Computing Word Classes Using Spectral Clustering

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Abstract

Clustering a lexicon of words is a well-studied problem in natural language processing (NLP). Word clusters are used to deal with sparse data in statistical language processing, as well as features for solving various NLP tasks (text categorization, question answering, named entity recognition and others).

Spectral clustering is a widely used technique in the field of image processing and speech recognition. However, it has scarcely been explored in the context of NLP; specifically, the method used in this work (Meila and Shi, 2001) has never been used to cluster a general word lexicon.

We apply spectral clustering to a lexicon of words, evaluating the resulting clusters by using them as features for solving two classical NLP tasks: semantic role labeling and dependency parsing. We compare performance with Brown clustering, a widely-used technique for word clustering, as well as with other clustering methods. We show that spectral clusters produce similar results to Brown clusters, and outperform other clustering methods. In addition, we quantify the overlap between spectral and Brown clusters, showing that each model captures some information which is uncaptured by the other.

1 Introduction

Word clusters (or word classes) are the result of dividing a lexicon of words into a pre-defined number of distinct groups, where the words comprising each group share some common characteristic. They have been well studied (Martin et al., 1998), in the context of part of speech (POS) induction (Christodoulopoulos et al., 2010), dealing with sparse data in statistical language processing (Brown et al., 1992), as well as for use as features in various NLP tasks such as text categorization (Bekkerman et al., 2003), question answering (Monttazi and Klakow, 2009), statistical parsing (Candito and Crabbé, 2009) named entity recognition (NER) (Miller et al., 2004), and others.

Brown Clustering (Brown et al., 1992), a hard hierarchical agglomerative clustering method, based on maximizing the quality of an induced class-based language model to make clustering decisions, is widely used to produce features for various NLP tasks (Monttazi and Klakow, 2009; Candito and Crabbé, 2009; Miller et al., 2004; Koo et al., 2008).

Spectral clustering is a clustering method which is broadly used for tasks such as image segmentation (Shi and Malik, 2000; Ng et al., 2002), speech recognition (Bach and Jordan, 2006) and topological mapping (Brunskill et al., 2007). It belongs to the family of dimensionality reduction algorithms and can actually be viewed as a weighted kernel K-means algorithm (Dhillon et al., 2004). However, it has scarcely been used for NLP tasks. Dhillon et al. (2011) use spectral methods for NER and chunking – but not for clustering – while Sun and Korhonen (2009) employ spectral clustering to improve verb clustering. Sedoc et al. (2017) used Signed Normalized Cut to produce word clusters; however, their motivation is focused on word-similarity (specifically, antonym-synonym relations), and they use mainly intrinsic evaluation methods. We aim to produce general spectral clusters and compare them to the widely-used Brown clusters using two NLP structured-prediction tasks for extrinsic evaluation.

While lacking the hierarchical nature of the Brown method, spectral clustering does possess a practical advantage over the former in that it produces – in addition to word clusters – a low-dimension represen-
tation \( v \in \mathbb{R}^m \) for each of the words in the lexicon as well as for each of the cluster centers. Beyond obvious theoretical interest, this representation may be quite useful for measuring word-word, word-cluster and even cluster-cluster distances, as well as allowing us to perform soft clustering, in which we assign each word a distribution over the clusters (instead of assigning a single cluster).

In this work we use the spectral clustering method presented in (Meila and Shi, 2001) over a general lexicon to produce word clusters. We proceed to show that when used as features for solving two classical yet complex NLP tasks – SRL and dependency parsing – spectral clusters produce very similar results to those produced by Brown clusters, despite lacking the firm language-modeling grounding inherent in the latter. Finally, we quantify the overlap (and difference) in the information contained in the two cluster sets, and show that combining them could potentially induce significantly better performance than using each on its own.

2 Background

2.1 Clustering Methods

Clustering deals with the problem of dividing set of \( n \) samples into \( k \) distinct clusters following a desired function measuring sample similarity. The definition of what makes two samples similar or related is not obvious or singular; most intuitively, the samples may reside in \( \mathbb{R}^n \) and the clusters should consist of samples that are similar following Euclidean distance, for example.

In the following sub-sections we will describe the best known clustering algorithm – K-means – which we use as a baseline, as well as spectral clustering – which is the focus of our work – and Brown clustering – which we compare to.

2.1.1 K-means

Dating back to six decades ago and published in (Lloyd, 1982), K-means is perhaps the most widely known clustering method. Given a set of points \( s_1, ..., s_n \) and a set of \( k \) initial means \( \mu_1^{(1)}, ..., \mu_k^{(1)} \) representing \( k \) clusters \( C_1^{(1)}, ..., C_k^{(1)} \), the algorithm iterates until convergence:

1. For each point, find the cluster with the closest mean and add the point to that cluster:

   \[
   (a) \quad i = \arg \min_{C_j^{(t)}} ||s_p - \mu_j^{(t)}||^2
   \]

   \[
   (b) \quad C_i^{(t)} \leftarrow s_p
   \]

2. Update clusters means:

   \[
   \mu_i^{(t+1)} = \frac{1}{|C_i^{(t)}|} \sum_{s_p \in C_i^{(t)}} s_p
   \]

   The means initialization may be random; different variants of K-means perform better with different types of initializations (Hamerly and Elkan, 2002).

   K-means performs best when the data is spatially separable, and uses Euclidean distance as metric. It also has a tendency to choose clusters occupying area of similar size on feature space.

2.1.2 Spectral Clustering

Spectral clustering refers to a group of algorithms that work on the spectral (eigenvector) decomposition of the samples’ affinity matrix, or similarity matrix. Each sample is represented by a vector; a metric is used in the resulting vector space to compute the distances between all point pairs – the affinity matrix. Following some mathematical manipulation on the affinity matrix, the eigenvectors are computed, and from them the clustering is derived, usually using the K-means algorithm. There are many flavors to this method, and we will describe here two of the main algorithms.

The first algorithm is introduced in (Ng et al., 2002):

Given a set of points \( S = s_1, ..., s_n \) in \( \mathbb{R} \) that we want to cluster into \( k \) subsets:

1. Form the affinity matrix \( A \in \mathbb{R}^{n \times n} \) defined by \( A_{ij} = \exp(-||s_i - s_j||^2/2\sigma^2) \) \( (A_{ii} = 0) \).
2. Define $D$ to be a diagonal matrix s.t. $D_{ii} = \sum_j A_{ij}$, and $L = D^{-1/2}AD^{-1/2}$.

3. Find the $k$ largest eigenvectors of $L$ (matching the largest $k$ eigenvalues), $x_1, \ldots, x_k$, and stack them in columns to form $X = [x_1 \ldots x_k] \in \mathbb{R}^{n \times k}$.

4. Derive $Y$ from $X$ by normalizing $X$’s rows to have unit length.

5. Cluster $Y$’s rows into $k$ clusters (using K-means, for example).

6. Assign point $s_i$ to cluster $j$ iff row $i$ of $Y$ was assigned to cluster $j$.

The second algorithm is presented by Shi and Malik (2000; Meila and Shi (2001) and takes a graph-theoretic approach to clustering. The data set is represented by a weighted undirected graph. Each point is represented by a vertex, and each pair of vertices are connected by a weighted edge, with the weight representing the similarity between the two respective points. In this setting, the problem of segmenting the data set into two groups is formulated as partitioning the graph into two groups of vertices where the similarity within each group is maximized while the similarity between the groups is minimized.

Given a weighted graph $G(V, E, w)$ and two subsets of the vertices in the graph $A, B \subseteq V$, Define:

$$w(A, B) = \sum_{u \in A, v \in B} w(u, v),$$

The normalized cut, a symmetric measure for the disassociation between the subsets, is defined to be:

$$Ncut(A, B) = \frac{w(A, B)}{w(A, V)} + \frac{w(A, B)}{w(B, V)}.$$

Along with a measure for association within each set:

$$Nassoc(A, B) = \frac{w(A, A)}{w(A, V)} + \frac{w(B, B)}{w(B, V)}.$$

Note that these two measures are related in the following way:

$$Ncut(A, B) = 2 - Nassoc(A, B).$$

Therefore, by minimizing the disassociation between groups we also maximize the association within each group.

Finding a cut that minimizes the $Ncut$ criterion is shown to be NP-hard, and the $Ncut$ algorithm is introduced, approximating in polynomial time the 2-way cut solution using the eigenvalues and eigenvectors of the affinity matrix. The algorithm is then used recursively to find a $k$-way partition of the graph, providing a clustering of the data set to $k$ groups.

A more efficient method, however, is presented in (Meila and Shi, 2001). After establishing a probabilistic theoretic foundation to the normalized cut framework by offering a random walk interpretation, they present the Modified $Ncut$ algorithm for a one-pass $k$-way segmentation:

1. Generate $D$ as defined in (Ng et al., 2002) (the first algorithm).

2. Generate $P = D^{-1}S$, $S$ being the similarity matrix, and find its eigenvalues and eigenvectors.

3. Discard the leading eigenvector and stack second-through-$k$ leading eigenvectors to form $X = [x_2 \ldots x_k]$.

4. Perform K-means (or equivalent) on rows of $X$ to find the clusters.

We chose to use the latter algorithm in our work, due to availability-of-code considerations.

Note that the last stage in the algorithm involves clustering the data points in the dense (due to the dimensionality reduction) vector space created by the computed eigenvectors. As mentioned in Section[1] this means that a by-product of this algorithm is a vector representation in this space for each of the clustered data points (in our case, lexicon words) as well as for the cluster centers. We discuss this further in Section[5].
2.1.3 Brown Clustering

Brown clustering (Brown et al., 1992) is a hard hierarchical agglomerative clustering method. It is based on the concept of maximizing the quality of an induced class-based language model, having originally been presented as a class-based n-gram model for natural language.

The algorithm generally follows the outline for hard hierarchical agglomerative clustering:

1. Start with a lexicon of types \( V \)
2. Sort \( V \) by corpus frequency, then put the top \( k \) types into clusters \( C_1, ..., C_k \)
3. Repeat \( |V| - k \) times:
   - Put the next type into a new cluster \( C_{k+1} \)
   - Merge the pair in the \( k + 1 \) clusters to receive clustering \( C \) that maximizes \( \text{Quality}(C) \)

The function measuring the quality of the clustering at each iteration – \( \text{Quality}(C) \) – was defined based on statistical language modeling reasoning. Given a corpus \( w_1, ..., w_n \), a first-order Hidden Markov Model is used (with the clusters \( C(w_1), ..., C(w_n) \) as the latent variables) to approximate the corpus probability:

\[
P(w_1, ..., w_n | C(w_1), ..., C(w_n)) \approx \prod_{i=1}^{n} P(w_i | C(w_i))P(C(w_i) | C(w_{i-1}))
\]

The quality function is defined to be:

\[
\text{Quality}(C) = \frac{1}{n} \log P(w_1, ..., w_n | C(w_1), ..., C(w_n))
\approx \frac{1}{n} \log \prod_{i=1}^{n} P(w_i | C(w_i))P(C(w_i) | C(w_{i-1}))
\]

Further decomposing into:

\[
= \sum_{1 \leq i,j \leq k} P(C_i, C_j) \log \frac{P(C_i, C_j)}{P(C_i)P(C_j)} + \sum_{w} P(w) \log P(w)
= I(C) - H(V)
\]

Which is the mutual information of the clustering minus the entropy of the vocabulary. The latter is constant over the clustering, meaning that the mutual information is maximized over the clustering in each iteration of the algorithm.

2.2 Evaluation

Conducting a comparison between two different clustering methods is not a trivial task. Viewing the clusters themselves may provide us with some insight. Some criteria exist for assessing the quality of the clusters, such as silhouette graphs (Rousseeuw, 1987), while others, such as the Variation of Information criterion (VI) (Meila, 2007) along with its normalized variant (Reichart and Rappoport, 2009), intend to compare two different clusterings over the same data set. These intrinsic methods, however, offer us very limited insight due to the unsupervised nature of our clustering task. We therefore turn to an extrinsic method of evaluation: using the clusters as features in a higher level task, and evaluate the clusters based on the performance in that task. We accomplish this by testing the clusters as features in two classical structured-prediction NLP tasks: semantic role labeling (SRL) and dependency parsing. As we wish to test the power of word clusters as features for these tasks, we use them exclusively without using any syntactic (POS tagging or dependency parses) or other semantic information.
2.2.1 Semantic Role Labeling

Semantic role labeling (SRL) is the task of detecting and labeling the different semantic arguments of a predicate in a sentence. The foundations for the task have been laid in (Gildea and Jurafsky, 2002), and had attracted much attention since. We use the PropBank annotated WSJ corpus, Palmer et al. (2005), which expands the Penn TreeBank with semantic annotations. These annotations were used as gold standard in the CoNLL-2005 shared task (Carreras and Marquez, 2005).

While most often the task is broken down to two parts — finding the arguments using syntactic features, then labeling them using semantic features (Marquez et al., 2008) — others try to solve the problem holistically, usually by training a statistical model such as Hidden Markov Model (HMM) or Conditional Random Fields (CRF), often solving some syntactic task jointly with SRL (Henderson et al., 2008). We adopt this approach in our work and use CRF for performing SRL.

Very few works attempt SRL without any syntactic features: Boxwell et al. (2011) report $F_1 = 0.44$ (although it is only reported for “completeness”). Recently, neural networks models have shown great promise in solving SRL. Boxwell et al. (2011), employing a unified neural network model to jointly learn POS tagging, chunking, NER and SRL, report $F_1 = 0.74$ on the SRL task. Zhou and Xu (2015), using a bi-directional long short-term memory (LSTM), report $F_1 = 81.07$, making their system state-of-the-art. These systems, however, are very complex and embody extensive fine-tuning in order to achieve the best possible results on SRL; our motivation, as previously stated, lies elsewhere.

2.2.2 Dependency Parsing

Dependency parsers build on the syntactic theory of Dependency Grammar. Proposed by Lucien Tesnière (1893-1954), the theory is based on relationships between words — between a head and a dependent. Starting from the verb, directed links connect all words of the sentence with links pointing from head to dependent, creating a directed rooted tree.

Dependency parsing is a major component of a large number of NLP applications. It is therefore one of the most well-studied tasks in NLP (McDonald et al., 2005; Nivre et al., 2007; Zhang and Nivre, 2011) and has been the focus of the CoNLL-2007 shared task (Nilsson et al., 2007). Extending classic syntactic features, Koo et al. (2008) use 4-6 bit prefixes of Brown clusters and full length clusters along with POS tags and word forms. They report an improvement in accuracy to 93.16% over a baseline of 92.02%. Bansal et al. (2014) challenge Brown clusters by using word embeddings and performing hierarchical clustering on them. They reach accuracy of 92.69%, same as their Brown baseline, but with considerably faster training time.

3 Experimental Setup

3.1 Representation

In order to employ K-means over the lexicon as well as compute an affinity matrix for spectral clustering, we first need to decide on a representation for the words in our lexicon. We choose to use a simple window-based count model, in which each word is represented by the number of times it appears within a window (with a pre-defined size) around a pre-defined set of descriptor words in some corpus.

More formally, given a corpus, we choose the set of $m$ most frequent words in the corpus to be our descriptor words, then each word $w$ in the corpus is represented as a vector in $\mathbb{R}^M$, where each coordinate denotes the number of times $w$ appears in the corpus in a window of a pre-determined size $W$ on the right size of the respective descriptor word. The result is a $N \times M$ matrix (where $N$ is the number of word in our lexicon) denoted $R$. We compute the exact same matrix using the left size of the descriptor word, and denote it $L$. These two matrices are finally concatenated to create a $N \times 2M$ sized matrix, in which each row contains the vector representation of a lexicon word, denoted the context matrix $C$.

In order to generate the context matrix we use the ukWaC corpus (Baroni et al., 2009), containing 2 billion words crawled from the .uk domain. Words are tokenized following the CoNLL format (separated ‘s, ‘nt, etc.), and the first word of a sentence is decapitalized.

For practical reasons, we artificially limit our lexicon size by choosing the top $N - 1$ most frequent words in the corpus, treating all other words as a special token "RARE" (making it the $N$-th word in our
For the experiments in this work we use $N = 12007$ and $M = 5000$ (the values were chosen empirically). In addition, we experiment with various values for the window size $W$ (2, 3 and 5).

In addition, for completeness, we perform an additional set of experiments using a state-of-the-art word embedding as our lexicon words representation. We chose to use word2vec, implementing the continuous skip-gram algorithm presented in Mikolov et al. (2013a) with negative sampling (Mikolov et al., 2013b).

### 3.2 Clustering

#### 3.2.1 Number of Clusters

We set the number of clusters on $k = 250$ throughout our experiments for all the clustering methods used in this work. This value was chosen empirically, having proven to provide the best results for both spectral and Brown clusters on the SRL task in preliminary experiments.

#### 3.2.2 Affinity Matrix

As noted in Section 2.1.2, the spectral clustering algorithm required an affinity matrix, representing the affinity between each pair of samples (lexicon words) as input. Rather than using the default Gaussian kernel (Ng et al., 2002), we follow Sun and Korhonen (2009) using the symmetrized skew-divergence for generating the affinity matrix. This approach was found to produce better results during preliminary testing. Given two vectors $v$ and $v'$, their skew-divergence is

$$d_{skew}(v, v') = D_{KL}(v' || a \cdot v + (1 - a) \cdot v')$$

Where $D_{KL}$ is the KL-divergence and $v$ is smoothed with $v'$ by parameter $a$ (we empirically choose $a = 0.999$). The symmetrized skew-divergence is then defined as:

$$d_{s-skew}(v, v') = \frac{1}{2}(d_{skew}(v, v') + d_{skew}(v', v))$$

Finally, the affinity matrix is computed using this measure. Given the $i$-th and $j$-th lexicon words ($i \neq j$), we compute their respective vector representations $v_i, v_j$. We then set

$$A_{ij} = A_{ji} = d_{s-skew}(v, v')$$

For each $i$, we set $A_{ii} = 0$.

### 3.3 Software

#### 3.3.1 Word Embeddings

In some of our experiments we utilize the word2vec embedding as word representation. We use the word embeddings available at [https://code.google.com/p/word2vec/](https://code.google.com/p/word2vec/). These embeddings were produced by a network which was trained on a partial Google News data set ($\sim$100 billion words), and generated 300-dimensional vectors.

#### 3.3.2 K-means

The K-means algorithm is employed as a baseline, as well as for the final stage in the spectral clustering algorithm (see Section 2.1.2). We use MATLAB’s implementation of Lloyd’s algorithm (Lloyd, 1982), with the K-MEANS++ algorithm for centroid initialization (Arthur and Vassilvitskii, 2007).

#### 3.3.3 Spectral Clustering

We use the spectral clustering package presented in [Cour et al., 2004](http://www.timotheecour.com/software/ncut/ncut.html), implementing the NCut spectral segmentation algorithm presented in [Shi and Malik, 2000](http://www.timotheecour.com/software/ncut/ncut.html), available at [http://www.timotheecour.com/software/ncut/ncut.html](http://www.timotheecour.com/software/ncut/ncut.html).
3.3.4 Brown Clustering
We use the implementation provided by Liang (2005), available at [https://github.com/percyliang/brown-cluster](https://github.com/percyliang/brown-cluster). The clusters were computed on the ukWaC corpus described in Section 3.1.

3.3.5 CRF
For the purpose of solving SRL we wanted a simple yet powerful, off-the-shelf learning algorithm. As stated in Section 2.2.1, we choose to use CRF, feeling it is a simple yet powerful enough tool that would allow us to place the focus on the features rather than the learning process. We use the CRF++ package, available at [https://taku910.github.io/crfpp/](https://taku910.github.io/crfpp/) implementing Lafferty et al. (2001)’s algorithm.

3.3.6 Dependency Parser
For our dependency parsing experiments we utilized the MSTParser, implementing the parser described in (McDonald et al., 2005) supplemented with second order features (McDonald and Pereira, 2006), available at [http://sourceforge.net/projects/mstparser/](http://sourceforge.net/projects/mstparser/).

4 Experiments & Results
4.1 SRL
Our first set of experiments is conducted on the PropBank annotated WSJ corpus (Palmer et al., 2005), which expands the Penn TreeBank with semantic annotations. We train a CRF model on sections 2-21 and test it on section 23. The features set for the CRF algorithm includes five surface features: whether the word contains a number, a hyphen or a capital letter, the position of the word w.r.t. current predicate and the word’s length in characters. In each experiment we add one of the following sets of features:

1. K-means clusters (using simple count model described in 3.1)
2. Spectral clusters (using simple count model described in 3.1)
3. Brown clusters
4. K-means clusters (using word2vec)
5. Spectral clusters (using word2vec)
6. POS tag
7. POS tag + dependency parent’s POS tag
8. POS tag + dependency parent’s POS tag + dependency grandparent’s POS tag

For each experiment we report per-argument precision, recall and $F_1$ score (Table 1)

Exercising the first 3 rows in Table 1 we note the very small difference in $F_1$ score between the spectral model and the Brown model (0.003), denoting virtually equal performance. This is a very interesting result, considering that Brown clustering is tailored for word clustering by incorporating a statistical language model in the target function used for the clustering itself, as opposed to spectral clustering which takes no lexical considerations into account during the clustering process. Both models improve over the baseline (K-means) by approximately 0.04.

Moving down in the table, we see that experimenting with the word2vec embedding as initial representation yields worse results than using the simple count model. This is also a surprising result given the amount of success word2vec is having in word similarity tasks (Baroni et al., 2014). We suspect

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1 Using a combination of all three window sizes, discussed in Section 3.1
2 Dependency parses were extracted from the CoNLL-2005 Shared Task annotated data set (Carreras and Márquez, 2005).
| Feature Set                              | precision | recall  | $F_1$  |
|-----------------------------------------|-----------|---------|--------|
| K-means (count model)                   | 0.620     | 0.504   | 0.555  |
| Spectral (count model)                  | 0.648     | 0.552   | 0.596  |
| Brown                                   | 0.663     | 0.547   | 0.599  |
| K-means (word2vec)                      | 0.628     | 0.471   | 0.539  |
| Spectral (word2vec)                     | 0.642     | 0.503   | 0.565  |
| POS                                     | 0.666     | 0.552   | 0.604  |
| POS + parent’s POS                      | 0.661     | 0.556   | 0.604  |
| POS + parent’s POS + grandparent’s POS  | 0.681     | 0.564   | 0.617  |

Table 1: Results for the SRL experiments.

| Feature Set                              | UAS      |
|-----------------------------------------|----------|
| Spectral (count model)                  | 0.867    |
| Brown                                   | 0.881    |
| POS                                     | 0.921    |

Table 2: Results for the dependency parsing experiments.

this is due to the fact that these embeddings are learned without any connection to the clustering task, possibly not retaining some clustering-related information due to their high density.

Moving on, we observe that using POS tags as features instead of word clusters does not significantly improve the results: merely by 0.005 over Brown clusters. Considering the POS tag of the word’s parent in the dependency tree does not improve the results, but considering the POS tag of the grandparent does improve performance by 0.013. All-in-all, we can improve over word clusters using syntactic information by the small amount of 0.018, but this requires second-order information from the dependency tree as well as POS tags, both which are generally expensive to manually produce.

4.2 Dependency Parsing

Our second set of experiments is conducted on the dependency annotations expansion to the WSJ corpus introduced in (Carreras and Marquez, 2005). We train the MSTParser on sections 2-21 of the data set and test the resulting model on section 23. We use several of the feature sets described in Section 4.1 to compare between the different clustering techniques.

For each experiment we report unlabeled attachment score (UAS). Results are shown in Table 2.

We can observe that these results are similar to the ones obtained for SRL. The difference between using Brown clusters to spectral ones is relatively small, though larger than in the SRL task (0.014 out of 0.881). Using POS improves performance here too, though by a larger margin (0.041). Over all, we observe the same phenomena, albeit on a slightly different scale.

4.3 Oracle Model

In order to analyze the amount of information overlap between spectral and Brown clusters (in the context of SRL and dependency parsing performance), we examine the performance of an oracle model.

Given a prediction task, along with two models trained for the task, an oracle model is a hypothetical model which is capable of determining which one of the two models will perform better on any given test sample. The decision is made per sample; for each sample, the oracle chooses the better model. As stated, this is obviously a hypothetical model, but it gives us the opportunity to estimate how overlapping our two models are. In the case of a complete overlap, the performance of the oracle model will not improve beyond those of the separate models; if the overlap is not complete, however, the amount of improvement achieved by the oracle may indicate the amount of difference between the models.

We perform the following procedure for the models learned using the spectral vs. the Brown clusters (feature sets 2 & 3 in 4.1). We go over the test samples one by one and check which model performed better, taking that model’s prediction to be the oracle’s prediction on that test sample. Finally, we eval-
Table 3: Results for oracle model analysis.

| Task                  | Spectral | Brown  | Oracle |
|-----------------------|----------|--------|--------|
| SRL ($F_1$)           | 0.596    | 0.599  | 0.664  |
| Dep. Parsing (UAS)    | 0.867    | 0.881  | 0.901  |

An oracle analysis reveals that spectral clusters and Brown clusters complement each other rather than completely overlap. The oracle model achieves a significant improvement of 12.5% in performance in the SRL task, with a more modest improvement of 2.3%-3.9% in the dependency parsing task. We further show that the two models agree merely on 61.4% of the test samples in the SRL test set, approximately evenly dividing the rest of samples between them (in terms of besting each other). This analysis reveals the complementary nature of these clustering methods, implying that each of them may be better suited for different cases in the same task.
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