
Approximate Knowledge Compilation by Online Collapsed Importance Sampling

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Abstract

We introduce collapsed compilation, a novel approximate inference algorithm for discrete probabilistic graphical models. It is a collapsed sampling algorithm that incrementally selects which variable to sample next based on the partial sample obtained so far. This online collapsing, together with knowledge compilation inference on the remaining variables, naturally exploits local structure and context-specific independence in the distribution. These properties are naturally exploited in exact inference, but are difficult to harness for approximate inference. Moreover, by having a partially compiled circuit available during sampling, collapsed compilation has access to a highly effective proposal distribution for importance sampling. Our experimental evaluation shows that collapsed compilation performs well on standard benchmarks. In particular, when the amount of exact inference is equally limited, collapsed compilation is competitive with the state of the art, and outperforms it on several benchmarks.

1 Introduction

Modern probabilistic inference algorithms for discrete graphical models are designed to exploit key properties of the distribution. In addition to classical conditional independence, they exploit local structure in the individual factors, determinism coming from logical constraints (Darwiche, 2009), and the context-specific independencies that arise in such distributions (Boutilier et al., 1996). The *knowledge compilation* approach in particular forms the basis for state-of-the-art probabilistic inference algorithms in a wide range of models, including Bayesian networks (Chavira & Darwiche, 2008), factor graphs (Choi et al., 2013), relational models (Chavira et al., 2006), probabilistic programs (Fierens et al., 2015), probabilistic databases (Van den Broeck & Suciu, 2017), and dynamic Bayesian networks (Vlasselaer et al., 2016). Based on logical reasoning techniques, knowledge compilation algorithms construct an *arithmetic circuit* representation of the distribution on which inference is guaranteed to be efficient (Darwiche, 2003). The inference algorithms listed above have one common limitation: they perform exact inference by compiling a worst-case exponentially-sized arithmetic circuit representation. Our goal in this paper is to upgrade these techniques to allow for *approximate probabilistic inference*, while still naturally exploiting the structure in the distribution. We aim to open up a new direction towards scaling up knowledge compilation to larger distributions.

When knowledge compilation produces circuits that are too large, a natural solution is to sample some random variables, and to do exact compilation on the smaller distribution over the remaining variables. This collapsed sampling approach suffers from two problems. First, collapsed sampling assumes that one can determine a priori which variables need to be sampled to make the distribution amenable to exact inference. When dealing with large amounts of context-specific independence, it is difficult to find such a set, because the independencies are a function of the particular values that

variables get instantiated to. Second, collapsed sampling assumes that one has access to a proposal distribution for the sampled variables, and the success of inference largely depends on the quality of this proposal. In practice, the user often needs to specify the proposal distribution, and it is difficult to automatically construct one which is general-purpose.

As our first contribution, Section 2 introduces *online collapsed importance sampling*, where the sampler chooses which variable to sample next based on the values sampled for previous variables. This algorithm is a solution to the first problem identified above. We show that the sampler corresponds to a classical collapsed importance sampler on an augmented graphical model and prove conditions for it to be asymptotically unbiased.

Section 3 describes our second contribution: a *collapsed compilation* algorithm that maintains a partially compiled arithmetic circuit during online collapsed sampling. This circuit serves as a highly-effective proposal distribution at each step of the algorithm. By setting a limit on the circuit size as we compile more factors into the model, we are able to sample exactly as many variables as needed to fit the arithmetic circuit in memory. Moreover, through online collapsing, the set of collapsed variables changes with every sample, exploiting different independencies in each case.

Finally, we experimentally validate the performance of collapsed compilation on standard benchmarks. We begin by empirically examining properties of collapsed compilation, to show the value of the proposal distribution and pick apart where performance improvements are coming from. Then, in a setting where the amount of exact inference is fixed, we find that collapsed compilation is competitive with state-of-the-art approximate inference algorithms, outperforming them on several benchmarks.

2 Online Collapsed Importance Sampling

We begin with a brief review of collapsed importance sampling, before motivating the need for dynamically selecting which variables to sample. We then demonstrate that we can select variables in an online fashion while maintaining the desired unbiasedness property, using an algorithm we call online collapsed importance sampling.

We denote random variables with uppercase letters (X), and their instantiation with lowercase letters (x). Bold letters denote sets (\mathbf{X}) and their instantiations (\mathbf{x}). We refer to Koller & Friedman (2009) for notation and formulae related to (collapsed) importance sampling.

2.1 Collapsed Importance Sampling

The basic principle behind collapsed sampling is that we can reduce the variance of an estimator by making part of the inference exact. That is, suppose we partition our variables into two sets: \mathbf{X}_p , and \mathbf{X}_d . In collapsed importance sampling, the distribution of variables in \mathbf{X}_p will be estimated via importance sampling, while those in \mathbf{X}_d will be estimated by computing exactly $P(\mathbf{X}_d|\mathbf{x}_p)$ for each sample \mathbf{x}_p . In particular, suppose we have some function $f(\mathbf{x})$ where \mathbf{x} is a complete instantiation of $\mathbf{X}_p \cup \mathbf{X}_d$, and a proposal distribution Q over \mathbf{X}_p . Then we estimate the expectation of f by

$$\hat{\mathbb{E}}(f) = \frac{\sum_{m=1}^M w[m] (\mathbb{E}_{P(\mathbf{X}_d|\mathbf{x}_p[m])} [f(\mathbf{x}_p[m], \mathbf{X}_d)])}{\sum_{m=1}^M w[m]}, \tag{1}$$

where we have drawn samples $\{\mathbf{x}_p[m]\}_{m=1}^M$ from a proposal distribution Q , and for each sample analytically computed the importance weights $w[m] = \frac{P(\mathbf{x}_p[m])}{Q(\mathbf{x}_p[m])}$, as well as $\mathbb{E}_{P(\mathbf{X}_d|\mathbf{x}_p[m])} [f(\mathbf{x}_p[m], \mathbf{X}_d)]$. Due to the properties of importance samplers, the estimator given by (1) is asymptotically unbiased. Moreover, if we can compute $P(\mathbf{x}_p[m])$ exactly rather than the unnormalized $\hat{P}(\mathbf{x}_p[m])$, then the estimator is unbiased (Tokdar & Kass, 2010).

2.2 Motivation

A critical decision that needs to be made when doing collapsed sampling is selecting a partition – that is, which variables go in \mathbf{X}_p and which go in \mathbf{X}_d . The choice of partition can have a large effect on the quality of the resulting estimator, and the process of choosing such a partition requires expert knowledge. Furthermore, selecting a partition a priori that works well is not always possible, as we

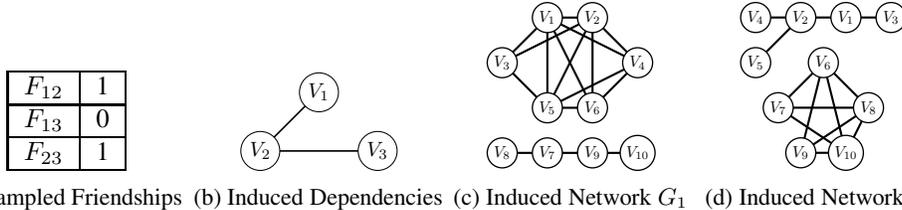


Figure 1: Different samples for \mathbf{F} can have a large effect on the resulting dependencies between \mathbf{V} .

will show in the following example. All of this raises the question whether it is possible to choose the partition on the fly for each sample, which we will discuss in Section 2.3.

Suppose we have a group of n people, denoted $1, \dots, n$. For every pair of people $(i, j), i < j$, there is a binary variable F_{ij} indicating whether i and j are friends. Additionally, we have features V_i for each person i , and $F_{ij} = 1$ (that is i, j are friends) implies that V_i and V_j are correlated. Suppose we are performing collapsed sampling on the joint distribution over \mathbf{F}, \mathbf{V} , and that we have already decided to place all friendship indicators F_{ij} in \mathbf{X}_p to be sampled. Next, we need to decide which variables in \mathbf{V} to include in \mathbf{X}_p for the remaining inference problem over \mathbf{X}_d to become tractable. Observe that given a sampled \mathbf{F} , due to the independence properties of \mathbf{V} relying on \mathbf{F} , a graphical model G is induced over \mathbf{V} (see Figures 1a,1b). Moreover, this graphical model can vary greatly between different samples of \mathbf{F} . For example, G_1 in Figure 1c densely connects $\{V_1, \dots, V_6\}$ making it difficult to perform exact inference. Thus, we will need to sample some variables from this set. However, exact inference over $\{V_7, \dots, V_{10}\}$ is easy. Conversely, G_2 in Figure 1d depicts the opposite scenario: $\{V_1, \dots, V_5\}$ forms a tree, which is easy for inference, whereas $\{V_6, \dots, V_{10}\}$ is now intractable. It is clearly impossible to choose a small subset of \mathbf{V} to sample that fits all cases, thus demonstrating a need for an online variable selection during collapsed sampling.

2.3 Algorithm

We now introduce our online collapsed importance sampling algorithm. It decides at sampling time which variables to sample and which to do exact inference on.

To gain an intuition, suppose we are in the standard collapsed importance sampling setting. Rather than sampling an instantiation \mathbf{x}_p jointly from Q , we can instead first sample $x_{p_1} \sim Q(X_{p_1})$, then $x_{p_2} \sim Q(X_{p_2}|x_{p_1})$, and so on using the chain rule of probability. In online collapsed importance sampling, rather than deciding X_{p_1}, X_{p_2}, \dots a priori, we select which variable we will use as X_{p_2} based on the previous sampled value x_{p_1} .

Definition 1. Let \mathbf{y} be an instantiation of $\mathbf{Y} \subset \mathbf{X}$. A variable selection policy π takes \mathbf{y} and either stops sampling or returns a distribution over $\mathbf{X} \setminus \mathbf{Y}$ that decides the next variable to sample.

For example, a naive policy could be to select a remaining variable uniformly at random. Once the policy π stops sampling, we are left with an instantiation \mathbf{x}_p and a set \mathbf{X}_d , where both are specific to the particular sample.

Algorithm 1 shows more precisely how online collapsed importance sampling generates a single sample, given a full set of variables \mathbf{X} , a variable selection policy π , and proposal distributions Q_{X_i} for any choice of X_i and \mathbf{x}_p . Note that \mathbf{x}_p is a set of variables together with their instantiations, while \mathbf{X}_d is just a set of variables. Notationally, for sample m we write these as \mathbf{x}_p^m and \mathbf{X}_d^m . The joint proposal $Q(\mathbf{x}_p)$ is left abstract for now (see Section 2.4.2 for a concrete instance). In general, it is induced by π and the individual local proposals Q_{X_i} .

Algorithm 1: Online Collapsed IS

Input: \mathbf{X} : The set of all variables,
 Q_{X_i} : Proposal distributions,
 π : Variable selection policy

Result: A sample $(\mathbf{X}_d^m, \mathbf{x}_p^m, w[m])$

```

1  $\mathbf{x}_p \leftarrow \{\}; \mathbf{X}_d \leftarrow \mathbf{X}$ 
2 while  $\pi$  does not stop do
3    $X_i \sim \pi(\mathbf{x}_p)$ 
4    $x_i \sim Q_{X_i}(X_i|\mathbf{x}_p)$ 
5    $\mathbf{x}_p \leftarrow \mathbf{x}_p \cup \{x_i\}$ 
6    $\mathbf{X}_d \leftarrow \mathbf{X}_d \setminus \{X_i\}$ 
7 return  $(\mathbf{X}_d, \mathbf{x}_p, \frac{\hat{P}(\mathbf{x}_p)}{Q(\mathbf{x}_p)})$ 

```

Definition 2. Given m samples $\{\mathbf{X}_d^m, \mathbf{x}_p^m, w[m]\}$ produced by online collapsed importance sampling, the online collapsed importance sampling estimator of f is

$$\hat{\mathbb{E}}(f) = \frac{\sum_{m=1}^M w[m] (\mathbb{E}_{P(\mathbf{X}_d^m | \mathbf{x}_p^m)} [f(\mathbf{x}_p^m, \mathbf{X}_d^m)])}{\sum_{m=1}^M w[m]}. \quad (2)$$

Note that the only difference compared to Equation 1 is that sets \mathbf{X}_p^m and \mathbf{X}_d^m vary with each sample.

2.4 Analysis

Our algorithm for online collapsed importance sampling raises two questions: does Equation 2 yield unbiased estimates, and how does one compute the proposal $Q(\mathbf{x}_p)$? We study both questions next.

2.4.1 Unbiasedness of Estimator

If we let π be a policy that always returns the same variables in the same order, then we recover classical offline collapsed importance sampling - and thus retain all of the properties of offline collapsed importance sampling. With this in mind, we present the augmented factor graph construction given in Figure 2, which will allow us to make a similar statement for any arbitrary policy π .



Figure 2: Online collapsed sampling corresponds to collapsed sampling on an augmented graph

Intuitively, X_i^* is a copy variable of X_i , where we are sampling X_i^* and computing X_i exactly. To make this possible without actually inferring the entire distribution exactly, we have f_{au} cause each S_i to act as an indicator for whether X_i^* and X_i are constrained to be equal. S_i can also be thought of as indicating whether or not we are sampling X_i in our original factor graph when doing online collapsed importance sampling. We are now ready to prove the following results.

Theorem 1. For any factor graph F and its augmented graph F_A , we have $\forall \mathbf{x}, \mathbb{E}_F(\mathbf{x}) = \mathbb{E}_{F_A}(\mathbf{x})$.

Theorem 2. Let F be a factor graph and let F_A be its augmented factor graph. Consider the collapsed importance sampling estimator (Eq. 1) with $\mathbf{X}_p = \mathbf{X}^* \cup \mathbf{S}$ and $\mathbf{X}_d = \mathbf{X}$. This estimator is equivalent to using the online collapsed importance sampling estimator (Eq. 2) on F .

Corollary 1. The estimator given by Equation 2 is asymptotically unbiased.

Proofs and the details of this construction can be found in Appendix A.

2.4.2 Computing the Proposal Distribution

Our next question is how to compute the global $Q(\mathbf{x}_p)$, given that we have each local $Q_{X_{p_i}}(X_{p_i} | \mathbf{x}_{p_{1:i-1}})$. Notice that since $Q_{X_{p_i}}$ is not a conditional from Q , but rather a generic distribution, this computation is not easy in general. In particular, considering $|\mathbf{X}_p| = n$ and our previous example of a uniformly random policy, then for any given instantiation \mathbf{x}_p , there are $n!$ different ways \mathbf{x}_p could be sampled from Q - one for each ordering that arrives at \mathbf{x}_p . In this case, computing $Q(\mathbf{x}_p)$ requires summing over exponentially many terms, which is undesirable. Instead, we restrict the variable selection policies we use to the following class, which makes computing $Q(\mathbf{x}_p)$ easy.

Definition 3. A deterministic variable selection policy $\pi(\mathbf{x}_p)$ is a function with a range of $\mathbf{X} \setminus \mathbf{X}_p$.

Theorem 3. For any sample \mathbf{x}_p and deterministic selection policy $\pi(\mathbf{x}_p)$, there is exactly one order in which the variables \mathbf{X}_p could have been sampled. Therefore, $Q(\mathbf{x}_p) = \prod_{i=1}^{|\mathbf{X}_p|} Q_{X_{p_i}}(x_{p_i} | \mathbf{x}_{p_{1:i-1}})$.

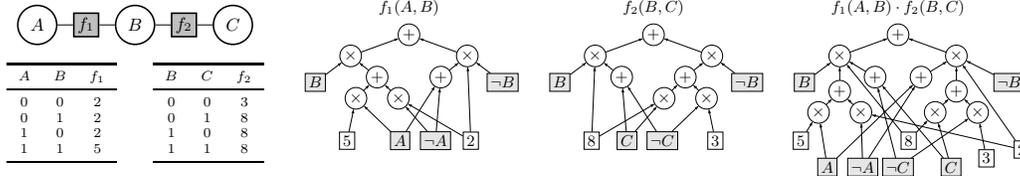


Figure 3: Multiplying Arithmetic Circuits: Factor graph and ACs for individual factors which multiply into a single AC for the joint distribution. Given an AC, inference is tractable by propagating inputs.

3 Collapsed Compilation

Online collapsed importance sampling presents us with a powerful technique for adapting to problems traditional collapsed importance sampling may struggle with. However, it also demands we solve several difficult tasks: one needs a good proposal distribution over any subset of variables, an efficient way of computing exactly the probability of a generic query, and an efficient way of finding the true probability of sampled variables. In this section, we introduce collapsed compilation, which tackles all three of these problems at once using techniques from knowledge compilation.

3.1 Knowledge Compilation Background

We begin with a short review of how to perform exact inference on a probabilistic graphical model using knowledge compilation to arithmetic circuits (ACs). Suppose we have a factor graph consisting of three binary variables A , B and C , and factors f_1, f_2 as depicted in Figure 3. Each of these factors, as well as the joint distribution of the product can be represented as an arithmetic circuit. These circuits have inputs corresponding to variable assignments (e.g., A and $\neg A$) or constants (e.g., 5). Internal nodes are sums or products. We can encode a complete assignment by setting the corresponding variable assignments to 1 and the opposing assignments to 0. Then, the root of the circuit evaluates the weight (unnormalized probability) of that world. Moreover, because of two important properties, we are also able to perform marginal inference. Product nodes are *decomposable*, meaning that their inputs are disjoint, having no variable inputs in common. Sum nodes are *deterministic*, meaning that for any given complete input assignment to the circuit, at most one of the sum’s inputs evaluates to a non-zero value. Now, by setting both assignments for the same variable to 1, we effectively marginalize out that variable. For example, by setting all inputs to 1, the arithmetic circuit evaluates to the partition function of the graphical model. We refer to Darwiche (2009) for further details on how to reason with arithmetic circuits.

In practice, arithmetic circuits are often compiled from graphical models through a logical task called weighted model counting, followed by Boolean circuit compilation techniques. We refer to Choi et al. (2013) and Chavira & Darwiche (2008) for details. As our compilation target, we will use the sentential decision diagram (SDD) (Darwiche, 2011). Moreover, given any two SDDs representing factors f_1, f_2 , and a variable X , we can efficiently compute the SDD representing the factor multiplication of f_1 and f_2 , as well as the result of conditioning f_1 on $X = x$. We call such operations **apply**, and they are the key to using knowledge compilation for doing online collapsed importance sampling. The result of multiplying two arithmetic circuits is depicted in Figure 3.

As a result of SDDs supporting the **apply** operations, we can directly compile graphical models to circuits in a bottom-up manner. Concretely, we start out by compiling each factor into a corresponding SDD representation using the encoding of Choi et al. (2013). Next, these SDDs are multiplied in order to obtain a representation for the entire model. As shown by Choi et al. (2013), this straightforward approach can be used to achieve state-of-the-art exact inference on probabilistic graphical models.

3.2 Algorithm

Now that we have proposed online collapsed importance sampling and given background on knowledge compilation, we are ready to introduce collapsed compilation, an algorithm that uses knowledge compilation to do online collapsed importance sampling. Collapsed compilation begins multi-

plying factors, and when the the SDD becomes too large, we sample and condition variables until it is sufficiently small again. At the end, the sampled variables form \mathbf{X}_p , and the variables remaining in the SDD form \mathbf{X}_d . All SDD **apply** operations are tractable (Van den Broeck & Darwiche, 2015).

Concretely, collapsed compilation repeatedly performs a few simple steps, following Algorithm 1:

1. Choose an order, and begin compiling factors until the size limit is reached
2. Select a variable X_i using the given policy π
3. Sample X_i according to its marginal in the current SDD for the partially compiled graph
4. Condition the SDD on the sampled value for X_i

The full technical description and implementation details can be found in Appendix B and C.

4 Experimental Evaluation

Data & Evaluation Criteria To empirically investigate collapsed compilation, we evaluate the performance of estimating a single marginal on a series of commonly used graphical models. Each model is followed in brackets by the count of nodes and factors.

From the 2014 UAI inference competition, we evaluate on linkage(1077,1077), Grids(100,300), DBN(40, 440), and Segmentation(228,845) problem instances. From the 2008 UAI inference competition, we use two semi-deterministic grid instances, 50-20(400, 400) and 75-26(676, 676). Here the first number indicates the number factor entries that are deterministic, and the second indicates the size of the grid. Finally, we generated a randomized frustrated Ising model on a 16x16 grid, frust16(256, 480). Beyond these seven benchmarks, we experimented on ten additional standard benchmarks. Because those were either too easy (showing no difference between collapsed compilation and the baselines), or similar to other benchmarks, we do not report on them here.

For evaluation, we run all sampling-based methods 5 times for 1 hour each. We report the median Hellinger distance across all runs, which for discrete distributions P and Q is given by $H(P, Q) = \frac{1}{\sqrt{2}} \sqrt{\sum_{i=1}^k (\sqrt{p_i} - \sqrt{q_i})^2}$. A * symbol means that the method performed exact inference.

Variable Selection Policies We will evaluate the policies detailed in Appendix C. The first policy RBVar explores the idea of picking the variable that least increases the Rao-Blackwell variance of the query (Darwiche, 2009). For a given query α , to select our next variable from \mathbf{X} , we use $\operatorname{argmin}_{X \in \mathbf{X}} \sum_X P(\alpha|X)^2 P(X)$. This can be computed in time linear in the size of the current SDD.

The next policy we look at is MinEnt, which selects the variable with the smallest entropy. Intuitively, this is selecting the variable for which sampling assumes the least amount of unknown information.

Finally, we examine a graph-based policy FD (FrontierDistance). At any given point in our compilation we have some frontier \mathcal{F} , which is the set of variables which have some but not all factors included in the current SDD. Then we select the variable in our current SDD which is, on the graph of our model, closest to the “center” induced by \mathcal{F} . That is, we use $\operatorname{argmin}_{X \in \mathbf{X}} \max_{F \in \mathcal{F}} \operatorname{distance}(X, F)$.

4.1 Understanding Collapsed Compilation

We begin our evaluation with experiments designed to shed some light on different components involved in collapsed compilation. First, we evaluate our choice in proposal distribution by comparison to marginal-based proposals. Then, we examine the effects of setting different size thresholds for compilation on the overall performance, as well as the sample count and quality.

Evaluating the Proposal Distribution Selecting an effective proposal distribution is key to successfully using importance sampling estimation (Tokdar & Kass, 2010). As discussed in Section 3, one requirement of online collapsed importance sampling is that we must provide a proposal distribution over any subset of variables. Alternatively, this means a proposal for any variable conditioned on any partial instantiation, which is challenging to provide for many traditional choices of proposal.

Table 1: Internal comparisons for collapsed compilation. Values represent Hellinger distances.

(a) Comparison of proposal distributions				(b) Comparison of size thresholds			
Policy	Dummy	True	SDD	Policy	10k	100k	1m
FD	2.37e-4	1.77e-4	3.72e-7	FD	7.33e-5	3.21e-7	7.53e-6
MinEnt	3.29e-4	1.31e-3	2.10e-8	MinEnt	1.44e-3	1.59e-5	8.07e-4
RBVar	5.81e-3	5.71e-3	7.34e-3	RBVar	2.96e-2	2.60e-2	8.81e-3

(c) Comparison of size thresholds (50 samples)				(d) Number of samples taken in 1 hour by size			
Policy	10k	100k	1m	Size Threshold	10k	100k	1m
FD	1.63e-3	5.08e-7	1.27e-6	Number of Samples	561.3	33.5	4.7
MinEnt	1.69e-2	1.84e-6	7.24e-6				
RBVar	1.94e-2	1.52e-1	3.07e-2				

To evaluate the quality of collapsed compilation’s proposal distribution, we compare it to using marginal-based proposals, and highlight the problem with such proposals. First, we compare to a dummy uniform proposal. Second, we compare to a proposal that uses the true marginals for each variable. Experiments on the 50-20 benchmark are shown in Table 1a.

Particularly with policies `FrontierDist` and `MinEnt`, the results underline the effectiveness of collapsed compilation’s proposal distribution over baselines. This is the effect of conditioning – even sampling from the true posterior marginals does not work very well, due to the missed correlation between variables. Since we are already conditioning for our partial exact inference, collapsed compilation’s proposal distribution is providing this improvement for very little added cost.

Choosing a Size Threshold A second requirement for collapsed compilation is to set a size threshold for the circuit being maintained. Setting the threshold to be infinity leaves us with exact inference which is in general intractable, while setting the threshold to zero leaves us with importance sampling using what is likely a poor proposal distribution (since we can only consider one factor at a time). Clearly, the optimal choice finds a trade-off between these two considerations.

Using benchmark 50-20 again, we compare the performance on three different settings for the circuit size threshold: 10,000, 100,000, and 1,000,000. Table 1b shows that generally, 100k gives the best performance, but the results are often similar. To further investigate this, Table 1c and Table 1d show performance with exactly 50 samples for each size, and number of samples per hour respectively. This is more informative as to why 100k gave the best performance - there is a massive difference in performance for a fixed number of samples between 10k and 100k or 1m. The gap between 100k and 1m is quite small, so as a result the increased samples for 100k leads to better performance. Intuitively, this is due to the nature of exact circuit compilation, where at a certain size point of compilation you enter an exponential regime. Ideally, we would like to stop compiling right before we reach that point. Thus, we proceed with 100k as our size-threshold setting for further experiments.

4.2 Memory-Constrained Comparison

In this section, we compare collapsed compilation to two related state-of-the-art methods: edge-deletion belief propagation (EDBP) (Choi & Darwiche, 2006), and IJGP-Samplesearch (SS) (Gogate & Dechter, 2011). Generally, for example in past UAI probabilistic inference competitions, comparing methods in this space involves a fixed amount of time and memory being given to each tool. The results are then directly compared to determine the empirically best performing algorithm. While this is certainly a useful metric, it is highly dependent on efficiency of implementation, and moreover does not provide as good of an understanding of the effects of being allowed to do more or less exact inference. To give more informative results, in addition to a time limit, we restrict our comparison at the algorithmic level, by controlling for the level of exact inference being performed.

Edge-Deletion Belief Propagation EDBP performs approximate inference by increasingly running more exact junction tree inference, and approximating the rest via belief propagation (Choi & Darwiche, 2006; Choi et al., 2005). To constrain EDBP, we limit the corresponding circuit size for the junction tree used. In our experiments we set these limits at 100,000 and 1,000,000.

Table 2: Hellinger distances across methods with internal treewidth and size bounds

Method	50-20	75-26	DBN	Grids	Segment	linkage	frust
EDBP-100k	2.19e-3	3.17e-5	6.39e-1	1.24e-3	1.63e-6	6.54e-8	4.73e-3
EDBP-1m	7.40e-7	2.21e-4	6.39e-1	1.98e-7	1.93e-7	5.98e-8	4.73e-3
SS-10	2.51e-2	2.22e-3	6.37e-1	3.10e-1	3.11e-7	4.93e-2	1.05e-2
SS-12	6.96e-3	1.02e-3	6.27e-1	2.48e-1	3.11e-7	1.10e-3	5.27e-4
SS-15	9.09e-6	1.09e-4	*	8.74e-4	3.11e-7	4.06e-6	6.23e-3
FD	9.77e-6	1.87e-3	1.24e-1	1.98e-4	6.00e-8	5.99e-6	5.96e-6
MinEnt	1.50e-5	3.29e-2	1.83e-2	3.61e-3	3.40e-7	6.16e-5	3.10e-2
RBVar	2.66e-2	4.39e-1	6.27e-3	1.20e-1	3.01e-7	2.02e-2	2.30e-3

IJGP-Samplesearch IJGP-Samplesearch is an importance sampler augmented with constraint satisfaction search (Gogate & Dechter, 2011, 2007). It uses iterative join graph propagation (Dechter et al., 2002) together with w -cutset sampling (Bidyuk & Dechter, 2007) to form a proposal, and then uses search to ensure that no samples are rejected. To constrain SS, we limit treewidth w at either 15, 12, or 10. For reference, a circuit of size 100,000 corresponds to a treewidth between 10 and 12.

Appendix D describes both baselines as well as the experimental setup in further detail.

4.2.1 Discussion

Table 2 shows the experimental results for this setting. Overall, we have found that when restricting all methods to only do a fixed amount of exact inference, collapsed compilation has similar performance to both samplesearch and EDBP. Furthermore, given a good choice of variable selection policy, it can often perform better. In particular, we highlight DBN, where we see that collapsed compilation with the RBVar or MinEnt policies is the only method that manages to achieve reasonable approximate inference. This follows the intuition discussed in Section 2.2: a good choice of a few variables in a densely connected model can lead to relatively easy exact inference for a large chunk of the model.

Another factor differentiating collapsed compilation from both EDBP and samplesearch is the lack of reliance on some type of belief propagation algorithm. Loopy belief propagation is a cornerstone of approximate inference in graphical models, but it is known to have problems converging to a good approximation on certain classes of models (Murphy et al., 1999). The problem instance frust16 is one such example - it is an Ising model with spins set up such that potentials can form loops, and the performance of both EDBP and samplesearch highlight these issues.

4.3 Probabilistic Program Inference

As an additional point of comparison, we introduce a new type of benchmark. We use the probabilistic logic programming language ProbLog (De Raedt & Kimmig, 2015) to model a graph with probabilistic edges, and then query for the probability of two nodes being connected. This problem presents a unique challenge, as every non-unary factor is deterministic.

Table 3 shows the results for this benchmark, with the underlying graph being a 12x12 grid. We see that EDBP struggles here due to the large number of deterministic factors, which stop belief propagation from converging in the allowed number of iterations. Samplesearch and collapsed compilation show similarly decent results, but interestingly they are not happening for the same reason. To contextualize this discussion, consider the stability of each method. Collapsed compilation draws far fewer samples than SS - some of this is made up for by how powerful collapsing is as a variance reduction technique, but it is indeed less stable than SS. For this particular instance, we found that while different runs for collapsed compilation tended to give different marginals fairly near the true value, SS consistently gave the same incorrect marginal. This suggests that if we ran each algorithm until convergence, collapsed compilation would tend toward the correct solution, while SS would not, and appears to have a bias on this benchmark.

Method	Prob12
EDBP-1m	3.18e-1
SS-15	3.87e-3
FD	1.50e-3

Table 3: Hellinger distances for ProbLog

5 Related Work and Conclusions

We have presented online collapsed importance sampling, an asymptotically unbiased estimator which allows for doing collapsed importance sampling without choosing which variables to collapse a priori. Using techniques from knowledge compilation, we developed collapsed compilation, an implementation of online collapsed importance sampling that draws its proposal distribution from partial compilations of the distribution, and naturally exploits structure in the distribution.

In related work, Lowd & Domingos (2010) study arithmetic circuits as a variational approximation of graphical models. Approximate compilation has been used for inference in probabilistic (logic) programs (Vlasselaer et al., 2015). Other approximate inference algorithms that exploit local structure include samplesearch and the family of universal hashing algorithms (Ermon et al., 2013; Chakraborty et al., 2014). Finally, collapsed compilation can be viewed as an approximate knowledge compilation method: each drawn sample presents a partial knowledge base along with the corresponding correction weight. This means that it can be used to approximate any query which can be performed efficiently on an SDD – for example the most probable explanation (MPE) query (Chan & Darwiche, 2006; Choi & Darwiche, 2017). We leave this as an interesting direction for future work.

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A Proof of Theorems

We begin with the formal definition of our augmented factor graph.

Definition 4. Suppose we have a discrete distribution represented by a factor graph F , with variables \mathbf{X} . Then we define the corresponding augmented factor graph F_A as follows:

- For each X_i , we introduce variables X_i^* and S_i
- For each S_i , add a unary factor such that S_i is uniform
- For each X_i^* , add a factor such that

$$P(X_i^*|S_i, X_i) = \begin{cases} 0.5, & \text{if } S_i = 0 \\ 1, & \text{if } S_i = 1, X_i = X_i^* \\ 0, & \text{otherwise} \end{cases} \quad (3)$$

A.1 Proof of Theorem 1

Proof. Consider some variable X_i together with its corresponding augmented variables X_i^* , S_i . Then examining the resulting graphical model, we see that X_i^* and S_i are only connected to the rest of the model via X_i . Due to the conditional independence properties, this means that we can first sum out S_i from $P(X_i^*|S_i, X_i)$, and then simply sum X_i^* . Thus, the result of any query over \mathbf{X} on F_A will be equivalent to the result on F , as we can simply sum out all additional variables, and end up at the same model. \square

A.2 Proof of Theorem 2

Proof. Our goal is to estimate

$$\begin{aligned} \mathbb{E}_{P(\xi)}[f(\xi)] &= \sum_{X, X^*, S} P(X, X^*, S) f(X) \\ &= \sum_{X, X^*, S} P(X) \prod_i P(S_i) P(X_i^*|X_i, S_i) \end{aligned}$$

We begin by introducing our proposal distribution $Q(X^*, S)$, and supposing we have samples $\{X^*[m], S[m]\}_{m=1}^M$ drawn from our proposal distribution Q . Now, substituting this in and continuing our previous equations gives us:

$$\begin{aligned} &\sum_{X^*[m], S[m]} \frac{1}{Q(X^*[m], S[m])} \\ &\left(\sum_X P(X) \prod_i P(S_i[m]) P(X_i^*[m]) \right) f(X) \end{aligned}$$

We can split our sum over X up into indices where $S_i[m] = 0$, and ones where $S_i[m] = 1$. For the sake of notation, we will refer to these sets of indices as I^C and I respectively:

$$\begin{aligned}
& \sum_{m=1}^M \frac{1}{Q(X^*[m], S[m])} \\
& \sum_{X_I} \prod_{i \in I} (P(S_i[m])P(X_i^*[m]|X_i, S_i[m])) \\
& \sum_{X_{I^C}} P(X) \prod_{i \in I^C} P(S_i[m])P(X_i^*[m]|X_i, S_i[m])f(X) \\
&= \sum_{m=1}^M \frac{1}{Q(X^*[m], S[m])} \\
& \sum_{X_I} \prod_{i \in I} (P(S_i[m])P(X_i^*[m]|X_i, S_i[m])) \\
& \sum_{X_{I^C}} P(X_I)P(X_{I^C}|X_I) \\
& \prod_{i \in I^C} (0.5)P(X_i^*[m]|X_i, S_i[m])f(X) \\
&= \sum_{m=1}^M \frac{1}{Q(X^*[m], S[m])} \\
& \sum_{X_I} P(X_I) \prod_{i^I} (P(S_i[m])P(X_i^*[m]|X_i, S_i[m])) \\
& \sum_{X_{I^C}} P(X_{I^C}|X_I)(0.5 * 0.5)^{|I^C|} f(X) \\
&= \sum_{m=1}^M \frac{(0.25)^{|I^C|}}{Q(X^*[m], S[m])} \sum_{X_I} P(X_I) \\
& \prod_{i \in I} (0.5)P(X_i^*[m]|X_i, S_i[m]) \sum_{X_{I^C}} P(X_{I^C}|X_I)f(X) \\
&= \sum_{m=1}^M \frac{(0.25)^{|I^C|}}{Q(X^*[m], S[m])} (0.5)^{|I|} \sum_{X_I} P(X_I) \\
& \prod_{i \in I} P(X_i^*[m]|X_i, S_i[m]) \mathbb{E}_{P(X_{I^C}|X_I)}[f(X)]
\end{aligned}$$

Now, observe that the term $\prod_{i \in I} P(X_i^*[m]|X_i, S_i[m]) = 1 \iff \forall i \text{ s.t. } S_i[m] = 1, X_i = X_i^*[m]$. Since there is only one setting of these indices that satisfies this (that is, letting $X_i = X_i^*[m]$ everywhere that $S_i[m] = 1$), this allows us to get rid of this sum, and our equation becomes:

$$\sum_{m=1}^M \frac{(0.25)^{|I^C|} (0.5)^{|I|} P_{X_I}(X_I^*[m]) \mathbb{E}_{P(X_{I^C}|X_I)}[f(X)]}{Q(X^*[m], S[m])}$$

Which is precisely what we wanted - this is the equation we would expect to use for our online collapsed importance sampler (once we adjust our proposal distribution for variables that are not actually sampled to correct for the 0.5s that are left). \square

A.3 Proof of Theorem 3

Proof. We will proceed by contradiction. Suppose we have two paths through our variables, $X_{i_1}^*, X_{i_2}^*, \dots, X_{i_n}^*$ and $X_{j_1}^*, X_{j_2}^*, \dots, X_{j_n}^*$ which can produce the same assignment to all variables. Now, there are two key facts we will observe and make use of:

1. The starting point for any path is fixed, that is $X_{i_1}^* = X_{j_1}^*$. Our heuristics are deterministic, and the memory limit remains constant, so the first variable to sampled is always the same.
2. Once an assignment is made to the current variable being sampled, the decision of which variable to sample is deterministic - again, because our heuristics must be deterministic. To put it another way, if $x_{i_k}^* = x_{j_k}^*$, then $X_{i_{k+1}}^* = X_{j_{k+1}}^*$.

Now, since path i and path j are different, there must be some first element k where $X_{i_k}^* \neq X_{j_k}^*$. By fact 1, $k > 1$. Also, observe that since k is the first such element, $X_{i_{k-1}}^* = X_{j_{k-1}}^*$. But since our 2 paths must give the same assignment to all variables, this means also that $x_{i_{k-1}}^* = x_{j_{k-1}}^*$, which means that $X_{i_k}^* = X_{j_k}^*$ by fact 2. This is a contradiction. □

B Collapsed Compilation: Algorithm Outline

Algorithm 2 describes Collapsed Compilation in more detail. Note that to avoid confusion, we hereafter refer to the processing of sampling $x_i \sim Q(X_i)$ as conditioning (since all future variables sampled are conditioned on x_i), and a single full run as a sample.

Algorithm 2: A single sample of Collapsed Compilation

Input: π : A variable selection policy computed using an SDD,
 M : A probabilistic graphical model,
 f : Function to estimate expectation for

Result: A single sample $\mathbf{X}_P^m[m]$, $\mathbb{E}_{P(\mathbf{X}_d^m | \mathbf{X}_P^m[m])}[f(\mathbf{x}_P^m[m], \mathbf{X}_d^m)]$, $w[m]$

```

1  $SDD \leftarrow True$ 
2  $\mathbf{X}_P \leftarrow \{\}$ 
3  $q \leftarrow 1$ 
4 while  $M$  is not compiled do
5    $SDD \leftarrow bottomUpCompileStep(SDD, M)$ 
6   while  $SDD$  is too large do
7      $X_j \leftarrow \pi(SDD, \mathbf{X}_P)$   $x_j \sim P_{SDD}(X_j)$ 
8      $q \leftarrow q * P_{SDD}(x_j)$ 
9      $\mathbf{X}_P \leftarrow \mathbf{X}_P \cup X_j = x_j$ 
10     $SDD \leftarrow SDD|x_j$ 
11 return  $\mathbf{X}_P$ ,  $\frac{WMC_f(SDD)}{WMC(SDD)}$ ,  $\frac{q}{WMC(SDD)}$ 
```

There are a few important things to take away here. First, if we at any point interrupt bottom-up compilation, what we will get is a complete compilation of some subset of the model. This means that on line 7, the proposal distribution P_{SDD} we are drawing from is the true distribution for X_j on some subset of M , conditioned on all previous variables.

Secondly, there are a few calls to a weighted model count function WMC on line 11. Recall that for an SDD representing a probability distribution $P(\mathbf{X})$, the weighted model count subject to a function f computes $\sum_{\mathbf{x}} P(\mathbf{x})f(\mathbf{x})$. Also, observe that the SDD we are left with when we finish compiling is representing the joint distribution $P(\mathbf{X}_d, \mathbf{X}_P = \mathbf{x}_P)$. Thus, observing that $P(\mathbf{X}_d, \mathbf{X}_P = \mathbf{x}_P) = \hat{P}(\mathbf{X}_d | \mathbf{X}_P = \mathbf{x}_P)$ we see that the weighted model count subject to f is actually $\sum_{\mathbf{x}_d} \hat{P}(\mathbf{X}_d = \mathbf{x}_d | \mathbf{X}_P = \mathbf{x}_P) f(\mathbf{x}_d, \mathbf{x}_P)$. But setting $f \equiv 1$ allows us to compute the normalization constant, meaning that $\frac{WMC_f(SDD)}{WMC(SDD)} = \mathbb{E}_{P(\mathbf{X}_d | \mathbf{X}_P = \mathbf{x}_P)}[f(\mathbf{x}_P, \mathbf{X}_d)]$.

C Collapsed Compilation: Algorithmic Details

There are many moving parts in this method, so in this section we will examine each in isolation.

C.1 Compilation Order

Once we have compiled an SDD for each factor in the graph, bottom up compilation allows us to choose in which order to multiply these SDDs. In our experiments, we look at two orders: `BFS` and `revBFS`. The first begins from the marginal query variable, and compiles outwards in a breadth first order. The second does the same, but in exactly the opposite order arriving at the query variable last.

C.2 Proposal Distribution

Given that we have decided to condition on a variable X_j , we decide its proposal distribution by computing the marginal probability of X_j in our currently compiled SDD. This can be done in time linear in the size of the circuit by computing the partial derivative of X_j with respect to the weighted model count in the current circuit (Darwiche, 2003).

C.3 Variable Selection

The manner in which we select the next variable to condition on - that is, our choice of `nextVar` in algorithm 2 - has a large effect on both the tractability and efficiency of our sampler. We will explore three policies, the first of which depends specifically on the marginal being queried for, while the other two do not. All of the policies we explore will satisfy Definition 3 - that is, they are all deterministic.

The first policy `RBVar` explores the idea of picking the variable which least increases the Rao-Blackwell variance of the query (Darwiche, 2009). For a given query α to select our next variable from \mathbf{X} we use $\operatorname{argmin}_{X \in \mathbf{X}} \sum_X P(\alpha|X)^2 P(X)$. This can be computed in time linear in the size of the current SDD.

The next policy we look at is `MinEnt`, which selects the variable with the smallest entropy. Intuitively, this is selecting the variable for which sampling it is assuming the least unknown information.

Finally, we examine a graph based policy `FrontierDist`. At any given point in our compilation we have some frontier \mathcal{F} , which is the set of variables which have some but not all factors included in the current SDD. Then we select the variable in our current SDD which is, on the graph of our model, closest to the “center” induced by \mathcal{F} . That is, we use $\operatorname{argmin}_{X \in \mathbf{X}} \max_{F \in \mathcal{F}} \operatorname{dist}(X, F)$.

In our experiments, policy `RBVar` is used with the compilation order `BFS`, while policies `MinEnt` and `FrontierDist` are used with `RevBFS`.

C.4 Determinism

A desirable property for samplers - particularly when there are a large number of deterministic relationships present in the model - is to be *rejection-free*. It is clear that in the presence of no deterministic factors (that is, no 0 entries in any factor), collapsed compilation will never reject samples. Here, we describe how this guarantee can be maintained in the presence of 0-valued factor entries.

Extracting a Logical Base Suppose we are given a factor over some variables X_1, X_2, \dots, X_n . Then a weight of 0 given to an assignment x_1, x_2, \dots, x_n indicates a logical statement. Specifically, we can say for certain over the entire model that $\neg(x_1 \wedge x_2, \dots, \wedge x_n)$. As a preprocessing step, we find all such factor entries in the graph, convert them each into the corresponding logical statement, and then take the conjunction of all of these. This forms the logical base for our model.

An Oracle for Determinism Once we have obtained this base for our model, we can precompile a circuit representing it. This allows us to make queries asking whether there is a non-zero probability that $X_j = x_j$, given all previous assignments \mathbf{X}_p (Darwiche, 1999). Thus, when we sample from

the marginal probability of X_j from our current SDD (our proposal distribution), we first renormalize this marginal to only include assignments which have a non-zero probability according to our determinism oracle. Of course, this is not always possible due to size constraints in the case where there is an enormous amount of determinism. For these cases we just run collapsed compilation as is - depending on the compilation order it will still tend to reject few samples.

Literal Entailment As a simple optimization, we can recognize any variables whose values are already deterministically chosen based on previously conditioned variables, and assign them as such in our SDD. Given a determinism oracle, deciding this can be done for all variables in the model in time linear in the size of the oracle (Darwiche, 2001).

D Experimental Details

D.1 Edge-Deletion Belief Propagation

Edge-deletion belief propagation (EDBP) is a method for doing approximate graphical model inference by using a combination of exact inference and belief propagation (Choi & Darwiche, 2006) (Choi et al., 2005). EDBP iteratively computes more and more of the model exactly using junction tree, at each step performing belief propagation to approximate the rest of the model. It can be viewed as the belief propagation analog of collapsed compilation, which makes it an interesting target for comparison. A major conceptual difference between the two is that while collapsed compilation is asymptotically unbiased and thus will give an accurate result given enough time, EDBP will tend to finish more quickly but has no way to improve once converged.

To capture a more direct comparison of the amount of exact inference being performed, we compare collapsed compilation to EDBP with the size of the junction tree used directly being limited, rather than the computer memory usage. In particular, we limit the size of the circuit corresponding to the junction tree to be similar to the sizes used for collapsed compilation. To this end, we use 100,000 and 1,000,000 as size limits for the junction tree, and otherwise run the algorithm as usual. Table 2 shows the results across all benchmarks. Keeping in mind that all policies for collapsed compilation use 100,00 as their size limit, collapsed compilation is comparable to EDBP. Both perform very well in linkage and Segment, and while collapsed compilation performs better on 50-20, EDBP does better on 75-26.

D.2 SampleSearch

IJGP-Samplesearch is an importance sampler augmented with constraint satisfaction search (Gogate & Dechter, 2011) (Gogate & Dechter, 2007). It uses iterative join graph propagation (Dechter et al., 2002) together with w -cutset sampling (Bidyuk & Dechter, 2007) to form a proposal, and then uses search to ensure that no samples are rejected.

Once again, we would like to control for the amount of exact inference being done directly at the algorithmic level, rather than via computer memory. For samplesearch, we do this by limiting w , which is the largest treewidth that is allowed when using collapsing to reduce variance. We run samplesearch with three different settings, limiting w to 15, 12, and 10 respectively. Table 2 shows the results of running our standard set of benchmarks with all of these settings. As a reference point, empirically a circuit size limit of 100,00 generally corresponds to a treewidth somewhere between 10 and 12. The results are generally similar to constraining the memory of EDBP, but with more constrained versions of samplesearch suffering more. For example, although linkage appears to be an easy instance in general, without a large enough w -cutset, samplesearch struggles compared to other methods.