

A Comparative Study of Parameter Estimation Methods for Statistical Natural Language Processing

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Abstract

This paper presents a comparative study of five parameter estimation algorithms on four NLP tasks. Three of the five algorithms are well-known in the computational linguistics community: Maximum Entropy (ME) estimation with L_2 regularization, the Averaged Perceptron (AP), and Boosting. We also investigate ME estimation with L_1 regularization using a novel optimization algorithm, and BLasso, which is a version of Boosting with Lasso (L_1) regularization. We first investigate all of our estimators on two re-ranking tasks: a parse selection task and a language model (LM) adaptation task. Then we apply the best of these estimators to two additional tasks involving conditional sequence models: a Conditional Markov Model (CMM) for part of speech tagging and a Conditional Random Field (CRF) for Chinese word segmentation. Our experiments show that across tasks, three of the estimators — ME estimation with L_1 or L_2 regularization, and AP — are in a near statistical tie for first place.

1 Introduction

Parameter estimation is fundamental to many statistical approaches to NLP. Because of the high-dimensional nature of natural language, it is often easy to generate an extremely large number of features. The challenge of parameter estimation is to find a combination of the typically noisy, redundant features that accurately predicts the target output variable and avoids overfitting. Intuitively, this can be achieved either by selecting a small number of highly-effective features and ignoring the others, or by averaging over a large number of weakly informative features. The first intuition motivates feature selection methods such as Boosting and BLasso (e.g., Collins 2000; Zhao and

Yu, 2004), which usually work best when many features are completely irrelevant. L_1 or Lasso regularization of linear models, introduced by Tibshirani (1996), embeds feature selection into regularization so that both an assessment of the reliability of a feature and the decision about whether to remove it are done in the same framework, and has generated a large amount of interest in the NLP community recently (e.g., Goodman 2003; Riezler and Vasserman 2004). If on the other hand most features are noisy but at least weakly correlated with the target, it may be reasonable to attempt to reduce noise by averaging over all of the features. ME estimators with L_2 regularization, which have been widely used in NLP tasks (e.g., Chen and Rosenfeld 2000; Charniak and Johnson 2005; Johnson et al. 1999), tend to produce models that have this property. In addition, the perceptron algorithm and its variants, e.g., the voted or averaged perceptron, is becoming increasingly popular due to their competitive performance, simplicity in implementation and low computational cost in training (e.g., Collins 2002).

While recent studies claim advantages for L_1 regularization, this study is the first of which we are aware to systematically compare it to a range of estimators on a diverse set of NLP tasks. Gao et al. (2006) showed that BLasso, due to its explicit use of L_1 regularization, outperformed Boosting in the LM adaptation task. Ng (2004) showed that for logistic regression, L_1 regularization outperforms L_2 regularization on artificial datasets which contain many completely irrelevant features. Goodman (2003) showed that in two out of three tasks, an ME estimator with a one-sided Laplacian prior (i.e., L_1 regularization with the constraint that all feature weights are positive) outperformed a comparable estimator using a Gaussian prior (i.e., L_2 regularization). Riezler and Vasserman (2004) showed that an L_1 -regularized ME estimator outperformed an L_2 -regularized estimator for ranking the parses of a stochastic unification-based grammar.

While these individual estimators are well described in the literature, little is known about the relative performance of these methods because the published results are generally not directly comparable. For example, in the parse re-ranking task, one cannot tell whether the L_2 -regularized ME approach used by Charniak and Johnson (2005) significantly outperforms the Boosting method by Collins (2000) because different feature sets and n -best parses were used in the evaluations of these methods.

This paper conducts a much-needed comparative study of these five parameter estimation algorithms on four NLP tasks: ME estimation with L_1 and L_2 regularization, the Averaged Perceptron (AP), Boosting, and BLasso, a version of Boosting with Lasso (L_1) regularization. We first investigate all of our estimators on two re-ranking tasks: a parse selection task and a language model adaptation task. Then we apply the best of these estimators to two additional tasks involving conditional sequence models: a CMM for POS tagging and a CRF for Chinese word segmentation. Our results show that ME estimation with L_2 regularization achieves the best performing estimators in all of the tasks, and AP achieves almost as well and requires much less training time. L_1 (Lasso) regularization also performs well and leads to sparser models.

2 Estimators

All the four NLP tasks studied in this paper are based on linear models (Collins 2000) which require learning a mapping from inputs $x \in X$ to outputs $y \in Y$. We are given:

- Training samples (x_i, y_i) for $i = 1 \dots N$,
- A procedure **GEN** to generate a set of candidates $GEN(x)$ for an input x ,
- A feature mapping $\Phi: X \times Y \mapsto \mathbb{R}^D$ to map each (x, y) to a vector of feature values, and
- A parameter vector $\mathbf{w} \in \mathbb{R}^D$, which assigns a real-valued weight to each feature.

For all models except the CMM sequence model for POS tagging, the components **GEN**, Φ and \mathbf{w} directly define a mapping from an input x to an output $F(x)$ as follows:

$$F(x) = \arg \max_{y \in GEN(x)} \Phi(x, y) \cdot \mathbf{w}. \quad (1)$$

In the CMM sequence classifier, locally normalized linear models to predict the tag of each word token are chained together to arrive at a probability esti-

mate for the entire tag sequence, resulting in a slightly different decision rule.

Linear models, though simple, can capture very complex dependencies because the features can be arbitrary functions of the input/output pair. For example, we can define a feature to be the log conditional probability of the output as estimated by some other model, which may in turn depend on arbitrarily complex interactions of ‘basic’ features. In practice, with an appropriate feature set, linear models achieve very good empirical results on various NLP tasks. The focus of this paper however is not on feature definition (which requires domain knowledge and varies from task to task), but on parameter estimation (which is generic across tasks). We assume we are given fixed feature templates from which a large number of features are generated. The task of the estimator is to use the training samples to choose a parameter vector \mathbf{w} , such that the mapping $F(x)$ is capable of correctly classifying unseen examples. We will describe the five estimators in our study individually.

2.1 ME estimation with L_2 regularization

Like many linear models, the ME estimator chooses \mathbf{w} to minimize the sum of the empirical loss on the training set and a regularization term:

$$\hat{\mathbf{w}} = \arg \min_{\mathbf{w}} \{L(\mathbf{w}) + R(\mathbf{w})\}. \quad (2)$$

In this case, the loss term $L(\mathbf{w})$ is the negative conditional log-likelihood of the training data,

$$L(\mathbf{w}) = -\sum_{i=1}^n \log P(y_i | x_i), \text{ where}$$

$$P(y|x) = \frac{\exp(\Phi(x, y) \cdot \mathbf{w})}{\sum_{y' \in GEN(x)} \exp(\Phi(x, y') \cdot \mathbf{w})}$$

and the regularizer term $R(\mathbf{w}) = \alpha \sum_j w_j^2$ is the weighted squared L_2 norm of the parameters. Here, α is a parameter that controls the amount of regularization, optimized on held-out data.

This is one of the most popular estimators, largely due to its appealing computational properties: both $L(\mathbf{w})$ and $R(\mathbf{w})$ are convex and differentiable, so gradient-based numerical algorithms can be used to find the global minimum efficiently.

In our experiments, we used the *limited memory quasi-Newton* algorithm (or L-BFGS, Nocedal and Wright 1999) to find the optimal \mathbf{w} because this method has been shown to be substantially faster than other methods such as Generalized Iterative Scaling (Malouf 2002).

Because for some sentences there are multiple best parses (i.e., parses with the same F-Score), we used the variant of ME estimator described in Riezler et al. (2002), where $L(\mathbf{w})$ is defined as the likelihood of the best parses $y \in Y(x)$ relative to the n -best parser output $\mathbf{GEN}(x)$, (i.e., $Y(x) \subseteq \mathbf{GEN}(x)$): $L(\mathbf{w}) = -\sum_{i=1}^n \log \sum_{y_i \in Y(x_i)} P(y_i | x_i)$.

We applied this variant in our experiments of parse re-ranking and LM adaptation, and found that on both tasks it leads to a significant improvement in performance for the L_2 -regularized ME estimator but not for the L_1 -regularized ME estimator.

2.2 ME estimation with L_1 regularization

This estimator also minimizes the negative conditional log-likelihood, but uses an L_1 (or Lasso) penalty. That is, $R(\mathbf{w})$ in Equation (2) is defined according to $R(\mathbf{w}) = \alpha \sum_j |w_j|$. L_1 regularization typically leads to sparse solutions in which many feature weights are exactly zero, so it is a natural candidate when feature selection is desirable. By contrast, L_2 regularization produces solutions in which most weights are small but non-zero.

Optimizing the L_1 -regularized objective function is challenging because its gradient is discontinuous whenever some parameter equals zero. Kazama and Tsujii (2003) described an estimation method that constructs an equivalent constrained optimization problem with twice the number of variables. However, we found that this method is impractically slow for large-scale NLP tasks. In this work we use the *orthant-wise limited-memory quasi-Newton* algorithm (OWL-QN), which is a modification of L-BFGS that allows it to effectively handle the discontinuity of the gradient (Andrew and Gao 2007). We provide here a high-level description of the algorithm.

A quasi-Newton method such as L-BFGS uses first order information at each iterate to build an approximation to the Hessian matrix, \mathbf{H} , thus modeling the local curvature of the function. At each step, a search direction is chosen by minimizing a quadratic approximation to the function:

$$Q(x) = \frac{1}{2}(x - x_0)' \mathbf{H}(x - x_0) + g_0'(x - x_0)$$

where x_0 is the current iterate, and g_0 is the function gradient at x_0 . If \mathbf{H} is positive definite, the minimizing value of x can be computed analytically according to: $x^* = x_0 - \mathbf{H}^{-1}g_0$.

L-BFGS maintains vectors of the change in gradient $g_k - g_{k-1}$ from the most recent iterations, and uses them to construct an estimate of the inverse Hessian \mathbf{H}^{-1} . Furthermore, it does so in such a way that $\mathbf{H}^{-1}g_0$ can be computed without expanding out the full matrix, which is typically unmanageably large. The computation requires a number of operations linear in the number of variables.

OWL-QN is based on the observation that when restricted to a single orthant, the L_1 regularizer is differentiable, and is in fact a linear function of \mathbf{w} . Thus, so long as each coordinate of any two consecutive search points does not pass through zero, $R(\mathbf{w})$ does not contribute at all to the curvature of the function on the segment joining them. Therefore, we can use L-BFGS to approximate the Hessian of $L(\mathbf{w})$ alone, and use it to build an approximation to the full regularized objective that is valid on a given orthant. To ensure that the next point is in the valid region, we project each point during the line search back onto the chosen orthant.¹ At each iteration, we choose the orthant containing the current point and into which the direction giving the greatest local rate of function decrease points.

This algorithm, although only a simple modification of L-BFGS, works quite well in practice. It typically reaches convergence in even fewer iterations than standard L-BFGS takes on the analogous L_2 -regularized objective (which translates to less training time, since the time per iteration is only negligibly higher, and total time is dominated by function evaluations). We describe OWL-QN more fully in (Andrew and Gao 2007). We also show that it is significantly faster than Kazama and Tsujii's algorithm for L_1 regularization and prove that it is guaranteed converge to a parameter vector that globally optimizes the L_1 -regularized objective.

2.3 Boosting

The Boosting algorithm we used is based on Collins (2000). It optimizes the pairwise *exponential loss* (ExpLoss) function (rather than the logarithmic loss optimized by ME). Given a training sample (x_i, y_i) , for each possible output $y_j \in \mathbf{GEN}(x_i)$, we

¹ This projection just entails zeroing-out any coordinates that change sign. Note that it is possible for a variable to change sign in two iterations, by moving from a negative value to zero, and on the next iteration moving from zero to a positive value.

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- 1 Set $w_0 = \operatorname{argmin}_w \operatorname{ExpLoss}(w)$; and $w_d = 0$ for $d=1 \dots D$
 - 2 Select a feature f_{k^*} which has largest estimated impact on reducing $\operatorname{ExpLoss}$ of Equation (3)
 - 3 Update $\lambda_{k^*} \leftarrow \lambda_{k^*} + \delta^*$, and return to Step 2
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Figure 1: The boosting algorithm

define the margin of the pair (y_i, y_j) with respect to \mathbf{w} as $M(y_i, y_j) = \Phi(x_i, y_i) \cdot \mathbf{w} - \Phi(x_i, y_j) \cdot \mathbf{w}$.

Then $\operatorname{ExpLoss}$ is defined as

$$\operatorname{ExpLoss}(\mathbf{w}) = \sum_i \sum_{y_j \in \operatorname{GEN}(x_i)} \exp(-M(y_i, y_j)) \quad (3)$$

Figure 1 summarizes the Boosting algorithm we used. It is an incremental feature selection procedure. After initialization, Steps 2 and 3 are repeated T times; at each iteration, a feature is chosen and its weight is updated as follows.

First, we define $\operatorname{Upd}(\mathbf{w}, k, \delta)$ as an updated model, with the same parameter values as w with the exception of w_k , which is incremented by δ :

$$\operatorname{Upd}(\mathbf{w}, k, \delta) = (w_1, \dots, w_k + \delta, \dots, w_D)$$

Then, Steps 2 and 3 in Figure 1 can be rewritten as Equations (4) and (5), respectively.

$$(k^*, \delta^*) = \operatorname{arg\,min}_{k, \delta} \operatorname{ExpLoss}(\operatorname{Upd}(\mathbf{w}, k, \delta)) \quad (4)$$

$$\mathbf{w}^t = \operatorname{Upd}(\mathbf{w}^{t-1}, k^*, \delta^*) \quad (5)$$

Because Boosting can overfit we update the weight of f_{k^*} by a small fixed step size ϵ , as in Equation (6), following the FSLR algorithm (Hastie et al. 2001).

$$\mathbf{w}^t = \operatorname{Upd}(\mathbf{w}^{t-1}, k^*, \epsilon \times \operatorname{sign}(\delta^*)) \quad (6)$$

By taking such small steps, Boosting imposes a kind of implicit regularization, and can closely approximate the effect of L_1 regularization in a local sense (Hastie et al. 2001). Empirically, smaller values of ϵ lead to smaller numbers of test errors.

2.4 Boosted Lasso

The Boosted Lasso (BLasso) algorithm was originally proposed in Zhao and Yu (2004), and was adapted for language modeling by Gao et al. (2006). BLasso can be viewed as a version of Boosting with L_1 regularization. It optimizes an L_1 -regularized $\operatorname{ExpLoss}$ function:

$$\operatorname{LassoLoss}(\mathbf{w}) = \operatorname{ExpLoss}(\mathbf{w}) + R(\mathbf{w}) \quad (7)$$

where $R(\mathbf{w}) = \alpha \sum_j |w_j|$.

BLasso also uses an incremental feature selection procedure to learn parameter vector \mathbf{w} , just as Boosting does. Due to the explicit use of the regu-

larization term $R(\mathbf{w})$, however, there are two major differences from Boosting.

At each iteration, BLasso takes either a *forward step* or a *backward step*. Similar to Boosting, at each forward step, a feature is selected and its weight is updated according to Eq. (8) and (9).

$$(k^*, \delta^*) = \operatorname{arg\,min}_{k, \delta = \pm \epsilon} \operatorname{ExpLoss}(\operatorname{Upd}(\mathbf{w}, k, \delta)) \quad (8)$$

$$\mathbf{w}^t = \operatorname{Upd}(\mathbf{w}^{t-1}, k^*, \epsilon \times \operatorname{sign}(\delta^*)) \quad (9)$$

There is a small but important difference between Equations (8) and (4). In Boosting, as shown in Equation (4), a feature is selected by its impact on reducing the loss with its optimal update δ^* . By contrast, in BLasso, as shown in Equation (8), rather than optimizing over δ for each feature, the loss is calculated with an update of either $+\epsilon$ or $-\epsilon$, i.e., grid search is used for feature weight estimation. We found in our experiments that this modification brings a consistent improvement.

The backward step is unique to BLasso. At each iteration, a feature is selected and the absolute value of its weight is reduced by ϵ if and only if it leads to a decrease of the LassoLoss, as shown in Equations (10) and (11), where θ is a tolerance parameter.

$$k^* = \operatorname{arg\,min}_{k: w_k \neq 0} \operatorname{ExpLoss}(\operatorname{Upd}(\mathbf{w}, k, -\epsilon \operatorname{sign}(w_k))) \quad (10)$$

$$\mathbf{w}^t = \operatorname{Upd}(\mathbf{w}^{t-1}, k^*, \operatorname{sign}(w_{k^*}) \times \epsilon) \quad (11)$$

if $\operatorname{LassoLoss}(\mathbf{w}^{t-1}, \alpha^{t-1}) - \operatorname{LassoLoss}(\mathbf{w}^t, \alpha^t) > \theta$

Figure 2 summarizes the BLasso algorithm we used. After initialization, Steps 4 and 5 are repeated T times; at each iteration, a feature is chosen and its weight is updated either backward or forward by a fixed amount ϵ . Notice that the value of α is adaptively chosen according to the reduction of $\operatorname{ExpLoss}$ during training. The algorithm starts with a large initial α , and then at each forward step the value of α decreases until $\operatorname{ExpLoss}$ stops decreasing. This is intuitively desirable: it is expected that most highly effective features are selected in early stages of training, so the reduction of $\operatorname{ExpLoss}$ at each step in early stages are more substantial than in later stages. These early steps coincide with the Boosting steps most of the time. In other words, the effect of backward steps is more visible at later stages. It can be proved that for a finite number of features and $\theta=0$, the BLasso algorithm shown in Figure 2 converges to the Lasso solution when $\epsilon \rightarrow 0$. See Gao et al. (2006) for implementation details, and Zhao and Yu (2004) for a theoretical justification for BLasso.

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- 1 Initialize w^0 : set $w_0 = \operatorname{argmin}_{w^0} \operatorname{ExpLoss}(w)$, and $w_d = 0$ for $d=1 \dots D$.
 - 2 Take a forward step according to Eq. (8) and (9), and the updated model is denoted by w^1
 - 3 Initialize $\alpha = (\operatorname{ExpLoss}(w^0) - \operatorname{ExpLoss}(w^1)) / \varepsilon$
 - 4 Take a backward step if and only if it leads to a decrease of LassoLoss according to Eq. (10) and (11), where $\theta = 0$; otherwise
 - 5 Take a forward step according to Eq. (8) and (9); update $\alpha = \min(\alpha, (\operatorname{ExpLoss}(w^{t-1}) - \operatorname{ExpLoss}(w^t)) / \varepsilon)$; and return to Step 4.
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Figure 2: The BLasso algorithm

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- 1 Set $w_0 = 1$ and $w_d = 0$ for $d=1 \dots D$
 - 2 For $t = 1 \dots T$ (T = the total number of iterations)
 - 3 For each training sample (x_i, y_i) , $i = 1 \dots N$
 - 4 Choose the best candidate z_i from $\operatorname{GEN}(x_i)$ using the current model w ,

$$z_i = \operatorname{argmax}_{z \in \operatorname{GEN}(x_i)} \Phi(x_i, z) \cdot w$$
 - 5 $w = w + \eta(\Phi(x_i, y_i) - \Phi(x_i, z_i))$, where η is the size of learning step, optimized on held-out data.
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Figure 3: The perceptron algorithm

2.5 Averaged Perceptron

The perceptron algorithm can be viewed as a form of incremental training procedure (e.g., using stochastic approximation) that optimizes a *minimum square error* (MSE) loss function (Mitchell, 1997). As shown in Figure 3, it starts with an initial parameter setting and updates it for each training example. In our experiments, we used the Averaged Perceptron algorithm of Freund and Schapire (1999), a variation that has been shown to be more effective than the standard algorithm (Collins 2002). Let $w^{t,i}$ be the parameter vector after the i^{th} training sample has been processed in pass t over the training data. The average parameters are defined as $\bar{w} = \frac{1}{TN} \sum_t \sum_i w^{t,i}$ where T is the number of epochs, and N is the number of training samples.

3 Evaluations

From the four tasks we consider, parsing and language model adaptation are both examples of re-ranking. In these tasks, we assume that we have been given a list of candidates $\operatorname{GEN}(x)$ for each training or test sample (x, y) , generated using a *baseline* model. Then, a linear model of the form in Equation (1) is used to discriminatively re-rank the candidate list using additional features which may or may not be included in the baseline model. Since

the mapping from x to y by the linear model may make use of arbitrary global features of the output and is performed “all at once”, we call such a linear model a *global model*.

In the other two tasks (i.e., Chinese word segmentation and POS tagging), there is no explicit enumeration of $\operatorname{GEN}(x)$. The mapping from x to y is determined by a sequence model which aggregates the decisions of local linear models via a dynamic program. In the CMM, the local linear models are trained independently, while in the CRF model, the local models are trained jointly. We call these two linear models *local models* because they dynamically combine the output of models that use only local features.

While it is straightforward to apply the five estimators to global models in the re-ranking framework, the application of some estimators to the local models is problematic. Boosting and BLasso are too computationally expensive to be applied to CRF training and we compared the other three better performing estimation methods for this model. The CMM is a probabilistic sequence model and the log-loss used by ME estimation is most natural for it; thus we limit the comparison to the two kinds of ME models for CMMs. Note that our goal is not to compare locally trained models to globally trained ones; for a study which focuses on this issue, see (Punyakanok et al. 2005).

In each task we compared the performance of different estimators using task-specific measures. We used the Wilcoxon signed rank test to test the statistical significance of the difference among the competing estimators. We also report other results such as number of non-zero features after estimation, number of training iterations, and computation time (in minutes of elapsed time on an XEONTM MP 3.6GHz machine).

3.1 Parse re-ranking

We follow the experimental paradigm of parse re-ranking outlined in Charniak and Johnson (2005), and fed the features extracted by their program to the five rerankers we developed. Each uses a linear model trained using one of the five estimators. These rerankers attempt to select the best parse y for a sentence x from the 50-best list of possible parses $\operatorname{GEN}(x)$ for the sentence. The linear model combines the log probability calculated by the Charniak (2000) parser as a feature with 1,219,272 additional features. We trained the fea-

	F-Score	# features	time (min)	# train iter
Baseline	0.8986			
ME/L2	0.9176	1,211,026	62	129
ME/L1	0.9165	19,121	37	174
AP	0.9164	939,248	2	8
Boosting	0.9131	6,714	495	92,600
BLasso	0.9133	8,085	239	56,500

Table 1: Performance summary of estimators on parsing re-ranking (ME/L2: ME with L_2 regularization; ME/L1: ME with L_1 regularization)

	ME/L2	ME/L1	AP	Boost	BLasso
ME/L2		>>	~	>>	>>
ME/L1	<<		~	>	~
AP	~	~		>>	>
Boost	<<	<	<<		~
BLasso	<<	~	<	~	

Table 2: Statistical significance test results (“>>” or “<<” means $P\text{-value} < 0.01$; > or < means $0.01 < P\text{-value} \leq 0.05$; “~” means $P\text{-value} > 0.05$)

ture weights w on Sections 2-19 of the Penn Treebank, adjusted the regularizer constant α to maximize the F-Score on Sections 20-21 of the Treebank, and evaluated the rerankers on Section 22. The results are presented in Tables 1² and 2, where **Baseline** results were obtained using the parser by Charniak (2000).

The ME estimation with L_2 regularization outperforms all of the other estimators significantly except for the AP, which performs almost as well and requires an order of magnitude less time in training. Boosting and BLasso are feature selection methods in nature, so they achieve the sparsest models, but at the cost of slightly lower performance and much longer training time. The L_1 -regularized ME estimator also produces a relatively sparse solution whereas the Averaged Perceptron and the L_2 -regularized ME estimator assign almost all features a non-zero weight.

3.2 Language model adaptation

Our experiments with LM adaptation are based on the work described in Gao et al. (2006). The various trained language models were evaluated according to their impact on Japanese text input accuracy, where input phonetic symbols x are mapped into a word string y . Performance of the application is measured in terms of character error

² The result of ME/L2 is better than that reported in Andrew and Gao (2007) due to the use of the variant of L_2 -regularized ME estimator, as described in Section 2.1.

	CER	# features	time (min)	#train iter
Baseline	10.24%			
MAP	7.98%			
ME/L2	6.99%	295,337	27	665
ME/L1	7.01%	53,342	25	864
AP	7.23%	167,591	6	56
Boost	7.54%	32,994	175	71,000
BLasso	7.20%	33,126	238	250,000

Table 3: Performance summary of estimators (lower is better) on language model adaptation

	ME/L2	ME/L1	AP	Boost	BLasso
ME/L2		~	>>	>>	>>
ME/L1	~		>>	>>	>>
AP	<<	<<		>>	~
Boost	<<	<<	<<		<<
BLasso	<<	<<	~	>>	

Table 4: Statistical significance test results.

rate (CER), which is the number of characters wrongly converted from x divided by the number of characters in the correct transcript.

Again we evaluated five linear rerankers, one for each estimator. These rerankers attempt to select the best conversions y for an input phonetic string x from a 100-best list $GEN(x)$ of possible conversions proposed by a baseline system. The linear model combines the log probability under a trigram language model as base feature and additional 865,190 word uni/bi-gram features. These uni/bi-gram features were already included in the trigram model which was trained on a *background* domain corpus (Nikkei Newspaper). But in the linear model their feature weights were trained discriminatively on an *adaptation* domain corpus (Encarta Encyclopedia). Thus, this forms a cross domain adaptation paradigm. This also implies that the portion of redundant features in this task could be much larger than that in the parse re-ranking task, especially because the background domain is reasonably similar to the adaptation domain.

We divided the Encarta corpus into three sets that do not overlap. A 72K-sentences set was used as training data, a 5K-sentence set as development data, and another 5K-sentence set as testing data. The results are presented in Tables 3 and 4, where **Baseline** is the word-based trigram model trained on background domain corpus, and **MAP** (maximum *a posteriori*) is a traditional model adaptation method, where the parameters of the background model are adjusted so as to maximize the likelihood of the adaptation data.

	Test F_1	# features	# train iter
ME/L2	0.9719	8,084,086	713
ME/L1	0.9713	317,146	201
AP	0.9703	1,965,719	162

Table 5. Performance summary of estimators on CWS

The results are more or less similar to those in the parsing task with one visible difference: L_1 regularization achieved relatively better performance in this task. For example, while in the parsing task ME with L_2 regularization significantly outperforms ME with L_1 regularization, their performance difference is not significant in this task. While in the parsing task the performance difference between BLasso and Boosting is not significant, BLasso outperforms Boosting significantly in this task. Considering that a much higher proportion of the features are redundant in this task than the parsing task, the results seem to corroborate the observation that L_1 regularization is robust to the presence of many redundant features.

3.3 Chinese word segmentation

Our third task is Chinese word segmentation (CWS). The goal of CWS is to determine the boundaries between words in a section of Chinese text. The model we used is the hybrid Markov/semi-Markov CRF described by Andrew (2006), which was shown to have state-of-the-art accuracy. We tested models trained with the various estimation methods on the Microsoft Research Asia corpus from the Second International Chinese Word Segmentation, and we used the same train/test split used in the competition. The model and experimental setup is identical with that of Andrew (2006) except for two differences. First, we extracted features from both positive and negative training examples, while Andrew (2006) uses only features that occur in some positive training example. Second, we used the last 4K sentences of the training data to select the weight of the regularizers and to determine when to stop perceptron training.

We compared three of the best performing estimation procedures on this task: ME with L_2 regularization, ME with L_1 regularization, and the Averaged Perceptron. In this case, ME refers to minimizing the negative log-probability of the correct segmentation, which is globally normalized, while the perceptron is trained using at each iteration the exact maximum-scoring segmentation with the

current weights. We observed the same pattern as in the other tasks: the three algorithms have nearly identical performance, while L_1 uses only 6% of the features, and the Averaged Perceptron requires significantly fewer training iterations. In this case, L_1 was also several times faster than L_2 . The results are summarized in Table 5.³

We note that all three algorithms performed slightly better than the model used by Andrew (2006), which also used L_2 regularization (96.84 F_1). We believe the difference is due to the use of features derived from negative training examples.

3.4 POS tagging

Finally we studied the impact of the regularization methods on a Maximum Entropy conditional Markov Model (MEMM, McCallum et al. 2000) for POS tagging. MEMMs decompose the conditional probability of a tag sequence given a word sequence as follows:

$$P(t_1 \dots t_n | w_1 \dots w_n) = \prod_{i=1}^n P(t_i | t_{i-1} \dots t_{i-k}, w_1 \dots w_n)$$

where the probability distributions for each tag given its context are ME models. Following previous work (Ratnaparkhi, 1996), we assume that the tag of a word is independent of the tags of all preceding words given the tags of the previous two words (i.e., $k=2$ in the equation above). The local models at each position include features of the current word, the previous word, the next word, and features of the previous two tags. In addition to lexical identity of the words, we used features of word suffixes, capitalization, and number/special character signatures of the words.

We used the standard splits of the Penn Treebank from the tagging literature (Toutanova et al. 2003) for training, development and test sets. The training set comprises Sections 0-18, the development set — Sections 19-21, and the test set — Sections 22-24. We compared training the ME models using L_1 and L_2 regularization. For each of the two types of regularization we selected the best value of the regularization constant using grid search to optimize the accuracy on the development set. We report final accuracy measures on the test set in Table 6.

The results on this task confirm the trends we have seen so far. There is almost no difference in

³ Only the L2 vs. AP comparison is significant at a 0.05 level according to the Wilcoxon signed rank test.

	Accuracy (%)	# features	# train iter
MEMM/ L_2	96.39	926,350	467
MEMM/ L_1	96.41	84,070	85

Table 6. Performance summary of estimators on POS tagging

accuracy of the two kinds of regularizations, and indeed the differences were not statistically significant. Estimation with L_1 regularization required considerably less time than estimation with L_2 , and resulted in a model which is more than ten times smaller.

4 Conclusions

We compared five of the most competitive parameter estimation methods on four NLP tasks employing a variety of models, and the results were remarkably consistent across tasks. Three of the methods — ME estimation with L_2 regularization, ME estimation with L_1 regularization, and the Averaged Perceptron — were nearly indistinguishable in terms of test set accuracy, with ME estimation with L_2 regularization perhaps enjoying a slight lead. Meanwhile, ME estimation with L_1 regularization achieves the same level of performance while at the same time producing sparse models, and the Averaged Perceptron provides an excellent compromise of high performance and fast training.

These results suggest that when deciding which type of parameter estimation to use on these or similar NLP tasks, one may choose any of these three popular methods and expect to achieve comparable performance. The choice of which to implement should come down to other considerations: if model sparsity is desired, choose ME estimation with L_1 regularization (or feature selection methods such as BLasso); if quick implementation and training is necessary, use the Averaged Perceptron; and ME estimation with L_2 regularization may be used if it is important to achieve the highest obtainable level of performance.

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