Hierarchical Non-Emitting Markov Models[1]

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Abstract

We describe a simple variant of the interpolated Markov model with non-emitting state transitions and prove that it is strictly more powerful than any Markov model. More importantly, the non-emitting model outperforms the classic interpolated model on natural language texts under a wide range of experimental conditions, with only a modest increase in computational requirements. The non-emitting model is also much less prone to overfitting.

Keywords: Markov model, interpolated Markov model, hidden Markov model, mixture modeling, non-emitting state transitions, state-conditional interpolation, statistical language model, discrete time series, Brown corpus, Wall Street Journal.

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1 Introduction

The Markov model has long been the core technology of statistical language modeling. Many other models have been proposed, but none has offered a better combination of predictive performance, computational efficiency, and ease of implementation. Here we add hierarchical non-emitting state transitions to the Markov model. Although the states in our model remain Markovian, the model itself is no longer Markovian because it can represent unbounded dependencies in the state order distribution. Consequently, the non-emitting Markov model is strictly more powerful than any Markov model, including the context model [19, 20, 26], the backoff model [5, 13], and the interpolated Markov model [12, 14]. More importantly, the non-emitting model consistently outperforms the best Markov models on natural language texts, under a wide range of experimental conditions. The non-emitting model is also nearly as computationally efficient and easy to implement as the interpolated Markov model.

The remainder of our report consists of five sections and one appendix. In section 2, motivate the fundamental problem of time series prediction, which is to combine the probabilities of events of different orders. Section 3 reviews the interpolated Markov model and briefly demonstrates the equivalence of interpolated models and basic Markov models of the same model order. Next, we introduce the hierarchical non-emitting Markov model in section 4, and prove that even a second order non-emitting model is strictly more powerful than any Markov model, of any model order. Section 5 provides efficient algorithms to optimize the parameters of a non-emitting model on data. In section 6, we report empirical results for the interpolated model and the non-emitting model on the Brown corpus and Wall Street Journal. Finally, in section 7 we conjecture that the non-emitting model excels empirically because it imposes a pseudo-Bayesian discipline on maximum likelihood techniques. Appendix A reviews the backoff model and explains how to construct a non-emitting backoff model that is strictly more powerful than any backoff model.

Our notation is as follows. Let A be a finite alphabet of distinct symbols, $|A| = k$, and let $x^T \in A^T$ denote an arbitrary string of length T over the alphabet A. Then x_i^j denotes the substring of x^T that begins at position i and ends at position j. For convenience, we abbreviate the unit length substring x_i^i as x_i and the length t prefix of x^T as x^t.

2 Time Series Prediction

A time series model must assign accurate probabilities to strings of unbounded length. Yet unbounded strings don't occur in recorded histories, which are always finite. Therefore, to estimate the probabilities of unbounded strings from a finite corpus, we must assume that each symbol in a given string depends only on a finite number of (equivalence classes of) contexts. The most widely

adopted independence assumption is the order n Markov assumption, which states that each symbol depends only on the immediately preceding n symbols, and is conditionally independent of the distant past.

$$\begin{aligned} p(x^T|T) &= \prod_{t=1}^{T} p(x_t|x^{t-1}) \\ &\approx \prod_{t=1}^{T} p(x_t|x_{t-n}^{t-1}) \end{aligned}$$

The simplest statistical model to incorporate an order n Markov assumption is the basic Markov model. A basic Markov model $\phi = \langle A, n, \delta_n \rangle$ consists of an alphabet A, a model order n, $n \geq 0$, and the state transition probabilities $\delta_n : A^n \times A \rightarrow [0, 1]$. With probability $\delta_n(y|x^n)$, a Markov model in the state x^n will emit the symbol y and transition to the state $x_2^n y$. Therefore, the probability $p_m(x_t|x^{t-1}, \phi)$ assigned by an order n basic Markov model ϕ to a symbol x^t in the history x^{t-1} depends only on the last n symbols of the history.

$$p_m(x_t|x^{t-1}, \phi) = \delta_n(x_t|x_{t-n}^{t-1}) \tag{1}$$

Since the Markov model contains only a finite number of parameters, it is in principle possible to estimate their values directly from data. All that remains is to choose the model order.

In real-world time series problems, the future depends on the entire past, even if only weakly. In order to more closely approximate a real-world source, we would like our model order to be as large as possible. Yet we have only a finite amount of training data from which to estimate our model parameters. An order n Markov model over an alphabet of k symbols has k^{n+1} events, while a corpus of length T has at most $T - n$ distinct events of order n. The exponential growth in events quickly exceeds the size of all available training data, and nearly all the higher-order events do not occur in the training data.

This tension between model complexity and data sparsity is fundamental to time series modeling. The probabilities of the lower order events can be more accurately estimated from the available training data, while the higher order events are better able to model complex real-world sources. An effective model, then, must include individual events of both higher and lower orders.

The two most widely-used techniques for combining individual events of varying orders are backoff and interpolation. In an interpolated model, the transition probabilities from lower and higher order states are combined stochastically using mixing parameters. In a backoff model, the event probabilities are combined according to a partial order which typically favors higher order events over lower order events. In section 3 and appendix A, we show that backoff models and interpolated models are formally equivalent to basic Markov models. Therefore, backoff and interpolation are simply parameter estimation schemes for basic Markov models.

3 Interpolation

Here we introduce the interpolated Markov model and explain why the interpolated model class is equivalent to the class of basic Markov models. In the next section 4, we introduce hierarchical non-emitting state transitions to the Markov model, and prove that the new non-emitting models are no longer Markovian even though their states are.

In the interpolated Markov model, the transition probabilities from states of different orders are combined using state-conditional mixing parameters. The mixing parameters smooth the transition probabilities from higher order states with those from lower order states [12]. Mixing the transition probabilities from states of different orders results in more accurate predictions than can be obtained from any fixed model order.

Formally, an interpolated Markov model $\phi = \langle A, n, \delta, \lambda \rangle$ consists of a finite alphabet A, a maximal model order n, the state transition probabilities $\delta = \delta_0 \ldots \delta_n$, $\delta_i : A^i \times A \to [0,1]$, and the state-conditional interpolation parameters $\lambda : A^n \times [0,n] \to [0,1]$. The state order is a hidden variable. The probability assigned by an interpolated model is a linear combination of the probabilities assigned by all the lower order Markov models.

$$p_c(y|x^n, \phi) = \sum_{i=0}^{n} \delta_i(y|x^i)\lambda(i|x^n) \tag{2}$$

An interpolated model is a valid probability model if every $\delta_i(\cdot|x^i)$ and every $\lambda(i|x^n)$ is valid. It is nonzero for all strings A^* if $\delta_0(\cdot)$ is strictly positive for all symbols A and no $\lambda(i|x^n)$ is unity when $\delta(\cdot|x^i)$ is zero for some symbol.

Estimating the $O(nk^n)$ state interpolation probabilities is considerably easier than estimating the $O(k^{n+1})$ state transition probabilities in an order n Markov model. To begin with, we set $\lambda(i|x^n)$ to 0 if the order i state x^i is novel. Now we need only to estimate the $O(nT)$ interpolation parameters that have been observed in the training data.

Nonetheless, there are still too many interpolation parameters to be accurately estimated. Further refinements are necessary to improve predictive performance. One refinement is to group similar parameters into equivalence classes and then constrain them to take the same values. This is called parameter tying. At one extreme, each state-conditional interpolation distribution is its own equivalence class. At the other extreme, all interpolation probabilities are tied together and we have the state-independent interpolated Markov model

$$p_c(y|x^n, \phi) = \sum_{i=0}^{n} \delta_i(y|x^i)\lambda_i \tag{3}$$

with only $n+1$ interpolation parameters. While parameter tying can improve performance, reducing state-conditional interpolation to state-independent interpolation results in poor performance.

A hierarchical parameterization of the full state-conditional interpolation is more effective. Let $\lambda_i : A^i \rightarrow [0,1]$ be the set of i^{th} order state interpolation parameters, where $\lambda_i(x^i)$ is the probability of using the i^{th}order state transition probability $\delta_i(\cdot|x^i)$, conditioned on the decision not to use any higher order state transition probability.

$$\lambda(i|x^n) = \lambda_i(x^n_{n+1-i}) \prod_{j=i+1}^{n} (1 - \lambda_j(x^n_{n+1-j}))$$

Then the probability $p_c(y|x^n, \phi)$ that the state x^n will emit the symbol y has a particularly simple form

$$\begin{aligned} p_c(y|x^i, \phi) &= \lambda_i(x^i)\delta_i(y|x^i) \\ &\quad +(1 - \lambda_i(x^i))p_c(y|x^i_2, \phi) \end{aligned} \tag{4}$$

where $\lambda_i(x^i) = 0$ for $i \geq n$, and therefore $p_c(x_t|x^{t-1}, \phi) = p_c(x_t|x^{t-1}_{t-n}, \phi)$, ie., the prediction depends only on the last n symbols of the history.

A quick glance at the form of (2) and (1) reveals the fundamental simplicity of the interpolated Markov model. Every interpolated model is equivalent to a basic Markov model of the same order, and every basic Markov model is an interpolated model of the same order. We may convert an interpolated model ϕ into a basic model ϕ' of the same model order n, simply by setting $\delta'_n(y|x^n)$ equal to $p_c(y|x^n, \phi)$ for all states $x^n \in A^n$ and symbols $y \in A$. Thus, the class interpolated Markov models is extensionally equivalent to the class of basic Markov models.

4 Non-Emitting Transitions

In the previous section, we explained how to combine events of varying orders using interpolation and backoff. Interpolation and backoff both use the probabilities of lower events to estimate the probabilities of higher order events. As a result, interpolated and backoff models are extensionally equivalent to each other and to basic Markov models of the same order. In this section, we explain how to combine events of varying orders using non-emitting state transitions.

The central idea is to allow actual non-emitting transitions between events of different orders. Unlike interpolation and backoff, non-emitting transitions are not merely an estimation method – they actually increase the expressive power of the model class. As a result, non-emitting models are strictly more powerful than the class of basic Markov models. The next section 5 provides efficient algorithms to evaluate the probability of a string according to a non-emitting model and to optimize the parameters of a non-emitting model on data.

A non-emitting mixture Markov model $\phi = \langle A, n, \delta, \lambda \rangle$ consists of a finite alphabet A, a maximal model order n, the emitting state transition probabilities $\delta_i : A^i \times A \rightarrow [0,1]$, and the non-emitting state transition probabilities λ_i :

$A^i \times [0,n] \rightarrow [0,1]$. The non-emitting model alternates between non-emitting and emitting transitions according to the λ and δ parameters, respectively. The parameter $\lambda(j|x^i)$ specifies the probability that the model will transition from the state x^i to the state x^j without emitting a symbol. The parameter $\delta_j(y|x^j)$ specifies the probability that the model will emit the symbol y from the state x^j and transition to the successor state $x^j y$. Then the probability $p_\epsilon(y^j|x^i,\phi)$ assigned to a string y^j in the state x^i has the form

$$p_\epsilon(y^j|x^i,\phi) = \sum_{l=0}^{i} \lambda(l|x^i)\delta_l(y_1|x^l)p_\epsilon(y_2^j|x^l y,\phi). \tag{5}$$

When the model order is sufficiently high, then a hierarchical parameterization of the non-emitting transition probabilities may improve performance. With probability $1-\lambda_i(x^i)$, a hierarchical non-emitting model will transition from the state x^i to the state x_2^i without emitting a symbol. With probability $\lambda_i(x^i)\delta_i(y|x^i)$, the model will transition from the state x^i to the state $x^i y$ and emit the symbol y.

Therefore, the probability $p_\epsilon(y^j|x^i,\phi)$ assigned to a string y^j in the history x^i by a hierarchical non-emitting model ϕ has the recursive form (6),

$$\begin{aligned} p_\epsilon(y^j|x^i,\phi) \quad &= \quad \lambda_i(x^i)\delta_i(y_1|x^i)p_\epsilon(y_2^j|x^i y_1,\phi) \\ &\quad +(1-\lambda_i(x^i))p_\epsilon(y^j|x_2^i,\phi) \end{aligned} \tag{6}$$

where $\lambda_i(x^i)=0$ for $i>n$ and $\lambda_0(\epsilon)=1$. Note that, unlike the basic Markov model, $p_\epsilon(x_t|x^{t-1},\phi) \neq p_\epsilon(x_t|x_{t-n}^{t-1},\phi)$ because the state order distribution of the non-emitting model depends on the prefix x^{i-n}. This simple fact will allow us to establish that there exists a non-emitting model that is not equivalent to any Markov model.

Lemma 4.1 states that there exists a non-emitting model ϕ that cannot be converted into an equivalent basic model of any order. There will always be a string x^T that distinguishes the non-emitting model ϕ from any given basic model ϕ' because the non-emitting model can encode unbounded dependencies in its state distribution.

Lemma 4.1 $\exists\phi\ \forall\phi'\ \exists x^T \in A^*\ [p_\epsilon(x^T|\phi,T) \neq p_m(x^T|\phi',T)]$

Proof. The idea of the proof is that our non-emitting model will encode the first symbol x_1 of the string x^T in its state distribution, for an unbounded distance. This will allow it to predict the last symbol x_T using its knowledge of the first symbol x_1. The basic model will only be able predict the last symbol x_T using the preceding n symbols, and therefore when T is greater than n, we can arrange for $p_\epsilon(x^T|\phi,T)$ to differ from any $p_m(x^T|\phi',T)$, simply by our choice of x_1.

The smallest non-emitting model capable of exhibiting the required behavior has order 2. Lower order non-emitting models are equivalent to interpolated

models of the same order, with the same parameters. The non-emitting transition probabilities λ and the interior of the string x_2^{T-1} will be chosen so that the non-emitting model is either in an order 2 state or an order 0 state, with no way to transition from one to the other. The first symbol x_1 will determine whether the non-emitting model goes to the order 2 state or stays in the order 0 state. No matter what probability the basic model assigns to the final symbol x_T, the non-emitting model can assign a different probability by the appropriate choice of x_1, $\delta_0(x_T)$, and $\delta_2(x_T|x_{T-2}^{T-1})$.

Consider the second order non-emitting model over a binary alphabet with $\lambda(0) = 1$, $\lambda(1) = 0$, and $\lambda(11) = 1$ on strings in $A1^*A$. When $x_1 = 0$, then x_2 will be predicted using the 1st order model $\delta_1(x_2|x_1)$, and all subsequent x_t will be predicted by the second order model $\delta_2(x_t|x_{t-2}^{t-1})$. When $x_1 = 1$, then all subsequent x_t will be predicted by the zeroth order model $\delta_0(x_t)$. Thus for all $t > p$, $p_\epsilon(x_t|x^{t-1}) \neq p_\epsilon(x_t|x_{t-p}^{t-1})$ for any fixed p, and no basic model is equivalent to this simple non-emitting model. $\square$

Every basic model is a non-emitting model, with the appropriate choice of non-emitting transition probabilities.

Lemma 4.2 $\forall\phi\ \exists\phi'\ \forall x^T \in A^*\ [p_\epsilon(x^T|\phi', T) = p_m(x^T|\phi, T)]$

Proof. A basic model $\phi = \langle A, n, \delta_n\rangle$ is equivalent to a non-emitting model $\phi' = \langle A, n, \delta', \lambda'\rangle$ where $\delta'_n = \delta_n$ and $\lambda'(n|x^n) = 1$ for all x^n. In the hierarchical parameterization, $\lambda'(x^n) = 1$ for all x^n. $\square$

Therefore, the class $\mathcal{P}_\epsilon$ of non-emitting Markov distributions is strictly more powerful than the class $\mathcal{P}_m$ of basic Markov distributions.

Theorem 1 $\mathcal{P}_m \subset \mathcal{P}_\epsilon$

Proof. $\mathcal{P}_m \neq \mathcal{P}_\epsilon$ by lemma 4.1 and $\mathcal{P}_m \subseteq \mathcal{P}_\epsilon$ by lemma 4.2. $\square$

Since interpolated models and backoff models are equivalent to basic Markov models, we have as a corollary that non-emitting Markov models are strictly more powerful than interpolated and backoff models. Note that non-emitting Markov models are considerably less powerful than the full class of stochastic finite state automata because their states are Markovian. For the same reason, non-emitting models are also less powerful than the full class of hidden Markov models.

Let us now turn to the algorithms required to evaluate the probability of a string according to a non-emitting mixture model and to optimize the non-emitting state transitions on a training corpus.

5 Estimation

Here we present an efficient expectation-maximization (EM) algorithm to optimize the parameters of a hierarchical non-emitting mixture model on data.

An EM algorithm iteratively maximizes the probability of the training data according to the model by computing the expectation of model parameters on the data and then updating the model parameters to maximize those expectations [2, 3, 6].

The non-emitting mixture model is sufficiently expressive that any maximum likelihood estimator will overfit its parameters to the training corpus. Unseen events will be assigned zero probability, and the overfit model will fail to accurately predict the future. The traditional solution to this problem for interpolated Markov models is cross-estimation [12]. Cross-estimation repeatedly partitions the training data into two blocks and optimizes the mixing parameters on one block after initializing the state transition parameters on the other block. We present a traditional cross-estimation algorithm for hierarchical non-emitting models.

We begin by partitioning the training corpus into a fixed set of blocks $\mathbf{B}$. Ideally our partition is linguistically meaningful and roughly uniform, but neither condition is essential. For example, we might divide a natural language text corpus on sentence, paragraph, or article boundaries. Next we call CROSS-ESTIMATE($\mathbf{B}$,ϕ) on our hierarchical non-emitting model ϕ.

```
CROSS-ESTIMATE(B,φ)
1.  Until convergence
2.      Initialize λ+, λ− to zero;
3.      For each block B_i in B
4.          Initialize δ using B − B_i;
5.          EXPECTATION-STEP(B_i,φ,λ+,λ−);
6.      MAXIMIZATION-STEP(φ,λ+,λ−);
7.  Initialize δ using B;
```

The variables $\lambda^+(x^i)$ and $\lambda^-(x^i)$ accumulate expectations for the non-emitting state transition parameter $\lambda(x^i)$. $\lambda^+(x^i)$ contains the expectation of emitting a symbol in state x^i, conditioned on being in state x^i, while $\lambda^-(x^i)$ contains the expectation of transitioning to x_2^i without emitting a symbol, conditioned on being in state x^i. Lines 3-5 enumerate all one-block partitions of the training corpus. The emitting state transitions δ are initialized to their maximum likelihood estimates on the larger block $\mathbf{B} - B_i$ and then the non-emitting state transitions λ are optimized on the smaller "withheld" block B_i.

The heart of the algorithm is the EXPECTATION-STEP() procedure, which calculates the expectation of the non-emitting transitions on the string x^b and then increments the λ^+, λ^- accumulators.

```
EXPECTATION-STEP(x^b,φ,λ+,λ−)
1.  α = FORWARD(x^b,φ);
2.  β = BACKWARD(x^b,φ);
3.  for t = b downto 1
```

4. for $i = 1$ upto $\min(n, t)$
5. $\lambda_t^-(i) = \alpha_t(i)(1 - \lambda_t(i))\beta_t(i-1)$;
6. $\lambda_{t-1}^+(i-1) += \alpha_{t-1}(i-1)\lambda_{t-1}(i-1)\delta_{t-1}(i-1)\beta_t(i)$;
7. if $(t > n)$ [$\lambda_{t-1}^+(n) += \alpha_{t-1}(n)\lambda_{t-1}(n)\delta_{t-1}(n)\beta_t(n)$;]

The forward variable $\alpha_t(i)$ contains the probability $p(x^t, o_t = i|\phi)$ that the model ϕ generated the prefix x^t and terminated in the order i state. The backward variable $\beta_t(i)$ contains the probability $p(x_{t+1}bT|x^t, o_t = i, \phi)$ that the model ϕ generated the suffix x_{t+1}^b given that it was in the order i state at time t. To simplify the notation, we define $\lambda_t(i)$ to be the probability $\lambda(x_{t+1-i}^t)$ of emitting a symbol from the i^{th} order state at time t, given that we are in that state. We also define $\delta_t(i)$ to be the probability $\delta_i(x_{t+1}|x_{t+1-i}^t)$ of the emitting transition from state x_{t+1-i}^t to state x_{t+1-i}^{t+1}.

The EXPECTATION-STEP() algorithm requires $O(nb)$ time and space for an order n non-emitting model on a string x^b of length b. A comparable interpolated model can take an expectation step in $O(nb)$ time and $O(1)$ space [1]. While the difference between $O(nb)$ and $O(1)$ space can be considerable, the additional space requirements of the non-emitting algorithm are small when compared to the cost of storing all the model parameters. An order n mixture model has $O(nT)$ parameters for a training corpus of size T, and the training corpus is typically an order of magnitude larger than the withheld block.

FORWARD(x^T,ϕ)
1. $\alpha_0(0) = 1$;
2. for $t = 1$ upto $T - 1$
3. for $i = \min(n-1, t)$ downto 0
4. $\alpha_t(i) += \alpha_t(i+1)(1 - \lambda_t(i+1))$;
5. $\alpha_{t+1}(i+1) := \alpha_t(i)\lambda_t(i)\delta_t(i)$;
6. if $(t \geq n)$ [$\alpha_{t+1}(n) += \alpha_t(n)\lambda_t(n)\delta_t(n)$;]
7. return(α);

BACKWARD(x^T,ϕ)
1. for $i = 0$ upto $\min(n-1, T-1)$;
2. $\beta_{T-1}(i) = \lambda_{T-1}(i)\delta_{T-1}(i)$;
3. if $(T > n)$ [$\beta_{T-1}(n) = \lambda_{T-1}(n)\delta_{T-1}(n)$;]
4. for $t = T - 1$ downto 1
5. for $i = 1$ upto $\min(n, t)$
6. $\beta_t(i) += (1 - \lambda_t(i))\beta_t(i-1)$;
7. $\beta_{t-1}(i-1) = \lambda_{t-1}(i-1)\delta_{t-1}(i-1)\beta_t(i)$;
8. if $(t > n)$ [$\beta_{t-1}(n) = \lambda_{t-1}(n)\delta_{t-1}(n)\beta_T(n)$;]
9. return(β);

The FORWARD() and BACKWARD() algorithms each require $O(nT)$ time and space. It is possible to evaluate the probability $p_\epsilon(x^T|\phi)$ of a string x^T according to an order n non-emitting model ϕ in $O(nT)$ time and $O(n)$ space. In contrast,

it is possible to evaluate the probability $p_c(x^T|\phi)$ according to an interpolated model in $O(nT)$ time and $O(1)$ space. Again, the small additional cost in space is negligible when compared to the cost of storing the model parameters.

Having done all the work in the expectation step, the maximization step is straightforward.

MAXIMIZATION-STEP(ϕ,λ^+,λ^-)
1. Forall states x^i in $A^{\leq n}$
2. $\quad\lambda(x^i) := \lambda^+(x^i)/(\lambda^+(x^i) + \lambda^-(x^i))$;

Line 2 reestimates each non-emitting state transition parameter $\lambda(x^i)$ as the expectation of emitting a symbol from that state divided by the expectation of being in that state. In order to ensure that no non-emitting state transition parameter $\lambda(x^i)$ is ever reestimated to 0 or 1, we typically initialize each accumulator to a small positive number (eg., 0.1) instead of zero.

When λ parameters are tied, then their λ^+ and λ^- expectations must be pooled before they are updated. Let $\tau(x^i)$ be the equivalence class of x^i under the tying scheme τ. For simplicity, imagine $\tau(x^i)$ to be an index. All algorithms in this section would use the tied parameter $\lambda(\tau(x^i))$ instead of the untied parameter $\lambda(x^i)$. The TIED-EXPECTATION-STEP() algorithm would increment the $\lambda^+(\tau(x^i))$ and $\lambda^-(\tau(x^i))$ accumulators, and the TIED-MAXIMIZATION-STEP() algorithm would be as follows.

TIED-MAXIMIZATION-STEP(ϕ,λ^+,λ^-,π)
1. Forall classes i in $\tau(A^{\leq n})$
2. $\quad\lambda(i) := \lambda^+(i)/(\lambda^+(i) + \lambda^-(i))$;

In some situations, cross-estimation may be approximated by forward-estimation. Like cross-estimation, forward-estimation initializes the δ parameters on one text block and optimizes the λ parameters on another block. Forward-estimation uses only a single text partition whereas cross-estimation uses all one-block text partitions. As result, forward-estimation is considerably faster than cross-estimation, both in the amount of time required per iteration and in the number of iterations until convergence. Unfortunately, it can lead to inferior results when there are too many mixing parameters.

FORWARD-ESTIMATE(B_δ,B_λ,ϕ)
1. Until convergence
2. $\quad$Initialize λ^+, λ^- to zero;
3. $\quad$Initialize δ using B_δ;
4. $\quad$EXPECTATION-STEP(B_λ,ϕ,λ^+,λ^-);
5. $\quad$MAXIMIZATION-STEP(ϕ,λ^+,λ^-);
6. Initialize δ using $B_\delta \cup B_\lambda$;

Implementation Note. Unless the corpus and the alphabet size are very small, then the $\alpha_t(i)$ and $\beta_t(i)$ values used in the EXPECTATION-STEP() procedure will exceed the representational range of double precision IEEE floating point numbers. When this happens, a floating point exception will occur and an alternate representation must be used for the probability values. The simplest approach is to use a logarithmic representation. Multiplication and division of probability values is straightforward in a logarithmic representation.

$$\begin{aligned} \log(x \cdot y) &= \log(x) + \log(y) \\ \log(x/y) &= \log(x) - \log(y) \end{aligned}$$

Addition of logarithmic probability values is more costly, and care must be taken to avoid underflow.

$$\log(x+y) = \begin{cases} \log(x) & \text{if } \log(y) - \log(x) \leq \Lambda \\ \log(x) + \log(1 + \exp(\log(y) - \log(x))) & \text{otherwise} \end{cases}$$

Here Λ is the smallest representable exponent, for example, -707.7 for IEEE double precision floating point numbers when the logarithms are natural (ie., base e). This test is necessary to avoid underflow in the call to exp().

While it is simple to implement, logarithmic arithmetic can be 15-50 times slower than straight probability arithmetic, depending on the speed of the floating point unit and the math library provided with the operating system. For this reason, our implementation used an extended exponent representation from the library of practical abstractions [24]. This `balanced_t` module provides single precision floating point numbers with 32 bit exponents. It is 1.5 to 3.0 times faster than the logarithmic representation, depending on the machine.

When computation time is at a premium, then the most effective solution is to periodically scale the probability values in the $\alpha_t(i)$ and $\beta_t(i)$ arrays to keep them in an acceptable range. Scaling is more difficult to implement than logarithmic arithmetic or `balanced_t` arithmetic, and it is inherently nonmodular.

6 Empirical Results

The ultimate measure of a statistical model is its predictive performance in the domain of interest. To take the true measure of non-emitting models for natural language texts, we evaluate their performance as character models on the Brown corpus [7] and as word models on the Wall Street Journal. Our results show that the non-emitting Markov model consistently gives better predictions than the traditional interpolated Markov model under equivalent experimental conditions. In all cases we compare non-emitting and interpolated models of identical model orders, with the same number of parameters. Note that the non-emitting bigram and the interpolated bigram are equivalent.

Corpus	*Alphabet*	*Size*	*Blocks*
Brown	90	6,004,032	21
WSJ 1989	20,293	6,219,350	22
WSJ 1987-89	20,092	42,373,513	152

All λ values were initialized uniformly to 0.5 and then optimized using cross-estimation on the first 90% of each corpus. The remaining 10% percent of each corpus was used to evaluate model performance. While this validation paradigm exposes the models to nonstationarity, it is simple to understand and easily reproduced.

We consider a single parameter tying scheme, in which all states with the same frequency and diversity are considered equivalent. The frequency $c(x^i)$ of a state is the number of times that the string x^i occurred in the training corpus. The diversity $q(x^i) \doteq |\{y : c(x^i y) > 0\}|$ of a state is the number of distinct symbols observed in the state. Experience with multinomial prediction suggests that frequency and diversity are necessary to accurately estimate the likelihood of novel symbols [21].

In related work [25], Thomas compares the performance of the interpolated and non-emitting models on the Brown corpus and Wall Street Journal with ten different parameter tying schemes. His experiments confirm that some parameter tying schemes improve model performance, although to a lesser degree when cross-estimation is used. The non-emitting model consistently outperformed the interpolated model on both corpora for all ten parameter tying schemes. Thomas shows that our frequency-diversity parameter tying scheme is one of the more effective parameter schemes.

6.1 Brown Corpus

Our first set of experiments were with character models on the Brown corpus [7]. The Brown corpus is an eclectic collection of English prose, containing 6,004,032 characters partitioned into 500 files. We performed 10 iterations of cross estimation on 21 blocks. Results are reported as per-character test message entropies (bits/char), $-\frac{1}{v}\log_2 p(y^v|v)$. The non-emitting model outperforms the interpolated model for all nontrivial model orders, particularly for larger model orders. The non-emitting model is considerably less prone to overfitting. After 10 EM iterations, the untied order 9 non-emitting model scores 1.996 bits/char while the untied order 9 interpolated model scores 2.334 bits/char. The untied non-emitting model even outperforms the *tied* interpolated model for all nontrivial model orders.

Model order	Interpolation untied	Interpolation tied	Non-Emitting untied	Non-Emitting tied
1	3.602	3.602	3.602	3.602
2	2.950	2.950	2.946	2.946
3	2.490	2.486	2.473	2.473
4	2.231	2.218	2.193	2.192
5	2.149	2.112	2.076	2.075
6	2.164	2.082	2.031	2.027
7	2.212	2.077	2.015	2.008
8	2.277	2.084	2.010	2.000
9	2.334	2.093	2.009	1.996

We also compared the performance of our techniques with two new interpolation schemes recently proposed by Potamianos and Jelinek [16]. Their DI-TD scheme uses hierarchical state-conditional interpolation $\lambda(x^i)$, variable-width frequency $\times$ order parameter tying, and "top-down optimization" on one withheld block. Their DI-BU scheme uses general state-conditional interpolation $\lambda(j|x^i)$, variable-width frequency $\times$ order parameter tying, and bottom-up optimization on one withheld block. The comparison is performed on a modified version of the Brown corpus, which they provided to us. This modified corpus eliminates the unusual punctuation of the original Brown corpus, reduces the alphabet size from 90 to 79, and separates distinct linguistic tokens with single spaces.

Corpus	*Alphabet*	*Size*	*Train*	*Test*	*Blocks*
Brown (std)	90	6,004,032	5,403,629	600,403	21
Brown (JHU)	79	6,093,662	5,607,270	486,392	21

Another difference between the Potamianos-Jelinek validation paradigm and ours lies in how the corpus is partitioned into training and testing blocks. In our experiments, the test block was the last 10% of the Brown corpus – the last 428 characters from br-n14.txt plus all files from br-n15.txt through br-r09.txt inclusive. In the Potamianos-Jelinek experiments, the test block consisted of complete sentences chosen uniformly from the entire (modified) Brown corpus.

To this comparison, we added the original interpolation schemes of Jelinek and Mercer [12] under 10 iterations of forward-estimation (DI-FE) and cross-estimation (DI-CE). Both models used hierarchical state-conditional interpolation $\lambda(x^i)$ and straight frequency $\times$ diversity parameter tying. We also added the hierarchical non-emitting model with straight frequency $\times$ diversity parameter tying, and 10 iterations of forward-estimation (NE-FE) and cross-optimization (NE-CE). The results are summarized in the following table as mean test message entropies (bits/char).

Model order	Interpolation				Non-Emitting	
	DI-TD	DI-BU	DI-FE	DI-CE	NE-FE	NE-CE
1	3.470	3.470	3.478	3.478	3.478	3.478
2	2.851	2.850	2.860	2.858	2.857	2.856
3	2.328	2.326	2.337	2.331	2.328	2.324
4	2.016	2.007	2.012	2.007	1.996	1.991
5	1.894	1.878	1.872	1.867	1.849	1.843
6	1.853	1.831	1.820	1.815	1.789	1.782
7	1.837	1.811	1.804	1.800	1.761	1.754
8	1.828	1.801	1.800	1.796	1.746	1.739
9	1.824	1.796	1.802	1.798	1.738	1.730

The non-emitting model consistently outperforms all interpolation schemes at all model orders above 2, by a significant margin. The original Jelinek-Mercer interpolation scheme also tends to outperform the two new DI-TD and DI-BU schemes at higher model orders, for both forward-estimation (DI-FE) and cross-estimation (DI-CE).

Note also that the best order 9 result in the Potamianos-Jelinek paradigm (1.730 bits/char) is considerably better than the best order 9 result in our validation paradigm (1.996 bits/char). We believe this is partially attributable to the reduced alphabet size of the modified corpus, and principally due to the difference in the two train-test partitions. The prediction problem posed by our paradigm is more difficult because the last 10% of the Brown files are appreciably different than the first 90% of the files.

6.2 WSJ 1989

The second set of experiments was on the 1989 Wall Street Journal corpus, which contains 6,219,350 words. Our vocabulary consisted of the 20,293 words that occurred at least 10 times in the entire WSJ 1989 corpus. All out-of-vocabulary words were mapped to a unique OOV symbol. We performed 10 iterations of cross estimation on 22 blocks. Following standard practice in the speech recognition community, results are reported as per-word test message perplexities $p(y^v|v)^{-\frac{1}{v}}$. The perplexity represents the effective alphabet size. Again, the non-emitting model outperforms the interpolated model for all nontrivial model orders, even without parameter tying.

Model order	Interpolation untied	Interpolation tied	Non-Emitting untied	Non-Emitting tied
1	175.2	174.9	175.2	174.9
2	123.7	122.8	119.6	119.0
3	121.3	119.0	111.9	111.1
4	123.0	117.2	110.6	109.5
5	124.5	116.3	110.4	109.0

6.3 WSJ 1987-89

The third set of experiments was on the 1987-89 Wall Street Journal corpus, which contains 42,373,513 words. Our vocabulary consisted of the 20,092 words that occurred at least 63 times in the entire WSJ 1987-89 corpus. Again, all out-of-vocabulary words were mapped to a unique OOV symbol. We performed 10 iterations of cross estimation on 152 blocks. Results are reported as test message perplexities. As with the WSJ 1989 corpus, the non-emitting model outperforms the interpolated model for all nontrivial model orders, even without parameter tying.

Model order	Interpolation untied	Interpolation tied	Non-Emitting untied	Non-Emitting tied
1	150.7	150.7	150.7	150.7
2	94.0	93.9	92.1	92.1
3	89.2	88.6	83.2	83.2

6.4 Posthoc Analysis

In order to understand the striking empirical advantage of the non-emitting model over the interpolated model, we conducted the following experiment. We induced order 9 interpolated and non-emitting models from the Brown corpus using forward estimation with no parameter tying. This configuration was chosen to maximize the performance difference between the two models. The resulting interpolated model predicts the Brown test corpus with 2.4480 bits/char while the resulting non-emitting model predicts the Brown test corpus with 2.1536 bits/char. Figure 1 visualizes the overwhelming advantage of the non-emitting model in this experiment.

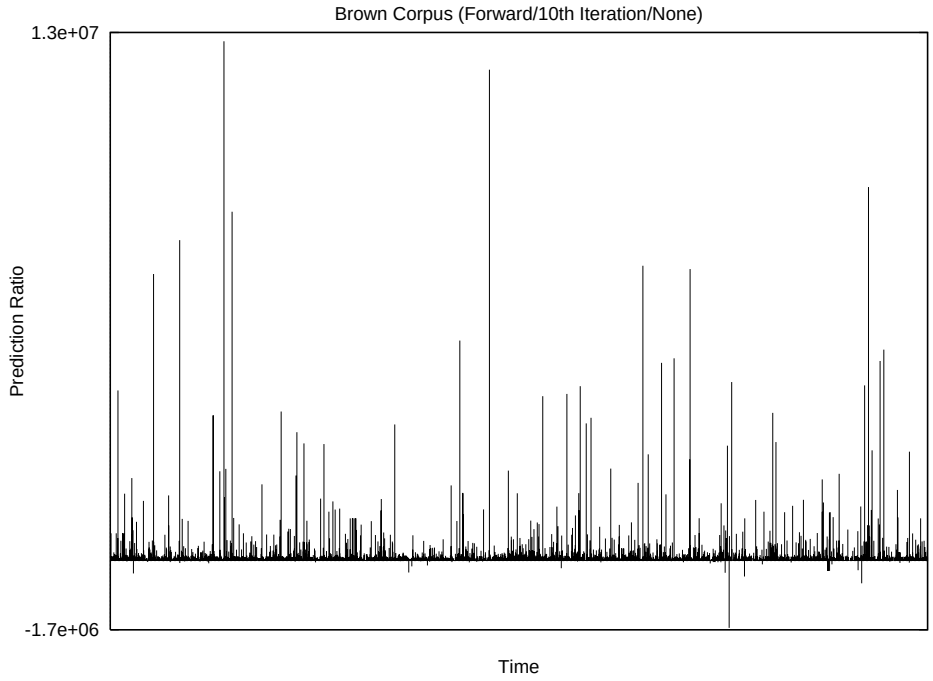

Figure 1: The relative performance of the non-emitting and interpolated models through the Brown test corpus on a character by character basis. The horizontal access represents time through the test corpus, proceeding from left to right. The vertical access represents the ratio $p_\epsilon(x_t|x^{t-1})/p_c(x_t|x^{t-1})$ when $p_\epsilon(x_t|x^{t-1}) > p_c(x_t|x^{t-1})$ and the ratio $p_c(x_t|x^{t-1})/p_\epsilon(x_t|x^{t-1})$ otherwise. Thus, each positive spike represents a win for the non-emitting model, and each negative spike represents a win for the interpolated model. The magnitude of the win is encoded by the magnitude of the spike. The largest win for the non-emitting model is a factor of 12,784,400 (23.61 bits). The largest win for the interpolated model is a factor of 1,654,650 (20.66 bits).

The following table shows the mean state order occupancy statistics for the two models on the Brown corpus.

Order	Interpolated	Non-Emitting
9	0.133	0.070
8	0.120	0.090
7	0.131	0.127
6	0.147	0.170
5	0.147	0.195
4	0.130	0.173
3	0.095	0.108
2	0.058	0.047
1	0.027	0.013
0	0.011	0.003
	5.639	5.357

As might be expected, the interpolated model spends more time than the non-emitting model in the higher order states (orders 7-9). It is arguably more surprising, however, that the interpolated model also spends more time in the lower order states (orders 0-2).

One of the smaller positive spikes in the figure 1 occurs when both models try to predict the space ␣ that follows the string `,␣but␣now␣Keith` in the Brown test corpus. Unfortunately, the string `␣Keith` does not occur in the training corpus. Nonetheless, the non-emitting model assigns 209 times more probability than the interpolated model to the event that a space will follow the string `␣Keith`. According to the non-emitting model, a space will follow the string `␣Keith` with probability 0.627. The interpolated model assigns probability 0.003 to the same event.

The reason is somewhat subtle. On the training corpus, the string `eith` is followed by the letter `e` with near certainty (0.9973). As a result, λ(`eith`) approaches unity in both the interpolated and non-emitting models. Since the model order 9 is sufficiently high, the interpolated model will use the `eith` state whenever it occurs and no higher order state is preferred (see figure 2).

The hierarchical non-emitting model has no such freedom (see figure 3). In order to reach the `eith` state, it must accurately predict every symbol in the string `eith`. Otherwise, it will be forced to a lower order state along the way. The transition to a lower order state occurs when the non-emitting model attempts to predict the symbol `t` from the state `ei`. Since `ei` is rarely followed by `t` in the training corpus (.0761), the non-emitting model is forced into the lower order state `i`, from which it is able to predict the symbol `t` with greater probability (.1172). As a result, the non-emitting model is never able to reach the `eith` state. Instead, it must predict the space ␣ after `␣Keith` using the state `ith`. This works quite well because `ith` is followed by ␣ with high probability in the training corpus (0.6136).

9	0.549	0.446					
8	0.275	0.223					
7	0.137	0.049					
6	0.037	0.082					
5	0.001	0.038					
4		0.067					1.000
3		0.050		0.376			
2		0.025	0.856	0.617	0.527	0.596	
1		0.015	0.142	0.004	0.415	0.297	
0		0.004	0.002	0.004	0.058	0.106	
	␣	K	e	i	t	h	␣
	0.998	0.000	0.273	0.006	0.093	0.226	0.003

Figure 2: State occupancy probabilities for the order 9 interpolated model on part of the Brown test corpus (2.4480 bits/char). The horizontal axis represents the position in the test string and the vertical access represents the hidden state order. The bottom column shows the conditional probability of the symbol, given the hidden state distribution. Thus the interpolated model is in the order 4 state `eith` with probability at least .9995 when predicting the final symbol, and it assigns probability 0.003 to this symbol.

9	0.166	0.121					
8	0.289	0.369					
7	0.340	0.449					
6	0.161	0.016					
5	0.035	0.017					
4	0.008	0.017					
3		0.008		0.472			0.734
2		0.002	0.939	0.527	0.848	0.749	0.187
1		0.001	0.061	0.001	0.118	0.207	0.003
0					0.033	0.044	0.075
	␣	K	e	i	t	h	␣
	0.977	0.000	0.277	0.006	0.081	0.232	0.627

Figure 3: State occupancy probabilities for the order 9 non-emitting model on part of the Brown test corpus (2.1536 bits/char). The horizontal axis represents the position in the test string and the vertical access represents the hidden state order. The bottom column shows the conditional probability of the symbol, given the hidden state distribution. Thus the non-emitting model is in the order 3 state `ith` with probability 0.734 when predicting the final symbol, and it assigns probability 0.627 to this symbol.

6.5 Posterior Tying

This posthoc analysis led John Lafferty (personal communication) to suggest that the interpolated model might be able to approximate the empirical performance of the non-emitting model with a suitable parameter tying scheme. According to the non-emitting model, two states should be considered equivalent if they are equally effective at predicting the future *and* they are equally well predicted by the model. A state is well-predicted if the string that it represents is assigned high probability, relative to the other states available at the time. A state provides strong predictions if the entropy of its emitting state transition probabilities is low.

The most effective way for the interpolated model to mimic the non-emitting model is to tie its states based on their expectations in the corresponding non-emitting model. In order to avoid implementing the non-emitting model, we may reasonably impose a uniform distribution on the non-emitting state transitions. And in order to avoid running the full EXPECTATION-STEP() algorithm, we may approximate the non-emitting state expectations by their forward expectations in $O(nT)$ time and $O(n)$ space.

A further simplification is to use the mean empirical posterior probability. The mean empirical posterior of a state is the empirical expectation $\delta[x^i]$ of the state divided by its frequency $c(x^i)$. The empirical expectation $\delta[x^i|y^T]$ of an i^{th} order state x^i in an order n mixture Markov model with respect to a string y^T is computed as follows

$$\delta[x^i|y^T] = \sum_{\{t:x^i=y_{t+1-i}^t\}} \delta(o_t = i|y^t),$$

with the empirical posterior

$$\delta(o = i|y^t) = \frac{\delta(y_{t+1-i}^t)}{\sum_{j=1}^n \delta(y_{t+1-j}^t)}.$$

Note that $\delta[x^i|y^T]$ may be calculated for all states in $O(nT)$ time using dynamic programming. The empirical posterior $\delta(o = i|y^t)$ of the i^{th} order state at time t could be weighted also by its predictive success $-\log \delta(y_{t+1}|y_{t+1-i}^t)$. A further refinement is to compute the mean empirical posterior on withheld data.

As a final step, these values must be quantized to a finite number of levels to construct the parameter tying scheme.

7 Conclusion

In this report, we propose a time series model that combines Markovian events of varying orders using stochastic non-emitting transitions. We prove that the resulting class of non-emitting Markov models is strictly more powerful than

the class of Markov models, including interpolated and backoff models. More importantly, our empirical investigation reveals that the non-emitting model consistently outperforms the strongest interpolated Markov models on natural language texts, with only a modest increase in computational requirements.

The expressive power of the non-emitting model comes from its ability to represent additional information in its state order distribution. To prove that the non-emitting model was strictly more powerful than any Markov model, we used the state order distribution to represent an unbounded dependency. In our posthoc analysis, we revealed how the model uses its hidden state order distribution to remember the short-term effectiveness of all available Markovian states.

The non-emitting model succeeds empirically because it imposes a pseudo-Bayesian discipline on maximum likelihood techniques. The interpolated model will favor a high-order state if it provides strong predictions on withheld data. The non-emitting model will favor a high-order state if the state provides strong predictions on withheld data *and it is well-predicted by the model*. In order to reach a high order state, the non-emitting model must assign high probability to each symbol in that state. Otherwise, the non-emitting model will be forced to transition to a lower order state at a previous time step and will not be able to reach the high order state. Thus, the state occupancies of the non-emitting model are influenced as much by their prior probabilities (pseudo-Bayes) as their past ability to predict the future (maximum likelihood).

Finally, we note the use of non-emitting transitions is a general modeling technique that may be employed in any time series model, for symbolic domains and for continuous domains.

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A Backoff

The backoff model is arguably the most widely used statistical language model, due in large part to its ease of implementation, computational efficiency, reasonable performance at lower model orders, and an influential paper [13]. Backoff models are also widely used in the data compression community, in large part due to their computational efficiency [5]. Here we review the backoff model, establish the equivalence of backoff models and basic Markov models, and then specify a class of non-emitting backoff models that is strictly more powerful than the class of traditional backoff models.

In a backoff model, event probabilities are combined according to a partial order. Typically, higher order events are preferred over lower order events. The event probabilities are rescaled as we move through the partial order so that the derived probability function is valid. The efficacy of the backoff model depends on the events that are included in the model, their individual probabilities, and the order in which they are combined.

Formally, a hierarchical backoff model $\theta = \langle A, E, \delta \rangle$ consists of an alphabet A, a dictionary E of selected state transitions, $E \subseteq A^* \times A$, and the state transition probabilities $\delta : E \to [0, 1]$. The state transition probabilities δ are extended to an unbounded domain by selecting the maximal suffix of the history that appears with the relevant symbol in the dictionary E of state transitions.

$$p_b(y|x^t, \theta) = \begin{cases} \delta(y|x^t) & \text{if } \langle x^t, y \rangle \in E \\ \eta(x^t) p_b(y|x_2^t, \theta) & \text{otherwise} \end{cases} \tag{7}$$

where $\eta(x^t)$ rescales the conditional probability distribution as we backoff from higher order events to lower order events

$$\eta(x^i) \doteq (1 - \delta(E(x^i)|x^i))/(1 - p_b(E(x^i)|x_2^i))$$

and $E(x^i)$ is the set of symbols available in the context x^i.

$$E(x^i) \doteq \{y : x^i y \in E\}$$

The rescalar $\eta(x^i)$ is computed directly form the transition probabilities $\delta(\cdot|x^i)$ in conjunction with the transition dictionary E. It is not a free parameter.

A hierarchical backoff model is a valid probability model if the dictionary E includes every 0^{th} order state transition – $\{\epsilon\} \times A \subset E$ – and every $\delta(E(x^i)|x^i)$ is a valid probability function. A backoff model is nonzero for all strings A^* if every $\delta(y|x^i)$ is nonzero and no $\delta(E(x^i)|x^i)$ is unity when $E(x^i) \subset A$.

In order to induce a hierarchical backoff model from data, we must select the state transition dictionary and estimate its probabilities. One simple – but highly effective – selection technique is to include every state transition whose frequency exceeds a fixed threshold, that may depend on the state order. More effective selection techniques require significant computational resources [22].

The state transition probabilities $\delta(y|x^t)$ are typically assigned by multinomial estimates, either as conditional events $y|x^i$ in the symbol alphabet A or as joint events x^iy in the string alphabet A^{i+1}. The most widely used multinomial estimates for statistical language modeling employ some form of discounting [5, 8, 9, 10, 11], although other estimators have also been shown to be effective [15, 18, 21].

A valid backoff model θ whose event dictionary E is a subset of A^{n+1} can be converted into an equivalent basic Markov model ϕ' of order n, simply by setting $\delta'_n(x_t|x_{t-n}^{t-1})$ equal to $p_b(x_t|x_{t-n}^{t-1}, \theta)$. Every basic model is a backoff model with a complete state transition dictionary. Consequently, the class of backoff models is extensionally equivalent to the class of basic Markov models.

The hierarchical non-emitting backoff model $\theta = \langle A, E, \delta \rangle$ has the same parameterization as the traditional backoff model. Unlike the traditional model, the backoff from the state x^i to its maximal proper suffix x_2^i is permanent in the non-emitting backoff model.

$$p_\epsilon(y^j|x^i, \theta) = \begin{cases} \delta(y_1|x^i)p_\epsilon(y_2^j|x^iy_1, \theta) & \text{if } \langle x^i, y_1 \rangle \in E \\ \eta(x^i)p_b(y^j|x_2^i, \phi) & \text{otherwise} \end{cases} \tag{8}$$

The rescalar $\eta(x^i)$ is identical in both version of the backoff model.

The class of non-emitting backoff models is strictly more powerful than the class of basic Markov models, by a similar argument as in lemma 4.1. Although the backoff model does not have any mixing parameters, we may use the presence or absence of a state transition $y|x^i$ in the dictionary E to control the hidden state order. Conversely, every order n backoff model can be converted into an equivalent non-emitting backoff model with a complete state transition dictionary $E = A^{n+1}$. Therefore, the class of non-emitting backoff models is strictly more powerful than the class of simple backoff models.