A Graph-based Semi-Supervised Learning for Question-Answering

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Abstract

We present a graph-based semi-supervised learning for the question-answering (OA) task for ranking candidate sentences. Using textual entailment analysis, we obtain entailment scores between a natural language question posed by the user and the candidate sentences returned from search engine. The textual entailment between two sentences is assessed via features representing high-level attributes of the entailment problem such as sentence structure matching, question-type named-entity matching based on a question-classifier, etc. We implement a semi-supervised learning (SSL) approach to demonstrate that utilization of more unlabeled data points can improve the answer-ranking task of QA. We create a graph for labeled and unlabeled data using match-scores of textual entailment features as similarity weights between data points. We apply a summarization method on the graph to make the computations feasible on large datasets. With a new representation of graph-based SSL on QA datasets using only a handful of features, and under limited amounts of labeled data, we show improvement in generalization performance over state-of-the-art QA models.

1 Introduction

Open domain natural language question answering (QA) is a process of automatically finding answers to questions searching collections of text files. There are intensive research in this area fostered by evaluation-based conferences, such as the Text REtrieval Conference (TREC) (Voorhees, 2004), etc. One of the focus of these research, as well as our work, is on factoid questions in English, whereby the answer is a short string that indicates a fact, usually a named entity.

A typical QA system has a pipeline structure starting from extraction of candidate sentences to ranking true answers. In order to improve QA systems' performance many research focus on different structures such as question processing (Huang et al., 2008), information retrieval (Clarke et al., 2006), information extraction (Saggion and Gaizauskas, 2006), textual entailment (TE) (Harabagiu and Hickl, 2006) for ranking, answer extraction, etc. Our QA system has a similar pipeline structure and implements a new TE module for information extraction phase of the QA task. TE is a task of determining if the truth of a text entails the truth of another text (hypothesis). Harabagui and Hickl (2006) has shown that using TE for filtering or ranking answers can enhance the accuracy of current QA systems, where the answer of a question must be entailed by the text that supports the correctness of this answer.

We derive information from pair of texts, i.e., question as hypothesis and candidate sentence as the text, potentially indicating containment of true answer, and cast the inference recognition as classification problem to determine if a question text follows candidate text. One of the challenges we face with is that we have very limited amount of labeled data, i.e., correctly labeled (true/false entailment) sentences. Recent research indicates that using labeled and unlabeled data in semi-supervised learning (SSL) environment, with an emphasis on graph-based methods, can improve the performance of information extraction from data for tasks such as question classification (Tri et al., 2006), web classification (Liu et al., 2006), relation extraction (Chen et al., 2006), passage-retrieval (Otterbacher et al., 2009), various natural language processing tasks such as partof-speech tagging, and named-entity recognition (Suzuki and Isozaki, 2008), word-sense disambiguation (Niu et al., 2005), etc.

We consider situations where there are much more unlabeled data, X_U , than labeled data, X_L , i.e., $n_L \ll n_U$. We construct a textual entailment (TE) module by extracting features from each paired question and answer sentence and designing a classifier with a novel yet feasible graphbased SSL method. The main contributions are:

- construction of a TE module to extract matching structures between question and answer sentences, i.e., q/a pairs. Our focus is on identifying good matching features from q/a pairs, concerning different sentence structures in section 2,

- representation of our linguistic system by a form of a special graph that uses TE scores in designing a novel affinity matrix in section 3,

- application of a graph-summarization method to enable learning from a very large unlabeled and rather small labeled data, which would not have been feasible for most sophisticated learning tools in section 4. Finally we demonstrate the results of experiments with real datasets in section 5.

2 Feature Extraction for Entailment

Implementation of different TE models has previously shown to improve the QA task using supervised learning methods (Harabagiu and Hickl, 2006). We present our recent work on the task of QA, wherein systems aim at determining if a text returned by a search engine contains the correct answer to the question posed by the user. The major categories of information extraction produced by our QA system characterizes features for our TE model based on analysis of q/a pairs. Here we give brief descriptions of only the major modules of our QA due to space limitations.

2.1 Pre-Processing for Feature Extraction

We build the following pre-processing modules for feature extraction to be applied prior to our textual entailment analysis.

Question-Type Classifier (QC): QC is the task of identifying the type of a given question among a predefined set of question types. The type of a question is used as a clue to narrow down the search space to extract the answer. We used our QC system presented in (Huang et al., 2008), which classifies each question into 6-coarse categories (i.e., *abbr., entity, human, location, number, description*) as well as 50-fine categories (i.e., *color, food, sport, manner*, etc.) with almost 90% accuracy. For instance, for question "How many states are there in US?", the question-type would be 'NUMBER' as course category, and 'Count' for the finer category, represented jointly as NUM:Count. The QC model is trained via support vector machines (SVM) (Vapnik, 1995) considering different features such as semantic headword feature based on variation of Collins rules, hypernym extraction via Lesk word disambiguation (Lesk, 1988), regular expressions for whword indicators, n-grams, word-shapes(capitals), etc. Extracted question-type is used in connection with our Named-Entity-Recognizer, to formulate question-type matching feature, explained next.

Named-Entity Recognizer (NER): This component identifies and classifies basic entities such as proper names of person, organization, product, location; time and numerical expressions such as year, day, month; various measurements such as weight, money, percentage; contact information like address, web-page, phone-number, etc. This is one of the fundamental layers of information extraction of our QA system. The NER module is based on a combination of user defined rules based on Lesk word disambiguation (Lesk, 1988), WordNet (Miller, 1995) lookups, and many userdefined dictionary lookups, e.g. renown places, people, job types, organization names, etc. During the NER extraction, we also employ phrase analysis based on our phrase utility extraction method using Standford dependency parser ((Klein and Manning, 2003)). We can categorize entities up to 6 coarse and 50 fine categories to match them with the NER types from QC module.

Phrase Identification(PI): Our PI module undertakes basic syntactic analysis (shallow parsing) and establishes simple, un-embedded linguistic structures such as noun-phrases (NN), basic prepositional phrases (PP) or verb groups (VG). In particular PI module is based on 56 different semantic structures identified in Standford dependency parser in order to extract meaningful compound words from sentences, e.g., "They heard <u>high pitched cries.</u>". Each phrase is identified with a head-word (*cries*) and modifiers (*high pitched*).

Questions in Affirmative Form: To derive linguistic information from pair of texts (statements), we parse the question and turn into affirmative form by replacing the *wh*-word with a placeholder and associating the question word with the question-type from the QC module. For example: "What is the capital of France?" is written in affirmative form as " $[X]_{LOC:City}$ is the capital of France_{LOC:Country}.". Here X is the answer text of LOC:City NER-type, that we seek.

Sentence Semantic Component Analysis: Using shallow semantics, we decode the underlying dependency trees that embody linguistic relationships such as head-subject (H-S), head-modifier (complement) (H-M), head-object (H-O), etc. For instance, the sentence "*Bank of America acquired Merrill Lynch in 2008.*" is partitioned as:

- Head (H): acquired
- Subject (S): Bank of America_[Human:group]
- Object (O): *Merrill Lynch*_[Human:group]
- Modifier (M): 2008_[Num:Date]

These are used as features to match components of questions like "Who purchased Merrill Lynch?".

Sentence Structure Analysis: In our question analysis, we observed that 98% of affirmed questions did not contain any object and they are also in *copula (linking) sentence* form that is, they are only formed by subject and information about the subject as: {*subject + linking-verb + subjectinfo.*}. Thus, we investigate such affirmed questions different than the rest and call them *copula* sentences and the rest as *non-copula* sentences. ¹ For instance our system recognizes affirmed question "*Fred Durst's group name is* [X]_{DESC:Def}". as *copula-sentence*, which consists of subject (underlined) and some information about it.

2.2 Features from Paired Sentence Analysis

We extract the TE features based on the above lexical, syntactic and semantic analysis of q/a pairs and cast the QA task as a classification problem. Among many syntactic and semantic features we considered, here we present only the major ones:

(1) (QTCF) Question-Type-Candidate Sentence NER match feature: Takes on the value '1' when the candidate sentence contains the fine NER of the question-type, '0.5' if it contains the coarse NER or '0' if no NER match is found.

(2) (QComp) Question component match features: The sentence component analysis is applied on both the affirmed question and the candidate sentence pairs to characterize their semantic components including subject(S), object(O), head (H) and modifiers(M). We match each semantic component of a question to the best matching component of a candidate sentence. For example for the given question, "When did Nixon die?", when the following candidate sentence, i.e., "Richard Nixon, 37th President of USA, passed away of stroke on April 22, 1994." is considered, we extract the following component match features:

- Head-Match: $die \rightarrow pass away$
- Subject-Match: Nixon→Richard Nixon
- Object-Match: -
- Modifier-Match: $[X] \rightarrow April 22, 1994$

In our experiments we observed that converted questions have at most one subject, head, object and a few modifiers. Thus, we used one feature for each and up to three for M-Match features. The feature values vary based on matching type, i.e., exact match, containment, synonym match, etc. For example, the S-Match feature will be "1.0" due to head-match of the noun-phrase.

(3) (LexSem) Lexico-Syntactic Alignment Features: They range from the ratio of consecutive word overlap between converted question (Q) and candidate sentence (S) including --Unigram/Bigram, selecting individual/pair of ad-

jacent tokens in Q matching with the S -Noun and verb counts in common, separately.

-When words don't match we attempt matching synonyms in WordNet for most common senses. -Verb match statistics using WordNet's *cause* and *entailment* relations.

As a result, each q/a pair is represented as a feature vector $x_i \in \Re^d$ characterizing the entailment information between them.

3 Graph Based Semi-Supervised Learning for Entailment Ranking

We formulate semi-supervised entailment rank scores as follows. Let each data point in $X = \{x_1, ..., x_n\}, x_i \in \mathbb{R}^d$ represents information about a question and candidate sentence pair and $Y = \{y_1, ..., y_n\}$ be their output labels. The labeled part of X is represented with $X_L = \{x_1, ..., x_l\}$ with associated labels $Y_L = \{y_1, ..., y_l\}^T$. For ease of presentation we concentrate on binary classification, where y_i can take on either of $\{-1, +1\}$ representing entailment or non-entailment. X has also unlabeled part, $X_U = \{x_1, ..., x_u\}$, i.e., $X = X_L \cup X_U$. The aim is to predict labels for X_U . There are also other testing points, X_{Te} , which has the same properties as X.

Each node V in graph g = (V, E) represents a feature vector, $x_i \in \Re^d$ of a q/a pair, characteriz-

¹One option would have been to leave out the non-copula questions and build the model for only copula questions.

ing their entailment relation information. When all components of a hypothesis (affirmative question) have high similarity with components of text (candidate sentence), then entailment score between them would be high. Another pair of q/a sentences with similar structures would also have high entailment scores as well. So similarity between two q/a pairs x_i, x_j , is represented with $w_{ij} \in \Re^{n \times n}$, i.e., edge weights, and is measured as:

$$w_{ij} = 1 - \sum_{q=1}^{d} \frac{|x_{iq} - x_{jq}|}{d}$$
(1)

As total entailment scores get closer, the larger their edge weights would be. Based on our sentence structure analysis in section 2, given dataset can be further separated into two, i.e., X_{cp} containing q/a pairs in which affirmed questions are copula-type, and X_{ncp} containing q/a pairs with non-copula-type affirmed questions. Since copula and non-copula sentences have different structures, e.g., copula sentences does not usually have objects, we used different sets of features for each type. Thus, we modify edge weights in (1) as follows:

$$\tilde{w}_{ij} = \begin{cases} 0 & x_i \in X_{cp}, x_j \in X_{ncp} \\ 1 - \sum_{q=1}^{d_{cp}} \frac{|x_{iq} - x_{jq}|}{d_{cp}} & x_i, x_j \in X_{cp} \\ 1 - \sum_{q=1}^{d_{ncp}} \frac{|x_{iq} - x_{jq}|}{d_{ncp}} & x_i, x_j \in X_{ncp} \end{cases}$$
(2)

The diagonal degree matrix **D** is defined for graph g by $\mathbf{D}=\sum_{j} \tilde{w}_{ij}$. In general graph-based SSL, a function over the graph is estimated such that it satisfies two conditions: 1) close to the observed labels , and 2) be smooth on the whole graph by:

$$\arg\min_{f} \sum_{i \subset L} (f_i - y_i)^2 + \lambda \sum_{i,j \in L \cup U} \tilde{w}_{ij} (f_i - f_j)^2$$
(3)

The second term is a regularizer to represent the label smoothness, $f^T \mathbf{L} f$, where $\mathbf{L} = \mathbf{D} - \mathbf{W}$ is the graph Laplacian. To satisfy the local and global consistency (Zhou et al., 2004), normalized combinatorial Laplacian is used such that the second term in (3) is replaced with normalized Laplacian, $\mathcal{L} = \mathbf{D}^{-1/2} \mathbf{L} \mathbf{D}^{-1/2}$, as follows:

$$\sum_{i,j\in L\cup U} w_{ij} \left(\frac{f_i}{\sqrt{d_i}} - \frac{f_j}{\sqrt{d_j}}\right)^2 = f^T \mathcal{L} f \qquad (4)$$

Setting gradient of loss function to zero, optimum f^* , where $Y = \{Y_L \cup Y_U\}$, $Y_U = \{y_{l+1}^n = 0\}$;

$$f^* = (\mathbb{1} + \lambda \, (\mathbb{1} - \mathcal{L}))^{-1} \, Y \tag{5}$$

Most graph-based SSLs are transductive, i.e., not easily expendable to new test points outside $L \cup U$. In (Delalleau et al., 2005) an induction scheme is proposed to classify a new point x_{Te} by

$$\hat{f}(x_{Te}) = \frac{\sum_{i \in L \cup U} w_{x_i} f_i}{\sum_{i \in L \cup U} w_{x_i}}$$
(6)

Thus, we use induction, where we can, to avoid re-construction of the graph for new test points.

4 Graph Summarization

Research on graph-based SSL algorithms point out their effectiveness on real applications, e.g., (Zhu et al., 2003), (Zhou and Schölkopf, 2004), (Sindhwani et al., 2007). However, there is still a need for fast and efficient SSL methods to deal with vast amount of data to extract useful information. It was shown in (Delalleau et al., 2006) that the convergence rate of the propagation algorithms of SSL methods is $O(kn^2)$, which mainly depends on the form of eigenvectors of the graph Laplacian (k is the number of nearest neighbors). As the weight matrix gets denser, meaning there will be more data points with connected weighted edges, the more it takes to learn the classifier function via graph. Thus, the question is, how can one reduce the data points so that weight matrix is sparse, and it takes less time to learn?

Our idea of summarization is to create *representative vertices* of data points that are very close to each other in terms of edge weights. Suffice to say that similar data points are likely to represent denser regions in the hyper-space and are likely to have same labels. If these points are close enough, we can characterize the boundaries of these group of similar data points with respect to graph and then capture their summary information by new *representative vertices*. We replace each data point within the boundary with their representative vertex, to form a summary graph.

4.1 Graph Summarization Algorithm

Let each selected dataset be denoted as $X^s = \{x_i^s\}$, i = 1...m, s = 1, ..., q, where m is the number of data points in the sample dataset and q is the number of sample datasets drawn from X. The labeled data points, i.e., X_L , are appended to each of these selected X^s datasets, $X^s = \{x_1^s, ..., x_{m-l}^s\} \cup X_L$. Using a separate learner, e.g., SVM (Vapnik, 1995), we obtain predicted outputs, $\hat{Y^s} = (\hat{y}_1^s, ..., \hat{y}_{m-l}^s)$ of X^s and append observed labels $\hat{Y^s} = \hat{Y^s} \cup Y_L$.



Figure 1: Graph Summarization. (a) Actual data point with predicted class labels, (b) magnified view of a single node (black) and its boundaries (c) calculated representative vertex, (d) summary dataset.

We define the weight W^s and degree D^s matrices of X^s using (1). Diagonal elements of D^s is converted into a column vector and is sorted to find the high degree vertices that are surrounded with large number of close neighbors.

The algorithm starts from the highest degree node $x_i^s \in X^s$, where initial neighbor nodes have assumably the same labels. This is shown in Figure 1-(b) with the inner square around the middle black node, corresponding high degree node. If its immediate k neighbors, dark blue colored nodes, have the same label, the algorithm continues to search for the secondary k neighbors, the light blue colored nodes, i.e., the neighbors of the neighbors, to find out if there are any opposite labeled nodes around. For instance, for the corresponding node (black) in Figure 1-(b) we can only go up to two neighbors, because in the third level, there are a few opposite labeled nodes, in red. This indicates boundary B_i^s for a corresponding node and unique nearest neighbors of same labels.

$$B_{i}^{s} = \left\{ x_{i}^{s} \cup \left\{ x_{j}^{s} \right\}_{j=1}^{nm} \right\}$$
(7)

In (7), nm denotes the maximum number of nodes of a B_i^s and $\forall x_j^s, x_{j'}^s \in B_i^s, y_j^s = y_{j'}^s = y_{B_i^s}^s$, where $y_{B_i^s}$ is the label of the selected boundary B_i^s .

We identify the edge weights w_{ij}^s between each node in the boundary B_i^s via (1), thus the boundary is connected. We calculate the weighted average of the vertices to obtain the representative summary node of B_i^s as shown in Figure 1-(c);

$$\overline{X}_{B_i}^s = \frac{\sum_{i \neq j=1}^{nm} \frac{1}{2} w_{ij}^s (x_i^s + x_j^s)}{\sum_{i \neq j=1}^{nm} w_{ij}^s}$$
(8)

The boundaries of some nodes may only contain themselves because their immediate neighbors may have opposite class labels. Similarly some may have only k + 1 nodes, meaning only immediate neighbor nodes have the same labels. For instance in Fig. 1 the boundary is drawn after the secondary neighbors are identified (dashed outer boundary). This is an important indication that some representative data points are better indicators of class labels than the others due to the fact that they represent a denser region of same labeled points. We represent this information with the local density constraints. Each new vertex is associated with a local density constraint, $0 \le \delta_i \le 1$, which is equal to the total number of neighboring nodes used to construct it. We use the normalized density constraints for ease of calculations. Thus, for a each sample summary dataset, a local density constraint vector is identified as $\delta^s = \{\delta_1^s, ..., \delta_{nb}^s\}^T$. The local density constraints become crucial for inference where summarized labeled data are used instead of overall dataset.

Algorithm 1	Graph Summa	ary of Large Dataset
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1: Given $X = \{x_1, ..., x_n\}, X = X_L \cup X_U$

2: Set $q \leftarrow \max$ number of subsets

3: for $s \leftarrow 1, ..., q$ do

Choose a random subset with repetitions 4:

- $X^{s} = \{x_{1}^{s}, ..., x_{m-l}^{s}, x_{m-l+1}, ..., x_{m}\}$ 5: 6: Summarize X^s to obtain \overline{X}^s in (9)
- 7: end for
- Obtain summary dataset $\overline{X} = \{\overline{X}^s\}_{s=1}^q = \{\overline{X}_i\}_{i=1}^p$ and local density constrains, $\delta = \{\delta_i\}_{i=1}^p$. 8:

After all data points are evaluated, the sample dataset X^s can now be represented with the summary representative vertices as

$$\overline{X}^{s} = \left\{ \overline{X}^{s}_{B_{1}}, ..., \overline{X}^{s}_{B_{nb}} \right\}.$$
(9)

and corresponding local density constraints as,

$$\delta^{s} = \{\delta_{1}^{s}, ..., \delta_{nb}^{s}\}^{T}, \ 0 < \delta_{i}^{s} \le 1$$
 (10)

The summarization algorithm is repeated for each random subset X^s , s = 1, ..., q of very large dataset $X = X_L \cup X_U$, see Algorithm 1. As a result q number of summary datasets \overline{X}^s each of which with nb labeled data points are combined to form a representative sample of X, $\overline{X} =$ $\{\overline{X}^s\}_{s=1}^q$ reducing the number of data from n to a much smaller number of data, $p = q * nb \ll n$. So the new summary of the X can be represented with $\overline{X} = {\{\overline{X}_i\}}_{i=1}^p$. For example, an original dataset with 1M data points can be divided up to q = 50 random samples of m = 5000data points each. Then using graph summarization each summarized dataset may be represented with $nb \cong 500$ data points. After merging summarized data, final summarized samples compile to $500 * 50 \cong 25K \ll 1M$ data points, reduced to 1/40 of its original size. Each representative data point in the summarized dataset \overline{X} is associated with a local density constraints, a p = q * nbdimensional row vector as $\delta = {\delta_i}_{i=1}^p$.

We can summarize a graph separately for different sentence structures, i.e., copula and noncopula sentences. Then representative data points from each summary dataset are merged to form final summary dataset. The Hybrid graph summary models in the experiments follow such approach.

4.2 Prediction of New Testing Dataset

Instead of using large dataset, we now use summary dataset with predicted labels, and local density constraints to learn the class labels of *nte* number of unseen data points, i.e., testing data points, $X_{Te} = \{x_1, ..., x_{nte}\}$. Using graph-based SSL method on the new representative dataset, $X' = \overline{X} \cup X_{Te}$, which is comprised of summarized dataset, $\overline{X} = \{\overline{X}_i\}_{i=1}^p$, as labeled data points, and the testing dataset, X_{Te} as unlabeled data points. Since we do not know estimated local density constraints of unlabeled data points, we use constants to construct local density constraint column vector for X' dataset as follows:

$$\delta' = \{1 + \delta_i\}_{i=1}^p \cup [1 \dots 1]^T \in \Re^{nte}$$
(11)

 $0 < \delta_i \leq 1$. To embed the local density constraints, the second term in (3) is replaced with the constrained normalized Laplacian, $\mathcal{L}^c = \delta^T \mathcal{L} \delta$,

$$\sum_{i,j\in L\cup T} w_{ij} \left(\frac{f_i}{\sqrt{\delta'_i * d_i}} - \frac{f_j}{\sqrt{\delta'_j * d_j}}\right)^2 = f^T \mathcal{L}^c f$$
(12)

If any testing vector has an edge between a labeled vector, then with the usage of the local density constraints, the edge weights will not not only be affected by that labeled node, but also how dense that node is within that part of the graph.

5 Experiments

We demonstrate the results from three sets of experiments to explore how our graph representation, which encodes textual entailment information, can be used to improve the performance of the QA systems. We show that as we increase the number of unlabeled data, with our graphsummarization, it is feasible to extract information that can improve the performance of QA models.

We performed experiments on a set of 1449 questions from TREC-99-03. Using the search engine², we retrieved around 5 top-ranked candidate sentences from a large newswire corpus for each question to compile around 7200 q/a pairs. We manually labeled each candidate sentence as true or false entailment depending on the containment of the true answer string and soundness of the entailment to compile quality training set. We also used a set of 340 QA-type sentence pairs from RTE02-03 and 195 pairs from RTE04 by converting the hypothesis sentences into question form to create additional set of q/a pairs. In total, we created *labeled* training dataset X_L of around 7600 q/a pairs . We evaluated the performance of graphbased QA system using a set of 202 questions from the TREC04 as testing dataset (Voorhees, 2003), (Prager et al., 2000). We retrieved around 20 candidate sentences for each of the 202 test questions and manually labeled each q/a pair as true/false entailment to compile 4037 test data.

To obtain more *unlabeled* training data X_U , we extracted around 100,000 document headlines from a large newswire corpus. Instead of matching headline and first sentence of the document as in (Harabagiu and Hickl, 2006), we followed a different approach. Using each headline as a query, we retrieved around 20 top-ranked sentences from search engine. For each headline, we picked the 1st and the 20th retrieved sentences. Our assumption is that the first retrieved sentence may have higher probability to entail the headline, whereas the last one may have lower probability. Each of these headline-candidate sentence pairs is used as additional unlabeled q/a pair. Since each head-

²http://lucene.apache.org/java/

Features	Model	MRR	Top1	Top5
Baseline	_	42.3%	32.7%	54.5%
QTCF	SVM	51.9%	44.6%	63.4%
	SSL	49.5%	43.1%	60.9%
LexSem	SVM	48.2%	40.6%	61.4%
	SSL	47.9%	40.1%	58.4%
QComp	SVM	54.2%	47.5%	64.3%
	SSL	51.9%	45.5%	62.4%

Table 1: MRR for different features and methods.

line represents a converted question, in order to extract the question-type feature, we use a matching NER-type between the headline and candidate sentence to set question-type NER match feature.

We applied pre-processing and feature extraction steps of section 2 to compile labeled and unlabeled training and labeled testing datasets. We use the rank scores obtained from the search engine as baseline of our system. We present the performance of the models using Mean Reciprocal Rank (MRR), top 1 (Top1) and top 5 prediction accuracies (Top5) as they are the most commonly used performance measures of QA systems (Voorhees, 2004). We performed manual iterative parameter optimization during training based on prediction accuracy to find the best k-nearest parameter for SSL, i.e., $k = \{3, 5, 10, 20, 50\}$, and best $C = \{10^{-2}, ..., 10^2\}$ and $\gamma = \{2^{-2}, ..., 2^3\}$ for RBF kernel SVM. Next we describe three different experiments and present individual results.

Graph summarization makes it feasible to execute SSL on very large unlabeled datasets, which was otherwise impossible. This paper has no assumptions on the performance of the method in comparison to other SSL methods.

Experiment 1. Here we test individual contribution of each set of features on our QA system. We applied SVM and our graph based SSL method with no summarization to learn models using labeled training and testing datasets. For SSL we used the training as labeled and testing as unlabeled dataset in transductive way to predict the entailment scores. The results are shown in Table 1. From section 2.2, QTCF represents question-type NER match feature, LexSem is the bundle of lexico-semantic features and QComp is the matching features of subject, head, object, and three complements. In comparison to the baseline, QComp have a significant effect on the accuracy of the QA system. In addition, QTCF has shown to improve the MRR performance by about 22%. Although the LexSem features have minimal semantic properties, they can improve MRR performance by 14%.

Experiment 2. To evaluate the performance of graph summarization we performed two separate experiments. In the first part, we randomly selected subsets of labeled training dataset $X_{I}^{i} \subset$ X_L with different sample sizes, $n_L^i = \{1\% * n_L,$ $5\% * n_L, 10\% * n_L, 25\% * n_L, 50\% * n_L,$ $100\% * n_L$, where n_L represents the sample size of X_L . At each random selection, the rest of the labeled dataset is hypothetically used as unlabeled data to verify the performance of our SSL using different sizes of labeled data. Table 2 reports the MRR performance of QA system on testing dataset using SVM and our graph-summary SSL (gSum SSL) method using the similarity function in (1). In the second part of the experiment, we applied graph summarization on copula and noncopula questions separately and merged obtained representative points to create labeled summary dataset. Then using similarity function in (2) we applied SSL on labeled summary and unlabeled testing via transduction. We call these models as Hybrid gSum SSL. To build SVM models in the same way, we separated the training dataset into two based on copula and non-copula questions, X_{cp}, X_{ncp} and re-run the SVM method separately. The testing dataset is divided into two accordingly. Predicted models from copula sentence datasets are applied on copula sentences of testing dataset and vice versa for non- copula sentences. The predicted scores are combined to measure overall performance of Hybrid SVM models. We repeated the experiments five times with different random samples and averaged the results.

Note from Table 2 that, when the number of labeled data is small $(n_L^i < 10\% * n_L)$, graph based SSL, gSum SSL, has a better performance compared to SVM. As the percentage of labeled points in training data increase, the SVM performance increases, however graph summary SSL is still comparable with SVM. On the other hand, when we build separate models for copula and non-copula questions with different features, the performance of the overall model significantly increases in both methods. Especially in Hybrid graph-Summary SSL, Hybrid gSum SSL, when the number of labeled data is small $(n_L^i < 25\% * n_L)$ performance improvement is better than rest

%	SVM		gSum SSL		Hybrid SVM		Hybrid gSum SSL					
#Labeled	MRR	Top1	Top5	MRR	Top1	Top5	MRR	Top1	Top5	MRR	Top1	Top5
1%	45.2	33.2	65.8	56.1	44.6	72.8	51.6	40.1	70.8	59.7	47.0	75.2
5%	56.5	45.1	73.0	57.3	46.0	73.7	54.2	40.6	72.3	60.3	48.5	76.7
10%	59.3	47.5	76.7	57.9	46.5	74.2	57.7	47.0	74.2	60.4	48.5	77.2
25%	59.8	49.0	78.7	58.4	45.0	79.2	61.4	49.5	78.2	60.6	49.0	76.7
50%	60.9	48.0	80.7	58.9	45.5	79.2	62.2	51.0	79.7	61.3	50.0	77.2
100%	63.5	55.4	77.7	59.7	47.5	79.7	67.6	58.0	82.2	61.9	51.5	78.2

Table 2: The MRR (%) results of graph-summary SSL (gSum SSL) and SVM as well as Hybrid gSum SSL and Hybrid SVM with different sizes of labeled data.

#Unlabeled	MRR	Top1	Top5
25K	62.1%	52.0%	76.7%
50K	62.5%	52.5%	77.2%
100K	63.3%	54.0%	77.2%

Table 3: The effect of number of unlabeled dataon MRR from Hybrid graph Summarization SSL.

of the models. As more labeled data is introduced, Hybrid SVM models' performance increase drastically, even outperforming the state-of-the art MRR performance on TREC04 datasets presented in (Shen and Klakow, 2006) i.e., MRR=67.0%, Top1=62.0%, Top5=74.0%. This is due to the fact that we establish two seperate entailment models for copula and non-copula q/a sentence pairs that enables extracting useful information and better representation of the specific data.

Experiment 3. Although SSL methods are capable of exploiting information from unlabeled data, learning becomes infeasible as the number of data points gets very large. There are various research on SLL to overcome the usage of large number of unlabeled dataset challenge (Delalleau et al., 2006). Our graph summarization method, Hybrid gsum SSL, has a different approach. which can summarize very large datasets into representative data points and embed the original spatial information of data points, namely local density constraints, within the SSL summarization schema. We demonstrate that as more labeled data is used, we would have a richer summary dataset with additional spatial information that would help to improve the the performance of the graph summary models. We gradually increase the number of unlabeled data samples as shown in Table 3 to demonstrate the effects on the performance of testing dataset. The results show

that the number of unlabeled data has positive effect on performance of graph summarization SSL.

6 Conclusions and Discussions

In this paper, we applied a graph-based SSL algorithm to improve the performance of QA task by exploiting unlabeled entailment relations between affirmed question and candidate sentence pairs. Our semantic and syntactic features for textual entailment analysis has individually shown to improve the performance of the QA compared to the baseline. We proposed a new graph representation for SSL that can represent textual entailment relations while embedding different question structures. We demonstrated that summarization on graph-based SSL can improve the QA task performance when more unlabeled data is used to learn the classifier model.

There are several directions to improve our work: (1) The results of our graph summarization on very large unlabeled data is slightly less than best SVM results. This is largely due to using headlines instead of affirmed questions, wherein headlines does not contain question-type and some of them are not in proper sentence form. This adversely effects the named entity match of questiontype and the candidate sentence named entities as well as semantic match component feature extraction. We will investigate experiment 3 by using real questions from different sources and construct different test datasets. (2) We will use other distance measures to better explain entailment between q/a pairs and compare with other semisupervised and transductive approaches.

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