new myopic strategy for a wide class of adaptive design of experiment (DOE) problems, where we wish to collect data in order to fulfil a given goal. Our approach, Myopic Posterior Sampling (MPS), which is inspired by the classical posterior sampling algorithm for multi-armed bandits, enables us to address a broad suite of DOE tasks where a practitioner may incorporate domain expertise about the system and specify her desired goal via a reward function. Empirically, this general-purpose strategy is competitive with more specialised methods in a wide array of synthetic and real world DOE tasks. More importantly, it enables addressing complex DOE goals where no existing method seems applicable. On the theoretical side, we leverage ideas from adaptive submodularity and reinforcement learning to derive conditions under which MPS achieves sublinear regret against natural benchmark policies.

1. Introduction

Many problems in adaptive decision-making under uncertainty fall into the design of experiments (DOE) framework, where one wishes to design a sequence of experiments and collect data so as to achieve a desired goal. For example, in electrolyte design for batteries, a chemist would like to conduct experiments that measure battery conductivity in order to identify an electrolyte design that maximises conductivity.

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ity. On a different day, she would like to experiment with different designs to learn how the viscosity of the electrolyte changes with the design. These two tasks, black-box optimisation and active learning, fall under the umbrella of DOE and are pervasive in industrial and scientific applications.

While several methods exist for specific DOE tasks, real world problems are broad and complex, and specialised methods have limited applicability. Continuing with the electrolyte example, the chemist can typically measure both conductivity and viscosity with a single experiment. Since such experiments are expensive, it is wasteful to first perform a set of experiments to optimise conductivity and then a fresh set to learn viscosity. Rather, it is desirable to design a single set of experiments that simultaneously achieves both goals. A second example is metallurgy, where one wishes to identify phase transitions in an alloy by carefully selecting a sequence of X-ray diffraction experiments (Bunn et al., 2016). Here and elsewhere, the goal of the experimenter is application specific and cannot be simply shoehorned into standard DOE formulations such as black-box optimisation, active learning, etc. In addition, domain knowledge about the problem may need to be considered in selecting experiments, as it may significantly reduce the number of experiments needed to achieve the desired goal.

To address these desiderata, we develop a general and flexible framework for goal oriented DOE, where a practitioner may specify her desired goal via a reward function λ . λ can depend on the data collected during the DOE process and unknown system characteristics, and hence cannot be directly computed by a decision-maker. We then develop an adaptive myopic strategy for DOE, inspired by posterior (Thompson) sampling for multi-armed bandits (Thompson, 1933), which uses results from past experiments to plan future experiments and achieve the goal, i.e. maximise λ . Our approach has two key advantages. First, our Bayesian formulation allows one to straightforwardly specify domain expertise. Moreover, modern tools for probabilistic programming enable pratitioners to apply a Bayesian algorithm such as ours in a fairly straightforward manner. Second, our myopic strategy is simple and computationally attractive in comparison with policies that engage in long-term planning. Nevertheless, borrowing ideas from submodular optimisation and reinforcement learning, we derive natural

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conditions under which our myopic policy is competitive with the globally optimal one. Our contributions are:

- We propose a flexible framework for DOE that allows a practitioner to describe their system via a probabilistic model, and specify their goal via a reward function. We derive a general purpose algorithm, Myopic Posterior Sampling (MPS), for this setting.
- 2. Empirically, we demonstrate MPS performs favourably in a variety of synthetic and real world DOE tasks. Despite its generality, MPS is competitive with specialised methods designed for particular settings. More importantly, it enables DOE in non-standard application-specific settings. Our implementation and experiments are available at github.com/kirthevasank/mps.
- 3. In our theoretical analysis, we explore conditions under which MPS, which learns about the system over time, is competitive with myopic and globally optimal strategies that have full knowledge of the system.

Related Work: The term DOE has been used to refer to different settings in the literature. Classically, the focus has been on learning an unknown system, and as such, the objective has been framed as maximising some notion of information gathered about the system. We will refer to these tasks as L-DOE problems to differentiate it from our setting, which subsumes L-DOE. Classical L-DOE focuses on discrete settings (Chernoff, 1972; Robbins, 1952) or linear models (Allen-Zhu et al., 2017; Fedorov, 1972). Recently, there have been several Bayesian approaches for L-DOE that adopt probabilistic programming in more complex models (Ouyang et al., 2016; Rainforth, 2017). However, L-DOE approaches may not be efficient or appropriate for an arbitrary user-specified reward λ . Moreover, many of these approaches are non-adaptive, aiming to find an optimal batch of experiments beforehand without incorporating feedback from completed experiments. While some do explore adaptive approaches for L-DOE, they aim for globally optimal policies (e.g. Rainforth (2017)), which can be computationally prohibitive, except in the most simple cases.

We focus on posterior sampling (PS) (Thompson, 1933) as the bandit algorithm, since it has proven to be quite general and admits a clean Bayesian analysis (Russo and Van Roy, 2016a). PS has been studied in a number of bandit settings (Gopalan et al., 2014; Kandasamy et al., 2018; Kawale et al., 2015), and some episodic RL problems (Gopalan and Mannor, 2015; Osband and Van Roy, 2014; Osband et al., 2013), where the agent is allowed to restart. In contrast, here we study PS on a single long trajectory with no restarts.

Myopic/greedy policies, while computationally simple, are known to be near-optimal for sequential decision making problems with *adaptive submodularity* (Golovin and Krause, 2011), which generalises submodularity (Nemhauser et al.,

1978) and formalises a diminishing returns property. Adaptive submodularity has been used for several adaptive DOE setups (Chen and Krause, 2013; Chen et al., 2014; 2017; Golovin et al., 2010). However, in these work, the reward only depends on the data collected and can be directly computed by the decision-maker. In our setting, this translates to the agent knowing the system characteristics. As such, these results are complementary to ours: adaptive submodularity controls the approximation error (the difference between myopic- and globally-optimal strategies, both of which know the system), while we control the estimation error (how close our policy which needs to learn about the system is to the myopic optimal policy that knows the system). As we show in Theorem 3, with adaptive submodularity, MPS can also compete with the globally optimal policy. In a similar vein, Frazier et al. (2008); Wang and Powell (2018) use knowledge gradient approaches for information collection tasks which are framed as myopic adaptive submodular set maximisation problems; but as before, the system is known to the decision-maker. Prior results for learning in submodular environments are episodic and allow restarts (Gabillon et al., 2013; 2014), which is unnatural in the DOE setup. In addition to the above, several papers have developed Bayesian methods for specific DOE applications such as black-box optimisation (Frazier, 2018), active search (Jiang et al., 2018), level set estimation (Gotovos et al., 2013a) and more (Kandasamy et al., 2015; Osborne et al., 2012).

Our theoretical analysis leverages ideas from reinforcement learning (RL) since at each round the agent makes a decision (what experiment to perform) with the goal of maximising a long-term reward. In that light, one goal of our work is to understand when myopic "bandit-like" strategies perform well in RL environments with long-term temporal dependencies. There are two main differences with prior work (Jaksch et al., 2010; Kearns and Singh, 2002; Liu and Brunskill, 2018; Osband and Van Roy, 2014; Strehl et al., 2009). First, we make no explicit assumptions about the complexity of the state and action space, instead placing assumptions on the reward structure and optimal policy, which is a better fit for our applications. Crucially, in our setup, the true reward is never revealed to the agent, and instead it receives sideobservations that provide information about an underlying parameter governing the environment. Secondly, our focus is on understanding when myopic strategies have reasonable performance rather than on achieving global optimality.

2. Set up and Method

Let Θ denote a parameter space, \mathcal{X} an action space, and \mathcal{Y} an outcome space. We consider a Bayesian setting where a *true parameter* $\theta_{\star} \in \Theta$ is drawn from a prior distribution ρ_0 . A decision maker repeatedly chooses an action $X \in \mathcal{X}$, conducts an experiment at X, and observes the outcome $Y_X \in \mathcal{Y}$. We assume Y_X is drawn from a *like*-

lihood $\mathbb{P}(\cdot|X,\theta_{\star})$, with known distributional form. This process proceeds for n rounds, resulting in a data sequence $D_n = \{(X_j, Y_{X_j})\}_{j=1}^n$, which is an ordered multi-set of action-observation pairs. Unlike, classical formalisms for DOE, we study a setting where we intend to achieve a desired goal, specified via a reward function $\lambda:\Theta\times\mathcal{D}\to\mathbb{R}$, that we wish to maximise. Here, \mathcal{D} denotes the set of all possible data sequences. In particular, after n rounds, we focus on the following two criteria, depending on the application:

(a)
$$\Lambda(\theta_{\star}, D_n) = \sum_{t=1}^{n} \lambda(\theta_{\star}, D_t)$$
 (b) $\lambda(\theta_{\star}, D_n)$, (1)

Here, $D_t = \{(X_j, Y_{X_j})\}_{j=1}^t$ denotes the *prefix* of length tof the data sequence D_n collected by the decision maker. The former notion is the cumulative sum of all rewards, while the latter corresponds to the reward once all experiments are complete. Since λ depends on the unknown true parameter θ_{\star} , the decision maker cannot compute the reward during the data collection process, and instead must infer the reward from observations in order to maximise it. This is a key distinction from existing work on reinforcement learning and sequential optimisation, and one of the new challenges in our setting.

Example 1. A motivating example is Bayesian active learning (Chen et al., 2017). Here, actions X correspond to data points while Y_X is the label and $\mathbb{P}(y|x,\theta)$ specifies an assumed discriminative model. We may set $\lambda(\theta, D_n) =$ $-\|\beta(\theta) - \hat{\beta}(D_n)\|_2^2$ where β is a parameter of interest and $\hat{\beta}$ is a predetermined estimator (e.g. via maximum likelihood). The true reward $\lambda(\theta_{\star}, D_n)$ is not available to the decision maker since it requires knowing $\beta(\theta_{\star})$.

Notation: For each $t \in \mathbb{N}$, let $\mathcal{D}_t = \{(X_j, Y_{X_j})\}_{j=1}^t$: $X_j \in \mathcal{X}, Y_{X_j} \in \mathcal{Y}$ denote the set of all data sequences of length t, so that $\mathcal{D} = \bigcup_{t \in \mathbb{N}} \mathcal{D}_t$. Let $D \uplus D'$ denote the concatenation of two sequences. $D \prec D'$ and $D' \succ D$ both equivalently denote that D is a prefix of D'. Given a data sequence D_t , we use $D_{t'}$ for t' < t to denote the prefix of the first t' action-observation pairs.

A policy for experiment design chooses a sequence of actions $\{X_j\}_{j\in\mathbb{N}}$ based on past actions and observations. In particular, for a randomised policy $\pi = {\pi_j}_{j \in \mathbb{N}}$, at time t, an action is drawn from $\pi_t(D_{t-1}) = \mathbb{P}(X_t \in \cdot | D_{t-1})$. Two policies that will appear frequently in the sequel are $\pi_{\rm M}^{\star}$ and $\pi_{\rm G}^{\star}$, both of which operate with knowledge of θ_{\star} . $\pi_{\rm M}^{\star}$ is the myopic optimal policy, which, from every data sequence D_t chooses the action X maximising the expected reward at the next step: $\mathbb{E}[\lambda(\theta_{\star}, D_t \uplus \{(X, Y_X)\})|\theta_{\star}, D_t].$ On the other hand π_G^{\star} is the non-myopic, globally optimal adaptive policy, which in state D_t with n-t steps to go chooses the action to maximise the expected long-term reward: $\mathbb{E}[\lambda(\theta_{\star}, D_t \uplus \{(X, Y_X)\} \uplus D_{t+2:n}) \mid \pi_{\mathbf{G}}^{\star}, \theta_{\star}, D_t]. \pi_{\mathbf{G}}^{\star}$

may depend on the time horizon n while π_{M}^{\star} does not.

Design of Experiments via Posterior Sampling

We present a simple and intuitive myopic strategy that aims to maximise λ based on the posterior of the data collected so far. For this, first define the expected look-ahead reward $\lambda^+: \Theta \times \mathcal{D} \times \mathcal{X} \to [0,1]$, such that $\lambda^+(\theta,D,x)$ is the expected reward at the next time step if $\theta \in \Theta$ were the true parameter, D was the current data sequence collected, and we were to take action $x \in \mathcal{X}$. Precisely,

$$\lambda^{+}(\theta, D, x) = \mathbb{E}_{Y_x \sim \mathbb{P}(Y|x,\theta)} \Big[\lambda \Big(\theta, D \uplus \{(x, Y_x)\} \Big) \Big]. \quad (2)$$

The proposed policy, presented in Algorithm 1, is called MPS (Myopic Posterior Sampling) and is denoted $\pi_{\rm M}^{\rm PS}$. At time step t, it first samples a parameter value θ from the posterior for θ_{\star} conditioned on the data, i.e. $\theta \sim \mathbb{P}(\theta_{\star}|D_{t-1})$. Then, it chooses the action X_t that is expected to maximise the reward λ by pretending that θ was the true parameter. It performs the experiment at X_t , collects the observation Y_{X_t} , and proceeds to the next time step.

Algorithm 1 MPS $(\pi_{\rm M}^{\rm PS})$

Require: Prior ρ_0 for θ_{\star} , Conditional $\mathbb{P}(Y|X,\theta)$.

- 1: $D_0 \leftarrow \emptyset$.
- 2: **for** $t = 1, 2, \dots$ **do**
- Sample $\theta \sim \rho_{t-1} \equiv \mathbb{P}(\theta_{\star}|D_{t-1})$.
- Choose $X_t = \operatorname{argmax}_{x \in \mathcal{X}} \lambda_{t-1}^+(\theta, D_{t-1}, x)$. $Y_{X_t} \leftarrow \text{conduct experiment at } X_t$.
- Set $D_t \leftarrow D_{t-1} \cup \{(X_t, Y_{X_t})\}.$ 6:

A natural question that may arise is the need to sample from the posterior ρ_{t-1} for θ_{\star} , instead of taking an expectation of λ^+ over ρ_{t-1} . In fact, many policies for non-adaptive L-DOE take an expectation over the posterior (Ouyang et al., 2016). However, in adaptive settings where θ_{\star} is unknown, taking the expectation may fail as it may not explore sufficiently. For example, in bandit problems, which is a special case of our setting (Section 3.5), this amounts to choosing the maximum of the posterior mean of the payoff function, which is known to fail spectacularly.

Computational considerations: It is worth pointing out some computational considerations in Algorithm 1. First, sampling from the posterior in step 3 might be difficult, especially in complex Bayesian models. Fortunately however, the field of Bayesian inference has made great strides in the recent past with the development of fast techniques for approximate inference methods such as MCMC or variational inference (Hensman et al., 2012; Neiswanger et al., 2015). Moreover, today we have efficient probabilistic programming tools (Bingham et al., 2018; Carpenter et al., 2017; Tran et al., 2017) that allow a practitioner to intuitively incorporate domain expertise via a prior and obtain the posterior given data. Secondly, the maximisation of the look ahead reward in step 4 can also be non-trivial, especially since it might involve empirically computing the expectation in (2). This is similar to existing work in Bayesian optimisation which assume access to such an optimisation oracle (Bull, 2011; Srinivas et al., 2010). That said, in many practical settings where experiments can cost significant time and money, these considerations are less critical.

Despite these concerns, it is worth mentioning that myopic strategies are still computationally far more attractive than policies which try to behave globally optimally. For example, extending MPS to a k step look-ahead might involve an optimisation over \mathcal{X}^k in step 4 of Algorithm 1 which might be impractical for large values of k. Moreover, in many problems where system characteristics θ_{\star} are known to the decision maker, myopic policies can be competitive with globally optimal policies (Golovin and Krause, 2011; Nemhauser et al., 1978; Wei et al., 2015). In Section 4, we identify conditions where $\pi_{\mathrm{M}}^{\mathrm{PS}}$ can be competitive with the globally optimal policy $\pi_{\mathrm{G}}^{\mathrm{c}}$ which knows θ_{\star} .

3. Examples & Experiments

We now describe some concrete examples of DOE problems that can be specified by a reward function λ and present experimental results. We compare $\pi_{\mathrm{M}}^{\mathrm{\scriptscriptstyle PS}}$ to random sampling (RAND), the myopically optimal policy π_{M}^{\star} which has access to θ_{\star} , and to specialised methods for the particular problem, when available. In the interest of aligning our experiments with our theoretical analysis, we compare methods on both criteria in (1), although in these applications, the final reward $\lambda(\theta_{\star}, D_n)$ is more relevant than the cumulative one $\Lambda(\theta_{\star}, D_n)$. In all cases, except Experiments 2 and 4 which have conjugate priors, We use variational inference (VI) in Edward (Tran et al., 2017) to approximate the posterior $\mathbb{P}(\theta_{\star}|D_t)$. While VI is known to underestimate the variance in practice, it worked well in our experiments. For better visualisation, we plot the negative reward in a semilog plot. We defer some experimental details to Appendix D.

High-level Takeaways: Despite being a quite general, $\pi_{\mathrm{M}}^{\mathrm{PS}}$ outperforms, or performs as well as, specialised methods. $\pi_{\mathrm{M}}^{\mathrm{PS}}$ is competitive, but typically worse than the non-realisable π_{M}^{\star} . Finally $\pi_{\mathrm{M}}^{\mathrm{PS}}$ enables effective DOE in complex settings where no prior methods seem applicable.

3.1. Active Learning

Problem: As described previously, we wish to learn some parameter $\beta_{\star} = \beta(\theta_{\star})$ which is a function of the true parameter θ_{\star} . Each time we query some $X \in \mathcal{X}$, we observe a label $Y \sim \mathbb{P}(Y|X,\theta_{\star})$. We conduct two synthetic exper-

iments in this setting. We use $\lambda(\theta_\star) \stackrel{\triangle}{=} - \|\beta_\star - \hat{\beta}(D_n)\|_2^2$ as the reward where $\hat{\beta}$ is a regularised maximum likelihood estimator. In addition to RAND and π_M^\star , we compare π_M^{PS} to ActiveSetSelect of Chaudhuri et al. (2015).

Experiment 1: We use the following parametric model: $Y_x|x,\theta \sim \mathcal{N}(f_\theta(x),\eta^2)$ where $f_\theta(x) = \frac{a}{1+e^{b(x-c)}}$ is a logistic function. The true parameter is $\theta_\star = (a,b,c,\eta^2)$ and our goal is to estimate $\beta_\star = (a,b,c)$. The MLE is computed via gradient ascent on the log likelihood. In our experiments, we used a=2.1,b=7,c=6 and $\eta^2=0.01$ as θ_\star . We used normal priors $\mathcal{N}(2,1),\mathcal{N}(5,3)$ and $\mathcal{N}(5,3)$ for a,b,c respectively and an inverse gamma $\mathrm{IG}(20,1)$ prior for η^2 . As the action space, we used $\mathcal{X}=[0,10]$. For variational inference, we used a normal approximation for the posterior for a,b,c and an inverse gamma approximation for η^2 . The results are given in the first column of Figure 1.

Experiment 2: In the second example, we use the following linear regression model: $Y_x|x,\theta \sim \mathcal{N}(f_\theta(x),0.01)$ where $f_\theta(x) = \sum_{i=1}^{16} \theta_{*i}\phi(x-c_i)$. Here, $\phi(v) = \frac{1}{\sqrt{0.2\pi}}e^{-5\|v\|_2^2}$ and the points c_1,\ldots,c_{16} were arranged in a 4×4 grid within $[0,1]^2$. We set $\theta_{*i}=g(c_i)$, with $g(v)=\sin(3.9\pi((v_1-0.1)^2+v_2+0.1))$. Our goal is to estimate $\beta_*=\theta_*$. As the action space, we used $\mathcal{X}=[0,1]^2$. The posterior for θ_* was calculated in closed form using a normal distribution $\mathcal{N}(0,I_{16})$ as the prior. The results are given in the second column of Figure 1.

Alternative Problem Formalism: A common formalism for parameter estimation in discriminative models (Chaudhuri et al., 2015; Frostig et al., 2015) is to maximise the expected likelihood of the data for a given sampling distribution Γ on \mathcal{X} . Here, one wishes to maximise $\lambda(\theta_\star, D_n) \stackrel{\Delta}{=} \mathbb{E}_{X \sim \Gamma, Y \sim \mathbb{P}(Y|X, \theta_\star)}[\log \mathbb{P}(Y|X, \hat{\theta})]$, where $\hat{\theta}$ is an estimator for θ obtained from D_n .

Experiments 3 & 4: We use the same models as in Experiment 1 & 2 but with the above reward function. We let Γ be the uniform distribution on the respective domains and $\hat{\theta}$ be the maximum likelihood estimator for θ . The results are given in the third and fourth columns of Figure 1.

3.2. Posterior Estimation & Active Regression

Problem: Consider estimating a non-parametric function $f_{\theta_{\star}}$, which is known to be uniformly smooth. An action $x \in \mathcal{X}$ queries $f_{\theta_{\star}}$, upon which we observe $Y_x = f_{\theta_{\star}}(x) + \epsilon$, where $\mathbb{E}[\epsilon] = 0$. If the goal is to learn $f_{\theta_{\star}}$ uniformly well in L^2 error, i.e. with reward $-\|f_{\theta_{\star}} - \hat{f}(D_n)\|^2$, adaptive techniques may not perform significantly better than non-adaptive ones (Willett et al., 2006). However, if our reward was $\lambda(\theta_{\star}, D_n) \stackrel{\triangle}{=} -\|\sigma(f_{\theta_{\star}}) - \sigma(\hat{f}(D_n))\|^2$ for some monotone super-linear transformation σ , then adaptive techniques may do better by requesting more evaluations at regions

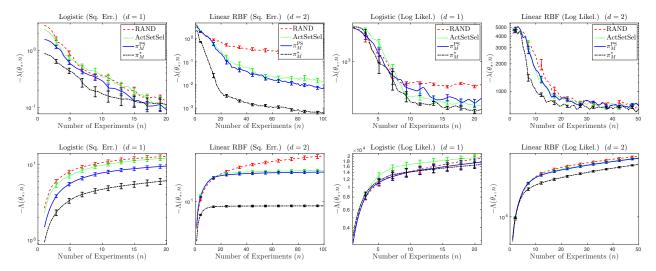


Figure 1: Results on the synthetic active learning experiments in Section 3.1. The title states the model and the dimensionality. In all figures, the x axis is the number of experiments n. In the top four figures, the y axis is the final negative reward $-\lambda(\theta_*, n)$ at the nth iteration. In the bottom figures, it is the corresponding negative cumulative reward $-\Lambda(\theta_*, n)$. Lower is better in both cases. The legend for all plots is given in the top left figure. All curves were averaged over 20 runs, and error bars indicate one standard error.

with high $f_{\theta_{\star}}$ value. This is because, $\lambda(\theta_{\star}, D_n)$ is more sensitive to such regions due to the transformation σ .

A particularly pertinent instance of this formulation arises in astrophysical applications where one wishes to estimate the posterior distribution of cosmological parameters, given some astronomical data Q (Parkinson et al., 2006). Here, an astrophysicist specifies a prior Ξ over the cosmological parameters $Z \in \mathcal{X}$, and the likelihood of the data for a given choice of the cosmological parameters $x \in \mathcal{X}$ is computed via an expensive astrophysical simulation. The prior and the likelihood gives rise to an unknown log joint density $f_{\theta_{\star}}$ defined on \mathcal{X} , and the goal is to estimate the the joint density $p(Z = x, Q) = \exp(f_{\theta_*}(x))$ so that we can perform posterior inference. Adopting assumptions from Kandasamy et al. (2015), we model $f_{\theta_{\star}}$ as a Gaussian process, which is reasonable since we expect a log density to be smoother than the density itself. As we wish to estimate the joint density, λ takes the above form with $\sigma = \exp$.

Experiment 5: We use data on Type I-a supernova from Davis et al (2007). We wish to estimate the posterior over the Hubble constant $H \in (60,80)$, the dark matter fraction $\Omega_M \in (0,1)$ and the dark energy fraction $\Omega_E \in (0,1)$, which constitute our three dimensional action space \mathcal{X} . The likelihood is computed via the Robertson-Walker metric. In addition to π_{M}^{\star} and RAND, we compare $\pi_{\mathrm{M}}^{\mathrm{PS}}$ to Gaussian process based exponentiated variance reduction (GP-EVR) (Kandasamy et al., 2015) designed for this setting. We evaluate the reward via numerical integration. The results are presented in the first column of Figure 2.

3.3. Level Set Estimation

Problem: In active level set estimation (LSE), one wishes to determine which regions of a space $\mathcal X$ fall above or below a given level set of an expensive to evaluate function $f_{\theta_{\star}}$. An experiment evaluates this function and returns $Y_x = f_{\theta_{\star}}(x) + \epsilon$, where $\mathbb{E}[\epsilon] = 0$. We adopt the setting of Gotovos et al. (2013a), where a method for LSE returns its predictions for being above/below the threshold on a pre-specified set of discrete points $\mathcal{X}' \subset \mathcal{X}$. The reward function λ is set to be average prediction accuracy.

Experiment 6: We used data on luminous red galaxies (LRGs) to compute the galaxy power spectrum of 9 cosmological parameters including the spatial curvature, cold dark matter density, and baryonic density. We wish to find regions of the cosmological parameter space, where the power spectrum is larger than a pre-specified threshold. Software and data were taken from Tegmark et al (2006). We compare $\pi_{\rm M}^{\rm PS}$ to random search, $\pi_{\rm M}^{\star}$, and the Gaussian process based level set estimation (GP-LSE) method of Gotovos et al. (2013a). Following Gotovos et al. (2013a), we model the power spectrum as a GP, and define the reward function as described above where \mathcal{X}' is a set of $\sim 20K$ points. The results are presented in the second column of Figure 2.

3.4. Combined and Customised Objectives

Problem: In many real world problems, one needs to design experiments with multiple goals. For example, an experiment might evaluate multiple objectives, and the task might be to optimise some of them, while learning the parameters for another. Classical methods specifically designed for active learning or optimisation may not be suitable in such

¹ One should not conflate the prior over \mathcal{X} specified with the astrophysics model, with prior over Θ assumed in our set up.

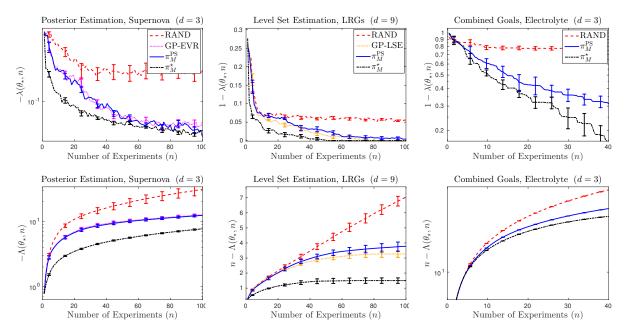


Figure 2: Results on the real experiments. The first column is for the posterior estimation problem in Sec. 3.2, the second column is for the level set estimation problem in Sec. 3.3, and the third column is for the combined objective problem in Sec. 3.2, In the top figures, the y axis is the negative reward $-\lambda(\theta_{\star}, D_n)$ and in the bottom figures, it is the negative cumulative reward $-\Lambda(\theta_{\star}, D_n)$ for the corresponding experiment. See caption under Figure 1 for more details.

settings. One advantage to the proposed framework is that it allows us to combine multiple goals in the form of a reward function. For instance, if an experiment measures two functions $f_{\theta_{\star},1}, f_{\theta_{\star},2}$ and we wish to learn f_1 while optimising f_2 , we can define the reward as $\lambda(\theta_{\star}, D_n) = -\|f_{\theta_{\star},1} - \hat{f}_1(D_n)\|^2 + \max_{X_t, t \leq n} \left(f_{\theta_{\star},2}(X_t) - \max_x f_{\theta_{\star},2}(x)\right)$. Here \hat{f}_1 is an estimate for $f_{\theta_{\star},1}$ obtained from the data, $\|\cdot\|$ is the L_2 norm and $\max_{X_t, t \leq n} f_{\theta_{\star},2}(X_t)$ is the maximum point of $f_{\theta_{\star},2}$ we have evaluated so far. Below, we demonstrate one such application.

Experiment 7: In battery electrolyte design, one tests an electrolyte composition under various physical conditions. On an experiment at $x \in \mathcal{X}$, we obtain measurements $Y_x = (Y_{x,\text{sol}}, Y_{x,\text{vis}}, Y_{x,\text{con}})$ which are noisy measurements of the solvation energy f_{sol} , the viscosity f_{vis} and the specific conductivity f_{con} . Our goal is to estimate f_{sol} and f_{vis} while optimising f_{con} . Hence,

$$\lambda(\theta_{\star}, D_n) = \alpha \Big(\max_{X_t, t \le n} f_{\text{con}}(X_t) - \max_{x \in \mathcal{X}} f_{\text{con}}(x) \Big)$$
$$-\beta \|f_{\text{sol}} - \hat{f}_{\text{sol}}(D_n)\|^2 - \gamma \|f_{\text{vis}} - \hat{f}_{\text{vis}}(D_n)\|^2,$$

where, the parameters α,β,γ were chosen so as to scale each objective and ensure that none of them dominate the reward. In our experiment, we use the dataset from Gering (2006). Our action space $\mathcal X$ is parametrised by the following three variables: $Q\in(0,1)$ measures the proportion of two solvents EC and EMC in the electrolyte, $S\in(0,3.5)$ is the molarity of the salt LiPF $_6$ and $T\in(-20,50)$ is the tempera-

ture in Celsius. We use the following prior which is based off a physical understanding of the interaction of these variables. $f_{\rm con}:\mathcal{X}\to\mathbb{R}$ is sampled from a Gaussian process (GP), $f_{\rm vis}(Q,S,T)=\exp(-aT)g_{\rm vis}(Q,S)$ where $g_{\rm vis}$ is sampled from a GP, and $f_{\rm sol}(Q,S,T)=b+\exp(cQ-dS-eT)$. We use inverse gamma priors for a,b,d,e and a normal prior for c. For variational inference, we used inverse gamma approximations for a,b,d,e, a normal approximation for c, and GP approximations for $f_{\rm con}$ and $g_{\rm vis}$. We use the posterior mean of $f_{\rm sol}$ and $f_{\rm vis}$ under this prior as the estimates $\hat{f}_{\rm sol},\hat{f}_{\rm vis}$. We present the results in the third column of Figure 2 where we compare RAND, $\pi_{\rm M}^{\rm PS}$ and $\pi_{\rm M}^{\star}$. This is an example of a customised DOE problem for which no prior method seems directly applicable.

3.5. Bandits & Bayesian Optimisation

Bandits and Bayesian optimisation are self-evident special cases of our formulation. Here, θ_{\star} specifies a function $f_{\theta_{\star}}: \mathcal{X} \to \mathbb{R}$. When we choose a point $X \in \mathcal{X}$ to evaluate the function, we observe $Y_X = f_{\theta_{\star}}(X) + \epsilon$ where $\mathbb{E}[\epsilon] = 0$. In the bandit framework, the reward is the instantaneous regret $\lambda(\theta_{\star}, D_n) = f_{\theta_{\star}}(X_n) - \max_{x \in \mathcal{X}} f_{\theta_{\star}}(x)$. In Bayesian optimisation, one is interested in simply finding a single value close to the optimum and hence $\lambda(\theta_{\star}, D_n) = \max_{t \leq n} f_{\theta_{\star}}(X_t) - \max_{x \in \mathcal{X}} f_{\theta_{\star}}(x)$. In either case, $\pi_{\mathrm{M}}^{\mathrm{PS}}$ reduces to the Thompson sampling procedure as $\arg\max_{x \in \mathcal{X}} \lambda^+(\theta_{\star}, D_{t-1}, x) = \arg\max_{x \in \mathcal{X}} f_{\theta}(x)$, where f_{θ} is a random function drawn from the posterior. Since

prior work has demonstrated that TS performs empirically well in several real world optimisation tasks (Chapelle and Li, 2011; Hernández-Lobato et al., 2017; Kandasamy et al., 2018), we omit experimental results for this example. One can also cast other variants of Bayesian optimisation, including multi-objective optimisation (Hernández-Lobato et al., 2016; Paria et al., 2018) and constrained optimization (Gardner et al., 2014), in our general formulation.

4. Theoretical Analysis

In this section we derive theoretical guarantees for $\pi_{\mathrm{M}}^{\mathrm{PS}}$. Our emphasis is on understanding conditions under which myopic algorithms which need to learn θ_{\star} can perform competitively with the myopic optimal and the globally optimal oracles π_{M}^{\star} , π_{G}^{\star} which know θ_{\star} (see Section 2). Going forward, to simplify the exposition, we will assume that λ is bounded, i.e. $\lambda:\Theta\times\mathcal{X}\to[0,1]$. Moreover, w.l.o.g, we will assume for all $\theta\in\Theta$, $\sup_{D\in\mathcal{D}}\lambda(\theta,D)=1$. This condition is for free since for any bounded reward λ' , we can set $\lambda(\theta,D)\stackrel{\triangle}{=} 1+\lambda'(\theta,D)-\sup_{D\in\mathcal{D}}\lambda'(\theta,D)$.

For criterion (a), we are interested in upper bounding $\mathbb{E}[\Lambda(\theta_\star,D_n)|D_n\sim\pi_{\mathrm{M}}^{\mathrm{PS}}]$ in terms of $\mathbb{E}[\Lambda(\theta_\star,D_n)|D_n\sim\pi_{\mathrm{M}}^{\star}]$, which yields a *cumulative regret bound*, and for criterion (b), we wish to bound $\mathbb{E}[\lambda(\theta_\star,D_n)|D_n\sim\pi_{\mathrm{M}}^{\mathrm{PS}}]$ in terms of the analogous quantities for π_{M}^{\star} , π_{G}^{\star} , which serves as a *final regret bound*. Note that a comparison with π_{G}^{\star} on (a) is meaningless since it might take low reward actions in the early stages in order to do well in the long run. In fact, our bounds for (a) will hold when $\lambda(\theta_\star,D)$ is an ordered multi-set function in D, but for (b) when $\lambda(\theta_\star,D)$ is a multi-set function, i.e. the ordering does not matter. Our bounds will hold in expectation over $\theta_\star\sim\rho_0$.

The following proposition shows that without further assumptions, a non-trivial regret bound is impossible. Such results are common in the RL literature, and necessitate several structural assumptions (Dann and Brunskill, 2015; Jaksch et al., 2010; Kearns and Singh, 2002). Its proof is given in Appendix B.4.

Proposition 1. For all policies π which do not know θ_{\star} , there exists a DOE problem where $\mathbb{E}_{\theta_{\star} \sim \rho_0}[\lambda(\theta_{\star}, D_n^{\star}) - \lambda(\theta_{\star}, D_n)|D_n^{\star} \sim \pi_{\mathrm{M}}^{\star}, D_n \sim \pi] \geq 1/2$ for all $n \geq 1$.

Motivated by this lower bound, we impose the following condition on the parameter space and reward structure, under which a policy can achieve sub-linear regret. For this, first note we can assume that, at all time steps, the observations $Y \sim \mathbb{P}(\cdot|x,\theta_\star)$ are generated for all $x \in \mathcal{X}$, but we only observe those for the chosen X_t . With this in consideration, let $\mathbb{E}_{Y,t+1:|\theta}$ denote expectation over all observations generated from time t+1 onwards when $\theta_\star = \theta$.

Condition 1. Let θ, θ' denote parameter values in Θ and $\pi_{\mathrm{M}}^{\theta}, \pi_{\mathrm{M}}^{\theta'}$ be the myopically optimal policies when $\theta_{\star} = \theta$,

and $\theta_{\star} = \theta'$ respectively. Let H denote a data sequence and D_n , D'_n be the data sequences collected by π^{θ}_M and $\pi^{\theta'}_M$ respectively when starting from H when $\theta_{\star} = \theta$ and $\theta_{\star} = \theta'$ respectively, i.e. the myopically optimal policies operate in their respective environments. Then, there exists sequences $\{\epsilon_n\}_{n\geq 1}$, $\{\tau_n\}_{n\geq 1}$ such that the following hold.

1. π_{M}^{θ} achieves asymptotically similar reward $\forall \theta \in \Theta$. That is.

$$\sup_{\theta,\theta' \in \Theta} \sup_{H \in \mathcal{D}} \left\{ \mathbb{E}_{Y,|H|+1:|\theta} \lambda(\theta, H \uplus D_n) - \mathbb{E}_{Y,|H|+1:|\theta'} \lambda(\theta', H \uplus D'_n) \right\} \leq \epsilon_n.$$

2. The rate of convergence is better than $\mathcal{O}(1/\sqrt{n})$. That is, letting $\sqrt{\tau_n} = 1 + \sum_{j=1}^n \epsilon_j$, we have $\tau_n \in o(n)$.

The condition states that when we execute π_{M}^{\star} , the myopically optimal policy which knows and depends on the value of θ_{\star} , from any prefix H, it achieves asymptotically similar λ for all values of θ_{\star} . It is worth emphasising that the condition involves executing $\pi_{\mathrm{M}}^{\theta}$ in the environment where $\theta_{\star} = \theta$. A condition of the above form seems necessary for any myopic algorithm π that does not know θ_{\star} for the following reason. Assume that the myopic π_{M}^{\star} can quickly achieve large λ value when $\theta_{\star} \in \Theta_q$ but is slow when $\theta_{\star} \in \Theta_b$. Since π does not know θ_{\star} it needs to hedge against the "bad" situation, i.e. $\theta_{\star} \in \Theta_b$. However, in doing so, it will necessarily perform poorly against $\pi_{\rm M}^{\star}$ when $\theta_{\star} \in \Theta_g$ as π_{M}^{\star} can quickly achieve large λ . Condition 1 prevents such situations. As we will see shortly, the regret for $\pi_{\mathrm{M}}^{\mathrm{PS}}$ will depend on au_n which dictates how differently π_{M}^{\star} can behave for different values of θ_{\star} . In particular, sublinearity of τ_n is necessary for sublinear regret with π_M^* .

In Appendix C we provide a more interpretable sufficient condition which implies Condition 1, and demonstrate that it is satisfied with $\tau_n \in \mathcal{O}(1)$ for bandit and black-box optimisation problems and $\tau_n \in \mathcal{O}(\log n)$ for an active learning problem. We also consider a setting where λ has "state-like" structure; under assumptions similar to standard assumptions in reinforcement learning with ergodic Markov decision processes, we are able to show that Condition 1 holds. Finally we mention that if Condition 1 holds for two reward functions λ_1, λ_2 , it is also true for the sum, $\lambda_1 + \lambda_2$, and the product, $\lambda_1 \cdot \lambda_2$, and can thus be applied to combined objective settings such as in Section 3.4.

Before stating the main theorem, we introduce the maximum information gain, Ψ_n , which captures the statistical difficulty of the learning problem.

$$\Psi_n = \max_{D_n \subset \mathcal{D}_n} I(\theta_\star; D_n). \tag{3}$$

Here $I(\cdot;\cdot)$ is the Shannon mutual information. Ψ_n measures the maximum information a set of n action-observation pairs can tell us about the true parameter θ_{\star} .

The quantity appears as a statistical complexity measure in many Bayesian adaptive data analysis settings (Gotovos et al., 2013b; Ma et al., 2015; Srinivas et al., 2010). Below, we list some examples of common models which demonstrate that Ψ_n is typically sublinear in n.

Example 2. We have the following bounds on Ψ_n for common models (Srinivas et al., 2010):

- 1. Finite sets: If Θ is finite, $\Psi_n \leq \log(|\Theta|)$ for all n.
- 2. Linear models: Let $\mathcal{X} \subset \mathbb{R}^d$, $\theta \in \mathbb{R}^d$, and $Y_x|x, \theta \sim \mathcal{N}(\theta^\top x, \eta^2)$. For a multi-variate Gaussian prior on θ_* , $\Psi_n \in \mathcal{O}(d\log(n))$.
- 3. Gaussian process: For a Gaussian process prior with RBF kernel over a compact domain $\mathcal{X} \subset \mathbb{R}^d$, and with Gaussian likelihood, we have $\Psi_n \in \mathcal{O}(\log(n)^{d+1})$.

We now state our main theorem for finite action spaces \mathcal{X} .

Theorem 2. Let X be finite and assume Condition 1 holds. Let τ_n be as defined in Condition 1. Then,

$$\mathbb{E}[\Lambda(\theta_\star, \pi_{\mathrm{M}}^\star) - \Lambda(\theta_\star, \pi_{\mathrm{M}}^{\scriptscriptstyle{\mathrm{PS}}})] \leq \sqrt{\frac{|\mathcal{X}| n \tau_n \Psi_n}{2}}.$$

Theorem 2 establishes a sublinear regret bound for $\pi_{\mathrm{M}}^{\mathrm{PS}}$ against π_{M}^{\star} when $\tau_n\Psi_n\in o(n)$. The $|\mathcal{X}|$ term captures the complexity of our action space, Ψ_n captures the complexity of the prior on θ_{\star} . The \sqrt{n} dependence is in agreement with prior results for Thompson sampling (Kaufmann et al., 2012; Russo and Van Roy, 2016b). Thus, under Condition 1, $\pi_{\mathrm{M}}^{\mathrm{PS}}$ is competitive with the myopic optimal policy π_{M}^{\star} , with average regret tending to 0.

We now compare $\pi_{\mathrm{M}}^{\mathrm{PS}}$ to the globally optimal policy π_{G}^{\star} , when λ is a multi-set function, i.e. the ordering in D_n does not matter. For this, we first introduce the notions of monotonicity and adaptive submodularity.

Condition 2. (Monotonicity and Adaptive Submodularity (Golovin and Krause, 2011)) Let \mathbb{E}_{Y_x} denote the expectation over the likelihood $Y_x \sim \mathbb{P}(\cdot|x,\theta_\star)$. The following two statements are true for all $\theta \in \Theta$, $D, D' \in \mathcal{D}$, $D \prec D'$, and $x \in \mathcal{X}$. λ is a monotone, meaning that $\mathbb{E}_{Y_x}[\lambda(\theta,D \uplus \{(x,Y_x)\})] \geq \lambda(\theta,D)$. Moreover, λ is adaptive submodular, meaning that,

$$\mathbb{E}_{Y_x}[\lambda(\theta, D \uplus \{(x, Y_x)\})] - \lambda(\theta_{\star}, D)$$

$$\geq \mathbb{E}_{Y_x}[\lambda(\theta, D' \uplus \{(x, Y_x)\})] - \lambda(\theta_{\star}, D').$$

Monotonicity states that adding more data increases the reward in expectation, while adaptive submodularity formalises a notion of diminishing returns. That is, performing the same action is more beneficial when we have less data. It is easy to see that some assumption is needed here, since

even in simple problems π_M^\star can be arbitrarily worse than π_G^\star . We now state the second main result of this paper.

Theorem 3. Assume that λ satisfies conditions 1 and 2. Let τ_n be as defined in Theorem 2. Then, for all $\gamma < 1$, we have

$$\mathbb{E}[\lambda(\theta_{\star}, D_{n})|D_{n} \sim \pi_{\mathrm{M}}^{\mathrm{PS}}] \geq (1 - \gamma)\mathbb{E}[\lambda(\theta_{\star}, D_{\gamma n}^{\star})|D_{\gamma n}^{\star} \sim \pi_{\mathrm{G}}^{\star}] - \sqrt{\frac{|\mathcal{X}|\tau_{n}\Psi_{n}}{2n}}.$$

The theorem states that $\pi_{\mathrm{M}}^{\mathrm{PS}}$ in n steps is guaranteed to perform up to a $1-\gamma$ factor as well as π_{G}^{\star} executed for $\gamma n < n$ steps, up to an additive $\sqrt{\tau_n \Psi_n/n}$ term. The result captures both approximation and estimation errors, in the sense that we are using a myopic policy to approximate a globally optimal one, and we are learning a good myopic policy from data. In comparison, prior works on adaptive submodular optimisation focus on approximation errors and typically achieve 1-1/e approximation ratios against the n steps of π_{G}^{\star} . Our bound is quantitatively worse, but focusing on a much more difficult task, and we view the results as complementary. Observe that an analogous bound holds against π_{M}^{\star} , since it is necessarily worse that π_{G}^{\star} .

Finally, we believe that the above results can be generalised to large or infinite action spaces under additional structure on λ . For example, when $\mathcal{X} \subset \mathbb{R}^d$, and the expected rewards are linear in the actions taken, we expect an $\mathcal{O}(d)$ dependence similar to linear bandit settings (Agrawal and Goyal, 2013; Russo and Van Roy, 2016b). Algorithm 1 can be applied as is when we can execute multiple experiments in parallel. We expect that similar results to Theorems 2 and 3 should hold, with mild dependence on the number of workers, using similar analyses to Kandasamy et al. (2018).

5. Conclusion

We study settings for adaptive goal oriented DOE problems in a Bayesian setting. Our formulation is quite general, allowing practitioners to incorporate domain knowledge via a probabilistic model, and specify their goal via a reward function that may depend on system characteristics. We focus on myopic policies due to their computational simplicity. Yet, our empirical results demonstrate that MPS has broad applicability, performing favourably with more specialised methods, and enabling complex DOE tasks where existing methods are not applicable. Our theoretical results establish conditions under which a myopic algorithm based on posterior sampling is competitive with myopic and globally optimal policies, both of which know the underlying system parameters. One interesting avenue for future theoretical work is to relax and/or find other conditions under which myopic strategies can do well. For instance, we believe that Condition 1 is stronger than necessary, and that it is sufficient if $\pi_{\rm M}^{\star}$ is able to do well in its own environment.

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