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Updating priors¶

In this notebook, I will show how it is possible to update the priors as new data becomes available. The example is a slightly modified version of the linear regression in the Getting started with PyMC3 notebook.

```
In [1]: %matplotlib inline
   import matplotlib.pyplot as plt
   import matplotlib as mpl
   import pymc3 as pm
   from pymc3 import Model, Normal, Slice
   from pymc3 import sample
   from pymc3 import traceplot
   from pymc3.distributions import Interpolated
   from theano import as_op
   import theano.tensor as tt
   import numpy as np
   from scipy import stats

plt.style.use('seaborn-darkgrid')
   print('Running on PyMC3 v{}'.format(pm.__version__))
```

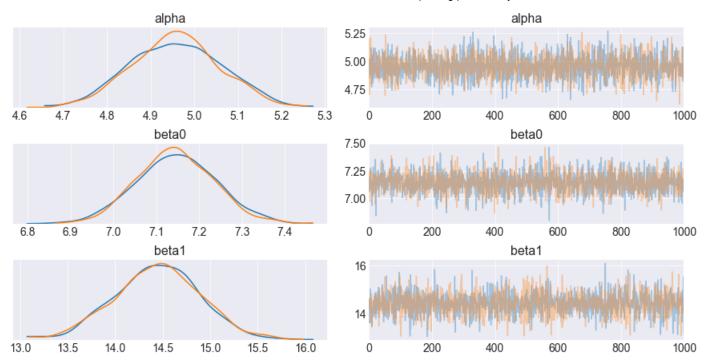
Running on PyMC3 v3.6

Generating data¶

Model specification¶

Our initial beliefs about the parameters are quite informative (sigma=1) and a bit off the true values.

```
In [3]: basic_model = Model()
        with basic_model:
            # Priors for unknown model parameters
            alpha = Normal('alpha', mu=0, sigma=1)
            beta0 = Normal('beta0', mu=12, sigma=1)
            beta1 = Normal('beta1', mu=18, sigma=1)
            # Expected value of outcome
            mu = alpha + beta0 * X1 + beta1 * X2
            # Likelihood (sampling distribution) of observations
            Y_obs = Normal('Y_obs', mu=mu, sigma=1, observed=Y)
            # draw 1000 posterior samples
            trace = sample(1000)
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt_diag...
        Multiprocess sampling (2 chains in 2 jobs)
        NUTS: [beta1, beta0, alpha]
        Sampling 2 chains: 100% 3000/3000 [00:01<00:00, 2271.43draws/s]
In [4]: traceplot(trace);
```



In order to update our beliefs about the parameters, we use the posterior distributions, which will be used as the prior distributions for the next inference. The data used for each inference iteration has to be independent from the previous iterations, otherwise the same (possibly wrong) belief is injected over and over in the system, amplifying the errors and misleading the inference. By ensuring the data is independent, the system should converge to the true parameter values.

Because we draw samples from the posterior distribution (shown on the right in the figure above), we need to estimate their probability density (shown on the left in the figure above). Kernel density estimation (KDE) is a way to achieve this, and we will use this technique here. In any case, it is an empirical distribution that cannot be expressed analytically. Fortunately PyMC3 provides a way to use custom distributions, via Interpolated class.

```
In [5]: def from_posterior(param, samples):
    smin, smax = np.min(samples), np.max(samples)
    width = smax - smin
    x = np.linspace(smin, smax, 100)
    y = stats.gaussian_kde(samples)(x)

# what was never sampled should have a small probability but not 0,
    # so we'll extend the domain and use linear approximation of density on it
    x = np.concatenate([[x[0] - 3 * width], x, [x[-1] + 3 * width]])
    y = np.concatenate([[0], y, [0]])
    return Interpolated(param, x, y)
```

Now we just need to generate more data and build our Bayesian model so that the prior distributions for the current iteration are the posterior distributions from the previous iteration. It is still possible to continue using NUTS sampling method because Interpolated class implements calculation of gradients that are necessary for Hamiltonian Monte Carlo samplers.

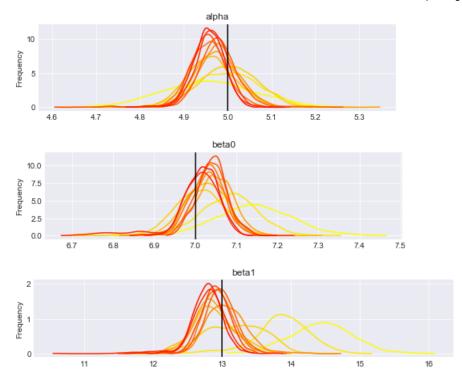
```
In [6]: traces = [trace]
In [7]: for _ in range(10):
            # generate more data
            X1 = np.random.randn(size)
            X2 = np.random.randn(size) * 0.2
            Y = alpha true + beta0 true * X1 + beta1 true * X2 + np.random.randn(size)
            model = Model()
            with model:
                # Priors are posteriors from previous iteration
                alpha = from_posterior('alpha', trace['alpha'])
                beta0 = from posterior('beta0', trace['beta0'])
                beta1 = from posterior('beta1', trace['beta1'])
                # Expected value of outcome
                mu = alpha + beta0 * X1 + beta1 * X2
                # Likelihood (sampling distribution) of observations
                Y_obs = Normal('Y_obs', mu=mu, sigma=1, observed=Y)
                # draw 10000 posterior samples
                trace = sample(1000)
                traces.append(trace)
        Auto-assigning NUTS sampler...
```

```
Initializing NUTS using jitter+adapt_diag...
Multiprocess sampling (2 chains in 2 jobs)
NUTS: [beta1, beta0, alpha]
Sampling 2 chains: 100% 3000/3000 [00:03<00:00, 988.14draws/s]
Auto-assigning NUTS sampler...
Initializing NUTS using jitter+adapt_diag...
Multiprocess sampling (2 chains in 2 jobs)
NUTS: [beta1, beta0, alpha]
Sampling 2 chains: 100% 3000/3000 [00:02<00:00, 1073.76draws/s]
Auto-assigning NUTS sampler...
Initializing NUTS using jitter+adapt_diag...
Multiprocess sampling (2 chains in 2 jobs)
NUTS: [beta1, beta0, alpha]
Sampling 2 chains: 100% 3000/3000 [00:02<00:00, 1155.98draws/s]
Auto-assigning NUTS sampler...
Initializing NUTS using jitter+adapt diag...
Multiprocess sampling (2 chains in 2 jobs)
NUTS: [beta1, beta0, alpha]
Sampling 2 chains: 100% 3000/3000 [00:03<00:00, 990.05draws/s]
Auto-assigning NUTS sampler...
Initializing NUTS using jitter+adapt diag...
Multiprocess sampling (2 chains in 2 jobs)
```

NUTS: [beta1, beta0, alpha]

```
Sampling 2 chains: 100%
                                 3000/3000 [00:02<00:00, 1117.98draws/s]
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt diag...
        Multiprocess sampling (2 chains in 2 jobs)
        NUTS: [beta1, beta0, alpha]
        Sampling 2 chains: 100% 3000/3000 [00:02<00:00, 1073.99draws/s]
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt diag...
        Multiprocess sampling (2 chains in 2 jobs)
        NUTS: [beta1, beta0, alpha]
        Sampling 2 chains: 100% 3000/3000 [00:02<00:00, 1073.40draws/s]
        The acceptance probability does not match the target. It is 0.8827037014950102, but should be close to 0.8. Try to increase the number of tuning steps.
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt diag...
        Multiprocess sampling (2 chains in 2 jobs)
        NUTS: [beta1, beta0, alpha]
        Sampling 2 chains: 100% 3000/3000 [00:02<00:00, 1045.30draws/s]
        The number of effective samples is smaller than 25% for some parameters.
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt diag...
        Multiprocess sampling (2 chains in 2 jobs)
        NUTS: [beta1, beta0, alpha]
        Sampling 2 chains: 100% 3000 3000 00:02<00:00, 887.64draws/s]
        The estimated number of effective samples is smaller than 200 for some parameters.
        Auto-assigning NUTS sampler...
        Initializing NUTS using jitter+adapt diag...
        Multiprocess sampling (2 chains in 2 jobs)
        NUTS: [beta1, beta0, alpha]
        Sampling 2 chains: 100% 3000/3000 [00:02<00:00, 1094.44draws/s]
In [8]: print('Posterior distributions after ' + str(len(traces)) + ' iterations.')
        cmap = mpl.cm.autumn
        for param in ['alpha', 'beta0', 'beta1']:
            plt.figure(figsize=(8, 2))
            for update i, trace in enumerate(traces):
               samples = trace[param]
               smin, smax = np.min(samples), np.max(samples)
               x = np.linspace(smin, smax, 100)
               v = stats.gaussian kde(samples)(x)
               plt.plot(x, y, color=cmap(1 - update_i / len(traces)))
            plt.axvline({'alpha': alpha true, 'beta0': beta0 true, 'beta1': beta1 true}[param], c='k')
            plt.ylabel('Frequency')
            plt.title(param)
        plt.tight layout();
        Posterior distributions after 11 iterations.
```

https://docs.pymc.io/notebooks/updating_priors.html



You can re-execute the last two cells to generate more updates.

What is interesting to note is that the posterior distributions for our parameters tend to get centered on their true value (vertical lines), and the distribution gets thiner and thiner. This means that we get more confident each time, and the (false) belief we had at the beginning gets flushed away by the new data we incorporate.



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