



Unique Qualifications And Capabilities

• Know-How

–Neuroscience

- Black, Wood, Wu, Method and System for automatic decoding of motor cortical activity, US Patent, 2005.
- Gasthaus, Wood, Görür, Teh, Dependent Dirichlet process spike sorting, 2009

–Machine Learning

- Neiswanger, Wood, Xing, The Dependent Dirichlet Process Mixture of Objects for Detection-free Tracking and Object Modeling, 2014
- Caron, Neiswanger, Wood, Doucet, Davy, Generalized Pólya Urn for Time-Varying Pitman-Yor Processes, In Submission 2014

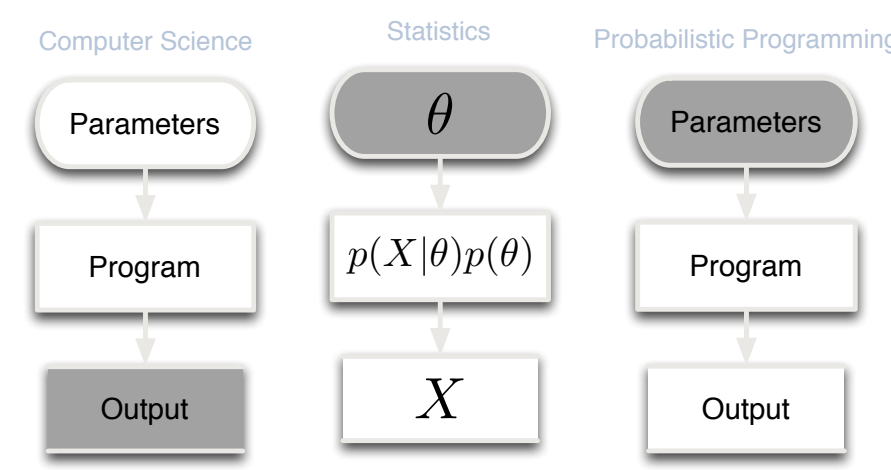
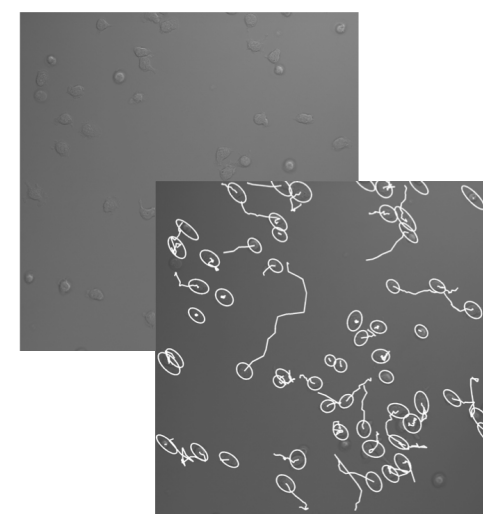
–Probabilistic Programming

- Wood, van de Meent, Mangsinghka, A New Approach to Probabilistic Programming Inference, AISTATS 2014
- Paige, Wood, A Compilation Target for Probabilistic Programming Languages, ICML 2014

• People

– Lab

– 3YP student stream



Research areas of interest

Anatomy

- Top-down biologically-parameterized, model-based regularization for tracking and structure discovery
- Efficient, massively parallel automatic inference

Function

- Automatic stochastic inversion of function simulators

Artificial Intelligence

- Expressive models and automatic inference via probabilistic programming

Probabilistic-C, a Probabilistic Programming Language @ICML 2014

Probabilistic C

Forward inference techniques such as sequential Monte Carlo [1] and particle Markov chain Monte Carlo [2] for probabilistic programming [3] can be implemented in any programming language by creative use of standardized operating system functionality including processes, forking, mutexes, and shared memory.

- Probabilistic C: a probabilistic programming intermediate representation language, which itself can be compiled to parallel machine code by standard compilers
- Standard C with two new directives: `observe` and `predict`
- Compiled programs automatically emit posterior samples of predicted quantities, conditioned on the observed data
- Freely incorporate existing libraries and C code, including black box simulators

Implementation

- The POSIX `fork` function causes a process to clone itself, creating a new process with an identical copy of the program execution state
 - Lazy copy-on-write procedure for duplicating memory contents: efficient
- Using `fork` we can branch and explore many downstream program execution paths
- Massively parallel: each downstream path is explored by an independent OS process
- Global synchronization is handled via shared memory segments

Usage

Simple example: iid draws from a Gaussian distribution with unknown mean.

$$\mu \sim \text{Normal}(1, 5)$$
$$y_1, y_2 \stackrel{iid}{\sim} \text{Normal}(\mu, 2)$$

Sample from prior:

```
#include "probabilistic.h"
int main(int argc, char **argv) {
    double var = 2;
    double mu = normal_rng(1, 5);
    printf("%f\n", normal_rng(mu, var));
    printf("%f\n", normal_rng(mu, var));
    printf("%f\n", mu);
    return 0;
}
```

Sample from posterior:

```
#include "probabilistic.h"
int main(int argc, char **argv) {
    double var = 2;
    double mu = normal_rng(1, 5);
    observe(normal_rng(8, mu, var));
    observe(normal_rng(8, mu, var));
    predict(&mu, &var, mu);
    return 0;
}
```

Another example: a hidden Markov model, with Gaussian emission distributions.

$$z_0 \sim \text{Discrete}([1/K, \dots, 1/K]) \quad z_n | z_{n-1} \sim \text{Discrete}(T_{z_{n-1}}) \quad y_n | z_n \sim \text{Normal}(\mu_{z_n}, \sigma^2)$$

Sample from prior:

```
#include "probabilistic.h"
#define K 3
#define N 11
/* Markov transition matrix */
static double T[K][K] = { { 0.1, 0.5, 0.4 },
                          { 0.2, 0.2, 0.6 },
                          { 0.15, 0.15, 0.7 } };
/* Prior distribution on initial state */
static double initial_state[K] = { 1.0/3, 1.0/3, 1.0/3 };
/* Per-state mean of Gaussian emission distribution */
static double state_mean[K] = { -1, 1, 0 };
/* Generative program for a HMM */
int main(int argc, char **argv) {
    int states[N];
    for (int n=0; n<N; n++) {
        states[n] = (n==0) ? discrete_rng(initial_state, K) :
                        discrete_rng(T[states[n-1]], K);
        if (n > 0) {
            printf("%d\n", states[n]);
            normal_rng(state_mean[states[n]], 1);
            printf("state=%d, %f, %f, states[n]\n",
                states[n], mu, var, states[n]);
        }
    }
    return 0;
}
```

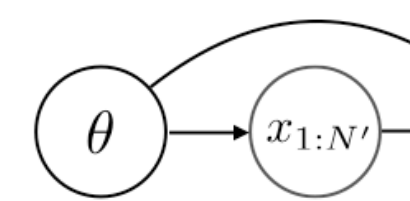
Sample from posterior:

```
#include "probabilistic.h"
#define K 3
#define N 11
/* Markov transition matrix */
static double T[K][K] = { { 0.1, 0.5, 0.4 },
                          { 0.2, 0.2, 0.6 },
                          { 0.15, 0.15, 0.7 } };
/* Observed data */
static double data[N] = { NAN, .9, .8, .1, 0, -.025,
                        -.5, -.2, -.1, 0, 0.13 };
/* Prior distribution on initial state */
static double initial_state[K] = { 1.0/3, 1.0/3, 1.0/3 };
/* Per-state mean of Gaussian emission distribution */
static double state_mean[K] = { -1, 1, 0 };
/* Generative program for a HMM */
int main(int argc, char **argv) {
    int states[N];
    for (int n=0; n<N; n++) {
        states[n] = (n==0) ? discrete_rng(initial_state, K) :
                        discrete_rng(T[states[n-1]], K);
        if (n > 0) {
            observe(normal_rng(data[n],
                               state_mean[states[n]], 1));
            predict(&states[n], &mu, &var, states[n]);
        }
    }
    return 0;
}
```

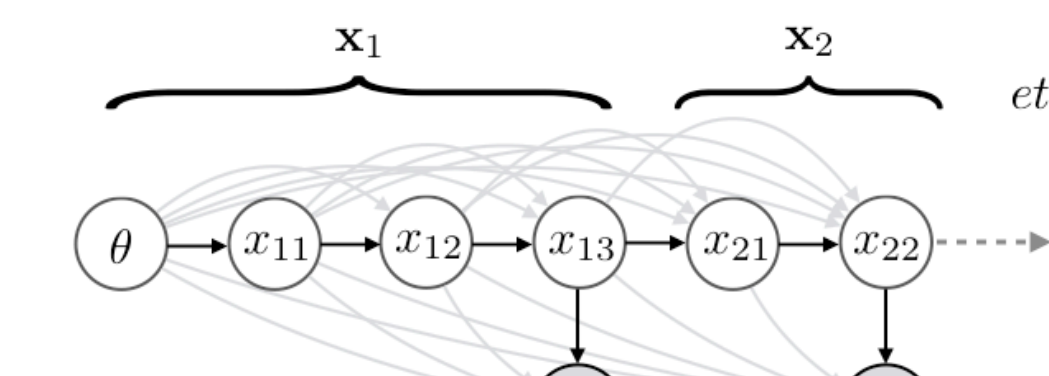
Inference

Define a probability distribution over program execution traces

- Enumerate all N observe statements, and the associated data points y_1, \dots, y_N
- Enumerate all N' random choices made during the execution of the program $x_1, \dots, x_{N'}$

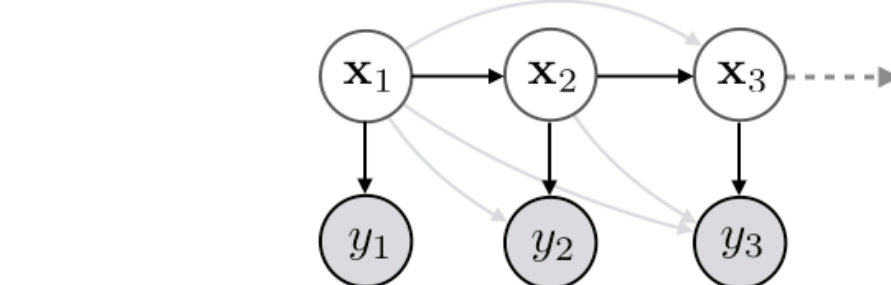


- Define a partition of the random choices such that each x_i contains all random choices made between observations y_{n-1} and y_n



- Run sequential Monte Carlo or particle MCMC algorithms over the resulting sequence of program execution states

$$p(y_{1:N}, x_{1:N}) = \prod_{n=1}^N g(y_n | x_{1:n}) f(x_n | x_{1:n-1})$$



- Note we only need to be able to simulate from $f(x_n | x_{1:n-1})$ and evaluate $g(y_n | x_{1:n})$

Algorithm Details

Algorithm 1 Parallel SMC program execution
Assume: N observations, L particles
launch L copies of the program
for $n = 1 \dots N$ do
wait until all L reach `observe` y_n (parallel)
update unnormalized weights w_n^i (serial)
if $ESS < \tau$ then (serial)
sample number of offspring $O_n^{i,j}$ (serial)
set weight $w_n^{i,j} = 1$ (serial)
for $j = 1 \dots L$ do (parallel)
fork or exit
else for
set all number of offspring $O_n^i = 1$ (serial)
end if
continue program execution (parallel)
end for
wait until L program traces terminate (barrier)
predict from L samples from $p(x_{1:N} | y_{1:N})$ (serial)

Algorithm 2 Parallel Particle Gibbs program execution
Assume: M iterations, N observations, L particles
for $m = 1 \dots M$ do
if $C \leftarrow L$ if $m = 1$, otherwise $L - 1$ (parallel)
launch S' copies of the program (barrier)
wait until all L' reach an observe (serial)
compute weights for all particles (serial)
if $m > 1$ then (serial)
signal num offspring to retained trace (serial)
end if
for $\ell = 1 \dots L'$ do (parallel)
spawn retain / branch process [Algo. 3] (barrier)
end for
wait until L particles finish branching (barrier)
continue program execution (serial)
end for
wait until L program traces terminate (barrier)
predict from L samples from $p(x_{1:N} | y_{1:N})$ (serial)
select and signal particles to retain (barrier)
wait until N processes are ready to branch (barrier)
continue to next iteration (parallel)
end for

Algorithm 3 Retain and Branch inner loop
Assume: input initial $C > 0$ children to spawn
is retained \leftarrow false
while $\tau == 0$ do
if $C = 0$ and not is retained then
discard this execution trace, exit
else ($C \geq 0$)
spawn N new children
end if
wait for signal which resets is retained
if is retained then
wait for signal which resets C
else
discard this execution trace, exit
end if
end while

References

- [1] Doucet, Arnaud, De Freitas, Nando, Gordon, Neil, et al. *Sequential Monte Carlo methods in practice*. Springer New York, 2001.
- [2] Andrieu, Christophe, Doucet, Arnaud, and Holenstein, Roman. Particle Markov chain Monte Carlo methods. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 72(3):269–342, 2010.
- [3] Goodman, Noah D., Mansinghka, Vikash K., Roy, Daniel M., Bonawitz, Keith, and Tenenbaum, Joshua B. Church: A language for generative models. In *Proceedings of the Twenty-Fourth Conference on Uncertainty in Artificial Intelligence (UAI2008)*, pp. 220–229, 2008.
- [4] Wingate, David, Stuhlmüller, Andreas, and Goodman, Noah D. Lightweight implementations of probabilistic programming languages via transformational compilation. In *Proceedings of the 14th international conference on Artificial Intelligence and Statistics*, pp. 131, 2011.
- [5] Wood, Frank, van de Meent, Jan Willem, and Mansinghka, Vikash. A new approach to probabilistic programming inference. In *Proceedings of the 17th International conference on Artificial Intelligence and Statistics*, 2014.

Capabilities sought and types of research supported

Anatomy

- Sought : data and bottom-up feature extraction
- Support : top-down model-based regularization

Function

- Sought : stochastic circuit simulators
- Support : automatic inversion

Artificial Intelligence

- Sought : ?
- Support : programs that write programs, etc.

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