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Machine learning inference of search engine heuristics

Part II Project

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Declaration of Originality

I, Karina Palyutina of St Catharine's College, being a candidate for Part II of the Computer Science Tripos, hereby declare that this dissertation and the work described in it are my own work, unaided except as may be specified below, and that the dissertation does not contain material that has already been used to any substantial extent for a comparable purpose.

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Date

Contents

In	Introduction 1							
1	\mathbf{Pre}	paration	3					
	1.1	Formulating the Goals	3					
		1.1.1 System Overview	3					
		1.1.2 Search Engine	5					
		1.1.3 Machine Learning	6					
	1.2	Development Strategy	11					
2	Imp	olementation	13					
	2.1	Data	13					
	2.2	Search Engine	13					
	2.3	Machine Learning	16					
		2.3.1 Naive Bayes	16					
		2.3.2 Support Vector Machine	17					
	2.4	Parser	20					
	2.5	Optimization	22					
3	Eva	Evaluation						
	3.1	The prototype	25					
	3.2	Bayes ranking: one to one comparison	25					
	3.3	SVR	25					
	3.4	Kernel function generality	25					
	3.5	Experiments with number of features	25					
	3.6	Score vs Rank	25					
4	Con	nclusion	27					
Bi	bliog	graphy	29					
\mathbf{A}	Pro	ject Proposal	31					



Introduction

This project is inspired by increasing importance of search engine rankings. To-day major search engines given a query return web pages in an order determined by secret algorithms. Such algorithms are believed to incorporate multiple unknown factors. For instance, Google claims to have over 200 unique factors that influence a position of a webpage in the search results relative to a query ². Only a handful of these factors are disclosed to the webmasters in the form of very general guidelines. Moreover, the Google algorithm in particular is updated frequently. However, most of the knowledge around the area amounts to speculation. Despite the fact that it is possible to pass a vast number of queries through the black box of any existing search engine, the immensity of the search space, and instability of such algorithms make them impossible to reverse engineer.

Machine learning is a natural approach to inferring the true algorithm from a subset of all possible observations. However, applying machine learning techniques to real search engines would be hardly effective, as the dynamic nature of the algorithms and the web as well as lack of meaningful feedback would prevent incremental improvement: when there are as many as 200 features in question, false assumptions made by a learner may have an unpredictable effect on its performance.

More generally, there are certain ambiguities associated with machine learning, which are 'problem-specific'. For example, it proves difficult to decide how much training data is necessary, as well as and selecting it to avoid over/under-fitting[1]. Similarly, it is not straightforward which machine learning technique is best for a particular problem.

This project is concerned with application of machine learning techniques to search engines. The aim of the project, in particular, is to explore how machine learning techniques can be used effectively to infer algorithms from search engines. To address the limitations imposed by existing search engines, part of the task is to develop a toy search engine that allows me to comtrol the nature and complexity of used heuristics. Such transparency addresses the problems stated

²http://www.google.com/competition/howgooglesearchworks.html

above and, more importantly, allows for useful evaluation of machine learning techniques by providing meaningful feedback.

Even though this study does not attempt to reverse engineer any existing heuristics, the results can be applied to such an ambitious task. Moreover, such a framework is potentially more general and can be used for a range of problems.

TODO: overview of the chapters here.

Chapter 1

Preparation

This chapter describes work that has been done before coding was started. In particular, it focuses on the reasoning behind the design of the system to be implemented. The first section is devoted to research undertaken to determine what can be done and how best to do it. The second section formulates the system requirements, namely formalizes everything that is developed in this project. The last section outlines the particulars of the software engineering approach to be adopted by this project.

1.1 Formulating the Goals

A particular difficulty in this project has been in planning what has to be done. Due to the exploratory nature of the project the course of action had to be predominantly determined by the outcome of a current tactic. Moreover, the unknowns originating from the machine learning further complicated matters.

1.1.1 System Overview

To achieve the goal of the project, a machine learning techniques comparison framework was necessary. In the Introduction I mentioned the benefit of having a transparent system as an object of learning. To further justify this decision, it is worth mentioning that generalisation using machine learning is different from most optimization problems in that the function that is being optimized is out of our reach, and all that is visible to the machine learner is the training error. Because our goal is not the correct classification of real data, but identifying the means to correct classification, it is important that informed choices are made towards improvement of the learner. Taking this into account, knowing the

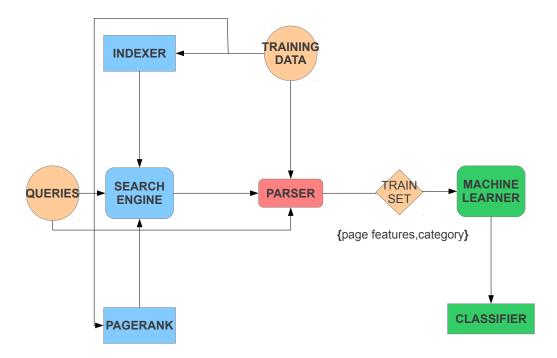


Figure 1.1: Overview of the system. Three major parts from left to right are search engine, parser and machine learner.

function that we want to learn and having direct control over it will guide the improvement of the search engine.

This argument motivates a system in three parts: a search engine, a machine learner and a parser to mediate between the two. Figure 1.1 illustrates the proposed learning system. Training data is a set of web pages set aside specifically for training purposes.

Choice of Programming Language

When choosing a programming language, main considerations reduced to library availability and simplicity. The project imposes no special requirements on the language, apart from, perhaps, library infrastructure for parsing web pages. Python is simple language with extensive library support. As for efficiency, all the mathematical operations in this project rely on python math libraries, which are implemented in C. I have not programmed in Python before the project, so a slight overhead was caused by having to learn a new language.

Data

Web pages used as Training and Test data are not required to have any special properties, but diversity and typicality are seen as advantageous. As for the size of the training data, Domingos [1] suggests that a primitive learner given more data performs better than a more complex one with less data. This, of course, is under certain assumptions of data quality, namely the assumption that the training data is a representative subset of all the possible data. Intuitively, provided there is no bias in data gathering, more data implies better generality. I have started with a training set spanning an order of a few thousands of pages, however, in practice, I found that there is no particular improvement beyond a thousand pages. **TODO:** Link to relevant part or example data here?

1.1.2 Search Engine

Next important decision regarded the search engine. Originally, I considered using open source existing engines, in particular, Lucene. Even though I could freely modify it for the purposes of the project, the complexity of it was superfluous. I saw writing a simple search engine as a more beneficial exercise, as developing it in the first place potentially gives an insight into the problem.

Functionally our search engine is a black box that takes a set of webpages and a set of queries and outputs an order. The order is determined by the features of the page, which together make up a score. The score is the function we want to infer using the ranking assigned by the search engine, however, we are only given the order as evidence. Machine learning paragraph below will address this issue in more details.

In general there are two aspects of information retrieval that have to be accounted for: precision and recall. Precision is the fraction of retrieved pages that are relevant to the query, whereas recall is the fraction of relevant documents that are retrieved. Even though both are important for a good search engine, but in practice, the web is very large, and so precision, or even precision at n^1) has become more prominent in defining a good search engine: very rarely the user actually browses returns that are not in the top few tens of returned pages. Therefore, modern search engines tend to focus on high precision at the expense of recall [4]. Therefore, we will concentrate primarily on precision, when designing a search engine.

¹'Precision at n' only evaluates precision with respect to n topmost returned pages.

PageRank

The PageRank algorithm was originally described in a research paper by Page and Brin[5]. It was first introduced as a way 'to measure the relative importance of web pages'. PageRank is interesting to include in this project not only because it is a defining feature of the Google search engine, but because it is a unique feature of its type, as it depends on the link structure of the whole web as opposed to other web features such as word count and alike.

The basic idea is to capture the link structure of the web and provide a ranking of pages that is robust against manipulation. To achieve this, the importance 'flows' forward: the backlinks share their importance with the pages they link to. Such a simplified ranking of a set of pages can be expressed as an assignment

$$R(u) = c \sum_{v \in B_u} \frac{R(v)}{N_v} \tag{1.1}$$

An intuitive justification for such ranking is by analogy with a citation network, we are likely to find highly cited pages more important. The equation is recursive, so iterating until convergence results in a steady state distribution, which corresponds to the PageRank vector. Together with this intuition the paper introduces the *Random Surfer Model*: we are interested in a steady state distribution of a random walk on the Web graph. At each step the surfer either follows a random link or 'teleports' to a random page. The paper justifies the approach of ignoring the dangling links, as it doesn't have a significant affect on the ranking.

The teleportation vector determines whether the PageRank is personalized. For the non-personalized version the teleportation vector holds equal probabilities for all pages in the Web, whereas in a personalized approach the probabilities are distributed according the knowledge of the surfer's previous activity. In this project we are only concerned with nonpersonalized ranking, which simplifies things a little.

1.1.3 Machine Learning

I have now covered the main peripheral decisions, but it is machine learning that constitutes the central part of the project. The field was completely new to me to start with, so research of different techniques was a big part of the preparation.

Machine learning is a vast field and very little is prescribed. The supervised learning approach, where the hidden function is inferred from the labelled training data, best fits our purposes for a number of reasons. Firstly, it applies well to the real-world problem of inferring the heuristics of the real search engines, as

we can select training data and obtain the labels simply by querying. Secondly, supervised learning is widely used and offers a variety of flexible techniques.

An interesting known problem with machine learning in general is referred to as the Curse of Dimensionality. In the real world scenario – suppose if we were inferring an existing search engine from the observed query responses – we could not be sure about which features are and which are not relevant to the hidden algorithm. Potentially, some features we might choose to use will increase the dimensionality, but would not actually be used by the real algorithm. Gathering more features can hurt, as it makes the learner infer nonexistent dependencies. Clearly within the scope of this project we will know at each point which features are or are not used. So observations can be made as to how using more features than the search engine will affect the classifier.

Another major issue that is common to all machine learning methods is over/under-fitting. These refer to the problem of finding balance between the generalized model and the training data at hand. This is yet another source of interest to this project, as we can observe the behaviour of the learner when certain control parameters are changed.

It is generally recommended that the simplest learners are tried first[1]. Of all learners Naive Bayesian is one of the most comprehensible. This in itself is a major advantage according to the Occam's razor principle, which finds ample application in machine learning. Hence, we start with describing the principles of the two machine learning techniques used - Naive Bayes and Support Vector Machines.

Naive Bayes

Naive Bayes is a probabilistic classifier based on the Bayes Theorem. The posterior probability $P(C|\vec{F})$ denotes the probability that a sample page with a feature vector $\vec{F} = (F_1, F_2, \dots, F_n)$ belongs to class C. The posterior probability is computed from the observable in the training data: the prior probability P(C) – the unconditional probability of a page belonging to the class C, the likelihood $P(\vec{F}|C)$ and the evidence $P(\vec{F})$:

$$P(C|\vec{F}) = \frac{P(C)P(\vec{F}|C)}{P(\vec{F})}$$
(1.2)

The simplicity of Bayesian approach owes to the conditional independence assumption: each F_i in \vec{F} is assumed to be independent of one another to get $P(\vec{F}|C) = P(F_1|C) * P(F_2|C) * \cdots * P(F_n|C)$. This leads to a concise classifier

definition:

$$\hat{C} = argmax_C P(C) \prod_{i=1}^n P(F_i|C)$$
(1.3)

where C is the result of classification of a page with feature vector F_1, F_2, \ldots, F_n .

In practice, the crude assumption rarely holds and is likely to be violated by our data, as we expect features of pages to be interdependent. However, it has been shown that Naive Bayes performs well under zero-one loss function in presence of dependencies[2]. This has a few implications for this project, particularly, on evaluation methods

As we have seen, Naive Bayes assigns probabilities to possible classifications in the process of classifying. Even though it generally performs well in classification tasks, these probability estimates are poor [3]. However, despite poor probability estimates, there exist several frameworks, which make use of Bayesian classification and achieve decent performance in ranking. For example, Zhang [7] experimentally found that Naive Bayes is locally optimal in ranking. The paper defines a classifier as locally optimal in ranking a positive example E if there is no negative example ranked after E and vice versa for a negative exaple. A classifier is global in ranking if it is locally optimal for all examples in the example set: in other words, it is optimal in pairwise ranking. It is particularly interesting that the paper discovered that Naive Bayes is globally optimal in ranking on linear functions that have been shown as not learnable by Naive Bayes². Another framework for ranking [6] is based on Placket-Luce model, which reconciles the concepts of score and rank. This framework is based on minimizing the Bayes risk over possible permutations. Existence of such frameworks suggest that Naive Bayes is an adequate choice for a prototype learner.

Although classification is a useful technique to try, it feels more natural to represent our score or rank as real numbers rather than classes. Regression is another approach to machine learning and the next learning technique we explore – Support Vector Regression – is non-probabilistic, to try something as distant from Naive Bayes as possible.

ϵ -Support Vector Regression

While the binary classification problem has as its goal the maximization of the margin between the classes, regression is concerned with fitting a hyperplane through the given training sequence.

²*M-of-N Concepts* and *Conjunctive Concepts* can't be learnt by Naive Bayes classifier but can be optimally ranked by it according to Zhang [7].

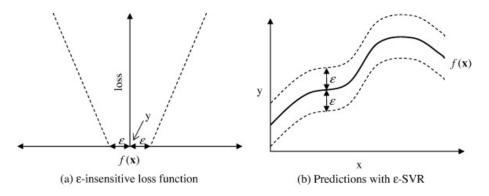


Figure 1.2: TODO: plot these yourself!

A great advantage of Support Vector machines is that they can perform nonlinear classification or regression by an implicit mapping of features into a higher dimensional space, so that the data is linearly separable. This mapping is referred to as the *Kernel Trick*. **TODO: something missing here** A training sequence is a set of training points $D = \{(\mathbf{x_1}, t_1), (\mathbf{x_2}, t_2), ..., (\mathbf{x_l}, t_l)\}$ where $\mathbf{x_i} \in R^n$ is a feature vector holding features of pages and $\mathbf{t_i} \in R$ is the corresponding ranking of each page.

In simple linear regression the aim is to minimize a regularized error function. We will be using an ϵ -insensitive error function (see Figure 1.2 (a)).

$$E_{\phi}(y((x) - t)) = \begin{cases} 0 & \text{if } |y(\mathbf{x}) - t)| < \epsilon \\ |y(\mathbf{x}) - t)| - \epsilon & \text{otherwise} \end{cases}$$

where $y((x) = \mathbf{w}^{\mathbf{T}} \phi(\mathbf{x}) + b$ is the hyperplane equation (and so $y(\mathbf{x})$ is the predicted output) and t_n is the target (true) output.

The regression tube then contains all the points for which $y(\mathbf{x_n}) - \epsilon \leq t_n \leq y(\mathbf{x_n}) + \epsilon$ as shown in Figure 1.2(b).

To allow variables to lie outside of the tube, slack variables $\xi_n \geq 0$ and $\xi_n^* \geq 0$ are introduced. The standard formulation of the error function for support vector regression (ref Vapnik 1998) can be written as follows:

$$E = C \sum_{n=1}^{N} (\xi_n + \xi_n^*) + \frac{1}{2} ||\mathbf{w}||^2$$
 (1.4)

E must be minimized subject to four constraints:

$$\xi_n \ge 0,\tag{1.5}$$

$$\xi_n^* \ge 0,\tag{1.6}$$

$$t_n \le y(\mathbf{x_n}) + \epsilon + \xi_n,\tag{1.7}$$

$$t_n \ge y(\mathbf{x_n}) - \epsilon - \xi_n^*,\tag{1.8}$$

This constraint problem can be transformed into its dual form by introducing Lagrange multipliers $a_n \geq 0, a_n^* \geq 0$. The dual problem involves maximizing

$$L(\mathbf{a}, \mathbf{a}^*) = -\frac{1}{2} \sum_{n=1}^{N} \sum_{m=1}^{N} (a_n - a_n^*) (a_m - a_m^*) K(\mathbf{x_n}, \mathbf{x_m})$$
(1.9)

$$-\epsilon \sum_{n=1}^{N} (a_n + a_n^*) + \sum_{n=1}^{N} t_n (a_n - a_n^*)$$

where $K(x_n, x_m)$ is the kernel function, t_n is the target output, subject to constraints

$$\sum_{n=1}^{N} (a_n - a_n^*) = 0, \tag{1.10}$$

$$0 \le a_n, \ a_n^* \le C, \ n = 1, ..., l$$
 (1.11)

Evaluation methodology

Two proposed techniques Naive Bayes and SVM regression are quite different, so comparing them is potentially erroneous. However, comparisons can be done within each method, as there is a lot of scope for variety of implementations in each. As a baseline for the Bayesian approach, a very primitive ranking model will be used. We will simply disregard the scoring function behind the rank and directly infer the ranking function instead. More sophisticated ranking models can then be compared to this basic performance. Figure 1.3 shows an evaluation framework for Naive Bayes. The Test Data is the data carefully set aside at the beginning that is never exposed to the learner.

For an SVM learner the baseline can be set to the performance using a linear kernel below.

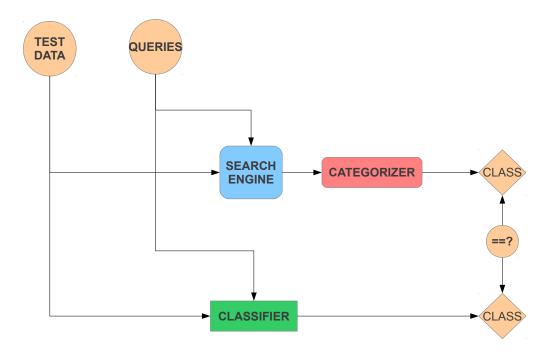


Figure 1.3: Evaluation system.

$$K(x, y) = x \cdot y$$

where \cdot denotes the dot product.

This is expected to be very high for linear heuristics but a lot lower for nonlinear ones.

1.2 Development Strategy

While the set up of Part II projects encourages waterfall-like development model, this project takes an iterative approach. The first iteration renders a prototype: a primitive search engine with a Naive Bayesian baseline classifier. The next iteration modifies each part of the system towards a more complex solution. Evaluation is performed at each iteration. Within each iteration the development follows the evolved waterfall model – the incremental build model. Each increment represents a functionally separate unit of the system: a search en-

gine, a learner, a parser and an assessment module. Increments are developed sequentially and regression testing is performed separately before integration.

The backup of the code is twofold: every time a substantial change is made a remote version control repository is updated to hold the newest version. Regularly both the code and the data used and obtained during evaluation are also backed up onto an external hard drive.

During the development of the learners a development pool of pages will be used, which must be disjoint with the Training or Test data. This ensures that no optimization is tailored to the data used for evaluation.

Requirements Analysis

The resulting system must comply with the requirements stated below.

- A search engine, which given a query must return a relevant subset of pages within 10 seconds in an order corresponding to a given heuristic, which can also be supplied to the search engine.
- Search engine must base ranking decisions on both dynamic (query dependent) and static (query independent) page features.
- A machine learner must be sufficiently isolated from the search engine through an encapsulating interface to avoid the possibility of undesirable interaction.
- Machine learners must be able to process thousands of pages in a reasonable time.
- The evaluation module must write results to persistent storage.

Chapter 2

Implementation

This section describes parts of the system that I have implemented. **TODO:** overview

2.1 Data

Development, Test and Training data for this project was implemented Separate directories..

2.2 Search Engine

The first basic block of the system – the search engine – can be logically split into two main parts: an indexer and a sorter. Because we concentrate on precision, as discussed in the Preparation chapter, we will assume for simplicity that all relevant documents are returned. This allowed me to an use existing library implementation of the indexer and focus on the sorter.

Indexer

The requirements on the indexer included flexibility, speed of indexing and retrieval, simplicity and usability. The 'Whoosh' python library provides all of these, so I used it to build an indexer. 'Whoosh' is an open source indexing library, so I had the option of modifying any part of it. It is built in pure Python, so it has a clean pythonic API. Its primary advantage is fast indexing and retrieval, although we are mostly concerned with retrieval speed, as indexing is done rarely. The predecessors of 'Whoosh' have served as the basis of well-known browsers

such as Lucene, so it is also a powerful indexing tool, should I have needed more sophistication.

I have defined a very simple schema for indexing. Perhaps, one notable detail is that 'Whoosh' can store timestamps with the index, which enabled me to provide both clean and incremental index methods. The incremental indexing relies on the timestamp stored with the index and compares it to the last-modified time provided by the file system. The user can specify whether indexing has to be done from scratch or updated to accommodate some document changes or document addition/deletion. I haven't originally expected to need an incremental indexing capability, but throughout the project it has permitted for a significant speedup.

Previously, I have defined the sorter as a logical unit that provides ranking to the retrieved documents. For the first prototype, the sorter returned the pages in the order of decreasing pagerank. Subsequently, the sorting has been decoupled from retrieval and is described in more detail in the Parser section.

PageRank

The PageRank vector is computed using matrix inversion. All matrix operations were performed with the help of the python numerical library 'numpy'. Take t to be the teleportation probability, s=1-t is the probability of following a random link, E is the teleportation probability: equiprobable transitions, as using non-personalized pagerank (see Preparation), E_i , j = 1/N for all i, j. G is a stochastic matrix holding the link structure of the data, such that

$$G_i, j = \begin{cases} 1/L & \text{there is a link from i to j and L = number of links from i} \\ 0 & \text{there is no link from i to j} \end{cases}$$

Then M is a stochastic matrix representing the web surfer activity, such that M_i , j is the probability of going from page i to page j,

$$M = s * G + t * E \tag{2.1}$$

In one step the new location of the surfer is described by the distribution Mp. We want to find a stationary distribution p, so must have

$$p = M * p \tag{2.2}$$

Substituting 2.1 into 2.2

$$p = (s * G + t * E) * p = s * G * p + t * E * p$$
(2.3)

Rearranging equation 2.3 gives

$$p * (I - s * G) = t * E * p \tag{2.4}$$

where I is the identity matrix

We can express E*p as P where P is a vector [1/N, 1/N, ..., 1/N].T, as members of p must sum to one. So computing pagerank amounts to

$$p = t * (I - s * G)^{-1} * P (2.5)$$

where $(I - s * G)^{-1}$ denotes a matrix inverse operation.

This solution is simple at the expense of being slow. Although computing inverse of a matrix is computationally expensive, we don't need to scale beyond a few thousands of pages. To avoid recomputation, I used python object serialization module - Pickle to store the pagerank vector for each directory. The resultant performance was actually very reasonable, the time spent computing pagerank was insignificantly small in comparison to the time spent crawling the directory.

TODO: Load class method TODO: byname array

Crawler

The matrix G, used for the pagerank computation, represents random link following activity. To obtain such a link structure each page has to be parsed, and all links recorded. Because our data is obtained from a single source page by a Wget spider, every page in a directory is guaranteed to be discovered by a spider.

The Crawler class recursively traverses the pages depth first starting with the seed page, the same as the seed page used for recursively downloading the pages from the web. To make sure each page is only explored once, a dictionary is used to hold pairs of absolute path, which uniquely identifies the page, and a numerical value corresponding to the timestamp when the page has been first discovered. **TODO:** link conversion and parsing Although every page has a unique path, the links to other pages are relative. Such links need to be normalized to maintain consistency. A page object is used to encapsulate path complexity: all link paths are converted to absolute paths before addition to the dictionary. All outbound links are stored with the page in a Set datastructure, such that no link is added more than once.

To produce the stochastic matrix G, we start with an empty NxN matrix, where N is the total number of pages. We assume that whenever a surfer encounters a dangling page – a page that has no outbound links – a teleportation step

Page	A
A	0
В	1/2
С	1/2
D	0

Table 2.1: Illustration of non-dangling pages: B and C share A's 'importance' equally.

occurs. Therefore, every dangling page links to every page in the pool including itself with equal probability 1/N. For non-dangling pages, all links are assumed equiprobable and all pages that are not linked to have probability of 0. So if page A links to pages B and C, but not itself or D, its row in G is described by the Table 2.1.

2.3 Machine Learning

In the course of this project two distinct machine learning algorithms have been used. The design goals for the implementations were as usual: speed and correctness. Another important aspect of the design is that the system and the machine learning modules must be sufficiently decoupled. This is an issue of scalability and code reuse. The implication on the machine learner implementations is that both must communicate with the system via the same interface. The rest of the section describes in detail how the machine learners have been implemented.

2.3.1 Naive Bayes

In Section 1.1 it has been shown that Naive Bayesian is a good learner to implement for the first prototype, so a very quick implementation was preferable to make sure the system can potentially function as intended. Due to its simplicity and popularity, Naive Bayesian is widely available in the libraries. Because the implementation of this module is straightforward, not central to the project, I decided to use one of many existing python implementations.

The *nltk* library implementation was particularly appealing as it offers a very concise interface. A classifier object is initialised by the **train** method on the **NaiveBayesClassifier** class. The format of the training set is defined as a list of tuples of featuresets and labels, e.g.

 $[(featureset_1, label_1), \dots, (featureset_N, label_N)].$ The train method simply computes the prior – the probability distribution over labels P(label) and

the likelihood – the conditional probability P(featureset = f|label) by simply counting and recording the relevant instances. The method outputs a NaiveBayesClassifier instance parametrized by the two probability distributions.

P(features) is not computed explicitly, instead, a normalizing denominator ?? is used.

$$\sum_{l \in labels} (P(l)P(featureset_1|l) \cdots P(featureset_N|l))$$
 (2.6)

The classify method on the NaiveBayesClassifier object takes exactly one featureset and returns a label which maximizes the posterior probability P(label—featureset) over all labels. Previously unseen features are ignored by the classifier, so that to avoid assigning zero probability to everything.

The only tangible difficulty with the implementation was the framework in which classification is done. The Parser was 'written around' the interface of the classifier. To batch classify everything in the test directory, the classifier object is passed to the constructor of the TestFeatureSetCollection class. Here the featuresets for the test directories are computed, classified and recorded for the evaluation stage later.

2.3.2 Support Vector Machine

In the Preparaion chapter we have looked at the theory of support vector machines as found in textbooks. This section continues from the theory and explains further transformations that were an essential part of the implementation as opposed to bookwork. **TODO:** rephrase

To implement a support vector machine one must solve a problem of optimizing a quadratic function ssubject to tlinear constraints – usually referred to as the *Quadratic Programming* (QP) problem. Therefore, the first implementation task was to convert our existing optimization problem into a generic QP form to make use of the available solvers.

The maximization problem 1.9 can be trivially expressed as a minimization problem (equation 2.7).

$$min_{\boldsymbol{\alpha},\boldsymbol{\alpha}^*} \frac{1}{2} (\boldsymbol{\alpha} - \boldsymbol{\alpha}^*)^T P(\boldsymbol{\alpha} - \boldsymbol{\alpha}^*) + \epsilon \sum_{i=1}^l (\alpha_i + \alpha_i^*) - \sum_{i=1}^l t_i (\alpha_i - \alpha_i^*)$$
 (2.7)

subject to constraints 2.8 and 2.9 below.

$$\mathbf{e}(\boldsymbol{\alpha} - \boldsymbol{\alpha}^*) = 0 \tag{2.8}$$

$$0 \le \alpha_i, \alpha_i^* \le C, \ i = 1, ..., l$$
 (2.9)

where $\mathbf{e} = [1, ..., 1]$, $P_{ij} = K(x_i, x_j)$, t_i is the target output, C > 0 and $\epsilon > 0$.

At this point in the implementation, for the first time, python did not seem like an ideal choice. *Cvxopt* is one of the few python libraries that implements a QP solver. The specification to the QP function is as follows: cvxopt.solvers.qp(P,q,G,h,A,b) solves a pair of primal and dual convex quadratic programs

$$\min \frac{1}{2}x^T P x + q^T x \tag{2.10}$$

subject to

$$Gx \le h \tag{2.11}$$

$$Ax = b (2.12)$$

Described in the next few pages are the transformations I devised to reconcile the minimization problem 2.7 and the library specification 2.10 and their respective constraints.

We take x to encode both α and α^* simultaneously, treating the upper half of x as α and the lower half as α^* :

$$x = \begin{bmatrix} \alpha \\ \alpha^* \end{bmatrix}$$

We will see later how this representation allows for elegant representation of the problem (2.7).

First, we express the first term in 2.10 to hold $(\alpha - \alpha^*)^T P(\alpha - \alpha^*)$. Take matrix P in equation 2.10 as

$$P = \begin{bmatrix} K & -K \\ -K & K \end{bmatrix}$$

where $K_{ij} = K(x_i, x_j)$ is the kernel.

Observe that now

$$x^T P x = \begin{bmatrix} \boldsymbol{\alpha} \\ \boldsymbol{\alpha}^* \end{bmatrix} \begin{bmatrix} K & -K \\ -K & K \end{bmatrix} \begin{bmatrix} \boldsymbol{\alpha} \\ \boldsymbol{\alpha}^* \end{bmatrix}$$

is equivalent to the first term of equation (1.9)

$$\sum_{n=1}^{N} \sum_{m=1}^{N} (a_n - a_n^*)(a_m - a_m^*) K(x_n, x_m)$$

Now expressing the remaining two terms of 2.7 by taking

$$q = \begin{bmatrix} \epsilon * \mathbf{e} - t_0 \\ \vdots \\ \epsilon * \mathbf{e} - t_{N-1} \\ \epsilon * \mathbf{e} + t_0 \\ \vdots \\ \epsilon * \mathbf{e} + t_{N-1} \end{bmatrix}$$

to achieve

$$q^T x = \epsilon \sum_{i=1}^{l} (\alpha_i + \alpha_i^*) - \sum_{i=1}^{l} t_i (\alpha_i - \alpha_i^*)$$

To encode constraint 2.9 consider the following pair of G and h.

The final constraint 2.8 is trivial to adapt by simply taking

$$A = [\overbrace{1, \dots, 1}^{N}, \overbrace{-1, \dots, -1}^{N}]$$

and

$$b = 0$$

Having worked out all the matrices, the rest of the implementation dealt simply with coding the matrices up. The Cvxopt library has its own matrix constructor, however, has generally limited functionality when it comes to matrix operations, so Numpy was used a lot for matrix manipulations. During the first implementation attempt, the solver gave mostly unintelligible error messages, so for the prototype SVM I actually used Matlab. Matlab's quadprog function had an identical specification to the Python solver I intended to use. Matrices in Matlab are a lot more straightforward to manipulate, which was also a good reason to first check the correctness of my matrices in Matlab.

Kernel functions

The first kernel function I used was linear

2.4 Parser

So far we have looked at the search engine. As a functional unit the search engine only retrieves the relevant pages. The Parser is an abstraction for a few classes which together perform a series of tasks involving page parsing. Search engine heuristics are also handled by the parsing module.

The primary function of the parser is to compute feature vectors for pages. At this point we abstract away from the web pages and html parsing: as far as the machine learner is concerned, a feature vector is a complete representation of a page. The infrastructure for feature vector computation is easily extensible to accommodate other related tasks, namely, training set and test set generation. To allow for multifunctionality, I have taken an object oriented approach to the design of the module.

The module operates in two modes: rank and score and handles both classification and regression. The high level specification is that we create two objects: a TestFeatureSetCollection and a LabelFeatureSetCollection objects and they encapsulate all the data, so can be passed around to the machine learners, as well as evaluation and plotting modules. These objects each operate within their own directories, to keep training and testing sufficiently separate.

Both category and rank are treated as page features and hence are part of the LabeledFeatureSet class state. The LabeledFeatureSetCollection class generates

2.4. PARSER 21

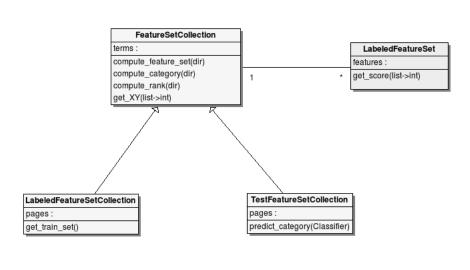


Figure 2.1: A UML class diagram describing operation of the Parser.

training sets from queries, whereas the TestFeatureSetCollection class computes predicted category given an instance of a classifier. Both, however, need to generate feature sets, one for the Training data and the other for the Test data. This common functionality is embodied in their abstract base class, FeatureSet-Collection. Figure 2.1 shows a UML class diagram illustrating the main structure of the module.

The rest of this section talks about the specifics of the implementation, in particular, the features used and how html parsing is done.

Rank mode. Conceptually a FeatureSetCollection is a dictionary of Labeled-FeatureSet objects indexed by both path to page and query term. After initialisation, all the features are computed and 'filled up' per query term per page.

Score mode.

Features. The requirements analysis does not prescribe the use of any particular features. However, it states that both dynamic (query dependent) and static(query independent) features must be used. PageRank is a dynamic feature that has quite special place among others: it takes into account all the pages in the pool and reflects on the structural hierarchy of the web pages. The parser

loads the pre-computed PageRank vector indexed by page name to extract the pagerank of any particular page.

An example of a dynamic feature used in the project is query term count: the number of times the query term occurs on the page. This is obtained directly from the html files using the BeautifulSoup library, just like for the link parsing in the crawler. The html is parsed into clean text and the count is computed on the string. A related but different feature – stem count – is the number of times the stem of the word occurs in the text. This is obtained using the PorterStemmer module in Natural Language Tool Kit (nltk) library.

Boolean features are a slightly different variety of feature, so was worth putting in. An example of this is a presence of an image on the page.

The features have been added incrementally as the project progressed. All other features are similar to the ones described above and are obtained using the same methods. **TODO:** enumerate features here

2.5 Optimization

One characteristic peripheral requirement for the system implemented in this project is speed. Even though it is not our direct goal to produce efficient implementations, optimization could not be overlooked, because significant amount of time is spent processing large quantities of data: indexing, pagerank and feature set computations all must complete in 'feasible' time, i.e. in the final implementation the longest computation takes order of minutes and processes a few thousand pages. Various optimizations have been used to achieve this.

Both pagerank vector and the index are precomputed and kept in persistent storage. Incremental indexing feature allows us to edit parts of the index as opposed to recomputing the whole index from scratch. These precomputations provide certain speedups, but were not enough.

Because the system is fairly complex and a lot of library code is used in places, it was hard to determine which code most affects the speed. 'Blind' attempts at optimization did not work well, which motivated the use of a profiling tool.

Profiling the first prototype of the complete system revealed a surprising fact: most time was spent in the library code parsing pages. To mitigate this issue I have tried using custom parsers instead. Apart from speed, robustness was another important consideration, as a failure of a parser increases compute time. Two of the most renowned python parsers are html5lib and lxml. Figure 2.2 below shows visual representation of time profiling of 3 different runs obtained using the

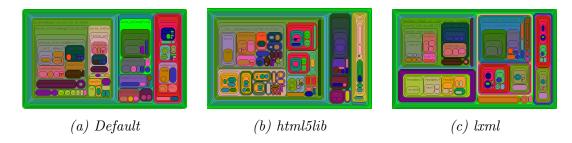


Figure 2.2: Visual profiles of three parsers from left to right: default python html parser, html5lib and lxml. The internal boxes are sized in proportion to the time spent in each function. In all profiles three distinct major compartments can be seen, the leftmost being the compartment of interest: time spent in parsing. Taking into consideration that the other modules are unaffected by changing the parser implementation, it can be observed that lxml (rightmost) is the fastest and html5lib is the slowest.

RunSnakeRun, each exploiting a different parser. Despite html5lib being quoted as the most robust/lenient, lxml was sufficiently faster to be preferable.

Another useful limitation was discovered due to profiling. The pagerank vector was loaded into memory every time a page was parsed: once for each page. This clearly is undesirable. I used caching to ensure that each of the two possible pagerank vectors (corresponding to the test and train directories) is loaded from memory exactly once. This is achieved via a double singleton class, which loads the vectors lazily.

Chapter 3

Evaluation

- 3.1 The prototype
- 3.2 Bayes ranking: one to one comparison
- 3.3 SVR
- 3.4 Kernel function generality
- 3.5 Experiments with number of features
- 3.6 Score vs Rank

Chapter 4

Conclusion

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Appendix A
 Project Proposal