### Switching to Sampling.. ...in order to Switch



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### What I'll talk about

- 1. Briefly compare stateless sampling vs. statefull sampling
- 2. Briefly discuss the general algorithm of mcmc
- 3. Briefly discuss the idea of sampling for inference
- 4. Discuss the switchpoint problem
- 5. Demo how to solve this via Pymc3
- 6. Mention some things to look out for in the future











This approach doesn't just work for  $\pi$ , it actually works for a lot of general cases too. Sampling might be a very useful to discover stochastic patterns in a system.

I'll demonstrate a simple numeric application.

```
nums = np.random.normal(3, 1, 10)
start_mu = 3
stepsize = 0.1
samples = []
for i in range(5000):
    new_mu = start_mu + np.random.normal(0, stepsize, 1)[0]
    old_loglik = np.prod([stats.norm.pdf(_, loc = start_mu) for _ in nums])
    new_loglik = np.prod([stats.norm.pdf(_, loc = new_mu) for _ in nums])
    if new_loglik > old_loglik:
        start_mu = new_mu
    else:
        if np.random.random() < new_loglik/old_loglik:
            start_mu = new_mu
    samples.append(start_mu)
```

nums = np.random.normal(3, 1, 5)



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nums = np.random.normal(3, 1, 10)



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nums = np.random.normal(3, 1, 100)



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# "Thats great Vincent, but what about applications?"

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There is a point in time when something change a system. We would like to find some algorithmic way of finding out when it happened and what the change was to the system.



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I'll first explain a simple example where the system has 5 parameters; au,  $\mu_1$ ,  $\sigma_1$ ,  $\mu_2$ ,  $\sigma_2$ .



Given a dataset we're interested in finding;

 $\mathbb{P}(s,\mu_1,\mu_2,\sigma_1,\sigma_2|D)$ 

Let's apply bayes rule.

 $\mathbb{P}(s,\mu_1,\mu_2,\sigma_1,\sigma_2|D) \propto \mathbb{P}(D|s,\mu_1,\mu_2,\sigma_1,\sigma_2)\mathbb{P}(s,\mu_1,\mu_2,\sigma_1,\sigma_2)$  $=\Pi_i \mathbb{P}(y_i|s,\mu_1,\mu_2,\sigma_1,\sigma_2)\mathbb{P}(s,\mu_1,\mu_2,\sigma_1,\sigma_2)$ 

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I know some information about  $\mathbb{P}(s, \mu_1, \mu_2, \sigma_1, \sigma_2)$ .

For example;  $\sigma_i > 0$ . Priors are useful!

The other part is the part that we model.

$$\mathbb{P}(y_i|s,\mu_1,\mu_2,\sigma_1,\sigma_2) = egin{cases} y_i \sim N(\mu_1,\sigma_1) \ y_i \sim N(\mu_2,\sigma_2) \end{cases}$$

So that means we have the building blocks for;

 $\mathbb{P}(s,\mu_1,\mu_2,\sigma_1,\sigma_2|D) \propto \Pi_i \mathbb{P}(y_i|s,\mu_1,\mu_2,\sigma_1,\sigma_2) \mathbb{P}(s,\mu_1,\mu_2,\sigma_1,\sigma_2)$ 

This begs the question, why would we prefer to use MCMC here?

I can come up with three reasons;

- Sampling is a very general framework. Can be parallized and take any shape.
- The search space in general isn't convex, so gradient methods lose their guarantee.
- By sampling, we get an impression of the distribution over all parameters; not just a mere MLE.

```
import pymc3 as pm
basic_model = pm.Model()
with basic_model:
   mu1 = pm.Normal('mu1', mu=0, sd=2)
   mu2 = pm.Normal('mu2', mu=0, sd=2)
    sigma1 = pm.HalfNormal('sigma1', sd=2)
    sigma2 = pm.HalfNormal('sigma2', sd=2)
    switchpoint = pm.DiscreteUniform('switchpoint', 0, time.max())
    tau_mu = pm.switch(time >= switchpoint, mu2, mu1)
```

```
tau_sigma = pm.switch(time >= switchpoint, sigma2, sigma1)
```

```
y = pm.Normal('y1', mu=tau_mu, sd=tau_sigma, observed=x)
trace = pm.sample(10000)
```

This last mentioned trace variable can be used for plotting.

= pm.traceplot(trace)

Let's look at the results for this dataset.



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Let's now change the problem.



```
adv_model = pm.Model()
with adv_model:
    mu1 = pm.Normal('mu1', mu=0, sd=4)
    mu2 = pm.Normal('mu2', mu=0, sd=4)
    sigma1 = pm.HalfNormal('sigma1', sd=4)
    sigma2 = pm.HalfNormal('sigma2', sd=4)
    switchpoint1 = pm.DiscreteUniform('switchpoint1', 0, time.max() - 1)
    switchpoint2 = pm.DiscreteUniform('switchpoint2', switchpoint1, time.max())
```

```
tau_mu1 = pm.switch(time >= switchpoint1, mu2, mu1)
tau_mu2 = pm.switch(time >= switchpoint2, mu1, tau_mu1)
tau_sigma1 = pm.switch(time >= switchpoint1, sigma2, sigma1)
tau_sigma2 = pm.switch(time >= switchpoint2, sigma1, tau_sigma1)
```

```
y = pm.Normal('y1', mu=tau_mu2, sd=tau_sigma2, observed=x)
adv_trace = pm.sample(5000)
```



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Combat the start of the trace by ignoring it.

\_ = pm.traceplot(trace[1000:])



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 $\left( \begin{array}{c} \\ \end{array} \right)$ 

### You can also retreive statistics.

 $rg\max_{\lambda} \mathbb{P}(D|\lambda)\mathbb{P}(\lambda)$ 

- > pm.find\_MAP(model=adv\_model)
- {'mu1': array(1.3253008730440974),
  - 'mu2': array(0.0),
  - 'sigma1\_log\_': array(0.8639975314985741),
  - 'sigma2\_log\_': array(1.3862943591402144),
  - 'switchpoint1': array(224),
  - 'switchpoint2': array(224)}

### The future

There's people doing this type of thing for neural networks.

 $\mathbb{P}(w|D) \propto \mathbb{P}(D|w)\mathbb{P}(w)$ 

Sampling this is crazy, so people sometimes resort to variational inference. It's an interesting thought and you can already start playing with it. <u>Here</u> is a blogpost from pymc3 contributor explaining how to use PyMC3 to train a neural network for the mnist dataset via lasagna.

### The future

Outside of Pymc3 it seems like <u>edward</u> is another contender for variational inference and probibalistic modelling.

If you're interested in just maximum likelihood; maybe check out pomegrenate. If allows you to make hidden markov models as well as general PGMs (though there currently is no support for continous variables for general PGM).

If you're interested in other ways to do sampling; maybe check out STAN or emcee. Maybe consider GP's!