

# Modular Probabilistic Models via Algebraic Effects

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Probabilistic programming languages (PPLs) allow programmers to construct statistical models and then simulate data or perform inference over them. Many PPLs restrict models to a particular instance of simulation or inference, limiting their reusability. In other PPLs, models are not readily composable. Using Haskell as the host language, we present an embedded domain specific language based on algebraic effects, where probabilistic models are modular, first-class, and reusable for both simulation and inference. We also demonstrate how simulation and inference can be expressed naturally as composable program transformations using algebraic effect handlers.

Additional Key Words and Phrases: Probabilistic programming, algebraic effects, modularity

## 1 INTRODUCTION

In statistics, a probabilistic model captures a real-world phenomenon as a set of relationships between random variables: the model's parameters, inputs, and outputs. By integrating such notions into general-purpose languages, probabilistic programming languages (PPLs) allow programmers to build and execute probabilistic models. For example, consider a simple linear regression model that assumes a linear relationship between input variables  $x$  and output variables  $y$ ; this can be represented using the standard mathematical notation shown on the left below. Using the language presented in this paper, the right-hand side shows how one could express the same model as a functional program.

$\mu \sim \text{Normal}(0, 3)$	<code>linRegr <math>x</math> = do</code>
$c \sim \text{Normal}(0, 2)$	<code><math>\mu \leftarrow \text{normal } 0 \ 3 \ \# \mu</math></code>
$\sigma \sim \text{Uniform}(1, 3)$	<code><math>c \leftarrow \text{normal } 0 \ 2 \ \# c</math></code>
$y \sim \text{Normal}(\mu * x + c, \sigma)$	<code><math>\sigma \leftarrow \text{uniform } 1 \ 3 \ \# \sigma</math></code>
	<code><math>y \leftarrow \text{normal } (\mu * x + c, \sigma) \ \# y</math></code>
	<code>return <math>y</math></code>

Both representations take an input  $x$  and specify the distributions which generate the model parameters  $\mu$ ,  $c$ , and  $\sigma$ ; the output  $y$  is then generated from the normal distribution using mean  $\mu * x + c$  and standard deviation  $\sigma$ . In the program representation, each primitive distribution is associated with a corresponding “observable variable”, indicated by the  $\#$  syntax; this is an optional argument, and its purpose will become clear shortly.

Given a probabilistic model, the programmer or data scientist will typically want to use it in at least two different ways. *Simulation* involves providing fixed values for the model parameters and inputs, to generate the resulting model outputs. Conversely, *inference* generally entails providing observed values for the model outputs and inputs, in an attempt to learn the model parameters.

For example, we might simulate from `linRegr` in our language as follows:

```
let xs = [ 0 .. 100 ]
    env = ( $\# \mu := [3]$ ) • ( $\# c := [0]$ ) • ( $\# \sigma := [1]$ ) • ( $\# y := []$ ) • nil
in map (simulate linRegr env) xs
```

First we declare a list of model inputs  $xs$  from 0 to 100. Then we define a “model environment” `env` which assigns values 3, 0, and 1 to observable variables  $\# \mu$ ,  $\# c$ , and  $\# \sigma$ . This expresses our intention to *observe* parameters  $\mu$ ,  $c$ , and  $\sigma$  — that is, to provide external data 3 as the value of random variable  $\mu$  whilst conditioning on the likelihood that  $\mu = 3$ , and similarly for  $c$  and  $\sigma$ . On the other hand no values are specified for  $\# y$  in `env`, expressing our intent to *sample* the model output  $y$  — that is, to draw a value from its probability distribution. We then use library function `simulate` to simulate a

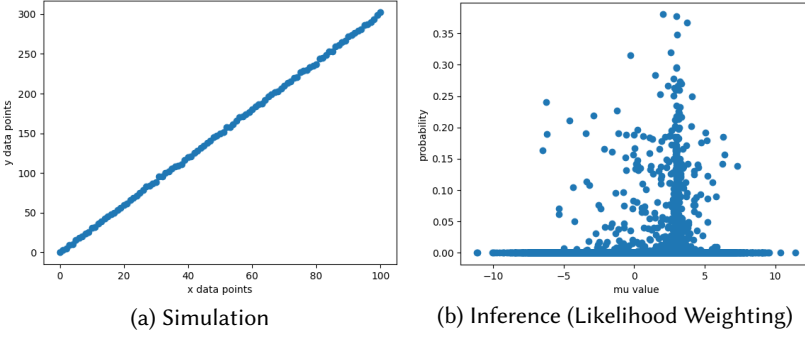


Fig. 1. Visualising Linear Regression

single output from the model for each data point in  $xs$  using the specified environment, producing the result visualised in Fig. 1a.

Alternatively, we can perform inference on `linRegr`, for example using the Likelihood Weighting algorithm [van de Meent et al. 2018], as follows:

```
let xs = [ 0 .. 100 ]
xys = [(x, env) | x ← xs, let env = (#μ := []) • (#c := []) • (#σ := []) • (#y := [3 * x]) • nil]
in map (lw 200 linRegr) xys
```

Here we define  $xys$  to pair each model input  $x$  with a model environment  $env$  that assigns the value  $3 * x$  to  $\#y$  but nothing to  $\#μ$ ,  $\#c$ , and  $\#σ$ . This expresses our intention to *observe*  $y$  but *sample*  $μ$ ,  $c$  and  $σ$ . We then use library function `lw` to perform 200 iterations of Likelihood Weighting for each pair of model input and environment, producing a trace of weighted parameters  $μ$ ,  $c$ , and  $σ$  whose distributions express the most likely parameter values to give rise to  $y$ . Fig. 1b visualises the likelihoods of samples for  $μ$ , where values around  $μ = 3$  clearly accumulate higher probabilities.

We refer to a model that can be used for both simulation and inference — where random variables can be switched between sample and observe modes without altering the model itself — as a *multimodal model*. While multimodal models have a clear benefit, letting the same model be interacted with for a variety of applications, few existing PPLs support them. Most frameworks, such as `MonadBayes` [Ścibior et al. 2018] and `Anglican` [Tolpin et al. 2016], instead require programmers to express models in terms of explicit sample and observe operations, which considerably limits their reusability. If the user wishes to interact with the “same” model in a new way, they have little choice but to reimplement it with a different configuration of sample and observe operations.

Indeed, the number of possible model interpretations extends far beyond the two general scenarios of simulation and inference, potentially including *any* combination of sample and observe operations that can be instantiated for a model’s random variables. Depending on available data and uncertainty about the model, it is common to explore the model’s output space by partially providing model parameters and randomly sampling the rest [Kline and Tamer 2016], or to alternate between which observable variables are being conditioned on [Moon 1996]. Ideally, all of these possible scenarios would be expressible with a single multimodal model definition, avoiding the need to define and separately maintain a different version of the model for each use case.

While some PPLs *do* support multimodal models, it is usually difficult or impossible to reuse existing models when creating new ones. `Stan` [Carpenter et al. 2017] and `WinBUGS` [Lunn et al. 2000] provide a bespoke language construct for models with its own distinctive semantics, but as well as lacking high-level programming features beyond those essential to model specification, model definitions are unable to reuse other model definitions. Languages like `Turing` [Ge et al. 2018] and `Gen` [Cusumano-Towner et al. 2019] take a different approach, supporting multimodal

models as macros that are compiled into functions; although they provide some support for compositionality, neither supports models as first-class values. These modularity limitations are especially significant for hierarchical modelling, where the goal is to explicitly define a composite model with independently defined sub-models [Gelman and Hill 2006].

In this paper, we present *Wasabaye*: a deeply embedded PPL in Haskell where probabilistic models are modular, first-class, and multimodal. Our solution uses algebraic effects [Plotkin and Power 2003] and handlers [Plotkin and Pretnar 2013], allowing models to be captured as syntax, and their semantics deferred to a choice of “model environment”. By embedding into a functional language, models can then naturally exist as (first-class) functions and leverage all the abstractions and features of the host [Elliott et al. 2003; Gibbons and Wu 2014].

Our approach uses two key type abstractions: polymorphic sums for expressing effect signatures, and extensible records for model environments. Both of these can in fact be subsumed by row polymorphism [Leijen 2005], and so any language with support for polymorphic rows, such as PureScript [Freeman 2017], OCaml [Leroy et al. 2020] or Links [Hillerström and Lindley 2018], or with a type system powerful enough to express something similar (such as Haskell), should be capable of capturing *Wasabaye*’s main features. The other type-level devices we use are ergonomic choices specific to Haskell, and are inessential to the main goals.

We begin by giving the necessary background and language overview in §2. Our contributions are then as follows:

- We demonstrate the features of our language via a realistic case study with real-world applications: the spread of disease during an epidemic (§3).
- We present an embedding technique that is novel in using algebraic effects to represent probabilistic models (§4), demonstrated with Haskell as the host language. To the best of our knowledge, ours is the first PPL to support models that are both multimodal and first-class.
- We provide a modular, type-safe mechanism for associating observed data to the random variables of a model, determining whether probabilistic operations should be interpreted as sample or observe (§5). The same mechanism is used to trace samples for plotting or debugging.
- We present a new approach to the compositional implementation of simulation and inference, using effect handlers to perform modular program transformations on models (§6). We illustrate the approach using Likelihood Weighting [van de Meent et al. 2018] and Metropolis Hastings [Wingate et al. 2011].
- We evaluate our language empirically, considering performance against two state-of-the-art PPLs, and language features supported across a range of modern PPLs (§7).

We discuss related work in more detail in §8. However, this is not the first time PPLs have been explored with Haskell as a host language. Erwig and Kollmansberger [2006] implement a probability monad for representing distributions in functional languages; Narayanan et al. [2016] use a tagless-final embedding [Kiselyov 2010] to encode inference algorithms as type class instances; Ścibior et al. [2018] use monad transformers [Liang et al. 1995] to demonstrate inference as effect composition. Our approach builds on the techniques offered by algebraic effects and extensible data.

The high-level notion of using interpreters to execute the effects of probabilistic models is an established technique in PPLs, and is similar in spirit to algebraic effects. Many PPLs accomplish this through context managers, coroutines, and continuation-passing style transformations [Bingham et al. 2019; Goodman and Stuhlmüller 2014; Tolpin et al. 2016]. However, these approaches fail to delineate between syntax and semantics, preventing models from being interpreted in a fully multimodal fashion. Moreover, the effect-interpreting mechanisms typically operate in weakly-typed, imperative settings, where effects are not associated with types and can occur unrestrictedly in a program. Algebraic effects have the potential to bring a type-safe, compositional discipline to

probabilistic programming. There is, however, little existing work in this area, and the topic has primarily remained a point of discussion [Moore and Gorinova 2018; Ścibior and Kammar 2015]. We present a novel design and implementation at the intersection of PPLs and algebraic effects.

In addition to the examples in this paper, our embedding has been tested with a range of models implemented in other PPLs, as well as well-known models such as those designed by Gelman and Hill [2006]; the full source code is freely available online.<sup>1</sup>

## 2 BACKGROUND AND LANGUAGE OVERVIEW

A probabilistic model, expressed using the  $\sim$  notation introduced in §1, describes how a set of random variables are distributed relative to some fixed input. If the model does not condition against any external data, the distribution it describes is the so-called *joint probability distribution*, giving the probabilities of all possible values that its random variables can assume. For example, the linear regression model in §1 describes the distribution  $\mathbb{P}(y, \mu, c, \sigma; x)$  – namely, the joint distribution over random variables  $y, \mu, c$ , and  $\sigma$  given fixed input  $x$  as a non-random parameter.

In real-world applications, we typically have known values for only a subset of these random variables, and are interested in how the other variables are distributed with respect to those known values. Consider providing known data  $\hat{y}$  for random variable  $y$  in linear regression; we say that we *condition on  $y$  having observed  $\hat{y}$* . Using the well-known chain rule for two random variables:

$$\mathbb{P}(X, Y) = \mathbb{P}(X \mid Y) \cdot \mathbb{P}(Y)$$

we can derive the resulting distribution as the product  $\mathbb{P}(\mu, c, \sigma \mid y = \hat{y}; x) \cdot \mathbb{P}(y = \hat{y})$ . The first component, called the *conditional distribution*, describes the probabilities of values for each random variable  $\mu, c, \sigma$  given  $y = \hat{y}$  and some input  $x$ ; the second component, called the *prior distribution*, gives the probability that  $y$  has value  $\hat{y}$ . Providing observed data to a probabilistic model can therefore be seen as specialising its joint distribution to some product of a conditional that is favourable to modelling and an associated prior.

### 2.1 Multimodal Models

The chain rule is a powerful tool that allows us to describe a jointly occurring set of events in terms of the variables we will provide data for, and then compute other variables of interest with respect to this; and there are of course as many ways to decompose a joint distribution as there are combinations of variables that can be conditioned on. Since statisticians often have a clear understanding of the variables they wish to learn and those they wish to condition against, models are in practice often specialised to specific conditional distributions (through the chain rule) and expressed as low-level algorithms that explicitly perform sampling and conditioning, such as in Ding et al. [2019]; Polson et al. [2013].

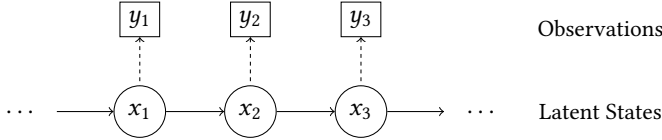
Most PPLs, such as WebPPL [Goodman and Stuhlmüller 2014] and Anglican [Tolpin et al. 2016], are then designed to support the direct translation of these low-level model specifications from paper to program via the operations `sample` and `observe`. These languages are useful for creating model instances tailored to specific situations, but the resulting models are not easy to experiment with. Tasks which should be straightforward, such as exploring random variable behaviours by isolating which ones are sampled from [Idreos et al. 2015] or selectively optimising model parameters [Yekutieli 2012], require alternative specialisations to be created by hand.

**2.1.1 Multimodal models via model environments.** With multimodal PPLs, the programmer specifies a single model which can be used to generate multiple specialisations, representing specific conditional distributions. Such languages require a mechanism for specifying observed data to random variables, determining whether they are to be sampled or observed. For example, Turing.jl

<sup>1</sup><https://github.com/min-nguyen/wasabay>

lets users choose whether to provide observed values as arguments when invoking a model, with omitting an argument defaulting to sampling [Ge et al. 2018]; in Pyro, users specify mappings between random variables and observed data via *context managers* that later constrain the values of runtime sampling operations [Bingham et al. 2019]. However, these solutions are dynamically typed with no guarantee that the named variables exist or are provided values of the correct type.

Our language supports multimodal models through a novel notion of *model environment*, which we explain in the context of a Hidden Markov Model (HMM) [Rabiner and Juang 1986]:



The idea of a HMM is that we have a series of latent states  $x_i$  which are related in some way to observations  $y_i$ . The HMM is then defined by two sub-models: a transition model ( $\rightarrow$ ) that determines how latent states  $x_i$  are transitioned between, and an observation model ( $\uparrow$ ) that determines how  $x_i$  is projected to an observation  $y_i$ . The objective is to learn about  $x_i$  given  $y_i$ .

A simple HMM expressed in typical statistical pseudocode is shown in Fig. 2a; we describe its corresponding implementation in our language in Fig. 2b:

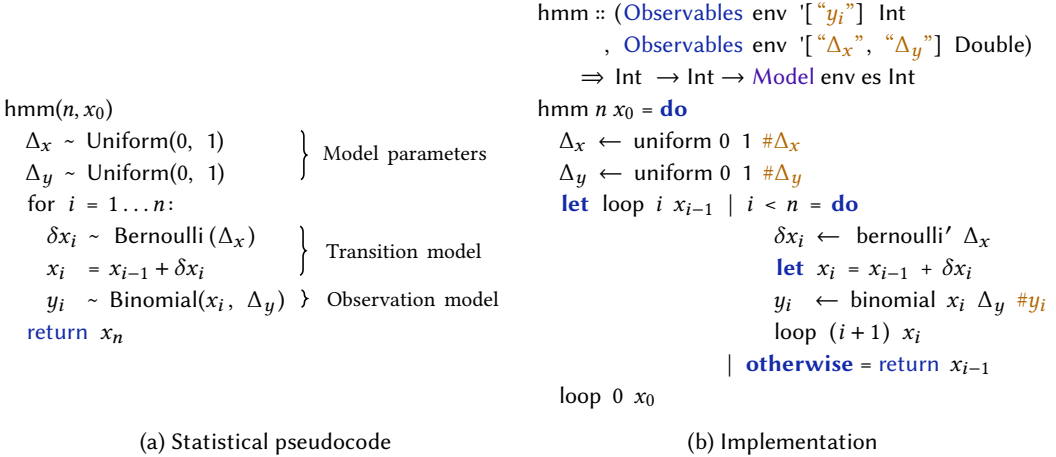


Fig. 2. Hidden Markov Model

The type of `hmm` says it is a function that takes two `Int`s as input and returns a `Model env es Int`, where `env` is the model environment, `es` is the effects which the model can invoke (detailed later in §4), and `Int` is the type of values the model generates. The constraint `Observables` states that `#yi :: Int`, `#Δx :: Double`, and `#Δy :: Double`, are *observable variables* in the model environment `env` which may be conditioned on later when the model is used.

The function `hmm` takes the HMM length  $n$  and initial latent state  $x_0$  as inputs, and specifies the transition and observation parameters  $\Delta_x$  and  $\Delta_y$  to be distributed uniformly. It then iterates over the  $n$  nodes, applying the transition and observation models at each step. The transition model computes latent state  $x_i$  from state  $x_{i-1}$  by adding a value  $\delta x_i$  generated from a Bernoulli distribution `bernoulli' Δx`. The observation model generates observation  $y_i$  from  $x_i$  via the distribution `binomial xi Δy`. The final latent state is returned at the end.

The hash syntax `#` is how the programmer associates variables in the `Observables` env constraint with specific primitive distributions. They do this to indicate that they may later be interested in providing observed values for these variables to condition on. When they execute a model, they must provide a concrete environment of type `env`, and the presence or absence of observed values in that environment (for a given variable  $x$ ) will determine whether the distribution tagged with `# $x$`  is to be interpreted as observe or sample. The distribution `bernoulli'`  $\Delta_x$  has no observable variable, indicating that it is not possible to condition on  $\delta x_i$ ; this makes sense because values of  $\delta x_i$  are latent, and so it is unlikely that we would ever want to provide data for them. (Primitive distributions like `bernoulli` come in primed variants that are always interpreted as sample.)

A model can then be interpreted as any of its conditioned forms by specifying an appropriate model environment. In our example, the HMM in its unspecialised form represents the joint distribution over its latent states  $x_i$ , observations  $y_i$ , and parameters  $\Delta_x, \Delta_y$  (given fixed input  $x_0$  as the first latent state):

$$\mathbb{P}(x_1 \dots x_n, y_1 \dots y_n, \Delta_x, \Delta_y; x_0)$$

and we can then simulate the HMM (with length  $n = 10$  and initial state  $x_0 = 0$ ) by providing values for `# $\Delta_x$`  and `# $\Delta_y$`  in an environment `env`:

```
let x0 = 0; n = 10;
    env = (# $\Delta_x$  := [0.5]) • (# $\Delta_y$  := [0.8]) • (# $y_i$  := []) • nil
in simulate (hmm n) env x0
```

This indicates that we want to observe 0.5 and 0.8 for  $\Delta_x$  and  $\Delta_y$ , and sample for each occurrence of  $y_i$  (because we provided no values for `# $y_i$` ); there are multiple occurrences of  $y_i$  at runtime, one for each  $i \in \{1 \dots n\}$ , thanks to the iterative structure of the HMM. By the chain rule, the probability density this expresses is:

$$\mathbb{P}(x_1 \dots x_{10}, y_1 \dots y_{10} \mid \Delta_x = 0.5, \Delta_y = 0.8; x_0 = 0) \cdot \mathbb{P}(\Delta_x = 0.5) \cdot \mathbb{P}(\Delta_y = 0.8)$$

In the case of inference, on the other hand, we provide an observation for each  $y_i$  and try to learn  $\Delta_x$  and  $\Delta_y$ . This is why (as the reader may already have noticed) a model environment provides a *list* of values for each observable variable, allowing for the situation where the observable variable has multiple dynamic occurrences. In this case we provide 10 observations, one for each  $i \in \{1 \dots n\}$ :

```
let x0 = 0; n = 10;
    env = (# $\Delta_x$  := []) • (# $\Delta_y$  := []) • (# $y_i$  := [0, 1, 1, 3, 4, 5, 5, 5, 6, 5]) • nil
in lw 100 (hmm n) (x0, env)
```

At runtime, the values associated with `# $y_i$`  in the model environment are used to condition against the occurrences of `# $y_i$`  that arise during execution, in the order in which they arise. By the chain rule, the probability density expressed by instantiating the model with this environment is:

$$\mathbb{P}(\Delta_x, \Delta_y, x_1 \dots x_{10} \mid y_1 = 0 \dots y_{10} = 5; x_0 = 0) \cdot \mathbb{P}(y_1 = 0) \dots \mathbb{P}(y_{10} = 5)$$

Although the type system ensures that model environments map observable variables to values of an appropriate type, it does not constrain the number of values that are provided. Should observed values run out for a particular variable, any remaining runtime occurrences of the variable will default to sample; any surplus of values is ignored. While this flexibility could certainly obscure programming errors, other PPLs (such as Turing.jl [Ge et al. 2018]) take a similar approach, and we also note that the correctness of inference is unaffected. We consider alternative designs in §8.2.



## 2.2 Modular, First-Class Models

Fig. 2a used a single procedure, written in statistical pseudocode, to express a Hidden Markov Model. Such notations are understood by most mathematicians and are widely used in statistical journals. Even when the model is complex, a monolithic style of presentation prevails, where models are defined from scratch each time rather than built out of reusable components. The design of PPLs such as Stan [Carpenter et al. 2017], PyMC3 [Salvatier et al. 2016] and Bugs [Lunn et al. 2000] reflect these non-modular conventions.

Programmers, on the other hand, recognise the importance of modularity to maintainability and reusability: they expect to be able to decompose models into meaningful parts. Fig. 3a shows how the programmer may imagine the same HMM as a composition of parts; our language can then support this treatment of models in Fig. 3b:

transModel( $\Delta_x, x_{i-1}$ )	transModel :: Double $\rightarrow$ Int $\rightarrow$ Model env es Int
$\delta x_i \sim \text{Bernoulli}(\Delta_x)$	transModel $\Delta_x \ x_{i-1} = \mathbf{do}$
return $x_{i-1} + \delta x_i$	$\delta x_i \leftarrow \text{bernoulli}' \ \Delta_x$
	return $x_{i-1} + \delta x_i$
obsModel( $\Delta_y, x_i$ )	obsModel :: (Observables env '[ $"y_i"$ ] Int)
$y_i \sim \text{Binomial}(x_i, \Delta_y)$	$\Rightarrow$ Double $\rightarrow$ Int $\rightarrow$ Model env es Int
return $y_i$	obsModel $\Delta_y \ x_i = \mathbf{do}$
	$y_i \leftarrow \text{binomial} \ x_i \ \Delta_y \ \#y_i$
	return $y_i$
hmmNode( $\Delta_x, \Delta_y, x_{i-1}$ )	hmmNode :: (Observables env '[ $"y_i"$ ] Int)
$x_i \sim \text{transModel}(\Delta_x, x_{i-1})$	$\Rightarrow$ Double $\rightarrow$ Double $\rightarrow$ Int $\rightarrow$ Model env es Int
obsModel( $\Delta_y, x_i$ )	hmmNode $\Delta_x \ \Delta_y \ x_{i-1} = \mathbf{do}$
return $x_i$	$x_i \leftarrow \text{transModel} \ \Delta_x \ x_{i-1}$
	obsModel $\Delta_y \ x_i$
	return $x_i$
hmm( $n, x_0$ )	hmm :: (Observables env '[ $"y_i"$ ] Int, Observables env '[ $"\Delta_x"$ , $"\Delta_y"$ ] Double)
$\Delta_x \sim \text{Uniform}(0, 1)$	$\Rightarrow$ Int $\rightarrow$ Int $\rightarrow$ Model env es Int
$\Delta_y \sim \text{Uniform}(0, 1)$	hmm $n \ x_0 = \mathbf{do}$
for $i = 1 \dots n$ ;	$\Delta_x \leftarrow \text{uniform} \ 0 \ 1 \ \#\Delta_x$
$x_i \sim \text{hmmNode}(\Delta_x, \Delta_y, x_{i-1})$	$\Delta_y \leftarrow \text{uniform} \ 0 \ 1 \ \#\Delta_y$
return $x_n$	foldl ( $\Rightarrow$ ) return (replicate $n \ (\text{hmmNode} \ \Delta_x \ \Delta_y)) \ x_0$
(a) Statistical Pseudocode	(b) Program Representation

Fig. 3. A Modular Hidden Markov Model

In Fig. 3b we define the transition and observation distributions as separate models transModel and obsModel. These are composed by hmmNode to define the behaviour of a single node, which is in turn used by hmm to create a chain of nodes of length  $n$ , using replicate and a fold of Kleisli composition ( $\Rightarrow$ ) to propagate each node's output to the next one in the chain.

( $\Rightarrow$ ) :: (a  $\rightarrow$  Model env es b)  $\rightarrow$  (b  $\rightarrow$  Model env es c)  $\rightarrow$  (a  $\rightarrow$  Model env es c)

The observable variables of hmm are now inherited from its sub-models: transModel has none (and so lacks an Observables constraint entirely), whereas obsModel declares  $\#y_i$  as its only observable variable.

As well as improving reusability, compositionality also allows programmers to organise models around the structure of the problem domain. For example the functions in Fig. 3b correspond in a straightforward way to the abstract components of a HMM: `transModel` makes it obvious that latent state  $x_i$  depends only on the previous state  $x_{i-1}$  (and  $\Delta_x$ ), and `obsModel` that observation  $y_i$  depends only on the state  $x_i$  that produced it (and  $\Delta_y$ ).

### 3 MODULAR, MULTIMODAL MODELS: A CASE STUDY

Before we turn to the details of our embedding approach in §4, we present a case study demonstrating our support for modular, first-class, multimodal models. §3.1 introduces our running example, the SIR (Susceptible-Infected-Recovered) model for the spread of disease [Liang and Li 2021]. §3.2 uses the SIR model to show how our language supports higher-order models which can be easily extended and adapted. §3.3 shows how a multimodal model can be used for both simulation and inference in the same application to facilitate Bayesian bootstrapping.

#### 3.1 The SIR Model

The SIR model predicts the spread of disease in a fixed population of size  $n$  partitioned into three groups:  $s$  for *susceptible to infection*,  $i$  for *infected*, or  $r$  for *recovered* (where  $s + i + r = n$ ). The model tracks how  $s$ ,  $i$ , and  $r$  vary over discrete time  $t$  measured in days. Because testing is both incomplete and unreliable, the true *sir* values for the population cannot be directly observed; however, we can observe the number of reported infections  $\xi$ . This problem is thus a good fit for a Hidden Markov Model (§2), where the *sir* values play the role of latent states of type `Popl`, and  $\xi$  as the observations of type `Reported`:

```
data Popl      = Popl { s :: Int, i :: Int, r :: Int }
type Reported = Int
```

We now show how our language can be used to implement the SIR model as a modular HMM, starting with the transition and observation models.

*SIR transition model.* The transition model describes how the *sir* values change over a single day; we model two specific dynamics. First, susceptible individuals  $s$  transition to infected  $i$  at a rate determined by the values of  $s$  and  $i$  and the contact rate  $\beta$  between the two groups. We use a binomial distribution to model each person in  $s$  having a  $1 - e^{-\beta i/n}$  probability of becoming infected, and update  $s$  and  $i$  accordingly:

```
transsi :: Double → Popl → Model env es Popl
transsi β (Popl s i r) = do
  let n = s + i + r
  δsi ← binomial' s (1 - exp ((-β * i) / n))
  return (Popl (s - δsi) (i + δsi) r)
```

Second, infected individuals  $i$  transition to the recovered group  $r$ , where a fixed fraction  $\gamma$  of people will recover in a given day. Again we use a binomial to model each person in  $i$  as having a  $1 - e^{-\gamma}$  probability of recovering, and use this to update  $i$  and  $r$ :

```
transir :: Double → Popl → Model env es Popl
transir γ (Popl s i r) = do
  δir ← binomial' i (1 - exp (-γ))
  return (Popl s (i - δir) (r + δir))
```



The overall transition model  $\text{trans}_{sir}$  is simply the sequential composition of  $\text{trans}_{si}$  and  $\text{trans}_{ir}$ . Given  $\beta$  and  $\gamma$  (aggregated into the type `TransParams`),  $\text{trans}_{sir}$  computes the changes from  $s$  to  $i$  and then  $i$  to  $r$  to yield the updated  $sir$  population over a single day:

```
data TransParams = TransParams {  $\beta$  :: Double,  $\gamma$  :: Double }

transsir :: TransParams → Popl → Model env es Popl
transsir (TransParams  $\beta$   $\gamma$ ) = transsi  $\gamma$  >=> transir  $\beta$ 
```

*SIR observation model.* For the observation model, we assume that the reported infections  $\xi$  depends only on the number of infected individuals  $i$ , of which a fixed fraction  $\rho$  will be reported. We use the Poisson distribution to model reports occurring with a mean rate of  $\rho * i$ :

```
type ObsParams = Double

obssir :: Observables env '[ $\xi$ ] Int ⇒ ObsParams → Popl → Model env es Reported
obssir  $\rho$  (Popl _  $i$  _) = do
   $\xi$  ← poisson ( $\rho * i$ ) # $\xi$ 
  return  $\xi$ 
```

Since we intend this as the observation model, we declare observable variable  $\# \xi :: \text{Int}$  in the `Observables` constraint and attach it to the Poisson distribution so we can condition on it later. (Binding the local name  $\xi$  and immediately returning it is technically redundant but emphasises the connection to the  $\sim$  notation used by statisticians.)

*HMM for the SIR model.* Now the transition and observation models can be combined into a HMM. We build on the modular design in Fig. 3, but go a step further by defining a HMM as a higher-order model that is parameterised by sub-models of type `TransModel` and `ObsModel`:

```
type TransModel env es ps lat = ps → lat → Model env es lat
type ObsModel env es ps lat obs = ps → lat → Model env es obs
```

Here `ps` represents the types of the model parameters, and `lat` and `obs` are the types of latent states and observations. The higher-order HMM is then defined as:

```
hmm :: Model env es ps1 → Model env es ps2
      → TransModel env es ps1 lat → ObsModel env es ps2 lat obs
      → Int → lat → Model env es lat

hmm transPrior obsPrior trans obs  $n$   $x_0$  = do
   $\theta$  ← transPrior
   $\phi$  ← obsPrior
  hmmNode  $x_{i-1}$  = do  $x_i$  ← trans  $\theta$   $x_{i-1}$ 
                   $y_i$  ← obs  $\phi$   $x_i$ 
                  return  $x_i$ 
  foldl (>=>) return (replicate  $n$  hmmNode)  $x_0$ 
```

The input models `transPrior` and `obsPrior` are first used to generate model parameters  $\theta$  and  $\phi$ , and `trans` and `obs` are arbitrary transition and observation models parameterised by  $\theta$  and  $\phi$  respectively. The last line creates a HMM of length  $n$  with initial state  $x_0$ .

Our SIR transition and observation parameters will be provided by models `transPriorsir` and `obsPriorsir` below, using primitive distributions `gamma` and `beta`; their observable variables  $\# \beta$ ,  $\# \gamma$ , and  $\# \rho$  will let us condition on those parameters later:

```

transPriorsir :: (Observables env '[ $\beta$ ,  $\gamma$ ] Double) ⇒ Model env es TransParams
transPriorsir = do
   $\beta$  ← gamma 2 1 # $\beta$ 
   $\gamma$  ← gamma 1 (1/8) # $\gamma$ 
  return (TransParams  $\beta$   $\gamma$ )

obsPriorsir :: (Observables env '[ $\rho$ ] Double) ⇒ Model env es ObsParams
obsPriorsir = do
   $\rho$  ← beta 2 7 # $\rho$ 
  return  $\rho$ 

```

We can now define the complete SIR model. From an initial population *sir* of susceptible, infected, and recovered individuals, *hmm<sub>sir</sub>* models the change in *sir* over *n* days given reported infections  $\xi$ :

```

hmmsir :: (Observables env '[ $\xi$ ] Int, Observables env '[ $\beta$ ,  $\gamma$ ,  $\rho$ ] Double)
  ⇒ Int → Popl → Model env es Popl
hmmsir = hmm transPriorsir obsPriorsir transsir obssir

```

We can simulate over this model, perhaps to explore some expected model behaviours, by specifying an input model environment *sim\_env<sub>in</sub>* of type *Env SIRenv* that provides specific values for  $\beta$ ,  $\gamma$ , and  $\rho$ , but provides no values for reported infections  $\xi$  (ensuring that we always sample for  $\xi$ ). Applying *simulate* to *hmm<sub>sir</sub>* 100 and input population *sir<sub>0</sub>* simulates the spread of the disease over 100 days.

```

type SIRenv = '[ $\beta$  := Double,  $\gamma$  := Double,  $\rho$  := Double,  $\xi$  := Int ]
simulateSIR :: IO (Popl, Env SIRenv)
simulateSIR = do
  let sim_envin = (# $\beta$  := [0.7]) • (# $\gamma$  := [0.009]) • (# $\rho$  := [0.3]) • (# $\xi$  := []) • nil
  sir0 = Popl { s = 762, i = 1, r = 0 }
  simulate (hmmsir 100) sim_envin sir0

```

This returns the final population *sir<sub>100</sub>* plus an *output* model environment *sim\_env<sub>out</sub>* mapping each observable variable to the values sampled for that variable during simulation. From this we can extract the reported infections  $\xi$ s:

```

do (sir100 :: Popl, sim_envout :: Env SIRenv) ← simulateSIR
let  $\xi$ s :: [Reported] = get # $\xi$  sim_envout
...

```

Fig. 4a shows a plot of these  $\xi$  values and their corresponding latent (population) states; but note that as it stands, the model provides no external access to the latent states shown in the plot (except the final one *sir<sub>100</sub>*). To generate this kind of output, it would be ideal if we could access those latent states without polluting the type signature of our HMM abstraction, which concisely corresponds to our statistical definition of a HMM. Fortunately, the setting of algebraic effects means users can orthogonally extend a model with a desired *effect*, such as *State*, *Reader*, or in this case a *Writer* effect for writing the intermediate latent states to a stream; this process is straightforward and shown later in §5.5.

### 3.2 Modular Extensions to the SIR Model

Although the SIR model is simplistic, realistic models may be uneconomical to run or too specific to be useful. When designing models, statisticians aim to strike a balance between complexity and precision, and modular models make it easier to incrementally explore this trade-off. We support this claim by showing how two possible extensions of the SIR model are made easy in our language; while these are by no means the most modular solutions possible, they should suffice to make our point.

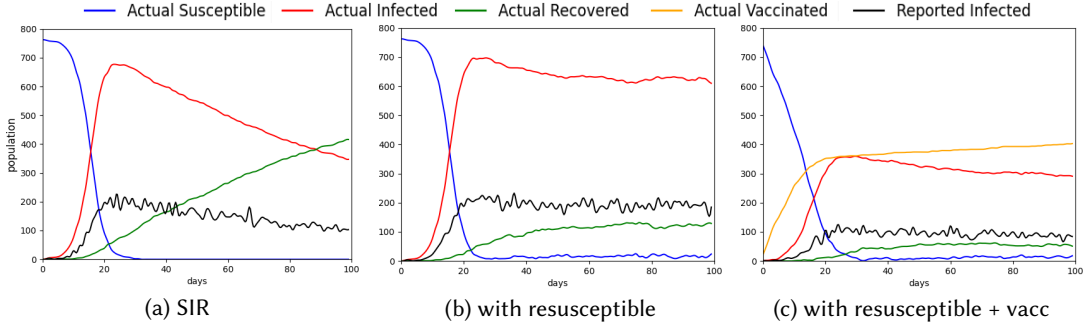


Fig. 4. SIR Hidden Markov Model Simulation

Suppose our disease does not confer long-lasting immunity, so that recovered individuals  $r$  transition back to being susceptible  $s$  after a period of time [Shi et al. 2008]. We can model this with a new transition behaviour:

```

data TransParams = TransParams {  $\beta :: \text{Double}$ ,  $\gamma :: \text{Double}$ ,  $\rho :: \text{Double}$ ,  $\eta :: \text{Double}$  }

transrs :: Double → Popl → Model env es Popl    transsir :: TransModel env es TransParams Popl
transrs  $\eta$  (Popl  $s$   $i$   $r$ ) = do                    transsir (TransParams  $\beta$   $\gamma$   $\eta$ ) =
   $\delta rs \leftarrow \text{binomial } r \ (1 - \exp(-\eta))$     transsi  $\beta$  >=> transir  $\gamma$  >=> transrs  $\eta$ 
  return (Popl ( $s + \delta rs$ )  $i$  ( $r - \delta rs$ ))

```

We need only modify **TransParams** to include a new parameter  $\eta$ ; define a new transition sub-model  $\text{trans}_{rs}$  (parameterised by  $\eta$ ) that stochastically moves individuals from recovered  $r$  to susceptible  $s$ ; and then adjust  $\text{trans}_{sir}$  to compose  $\text{trans}_{rs}$  with our existing transition behaviours. A simulation of the resulting system is shown in Fig. 4b.

Now consider adding a variant where susceptible individuals  $s$  can become vaccinated  $v$  [Ameen et al. 2020]. This involves adding a new sub-population  $v$  to the latent state:

```

data Popl = Popl {  $s :: \text{Int}$ ,  $i :: \text{Int}$ ,  $r :: \text{Int}$ ,  $v :: \text{Int}$  }
data TransParams = TransParams {  $\beta :: \text{Double}$ ,  $\gamma :: \text{Double}$ ,  $\rho :: \text{Double}$ ,  $\eta :: \text{Double}$ ,  $\omega :: \text{Double}$  }

transsv :: Double → Popl → Model env es Popl    transsir :: TransModel env es TransParams Popl
transsv  $\omega$  (Popl  $s$   $i$   $r$   $v$ ) = do                    transsir (TransParams  $\beta$   $\gamma$   $\eta$   $\omega$ ) =
   $\delta sv \leftarrow \text{binomial } s \ (1 - \exp(-\omega))$     transsi  $\beta$  >=> transir  $\gamma$  >=>
  return (Popl ( $s - \delta sv$ )  $i$   $r$  ( $v + \delta sv$ ))    transrs  $\eta$  >=> transsv  $\omega$ 

```

We add field  $v$  to **Popl** representing vaccinated individuals, and add  $\omega$  to **TransParams**, determining the rate at which  $s$  individuals transition to  $v$ . The new behaviour is expressed by  $\text{trans}_{sv}$  and then composed into a new transition model  $\text{trans}_{sir}$ . A simulation of this is shown in Fig. 4c.

### 3.3 Exploring Multimodality in the SIR Model

We now show how our support for higher-order, modular models is complemented by the flexibility of multimodal models. Suppose the goal were to infer SIR model parameters  $\beta, \gamma, \rho$  given data on reported infections  $\xi$ . Ideally we would have a real dataset of reported infections to condition on. But what if our dataset were sparse, or if we were interested in quick hypothesis testing? A common option is to use simulated data as observed data, a method called Bayesian bootstrapping [Fushiki 2010]. This task is made simple with multimodal models, because we can take the outputs from simulation over a model and plug them into an environment that specifies inference over the same model:

```

inferSIR :: Env SIRenv → IO (Env SIRenv)
inferSIR sim_env_out = do
  let ξs      = get #ξ sim_env_out
      mh_env_in = (#β := []) • (#γ := [0.0085]) • (#ρ := []) • (#ξ := ξs) • nil
      sir0     = Popl { s = 762, i = 1, r = 0 }
      mh 50000 (hmmsir 100) (sir0, mh_env_in)

```

Here we take the output model environment `sim_env_out` produced by `simulateSIR`, and use its  $\xi$  values to define an input model environment `mh_env_in` that conditions against  $\# \xi$ . Moreover, suppose we already have some confidence about a particular model parameter, such as the recovery rate  $\gamma$ ; for efficiency, we can avoid inference on  $\gamma$  by setting  $\# \gamma$  to an estimate 0.0085 and sampling only for the remaining parameters  $\# \beta$  and  $\# \rho$ . We then run Metropolis Hastings [Wingate et al. 2011] for 50,000 iterations, which returns an output environment (also of type `Env SIRenv`) containing the values sampled from all iterations.

The inferred posterior distributions for  $\beta$  and  $\rho$  can then be obtained simply by extracting them from that environment:

```

do mh_env_out ← inferSIR ξs
  let βs = get #β mh_env_out
      ρs = get #ρ mh_env_out
  ...

```

These are visualised in Fig. 5. Values around  $\beta = 0.7$  and  $\rho = 0.3$  occur more frequently, because these were the parameter values provided in `simulateSIR` in §3.1.

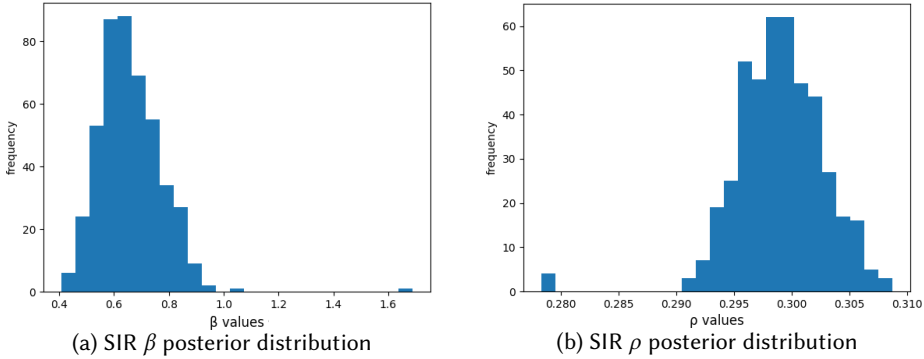


Fig. 5. SIR Inference (Metropolis Hastings)

#### 4 A SYNTACTIC EMBEDDING OF MULTIMODAL MODELS

To support multimodality, a key requirement is that models have a purely syntactic representation which can be assigned a semantics at a later stage. This will allow us to defer the interpretation of primitive distributions as either sample or observe operations until we know the variables we want to condition on. An embedding technique that supports this is an *algebraic effect* embedding [Plotkin and Power 2003]. In this approach, an effectful program which computes a value of type  $a$  has the following type:

```
prog :: Prog es a
```

The parameter  $es$  is called an *effect signature*, in our embedding represented as a type-level list, containing the computational effects that the program may perform. Such programs are syntax

trees whose nodes contain operation calls belonging to effect types  $e \in es$  and whose leaves contain pure values of type  $a$ .

Programs in this form can be interpreted by *algebraic effect handlers* [Plotkin and Pretnar 2013] which handle specific effects in the program:

```
handlee :: Prog (e : es) a → Prog es b
```

Such a handler assigns partial meaning to a program by interpreting all operations  $op$  of type  $e$ , discharging  $e$  from the effect signature, and transforming the return value of type  $a$  into a value of type  $b$ . Effect handlers are modular building blocks that can be composed into various program interpretations.

When we use this approach to represent a probabilistic computation, nodes of the syntax tree will contain calls to probabilistic operations, and inference and simulation will be implemented as effect handlers that interpret those operation calls. Mechanisms (such as sample tracing or particle simulation) specific to particular simulation or inference algorithms can be supported via effects that inject additional computational structure into the model, such as statefulness or non-determinism.

In the rest of this section (§4), we describe our syntactic representation of multimodal models; providing semantics to this embedding via effect handlers will follow in §5.

#### 4.1 An Infrastructure for Algebraic Effects

There are many ways to represent  $\text{Prog } es \ a$ , programs indexed by arbitrary effects; see Hillerström [2015] for a useful summary. Our embedding uses a simple version of the *freer monad* [Kiselyov and Ishii 2015]:

```
data Prog es a where
  Val :: a → Prog es a
  Op  :: forall x. EffectSum es x → (x → Prog es a) → Prog es a
```

Here, a term of type  $\text{Prog } es \ a$  is a syntax tree whose leaves,  $\text{Val } x$ , contain a pure value  $x$  of type  $a$ , and whose nodes,  $\text{Op } op \ k$ , contain an *operation*  $op$  of the abstract datatype  $\text{EffectSum } es \ x$ :

```
data EffectSum (es :: [Type → Type]) (x :: Type) where ...
```

A value of type  $\text{EffectSum } es \ x$  represents a single operation of type  $e \ x$ , where  $e$  is some effect type constructor that occurs in the type-level list  $es$ , and  $x$  is the type of value the operation produces. The freer monad thus supports multiple effects directly via the constructor  $\text{Op}$ , avoiding the need for explicit coproducts (cf. Swierstra [2008]). Effect sums cannot be constructed directly; their implementation is kept abstract, and the following type class,  $\text{Member}$ , is instead provided for working with them:

```
class FindElem e es ⇒ Member e es where
  inj :: e x → EffectSum es x
  prj :: EffectSum es x → Maybe (e x)
```

The constraint  $\text{Member } e \ es$  asserts that if we can determine  $e$ 's position in  $es$  via the type class  $\text{FindElem } e \ es$  (omitted), then we can safely inject and project an effectful operation, of type  $e \ x$ , into and out of  $\text{EffectSum } es \ x$ .

The second argument of the  $\text{Op}$  constructor is a continuation  $k$  of type  $x \rightarrow \text{Prog } es \ a$  that takes the result of the operation and constructs the remainder of the syntax tree. This encoding means effect types  $e$  need not be functors, in contrast to approaches based on the free monad. In turn this means the type of distributions (§4.2.1 below) can be expressed as a GADT, and is in part why we

prefer Kiselyov and Ishii's approach to alternatives such as Wu and Schrijvers [2015] and Kiselyov et al. [2013]. Other details of their design specific to performance are omitted for simplicity.

## 4.2 Multimodal Models as Effectful Programs

We now define a probabilistic model, or simply *model*, to be an effectful program of type `Prog es a` where `es` includes at least two specific effects: `Dist` and `ObsReader env`. The *distribution* effect `Dist` allows the model to make use of primitive distributions such as normal and uniform distributions; the *observable-reader* effect `ObsReader env` allows the model to read and update the values of observable variables in a model environment of type `env`.

```
newtype Model env es a =
  Model { runModel :: (Member Dist es, Member (ObsReader env) es) => Prog es a }
```

The `Member` constraints specify that `es` contains `Dist` and `ObsReader env`. While these two effects suffice for model specification, others may be useful for model execution, and this type allows the model to remain polymorphic in any such additional effects.

**4.2.1 Effects for distributions.** The core computational effect of a probabilistic model is the distribution effect `Dist`, allowing models to be formulated in terms of primitive probability distributions. The constructors represent the operations of the effect type, and thus correspond to various primitive distributions. We present a representative selection below:

```
data Dist a where
  Normal   :: Double -> Double -> Maybe Double -> Dist Double
  Uniform  :: Double -> Double -> Maybe Double -> Dist Double
  Bernoulli :: Double -> Maybe Bool -> Dist Bool
  Discrete :: [(a, Double)] -> Maybe a -> Dist a
  ...
```

Every use of a primitive distribution must, at runtime, be interpretable as either sample or observe, depending on the availability of an observed value. Each `Dist` operation therefore takes an additional parameter of type `Maybe a` (where `a` is the base type of the distribution) indicating the presence or absence of a value to condition on. This detail is hidden from the user, but is used internally to determine how operation calls are to be interpreted, as we discuss next.

**4.2.2 Effects for reading observable variables.** To support conditioning on observable variables via model environments, we require an *observable-reader* effect, `ObsReader env`, with an operation that can read and update values of observable variables in `env`. This is similar in spirit to the well-known `Reader` type.

```
data ObsReader env a where
  Ask :: Observable env x a => ObsVar x -> ObsReader env (Maybe a)
```

The single operation `Ask` of `ObsReader env` has the constraint `Observable env x a`. This associates a *list* of values of type `a` with observable variable `x` in any environment of type `env`, supporting conditioning on multiple dynamic instances of the same observable variable. The `ObsVar x` argument represents the observable variable name (specified using `#`). The result is then of type `Maybe a` indicating the presence or absence of a value to condition on, suitable for passing directly to one of the `Dist` constructors above. We defer the concrete implementations of these types to §5.1.

However, the user is not expected to specify `ObsReader` effects directly. Rather, we provide an interface of *smart constructors* [Swierstra 2008] which manage the requests to read from observable variables. For example:

```

normal :: (Observable env x Double) ⇒ Double → Double → ObsVar x → Model env es Double
normal mu sigma x' = Model (do maybe_v ← call (Ask x'); call (Normal mu sigma maybe_v))

```

Recall that, in addition to the usual normal distribution parameters  $\mu$  and  $\sigma$ , the constructor `Normal` of `Dist` also expects an argument of type `Maybe Double` representing the presence or absence of an observed value. The smart constructor `normal` has a similar signature, but with an observable variable name  $x'$  in place of the `Maybe Double`, thereby associating the distribution with a random variable (rather than particular value) to condition on. The role of the smart constructor is to insert an `Ask` operation into the program which retrieves the corresponding value `maybe_v` of type `Maybe Double` from the model environment, which is then used to call the `Normal` operation.

The helper function `call` simplifies the construction of operation calls, taking care of the injection into `EffectSum es x` and supplying the leaf continuation `Val`:

```

call :: Member es ⇒ e x → Prog es x
call op = Op (inj op) Val

```

A primed variant of each smart constructor is also provided, using `Nothing` as the observed value, for the common case when a primitive distribution does not need to be conditioned on:

```

normal' :: Double → Double → Model env es Double
normal' mu sigma = Model (call (Normal mu sigma Nothing))

```

As an illustration of how the smart constructors work, consider the simple model in Fig. 6 which generates a bias  $p$  from a uniform distribution and uses it to parameterise a Bernoulli distribution, determining whether the outcome of a coin flip  $y$  is more likely to be heads (`True`) or tails (`False`). Fig. 6a shows how the user might write the program (omitting the type signature); Fig. 6b shows the equivalent program written without smart constructors.

<pre> coinFlip = do   p ← uniform 0 1 #p   y ← bernoulli p #y   return y </pre>	<pre> coinFlip = do   maybe_p ← call (Ask #p)   p       ← call (Uniform 0 1 maybe_p)   maybe_y ← call (Ask #y)   y       ← call (Bernoulli p maybe_y)   return y </pre>
(a) User code, with smart constructors	(b) Without smart constructors

Fig. 6. Behaviour of smart constructors

## 5 INTERPRETING MULTIMODAL MODELS

We now turn to using effect handlers to assign semantics to models. Interpreting a multimodal model has two stages: specialising the model into the conditional form determined by a model environment, and then executing the resulting specialised model using a particular simulation or inference algorithm.

The first of these stages is the focus of this section. We start with the implementation of model environments (§5.1), and then define effect handlers for reading observed values (§5.2) and interpreting primitive distributions in response to whether observed values have been provided (§5.3). Effect handlers for simulation and inference are the topic of §6.

### 5.1 Model Environments

Model environments allow the user to assign values to observable variables which the model can then read from. For the sake of compositionality, models should only need to mention the observable



variables they make use of, and be polymorphic in the rest. Moreover, a given observable variable may bind multiple successive values at runtime – one for each time the variable is evaluated.

These two design constraints suggest a representation of model environments as extensible records, where the fields are observable variable names and the values are *lists* of observed values, as expressed by the datatype `Env`:

```
data Env (env :: [Assign Symbol Type]) where
  ENil    :: Env '[]
  ECons   :: [a] → Env env → Env ((x := a) : env)

data Assign x a = x := a
```

The type parameter `env` represents the type of the model environment as a type-level list of pairs `Assign x a`, associating type-level variable names `x` of kind `Symbol` with value types `a` of kind `Type`; this tracks the variables in the environment and their corresponding types. The constructor `ENil` is the empty environment, and the constructor `ECons` takes a list of values of type `a` and an environment of type `env` and prepends a new entry for `x`, producing an environment of type `(x := a) : env`.

The observable variable names in `Env`, being type-level strings of kind `Symbol`, have no value representation; to use them as record fields at the value-level, we give the singleton datatype `ObsVar`:

```
data ObsVar (x :: Symbol) where
  ObsVar :: KnownSymbol x ⇒ ObsVar x
```

This acts as a container for `Symbols`, storing them as a phantom parameter `x`. String values can be neatly promoted to such containers by deriving an instance of the `IsLabel` class, using Haskell's `OverloadedLabels` language extension:

```
instance (KnownSymbol x, x ~ x') ⇒ IsLabel x (ObsVar x') where
```

This enables values of type `ObsVar` to be created using the `#` syntax, so that for example value `#foo` has type `ObsVar "foo"`. Model environments are then constructed using the following interface, which provides `nil` for the empty environment, and an infix cons-like operator `(•)` which makes use of the `:=` notation at the value-level:

```
(•) :: Assign (ObsVar x) [a] → Env env → Env ((x := a) : env)
nil :: Env '[]
```

Finally, constraining a polymorphic model environment is done via the type class `Observable`, which provides type-safe access and updates to observable variables:

```
class (FindElem x env, LookupType x env ~ a) ⇒ Observable env x a
  get :: ObsVar x → Env env → [a]
  set :: ObsVar x → [a] → Env env → Env env
```

```
type family LookupType x env where
  LookupType x ((x := a) : env) = a
  LookupType x ((x' := a) : env) = LookupType x env
```

The methods `get` and `set` use the type class `FindElem` (used in §4.1) to find the position of a variable `x` in `env`, and the type family `LookupType` to retrieve its type `a`. Many observable variables of the same type can be specified with the type family `Observables` `env xs a` below, which returns a nested tuple of constraints `Observable env x a` for each variable `x` in `xs`:

```
type family Observables env (xs :: [Symbol]) a :: Constraint where
  Observables env (x : xs) a = (Observable env x a, Observables env xs a)
  Observables env '[] v = ()
```

Although other designs are possible, using lists to represent observations for random variables is justified by the fact that, for correctness, general-purpose inference algorithms must compute the same distribution on traces (sequences of sampled or observed values) [Tolpin et al. 2016].

## 5.2 Handling Reading of Observable Variables

The first step in specialising a model to a particular model environment is to handle the Ask operations of **ObsReader**, representing environment read requests.

In the setting of algebraic effects, denoting that an effect  $e$  has been handled is done by *discharging* it from the front of an effect signature  $e : es$ . For this, we introduce a helper function `discharge` which pattern-matches an operation of type **EffectSum** ( $e : es$ )  $a$  to determine whether it inhabits the leftmost component  $e$  of the sum:

```
discharge :: EffectSum (e : es) a → Either (EffectSum es a) (e a)
```

If the operation indeed belongs to  $e$ , it is returned as `Right op` where  $op$  has type  $e a$ . Otherwise the operation belongs to an effect in  $es$  and it is returned as `Left op` where  $op$  has type **EffectSum**  $es a$ , discharging  $e$  from the effect signature.

We then define a handler which, given a model environment  $env$  and a program with effect signature (**ObsReader**  $env : es$ ), discharges the **ObsReader**  $env$  effect by interpreting Ask operations:

```
handleRead :: Env env → Prog (ObsReader env : es) a → Prog es a
handleRead env (Op op k) = case discharge op of
  Right (Ask x) → let vs      = get x env
                   maybe_v = safeHead vs
                   env'    = set x (safeTail vs) env
                   in handleRead env' (k maybe_v)
  Left op'      → Op op' (handleRead env . k)
handleRead env (Val x) = return x
```

There are three cases. On matching an operation as `Right (Ask x)` containing a request to read from  $x$ , the list of values  $vs$  associated with  $x$  is looked up in  $env$ . If  $vs$  has a head element, contained in  $maybe_v$ , that value becomes the current observation of  $x$  and is removed from  $env$ ; this ensures that no observed value is conditioned on more than once during an execution, and that the order in which observations are consumed matches the execution order. Otherwise there is no observation of  $x$  and  $env$  is unchanged. The resulting  $maybe_v$  is provided to the continuation  $k$  to construct the remainder of the program, which is then recursively handled by `handleRead` using the updated environment.

If we match an operation as `Left op'`, then  $op'$  does not belong to **ObsReader**. The operation is left intact and the remainder of the program is handled via `(handleRead env . k)`. Lastly, reaching the non-operation `Val x` will simply return value  $x$  as the program's output.

In the handler definitions that follow, we omit the `Left op'` and `Val x` clauses when their implementation follows the pattern above. For an overview of effect handlers we refer the reader to Kiselyov and Ishii [2015].

## 5.3 Handling Distributions

The second step of model specialisation is the handler for the **Dist** effect, which interprets a primitive distribution call as a sampling or observing operation depending on the presence or absence of an observed value. This requires two new effects, **Sample** and **Observe**:

**data Sample a where**

`Sample :: Dist a → Sample a`

**data Observe a where**

`Observe :: Dist a → a → Observe a`

Each has a single operation. Sample takes a distribution to sample from, whereas Observe takes a distribution and an observed value. The distribution handler is then given by:

```

handleDist :: (Member Sample es, Member Observe es) ⇒ Prog (Dist : es) a → Prog es a
handleDist (Op op k) = case discharge op of
  Right d → case getObs d of Just v → (do x ← call (Observe d v)
                                         handleDist (k x))
                                         Nothing → (do x ← call (Sample d)
                                                         handleDist (k x))
  getObs :: Dist a → Maybe a

```

The accessor function `getObs` retrieves the optional observation associated with a primitive distribution (§4.2.1). On encountering a distribution `d`, the handler uses `getObs` to try to retrieve its observed value; if there is such a value `v`, we call a corresponding `Observe` operation, and otherwise we call `Sample`.

#### 5.4 Specialising Multimodal Models

Together, these two handlers specialise a multimodal model into the conditional form determined by a particular model environment. Their net effect on the type of the model is illustrated by the following composite handler:

```

handlecore :: (Member Observe es, Member Sample es)
  ⇒ Env env → Model env (ObsReader env : Dist : es) a → Prog es a
handlecore env = handleDist . (handleRead env) . runModel

```

This handles both `Dist` and `ObsReader`, replacing all primitive distributions with explicit calls to `Sample` and `Observe`. As an example, consider the `coinFlip` model presented earlier in Fig. 6b. Initially applying `handleRead` using the example environment  $((\#p := [0.5]) \cdot (\#y := []) \cdot \text{nil})$  would produce `coinFlip'` in Fig. 7a as an intermediate program, in which all distributions calls have been parameterised by either a concrete observed value or `Nothing`; the updated environment would be  $((\#p := []) \cdot (\#y := []) \cdot \text{nil})$  where `#p` is fully consumed and `#y` is unchanged. Then applying `handleDist` to `coinFlip'` would yield Fig. 7b; observed values in primitive distributions are retained, although they are rendered redundant by the information in the `Observe` constructor.

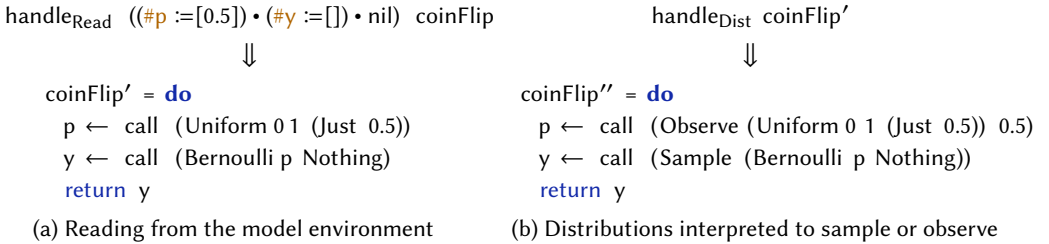


Fig. 7. `coinFlip` model: handling effects

This initial pipeline of effect handling turns a multimodal model into a form suitable for further specialisation by a simulation or inference algorithm, which we show in §6.

#### 5.5 Extending Models with Additional Effects

When building models, users are not restricted to using only the two base effects `Dist` and `ObsReader`: the effect signature `es` in `Model env es a` can be easily extended with an arbitrary desired effect `e`, by first constraining the model with `Member e es`, and then handling `e` with a corresponding handler.

As an example, we revisit the SIR model in §3.1 where Fig. 4 plotted *all* intermediate values `Popl` of susceptible, infected, and recovered individuals over  $n$  days, despite our implementation `hmmsir` only returning *final* one:

```
hmmsir :: (Observables env '[ "ξ" ] Int, Observables env '[ "β", "γ", "ρ" ] Double)
        ⇒ Int → Popl → Model env es Popl
hmmsir n = hmm transPriorsir obsPriorsir transsir obssir n
```

To record all *sir* values produced by `hmmsir`, we introduce the constraint `Member (Writer w) es` to require `es` to contain the effect `Writer w`, representing computations that produce a stream of data of type `w`; here we choose `w` to be a list of *sir* values `[Popl]`. The transition model `transsir` can then use the `Writer` operation `tell` to concatenate each new *sir* value to the existing trace of values:

```
tell      :: Member (Writer w) es ⇒ w → Model env es ()
transsir :: Member (Writer [Popl]) es ⇒ TransModel env es TransParams Popl
transsir (TransParams β γ) sir = do
  sir' ← (transsi β >=> transir γ) sir
  tell [sir']
  return sir'
```

Models with user-specified effects can then be easily reduced into a form suitable for specialisation under a model environment (§5.4), by handling those effects beforehand with a suitable handler:

```
handleWriter :: Monoid w ⇒ Model env (Writer w : es) a → Model env es (a, w)
hmm'sir :: (Observables env '[ "ξ" ] Int, Observables env '[ "β", "γ", "ρ" ] Double)
        ⇒ Int → Popl → Model env es (Popl, [Popl])
hmm'sir n = handleWriter . hmmsir n
```

Here, composing `hmmsir` with `handleWriter` interprets the `tell` operations arising from the transition model, producing a new SIR model `hmm'sir` that returns the trace of *sir* values as an additional output of type `[Popl]`.

## 6 SIMULATION AND INFERENCE AS EFFECT HANDLERS

Simulation or inference over a model is expressed through the semantics we assign to its `Sample` and `Observe` effects; in principle we could therefore define model execution in terms of two handlers, one for each of those effects. However, many algorithmic approaches such as Monte-Carlo methods [Chong et al. 2010] often rely on common mechanisms, such as sample tracing, probability mapping, and model reparameterisation [Robert and Titterton 1998]. By identifying some of these mechanisms as basic building blocks, the possibility arises of composing them in novel ways, giving rise to many useful variants of algorithms [Ścibior et al. 2018]. This motivates a compositional approach in which aspects of model execution can be defined and extended modularly.

In this section, we use effect handlers to define a series of composable program transformations for probabilistic programs, which iteratively refines a model by installing new effects around existing ones. The result is an interpretation of the model in the context of specific algorithms for simulation in §6.1 and inference in §6.2.

### 6.1 Simulation as Effect Handlers

Simulation can be considered the most basic form of model execution. It runs the provided model as a generative process, using observed data from the model environment when available and otherwise drawing new samples. At the end, it returns a model output, plus a *sample trace* uniquely

identifying each *runtime* `Sample` operation with its sampled value; this becomes especially pertinent (§6.2.2) for the correctness and implementation of generic inference algorithms [Tolpin et al. 2016].

To support sample traces, each `Sample` occurrence thus needs a unique dynamic address  $\alpha$  assigned; in our actual embedding, this feature is implemented by `handleDist`. For simplicity, we omit these low-level details, and assume operations `Sample` and `Observe` are now parameterised by an address  $\alpha$  of abstract type `Addr`. The type of sample traces, `STrace`, is then a map from addresses  $\alpha$  to values of abstract type `PrimVal` that primitive distributions can generate (concretely, `PrimVal` is an open sum, but one can avoid this using dependent maps):

```
type STrace = Map Addr PrimVal
```

Updating this sample trace is to be performed by the well-known `State` effect:

```
data State s a where
  Modify :: (s → s) → State s ()

handleState :: s → Prog (State s : es) a → Prog es (a, s)
```

Its operation `Modify` takes a function of type  $s \rightarrow s$  and applies this to state  $s$ . Its handler `handleState`, given an initial state, additionally returns the final state that results from handling a program.

We implement the tracing of samples as the *program transformation* `traceSamples` below, which is a handler that installs a runtime `State STrace` effect after each `Sample` operation:

```
traceSamples :: (Member Sample es, Member (State STrace) es) ⇒ Prog es a → Prog es a
traceSamples (Op op k) = case prj op of
  Just (Sample d α) → Op op (λx → do call (Modify (Map.insert α x))
                                     traceSamples (k x))
```

Notice that we apply `prj` (§4.1) to `op`, rather than `discharge`; this is because we do not intend to handle any effects, and so we place no constraints on the order of the effect signature `es`. Instead, we leave the `Sample` operation unhandled as `op` in the program, and construct a new continuation: this takes the future output  $x$  from `Sample` and calls a `Modify` operation to store  $x$  at address  $\alpha$  in the sample trace, before continuing with the original continuation  $k$ .

We can now define the handlers for `Observe` and `Sample` orthogonally from this transformation. For `Observe` operations, `handleObs` performs no conditioning side-effects, and simply needs to return the observed value  $y$  to its continuation  $k$ :

```
handleObs :: Prog (Observe : es) a → Prog es a
handleObs (Op op k) = case discharge op of
  Right (Observe d y _) → handleObs (k y)
```

For `Sample` operations, `handleSamp` takes the provided distribution  $d$  and applies the function `sampleIO`, defined using an external statistics library; the generated value  $v$  is then passed to  $k$ :

```
handleSamp :: Prog '[Sample] a → IO (a, STrace)
handleSamp (Op op k) = case discharge op of
  Right (Sample d _) → do v ← sampleIO d
                          handleSamp (k v)

sampleIO :: Dist a → IO a
```

Running this dispatches the final effect in `Prog` to produce an `IO` effect, hence it is always executed as the last handler where only `Sample` operations can occur in the program.

The complete definition for simulation is then given by the handler composition `runSimulate`:

```

runSimulate :: es ~ '[ObsReader env, Dist, State STrace, Observe, Sample]
              ⇒ Env env → Model env es a → IO (a, STrace)
runSimulate env = handleSamp . handleObs . (handleState Map.empty) . traceSamples . (handleCore env)

```

Above depicts how the logic of model execution can be decomposed into a modular and transparent system. The concrete effects in `es`, specified by the type coercion  $\sim$ , are kept abstract to the user interested in simply building and using models; simultaneously, the compositional nature of handlers makes it easy for programmers who wish to implement and extend new forms of model execution, as demonstrated next in §6.2.

As a last remark, the reader may notice that the top-level function `simulate` (as seen in §3.1), which uses `runSimulate`, will differ slightly in its type signature, which we give now:

```

simulate :: es ~ '[ObsReader env, Dist, State STrace, Observe, Sample]
           ⇒ (x → Model env es a) → Env env → x → IO (a, Env env)

```

This allows for better composition when simulating a model over many inputs  $x$ . Then for type-safe user-access to an `STrace` structure, this is reified into an *output* environment of type `Env env`; for this, we use simple type-level programming to extract values from the trace whose addresses  $\alpha$  (in the full implementation) are indexed by an observable variable name from the input environment `env`. A similar approach is taken for Likelihood Weighting (`lw`) and Metropolis Hastings (`mh`).

## 6.2 Inference as Effect Handlers

Approximative Bayesian inference attempts to learn the posterior distribution of a model's parameters given some observed data. We reuse the ideas introduced in §6.1 to implement Likelihood Weighting and Metropolis Hastings as inference algorithms.

**6.2.1 Likelihood Weighting (LW).** If one uses simulation as a process for randomly proposing model parameters, the LW algorithm [van de Meent et al. 2018] then assigns these proposals a *weight*, that is, the total likelihood of them having generated some specified observed data.

Its implementation is in fact extremely simple, as most of the work has been done when implementing `runSimulate`. The only change is to how the `Observe` effect is interpreted:

```

handleObsLW :: Double → Prog (Observe : es) a → Prog es (a, Double)
handleObsLW lp (Op op k) = case discharge op of
  Right (Observe d y _) → handleObsLW (lp + logProb d y) (k y)
handleObsLW lp (Val x)   = return (x, lp)

logProb :: Dist a → a → Double

```

The function `handleObsLW` is now parameterised by a log probability `lp`, where at each operation `Observe d y α`, it computes and adds to `lp` the log probability of distribution `d` having generated observed value `y`. The total log probability is returned upon reaching `Val x`.

The complete definition for a single iteration of LW is given as `runLW`:

```

runLW :: es ~ '[ObsReader env, Dist, State STrace, Observe, Sample]
        ⇒ Env env → Model env es a → IO ((a, STrace), Double)
runLW env = handleSamp . (handleObsLW 0) . (handleState Map.empty) . traceSamples . (handleCore env)

```

This is identical to simulation but we now use `handleObsLW` instead of `handleObs`; running this returns the log-likelihood of the values in `STrace` giving rise to the observed data in provided environment `env`. Similar to `runSimulate` and `simulate`, `runLW` is called via a top-level function `lw` (omitted), but is instead performed iteratively to produce a trace of weighted parameter proposals (Fig. 1b).

Likelihood Weighting, however, becomes ineffective as the number of random variables sampled from increases: as values are freshly generated for *all* Sample operations, achieving a high likelihood means sampling an entire set of likely proposals. Our next, final example offers a solution to this:

**6.2.2 Metropolis Hastings (MH).** MH [Wingate et al. 2011] supports *incremental* parameter proposals by randomly choosing a “proposal address”  $\alpha_0$  at the start of each iteration; this denotes the address from which a new sample is to be drawn, where all other addresses are to instead reuse old samples from previous iterations. At the end of an iteration, we decide whether to accept the new sample by comparing the *individual* log probabilities of each probabilistic operation.

To support this, we take a similar approach to the one shown for traceSamples: we define the type `LPTrace` to map addresses of probabilistic operations to their log probabilities, and then the program transformation `traceLPs` to update this as a state.

```
type LPTrace = Map Addr Double
```

```
traceLPs :: (Member Observe es, Member (State LPTrace) es) => Prog es a -> Prog es a
traceLPs (Op op k) = case prj op of
  Just (Observe d y  $\alpha$ ) -> Op op ( $\lambda x \rightarrow$  do call (Modify (Map.insert  $\alpha$  (logProb d y)))
                                     traceLP (k x))
```

On matching against `Observe d y  $\alpha$` , the above installs a `State LPTrace` effect by defining a new continuation that stores the log probability of `d` having generated `y`. The same is done for `Sample` operations, but we compute the log probability of the values sampled.

All of the conditioning is in fact taken care of by `traceLPs`, so the handler for `Observe` operations only needs to return any observed values to their continuations – its implementation is therefore the same as for simulation (§6.1). What is left to define is how `Sample` is interpreted:

```
handleSampMH :: STTrace -> Addr -> Prog '[Sample] a -> IO a
handleSampMH sTrace  $\alpha_0$  (Op op k) = case discharge op of
  Right (Sample d  $\alpha$ ) -> do x <- lookupSample sTrace d  $\alpha$   $\alpha_0$ 
                           handleSampMH  $\alpha_0$  sTrace (k x)
```

```
lookupSample :: STTrace -> Dist a -> Addr -> Addr -> IO a
```

The handler `handleSampMH` takes the sample trace `sTrace` of the previous MH iteration, and an address  $\alpha_0$  denoting the proposed sample site. On matching against an operation `Sample d  $\alpha$` , the function `lookupSample` will generate a new sample if  $\alpha$  matches  $\alpha_0$ , or if no previous MH iterations have yet generated a sample for it; otherwise, it reuses the old sample value of  $\alpha$  found in `sTrace`.

The final definition for a single iteration of MH is given as `runMH`:

```
runMH :: es ~ '[ObsReader env, Dist, State STTrace, State LPTrace, Observe, Sample]
      => Env env -> STTrace -> Addr -> Model env es a -> IO ((a, STTrace), LPTrace)
runMH env sTrace  $\alpha_0$  = (handleSampMH sTrace  $\alpha_0$ ) . handleObs
                      . (handleState Map.empty) . (handleState Map.empty)
                      . traceLPs . traceSamples . (handleCore env)
```

This reuses many of the building blocks of simulation, and extends these with the transformation `traceLPs` and handler for `State LPTrace`; the `handleSampMH` variant is then used instead of `handleSamp`.

Although `runMH` adheres to the MH algorithm in terms of how it samples and conditions, it does not decide the proposal address  $\alpha_0$  at the start of an MH iteration, nor whether newly proposed parameters are accepted at the end of an iteration. We keep this logic distinct from effect handlers, and instead implement it in a wrapper function `mh` (omitted) which folds over iterations of `runMH`, propagating information from the previous MH iteration to support decision making in the next; the end result generates a trace of accepted parameter proposals (Fig. 5).



## 7 EVALUATION

### 7.1 Quantitative Evaluation

We compare our language's performance, which for simplicity is implemented using `freer-simple`<sup>2</sup>, against two state-of-the-art PPLs. First, `MonadBayes` [Ścibior et al. 2018], a Haskell embedded language which uses a monad transformer library (`mtl`)<sup>3</sup> approach as an effect system for PPLs, but does not support multimodal models. Second, `Turing` [Ge et al. 2018], which achieves multimodal models via macro compilation in Julia as a host language.

Our evaluation strategy uses a set of popular benchmarks [Kulkarni et al. 2020], comparing simulation (SIM), Likelihood Weighting (LW), and Metropolis Hastings (MH) as forms of execution algorithms. We apply these to the following example models: linear regression, hidden Markov model, and latent dirichlet allocation. These benchmarks, given fully in Appendix A, are performed on an AMD Ryzen 5 1600 Six-Core Processor with 16GB of RAM.

Across all models, our performance scales linearly with the number of samples each algorithm generates; this remains mostly the case when varying the dataset size to models, except for inference on hidden Markov model where we begin to scale quadratically.

Against `Turing`, we are on average 4.0x and 1.2x faster for SIM and LW. As our benchmarks do not consider the overhead from `Turing`'s macro-compilation stage for specialising multimodal models, this may indicate our runtime approach to model specialisation does not notably affect performance. `MonadBayes` is on average 3.3x faster than us for SIM and has close to constant-time performance for LW, making it difficult to compare the latter. In contrast to us, they choose to not store and update sample traces for SIM and LW; this effectively means their performance is impacted only by the cost of sampling operations (of which there are naturally fewer for LW), whereas we incur overhead from additional `State` operations. Another factor we heed is the performative difference between `freer-simple` and `mtl` for large, non-synthetic programs, which still requires investigation.

For MH in particular, `Turing` and `MonadBayes` are on average 3.1x and 1.8x faster. Our implementation is naive in that we traverse the whole sample trace for MH updates despite a fixed number of variables being updated – this is straightforward to amend [Wingate et al. 2011]. We also remark that `Turing` and `MonadBayes` only perform MH updates to *one* variable per iteration, and do not consider the dependent variables which consequently need updating [Kiselyov 2016b]; our implementation does account for this and hence incurs costs for updating groups of dependencies, which can be larger for models with more complex dependency graphs.

We find our language competitively performant despite not yet having explored potential optimisation, and expect to benefit greatly from off-the-shelf Haskell techniques such as inlining and more efficient data structures. Exploring performance of effect handlers with PPLs in general is an important and challenging topic we plan to investigate separately, including techniques specific to algebraic effects, such as the codensity monad [Voigtländer 2008], and alternative representations of effectful programs [Maguire 2019; Wu and Schrijvers 2015]. We also observe that *all* of our handlers are applied on each model execution, whereas we suspect partially pre-evaluated models can be executed by reconsidering the order of handling; this may lead to substantial performance gains.

### 7.2 Qualitative Evaluation

Finally, we compare our supported features across a larger range of modern PPLs in Table 1. To the best of our knowledge, our language is the first to fully support both multimodal and higher-order models. Models in Gen [Cusumano-Towner et al. 2019], denoted with the special `@gen` syntax,

<sup>2</sup><https://github.com/lexi-lambda/freer-simple>

<sup>3</sup><https://github.com/haskell/mtl>

Table 1. Comparison of PPLs in terms of supported features for models, where ● is full support, ◐ is partial support, and ○ is no support. Modular models are those that can be defined in terms of other models.

Supported model features	Wasabaye	Gen	Turing	Stan	Pyro	MonadBayes	Anglican	WebPPL
Multimodal	●	●	●	●	◐	○	○	○
Modular	●	●	●	○	●	●	●	●
Higher-order	●	◐	○	○	●	●	◐	●
Type-safe	●	○	○	●	○	●	○	○

have support for higher-order interactions with other @gen terms; however, applying a standard higher-order function to a model, at least without programmer intervention, will escape the tracing of the model’s computation. In Pyro [Bingham et al. 2019], models are Python functions and as such are first-class, but support for multimodal models is limited: random variables may have only *one* observed value at runtime, and so Pyro expects this value to be a matrix where all contained data is conditioned on simultaneously. This approach cannot be used for models with sequential behaviour such as HMMs.

As far as we know, our language is also the first general-purpose PPL with multimodal models in a statically typed paradigm. Stan [Carpenter et al. 2017] is type-safe but special-purpose. Other general-purpose PPLs with multimodality are dynamically typed (Pyro) or use just-in-time compilation (Turing, Gen), and so do not guarantee that the observed values assigned to random variables are correctly typed, or that these variables exist; however, some progress has been made towards a type-safe version of Gen [Lew et al. 2019]. Although MonadBayes is statically typed, we are not aware of any attempts to extend its mtl approach to support multimodal models.

We note that as our language is experimental, the range of model execution algorithms we present so far is small. Also, whilst we provide basic utilities such as IO debugging and type-safe interfacing with sample traces, richer PPL features such as run-time inference diagnostics have not yet been implemented. We believe our language’s infrastructure, having embedded into an algebraic effect setting, can readily support extensions in both of these areas.

## 8 CONCLUSION

Probabilistic programming is an active research topic, with applications ranging from artificial intelligence [Tran et al. 2017] to market research [Letham et al. 2016]. Many existing PPLs are either too low-level to capture multimodal descriptions of models, or too rigid to allow models to be easily reused. In this paper we used an algebraic effects embedding to implement a PPL where multimodal models are first-class citizens that can be modularly defined and easily combined.

### 8.1 Related Work

**8.1.1 Effect abstractions for PPLs.** Although the design pattern of using interpreters to execute models is common in PPLs [Carpenter et al. 2017; Salvatier et al. 2016; Tolpin et al. 2016], viewing these interpreters formally as algebraic effect handlers has little existing literature. Ścibior and Kammar [2015] first demonstrate in Haskell how rejection sampling can be abstracted into handlers. Moore and Gorinova [2018] take inspiration from effect handlers and use Python *context managers* for sample tracing and conditioning; we distinguish these from algebraic effects, mainly because the operations are methods rather than syntax, handlers are coroutines invoked *by* programs, and operations and handlers are not associated with effects.

A notable alternative effect system is the *monad transformer library* (mtl) [Liang et al. 1995] approach adopted in MonadBayes [Ścibior et al. 2018]. The observation behind MonadBayes is that inference algorithms can be decomposed into monadic building blocks which together form a monad

transformer stack. This is comparable to our method, in that it also uses program transformations to augment models with additional inference-specific behaviours, but whereas we transform a syntactic representation of models via handler composition, [Ścibior et al.](#) directly execute their models via composition of monadic functions. Our handlers always produce intermediate programs of type `Prog es`, and are therefore compositional; however, making handlers compose for arbitrary effects requires additional infrastructure to appropriately forward the effects in a way that is compatible with monad transformers [[Schrijvers et al. 2019](#)].

For constructing models, we find the `mtl` approach of `MonadBayes` offers a similar user experience to algebraic effects: in particular both support abstract effect signatures for model components, allowing effect constraints to be propagated bottom-up through the program. For implementing inference algorithms, the relative advantages of monads vs. algebraic effects are more apparent. Monads appear to be more lightweight building blocks when making incremental extensions to inference, allowing for finer-grained composition. On the other hand, working with monads can be tricky when building sophisticated algorithms: the `mtl` design pattern of `MonadBayes` is to have several monads invoke their own side-effects when sampling or observing, arranged by having each monad “lift” the computation up through the transformer stack. To invoke only a selection of particular monads requires the programmer to correctly position a monad in the stack and have it terminate the sequence of lifts early. Moreover as `mtl` performs all lifts implicitly via type-class instances, this tends to obscure the current monadic context. By contrast, expressing such logic in a single effect handler can yield a more cohesive solution. Monads also come with pre-defined semantics, whereas in inference it is often advantageous to override certain behaviours e.g. static vs. dynamic tracing of random choices [[Ścibior et al. 2018](#)]. To achieve this, the same monad along with *all* of its operations and type-class instances must be redefined afresh in separate modules; this results in a proliferation of redundant boilerplate, and can make arduous the task of parameterising algorithms by new semantics.

**8.1.2 Embedding PPLs.** Tagless-final shallow embedding, as conceived by [Kiselyov \[2010\]](#), has often been used to correspond PPL syntax to type class methods [[Kiselyov 2016a](#); [Kiselyov and Shan 2009](#); [Narayanan et al. 2016](#)], where type class instances correspond to interpretations of programs under particular inference algorithms. This is well suited for applying a semantic domain uniformly over a program, but we found it difficult to iterate transformations (such as interpretations to sample and observe) over models to implement inference compositionally.

Free monads have also seen use, where [Ścibior et al. \[2015\]](#) embed primitive and conditional distributions as an intermediate free monad representation. This bears an initial resemblance to us, but the semantics are instead provided using type classes, and their direct encoding of conditional distributions means models are not multimodal. Later work by [Ścibior et al. \[2018\]](#) more closely coincides with our approach, where their implementation of Metropolis-Hastings uses free monad transformers [[Schrijvers et al. 2019](#)] to encode sampling operations as syntax; this allows models to be executed so that sampling can either invoke an IO effect or reuse previous samples.

**8.1.3 Observed data for multimodal models.** In terms of specifying observed data to multimodal models, `Pyro` [[Bingham et al. 2019](#)] and `Gen` [[Cusumano-Towner et al. 2019](#)] take the most similar approach to us initially: letting users provide mappings between variable names and observed values to a top-level function. `Pyro` then instead wraps models in a runtime context manager that constrains the values of variables to observed data, whereas `Gen` uses the data to macro-expand models into either a static computation graph or standard function. To support *many* observed values for the same variable, `Pyro` relies on these to be presented as a single matrix to be conditioned on together (§7.2), and `Gen` requires values to be individually mapped to *runtime* occurrences of a variable; our approach instead allows traces (lists) of values to be assigned to the same variable.

Special-purpose PPLs such as Stan [Carpenter et al. 2017] and WinBUGS [Lunn et al. 2000] specify models and observed data separately via bespoke language constructs, where common variable names are consolidated during compilation to a different target language.

Although our specific approach to model environments appears to be novel, closely related is work by Lew et al. [2019]. They investigate row polymorphism for assigning types to names of random choices when tracing probabilistic computations; we hope to build on this idea when formalising our language, in particular, for characterising the execution space of a multimodal model under a certain model environment.

## 8.2 Future Work

Our implementation currently responds to running out of observed values in the model environment by switching to sampling, as is typical in PPLs [Bingham et al. 2019; Ge et al. 2018]. However, it may be possible in some settings to use type-level naturals to statically constrain the number of observations required. Alternatively, a dynamic check to signal when too many or too few observations are provided would be easy to implement.

We are also investigating how naming conflicts between observable variables should be resolved when combining models. This is not considered an issue by most existing PPLs; many require programmers to uniquely name each dynamic random variable instance [Cusumano-Towner et al. 2019; Salvatier et al. 2016], perhaps failing at runtime if this condition is not met [Bingham et al. 2019]. Our language is type-safe in allowing model environments to have typed, orthogonally combinable variables (via constraint kinds), but for additional modularity, we are also considering a renaming mechanism for rebinding observable variables when name clashes arise.

A related topic is when the programmer statically refers to the same observable variable more than once in a model:

```
do x1 ← normal 0 1 #x
  x2 ← normal 0 2 #x
return (x1 + x2)
```

Technically, this is an invalid use of a random variable, resulting in an ill-formed model where `#x` is (confusingly) distributed according to two different distributions. For now, programmers must take care not to misuse observable variables in this way; a solution we intend to explore is an affine type system for model contexts which will disallow multiple static uses of the same observable variable. We are also working on a formalisation of our language and embedding technique.

Lastly, we aim to explore how using effect handlers as program transformations can help with the implementation of sophisticated, compositional inference algorithms such as SMC<sup>2</sup> [Doucet et al. 2001] and PMMH [Chopin 2002]; we anticipate needing to reason about interactions with new effects, for example with non-determinism and exceptions, and whether such effects distribute over [Wu et al. 2014] and commute with each other [Gibbons and Hinze 2011] in a probabilistic setting. For scaling up to advanced techniques that require gradient information, e.g. HMC and NUTS [Hoffman et al. 2014], models also need to be differentiable. This is generally done in PPLs via *automatic differentiation* (AD), a procedure for instantiating standard programs as differentiable functions. We believe it possible for our language's infrastructure to support this as a new effect, perhaps by building on top of recent work by Sigal [2021] who show it indeed possible to express AD through effect handlers. A simpler option may be to handle a model into a form suitable for use in an existing AD library [Kmett et al. 2021]. In either case, we would need to investigate the set of valid operations or syntactic forms that enable an AD interpretation of multimodal models in our language.

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## A PERFORMANCE BENCHMARKS

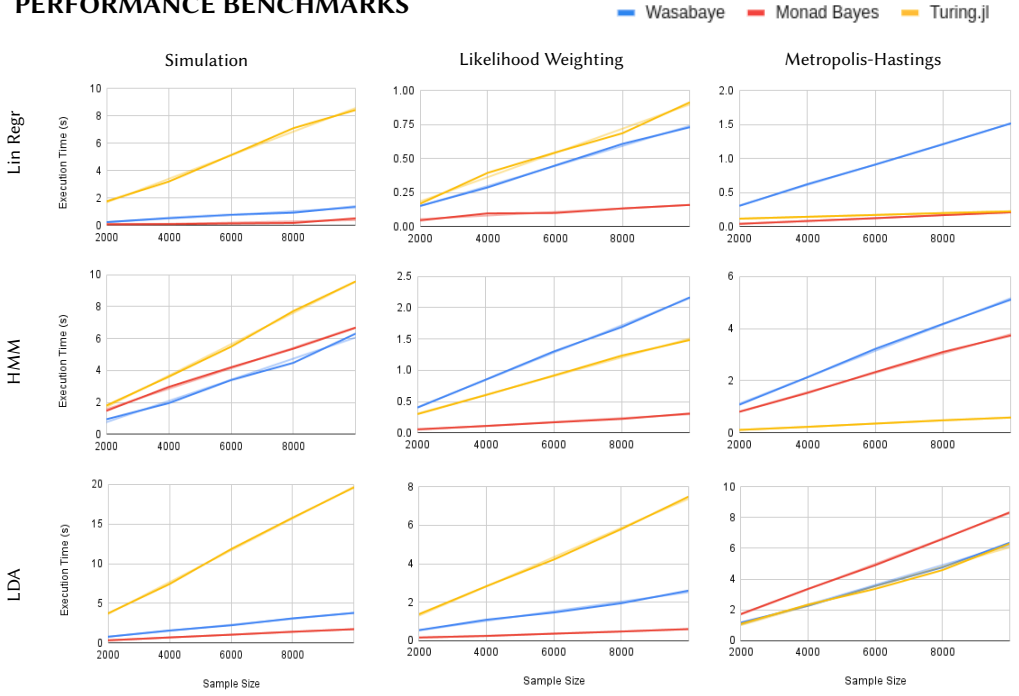


Fig. 8. Execution times of algorithms (seen at top) applied to models (seen on left) with respect to the number of samples each algorithm generates (i.e. number of algorithm iterations).

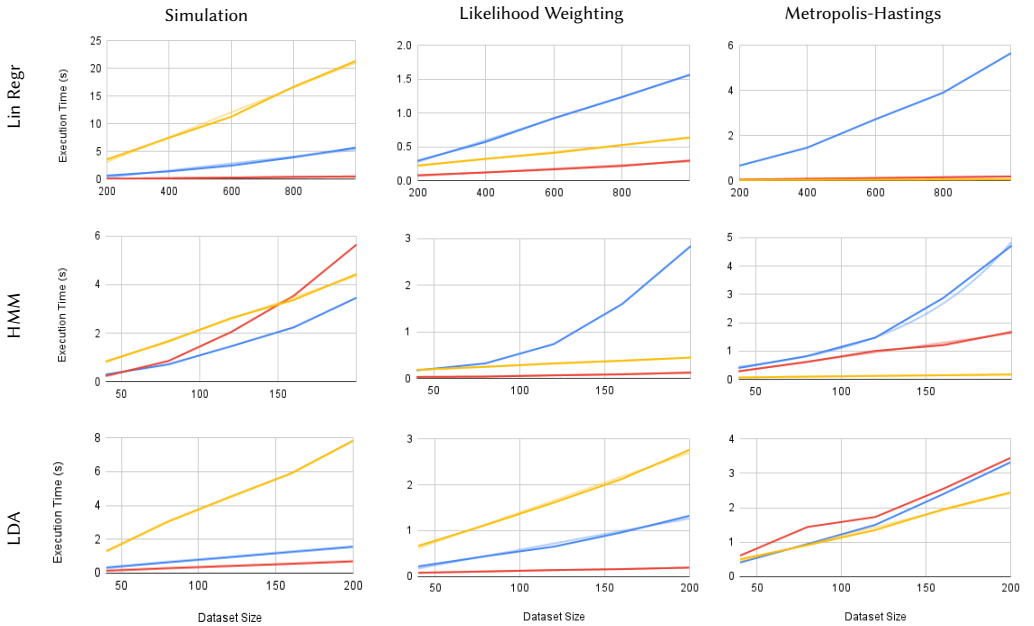


Fig. 9. Execution times of algorithms (at top) applied to models (on left) with respect to dataset size provided to models. The data we vary over is: the number of data points in linear regression, the number of Markov nodes in the hidden Markov model, and the length of text document for latent dirichlet allocation.