E6893 Big Data Analytics Lecture 5:

*Big Data Analytics Algorithms -- II*

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Review — Key Components of Mahout

Collaborative Filtering
- User-Based Collaborative Filtering - single machine
- Item-Based Collaborative Filtering - single machine / MapReduce
- Matrix Factorization with Alternating Least Squares - single machine / MapReduce
- Matrix Factorization with Alternating Least Squares on Implicit Feedback - single machine / MapReduce
- Weighted Matrix Factorization, SVD++, Parallel SGD - single machine

Classification
- Logistic Regression - trained via SGD - single machine
- Naive Bayes/ Complementary Naive Bayes - MapReduce
- Random Forest - MapReduce
- Hidden Markov Models - single machine
- MultiLayer Perceptron - single machine

Clustering
- Canopy Clustering - single machine / MapReduce (deprecated, will be removed once Streaming k-Means is stable enough)
- k-Means Clustering - single machine / MapReduce
- Fuzzy k-Means - single machine / MapReduce
- Streaming k-Means - single machine / MapReduce
- Spectral Clustering - MapReduce
Clustering a collection involves three things:

- *An algorithm*—This is the method used to group the books together.
- *A notion of both similarity and dissimilarity*—In the previous discussion, we relied on your assessment of which books belonged in an existing stack and which should start a new one.
- *A stopping condition*—In the library example, this might be the point beyond which books can’t be stacked anymore, or when the stacks are already quite dissimilar.
Clustering — on feature plane
Clustering example

\[(1, 1)\]
\[(2, 1)\]
\[(1, 2)\]
\[(2, 2)\]
\[(3, 3)\]
\[(8, 8)\]
\[(8, 9)\]
\[(9, 8)\]
\[(9, 9)\]

(0, 0)

\[x\text{-axis}\]

\[y\text{-axis}\]
Steps on clustering

1. Generate Vectors from input data
2. Write Vectors to input directory
3. Write initial cluster centers
4. Run clustering job
5. Read clusters from output directory
K-mean clustering

K-means clustering in action. Starting with three random points as centroids (top left), the map stage (top right) assigns each point to the cluster nearest to it. In the reduce stage (bottom left), the associated points are averaged out to produce the new location of the centroid, leaving you with the final configuration (bottom right). After each iteration, the final configuration is fed back into the same loop until the centroids come to rest at their final positions.
Making initial cluster centers
Parameters to Mahout k-mean clustering algorithm

- The `SequenceFile` containing the input vectors.
- The `SequenceFile` containing the initial cluster centers. In this case, we seed two clusters, so there are two centers.
- The similarity measure to be used. We use `EuclideanDistanceMeasure` as the measure of similarity here, and we explore other similarity measures later in this chapter.
- The convergence threshold. If in a particular iteration the centers of the clusters don’t change beyond this threshold, no further iterations are done.
- The number of iterations to be done.
- The vector implementation used in the input files.
HelloWorld clustering scenario

```java
public static final double[][] points = {
    {1, 1}, {2, 1}, {1, 2},
    {2, 2}, {3, 3}, {8, 8},
    {9, 8}, {8, 9}, {9, 9}};

public static void writePointsToFile(List<Vector> points,
    String fileName,
    FileSystem fs,
    Configuration conf) throws IOException {
    Path path = new Path(fileName);
    SequenceFile.Writer writer = new SequenceFile.Writer(fs, conf,
        path, LongWritable.class, VectorWritable.class);
    long recNum = 0;
    VectorWritable vec = new VectorWritable();
    for (Vector point : points) {
        vec.set(point);
        writer.append(new LongWritable(recNum++), vec);
    }
    writer.close();
}

public static List<Vector> getPoints(double[][] raw) {
    List<Vector> points = new ArrayList<Vector>();
    for (int i = 0; i < raw.length; i++) {
        double[] fr = raw[i];
        Vector vec = new RandomAccessSparseVector(fr.length);
        vec.assign(fr);
        points.add(vec);
    }
    return points;
}
```
public static void main(String args[]) throws Exception {
    int k = 2;
    List<Vector> vectors = getPoints(points);
    File testData = new File("testdata");
    if (!testData.exists()) {
        testData.mkdirs();
    }
    testData = new File("testdata/points");
    if (!testData.exists()) {
        testData.mkdirs();
    }
    Configuration conf = new Configuration();
    FileSystem fs = FileSystem.get(conf);
    writePointsToFile(vectors,
                      "testdata/points/file1", fs, conf);
    Path path = new Path("testdata/clusters/part-00000");
    SequenceFile.Writer writer
        = new SequenceFile.Writer(
                                fs, conf, path, Text.class, Cluster.class);
    for (int i = 0; i < k; i++) {
        Vector vec = vectors.get(i);
        Cluster cluster = new Cluster(
            vec, i, new EuclideanDistanceMeasure());
        writer.append(new Text(cluster.getIdentifier()), cluster);
    }
    writer.close();
KMeansDriver.run(conf, new Path("testdata/points"),
    new Path("testdata/clusters"),
    new Path("output"), new EuclideanDistanceMeasure(),
    0.001, 10, true, false);

SequenceFile.Reader reader
    = new SequenceFile.Reader(fs,
        new Path("output/" + Cluster.CLUSTERED_POINTS_DIR
            + "/part-m-00000"), conf);

IntWritable key = new IntWritable();
WeightedVectorWritable value = new WeightedVectorWritable();
while (reader.next(key, value)) {
    System.out.println(
        value.toString() + " belongs to cluster "
        + key.toString());
}
reader.close();
1.0: [1.000, 1.000] belongs to cluster 0
1.0: [2.000, 1.000] belongs to cluster 0
1.0: [1.000, 2.000] belongs to cluster 0
1.0: [2.000, 2.000] belongs to cluster 0
1.0: [3.000, 3.000] belongs to cluster 0
1.0: [8.000, 8.000] belongs to cluster 1
1.0: [9.000, 8.000] belongs to cluster 1
1.0: [8.000, 9.000] belongs to cluster 1
1.0: [9.000, 9.000] belongs to cluster 1
Testing difference distance measures

Euclidean distance measure

\[ d = \sqrt{(a_1 - b_1)^2 + (a_2 - b_2)^2 + \ldots + (a_n - b_n)^2} \]

Squared Euclidean distance measure

\[ d = (a_1 - b_1)^2 + (a_2 - b_2)^2 + \ldots + (a_n - b_n)^2 \]

Manhattan distance measure

\[ d = |a_1 - b_1| + |a_2 - b_2| + \ldots + |a_n - b_n| \]
**Manhattan distance measure**

\[ d = |a_1 - b_1| + |a_2 - b_2| + \ldots + |a_n - b_n| \]

**Cosine distance measure**

\[ d = 1 - \frac{(a_1 b_1 + a_2 b_2 + \ldots + a_n b_n)}{\left(\sqrt{a_1^2 + a_2^2 + \ldots + a_n^2}\right)\sqrt{b_1^2 + b_2^2 + \ldots + b_n^2}} \]
Tanimoto distance and weighted distance

**Tanimoto distance measure**

\[
d = 1 - \frac{a_1 b_1 + a_2 b_2 + \ldots + a_n b_n}{\sqrt{a_1^2 + a_2^2 + \ldots + a_n^2} + \sqrt{b_1^2 + b_2^2 + \ldots + b_n^2} - (a_1 b_1 + a_2 b_2 + \ldots + a_n b_n)}
\]

**Weighted distance measure**

Mahout also provides a `WeightedDistanceMeasure` class, and implementations of Euclidean and Manhattan distance measures that use it. A weighted distance measure is an advanced feature in Mahout that allows you to give weights to different dimensions in order to either increase or decrease the effect of a dimension.
## Results comparison

<table>
<thead>
<tr>
<th>Distance measure</th>
<th>Number of iterations</th>
<th>Vectors in cluster 0</th>
<th>Vectors in cluster 1</th>
</tr>
</thead>
<tbody>
<tr>
<td>EuclideanDistanceMeasure</td>
<td>3</td>
<td>0, 1, 2, 3, 4</td>
<td>5, 6, 7, 8</td>
</tr>
<tr>
<td>SquaredEuclideanDistanceMeasure</td>
<td>5</td>
<td>0, 1, 2, 3, 4</td>
<td>5, 6, 7, 8</td>
</tr>
<tr>
<td>ManhattanDistanceMeasure</td>
<td>3</td>
<td>0, 1, 2, 3, 4</td>
<td>5, 6, 7, 8</td>
</tr>
<tr>
<td>CosineDistanceMeasure</td>
<td>1</td>
<td>1</td>
<td>0, 2, 3, 4, 5, 6, 7, 8</td>
</tr>
<tr>
<td>TanimotoDistanceMeasure</td>
<td>3</td>
<td>0, 1, 2, 3, 4</td>
<td>5, 6, 7, 8</td>
</tr>
</tbody>
</table>

![Diagram](image)
Data preparation in Mahout — vectors

In Mahout, vectors are implemented as three different classes, each of which is optimized for different scenarios: DenseVector, RandomAccessSparseVector, and SequentialAccessSparseVector.

- **DenseVector** can be thought of as an array of doubles, whose size is the number of features in the data. Because all the entries in the array are preallocated regardless of whether the value is 0 or not, we call it *dense*.

- **RandomAccessSparseVector** is implemented as a HashMap between an integer and a double, where only nonzero valued features are allocated. Hence, they’re called as *SparseVectors*.

- **SequentialAccessSparseVector** is implemented as two parallel arrays, one of integers and the other of doubles. Only nonzero valued entries are kept in it. Unlike the **RandomAccessSparseVector**, which is optimized for random access, this one is optimized for linear reading.
vectorization example

0: weight
1: color
2: size

[0 => 100 gram, 1 => red, 2 => small]

<table>
<thead>
<tr>
<th>Apple</th>
<th>Weight (kg) (0)</th>
<th>Color (1)</th>
<th>Size (2)</th>
<th>Vector</th>
</tr>
</thead>
<tbody>
<tr>
<td>Small, round, green</td>
<td>0.11</td>
<td>510</td>
<td>1</td>
<td>[0.11, 510, 1]</td>
</tr>
<tr>
<td>Large, oval, red</td>
<td>0.23</td>
<td>650</td>
<td>3</td>
<td>[0.23, 650, 3]</td>
</tr>
<tr>
<td>Small, elongated, red</td>
<td>0.09</td>
<td>630</td>
<td>1</td>
<td>[0.09, 630, 1]</td>
</tr>
<tr>
<td>Large, round, yellow</td>
<td>0.25</td>
<td>590</td>
<td>3</td>
<td>[0.25, 590, 3]</td>
</tr>
<tr>
<td>Medium, oval, green</td>
<td>0.18</td>
<td>520</td>
<td>2</td>
<td>[0.18, 520, 2]</td>
</tr>
</tbody>
</table>
public static void main(String args[]) throws Exception {
    List<NamedVector> apples = new ArrayList<NamedVector>();

    NamedVector apple;
    apple = new NamedVector(
        new DenseVector(new double[] {0.11, 510, 1}),
        "Small round green apple");
    apples.add(apple);
    apple = new NamedVector(
        new DenseVector(new double[] {0.23, 650, 3}),
        "Large oval red apple");
    apples.add(apple);
    apple = new NamedVector(
        new DenseVector(new double[] {0.09, 630, 1}),
        "Small elongated red apple");
    apples.add(apple);
    apple = new NamedVector(
        new DenseVector(new double[] {0.25, 590, 3}),
        "Large round yellow apple");
    apples.add(apple);
    apple = new NamedVector(
        new DenseVector(new double[] {0.18, 520, 2}),
        "Medium oval green apple");
Configuration conf = new Configuration();
FileSystem fs = FileSystem.get(conf);

Path path = new Path("appledata/apples");
SequenceFile.Writer writer = new SequenceFile.Writer(fs, conf,  
        path, Text.class, VectorWritable.class);
VectorWritable vec = new VectorWritable();
for (NamedVector vector : apples) {  
    vec.set(vector);
    writer.append(new Text(vector.getName()), vec);  
}  
writer.close();

SequenceFile.Reader reader = new SequenceFile.Reader(fs,  
        new Path("appledata/apples"), conf);

Text key = new Text();
VectorWritable value = new VectorWritable();
while (reader.next(key, value)) {
    System.out.println(key.toString() + " "  
        + value.get().asFormatString());
}  
reader.close();
Vectorization of text

Vector Space Model: Term Frequency (TF)

For example, if the word *horse* is assigned to the 39,905\textsuperscript{th} index of the vector, the word *horse* will correspond to the 39,905\textsuperscript{th} dimension of document vectors. A document’s vectorized form merely consists, then, of the number of times each word occurs in the document, and that value is stored in the vector along that word’s dimension. The dimension of these document vectors can be very large.

Stop Words: \textit{a, an, the, who, what, are, is, was, and so on.}

Stemming:

A stemmer for English, for example, should identify the \textit{string} "cats" (and possibly "catlike", "catty" etc.) as based on the root "cat", and "stemmer", "stemming", "stemmed" as based on "stem". A stemming algorithm reduces the words "fishing", "fished", and "fisher" to the root word, "fish". On the other hand, "argue", "argued", "argues", "arguing", and "argus" reduce to the stem "argu" (illustrating the case where the stem is not itself a word or root) but "argument" and "arguments" reduce to the stem "argument".
Most Popular Stemming algorithms

Lookup algorithms

A simple stemmer looks up the inflected form in a lookup table. The advantages of this approach is that it is simple, fast, and easily handles exceptions. The disadvantages are that all inflected forms must be explicitly listed in the table: new or unfamiliar words are not handled, even if they are perfectly regular (e.g. iPads ~ iPad), and the table may be large. For languages with simple morphology, like English, table sizes are modest, but

Suffix-stripping algorithms

Suffix stripping algorithms do not rely on a lookup table that consists of inflected forms and root form relations. Instead, a typically smaller list of "rules" is stored which provides a path for the algorithm, given an input word form, to find its root form. Some examples of the rules include:

- if the word ends in 'ed', remove the 'ed'
- if the word ends in 'ing', remove the 'ing'
- if the word ends in 'ly', remove the 'ly'
Term Frequency — Inverse Document Frequency (TF-IDF)

The value of word is reduced more if it is used frequently across all the documents in the dataset.

To calculate the inverse document frequency, the document frequency (DF) for each word is first calculated. Document frequency is the number of documents the word occurs in. The number of times a word occurs in a document isn’t counted in document frequency. Then, the inverse document frequency or IDF$_i$ for a word, $w_i$, is

$$IDF_i = \frac{1}{DF_i}$$

$$W_i = TF_i \cdot IDF_i = TF_i \cdot \frac{N}{DF_i}$$

or

$$W_i = TF_i \cdot \log \frac{N}{DF_i}$$
It was the best of time. it was the worst of times.  

\[ \Rightarrow \]

bigram

It was  
was the  
the best  
best of  
of times  
times it  
it was  
was the  
the worst  
worst of  
of times

Mahout provides a log-likelihood test to reduce the dimensions of n-grams
Examples — using a news corpus

Reuters-21578 dataset: 22 files, each one has 1000 documents except the last one.

http://www.daviddlewis.com/resources/testcollections/reuters21578/

Extraction code:

```java
mvn -e -q exec:java
-Dexec.mainClass="org.apache.lucene.benchmark.utils.ExtractReuters"
-Dexec.args="reuters/ reuters-extracted/"
```

Using the extracted folder, run the `SequenceFileFromDirectory` class. You can use the launcher script from the Mahout root directory to do the same:

```bash
bin/mahout seqdirectory -c UTF-8
-i examples/reuters-extracted/ -o reuters-seqfiles
```

This will write the Reuters articles in the `SequenceFile` format. Now the only step left is to convert this data to vectors. To do that, run the `SparseVectorsFromSequenceFiles` class using the Mahout launcher script:

```bash
bin/mahout seq2sparse -i reuters-seqfiles/ -o reuters-vectors -ow
```
<table>
<thead>
<tr>
<th>Option</th>
<th>Flag</th>
<th>Description</th>
<th>Default value</th>
</tr>
</thead>
<tbody>
<tr>
<td>Overwrite</td>
<td>-ow</td>
<td>If set, the output folder is overwritten. If not set, the output folder is created if the folder doesn’t exist. If the output folder does exist, the job fails and an error is thrown. Default is unset.</td>
<td>N/A</td>
</tr>
<tr>
<td>Lucene analyzer name</td>
<td>-a</td>
<td>The class name of the analyzer to use.</td>
<td><code>org.apache.lucene.analyses.standard.StandardAnalyzer</code></td>
</tr>
<tr>
<td>Chunk size</td>
<td>-chunk</td>
<td>The chunk size in MB. For large document collections (sizes in GBs and TBs), you won’t be able to load the entire dictionary into memory during vectorization, so you can split the dictionary into chunks of the specified size and perform the vectorization in multiple stages. It’s recommended you keep this size to 80 percent of the Java heap size of the Hadoop child nodes to prevent the vectorizer from hitting the heap limit.</td>
<td>100</td>
</tr>
<tr>
<td>Weighting</td>
<td>-wt</td>
<td>The weighting scheme to use: tf for term-frequency based weighting and tfidf for TF-IDF based weighting.</td>
<td>tfidf</td>
</tr>
<tr>
<td>Minimum support</td>
<td>-s</td>
<td>The minimum frequency of the term in the entire collection to be considered as a part of the dictionary file. Terms with lesser frequency are ignored.</td>
<td>2</td>
</tr>
</tbody>
</table>
## Mahout dictionary-based vectorizer — II

<table>
<thead>
<tr>
<th>Option</th>
<th>Flag</th>
<th>Description</th>
<th>Default value</th>
</tr>
</thead>
<tbody>
<tr>
<td>Minimum document frequency (int)</td>
<td>-md</td>
<td>The minimum number of documents the term should occur in to be considered a part of the dictionary file. Any term with lesser frequency is ignored.</td>
<td>1</td>
</tr>
<tr>
<td>Max document frequency percentage (int)</td>
<td>-x</td>
<td>The maximum number of documents the term should occur in to be considered a part of the dictionary file. This is a mechanism to prune out high frequency terms (stop-words). Any word that occurs in more than the specified percentage of documents is ignored.</td>
<td>99</td>
</tr>
<tr>
<td>N-gram size (int)</td>
<td>-ng</td>
<td>The maximum size of n-grams to be selected from the collection of documents.</td>
<td>1</td>
</tr>
<tr>
<td>Option</td>
<td>Flag</td>
<td>Description</td>
<td>Default value</td>
</tr>
<tr>
<td>---------------------------------------------</td>
<td>------</td>
<td>-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------</td>
<td>---------------</td>
</tr>
<tr>
<td>Minimum log-likelihood ratio (LLR) (float)</td>
<td>-ml</td>
<td>This flag works only when $n$-gram size is greater than 1. Very significant $n$-grams have large scores, such as 1000; less significant ones have lower scores. Although there's no specific method for choosing this value, the rule of thumb is that $n$-grams with a LLR value less than 1.0 are irrelevant.</td>
<td>1.0</td>
</tr>
<tr>
<td>Normalization (float)</td>
<td>-n</td>
<td>The normalization value to use in the $L_p$ space. A detailed explanation of normalization is given in section 8.4. The default scheme is to not normalize the weights.</td>
<td>0</td>
</tr>
<tr>
<td>Number of reducers (int)</td>
<td>-nr</td>
<td>The number of reducer tasks to execute in parallel. This flag is useful when running a dictionary vectorizer on a Hadoop cluster. Setting this to the maximum number of nodes in the cluster gives maximum performance. Setting this value higher than the number of cluster nodes leads to a slight decrease in performance. For more details, read the Hadoop documentation on setting the optimum number of reducers.</td>
<td>1</td>
</tr>
<tr>
<td>Create sequential access sparse vectors (bool)</td>
<td>-seq</td>
<td>If set, the output vectors are created as <code>SequentialAccessSparseVectors</code>. By default the dictionary vectorizer generates <code>RandomAccessSparseVectors</code>. The former gives higher performance on certain algorithms like k-means and SVD due to the sequential nature of vector operations. By default the flag is unset.</td>
<td>N/A</td>
</tr>
</tbody>
</table>
Outputs & Steps

1. Tokenization using Lucene StandardAnalyzer
2. n-gram generation step
3. converts the tokenized documents into vectors using TF
4. count DF and then create TF-IDF
A practical setting of flags

- `-a`—Use `org.apache.lucene.analysis.WhitespaceAnalyzer` to tokenize words based on the whitespace characters between them.
- `-chunk`—Use a chunk size of 200 MB. This value won't produce any effect on the Reuters data, because the dictionary sizes are usually in the 1 MB range.
- `-wt`—Use the `tfidf` weighting method.
- `-s`—Use a minimum support value of 5.
- `-md`—Use a minimum document frequency value of 3.
- `-x`—Use a maximum document frequency percentage of 90 percent to aggressively prune away high-frequency words.
- `-ng`—Use an `n`-gram size of 2 to generate both unigrams and bigrams.
- `-ml`—Use a minimum log-likelihood ratio (LLR) value of 50 to keep only very significant bigrams.
- `-seq`—Set the `SequentialAccessSparseVectors` flag.

Run the vectorizer using the preceding options in the Mahout launcher script:

```bash
bin/mahout seq2sparse -i reuters-seqfiles/ -o reuters-vectors-bigram -ow
-a org.apache.lucene.analysis.WhitespaceAnalyzer
-chunk 200 -wt tfidf -s 5 -md 3 -x 90 -ng 2 -ml 50 -seq
```
Some documents may pop up showing they are similar to all the other documents because it is large. ==> Normalization can help.

In Mahout, normalization uses what is known in statistics as a $p$-norm. For example, the $p$-norm of a 3-dimensional vector, $[x, y, z]$, is

$$
\frac{x}{(|x|^p + |y|^p + |z|^p)^{1/p}}, \quad \frac{y}{(|x|^p + |y|^p + |z|^p)^{1/p}}, \quad \frac{z}{(|x|^p + |y|^p + |z|^p)^{1/p}}
$$
Clustering methods provided by Mahout

- K-means clustering
- Centroid generation using canopy clustering
- Fuzzy k-means clustering and Dirichlet clustering
- Topic modeling using latent Dirichlet allocation as a variant of clustering
K-means clustering in action. Starting with three random points as centroids (top left), the map stage (top right) assigns each point to the cluster nearest to it. In the reduce stage (bottom left), the associated points are averaged out to produce the new location of the centroid, leaving you with the final configuration (bottom right). After each iteration, the final configuration is fed back into the same loop until the centroids come to rest at their final positions.
Hadoop k-mean clustering jobs

In Mahout, the MapReduce version of the k-means algorithm is instantiated using the KMeansDriver class. The class has just a single entry point—the runJob method.

- The Hadoop configuration.
- The SequenceFile containing the input Vectors.
- The SequenceFile containing the initial Cluster centers.
- The similarity measure to be used. We’ll use EuclideanDistanceMeasure as the measure of similarity and experiment with the others later.
- The convergenceThreshold. If in an iteration, the centroids don’t move more than this distance, no further iterations are done and clustering stops.
- The number of iterations to be done. This is a hard limit; the clustering stops if this threshold is reached.
K-mean clustering running as MapReduce job

SequenceFile containing Vectors

chunk-0

chunk-0

...

chunk-n

Each Mapper reads in the centroid vector at startup

Parallel Mappers

Each Reducer gets the partial sums of all points of a cluster from each Mapper and recomputes the centroids

Parallel Reducers

Each Mapper computes nearest centroid for a vector

Repeat until converged

K-Means in MapReduce
Hadoop k-mean clustering code

KmeansDriver.runJob(hadoopConf,
    inputVectorFilesDirPath, clusterCenterFilesDirPath,
    outputDir, new EuclideanDistanceMeasure(),
    convergenceThreshold, numIterations, true, false);

Mahout reads and writes data using the Hadoop FileSystem class. This provides seamless access to both the local filesystem (via java.io) and distributed filesystems like HDFS and S3FS (using internal Hadoop classes). This way, the same code that works on the local system will also work on the Hadoop filesystem on the cluster, provided the paths to the Hadoop configuration files are correctly set in the environment variables. In Mahout, the bin/mahout shell script finds the Hadoop configuration files automatically from the $HADOOP_CONF environment variable.

$ bin/mahout kmeans -i reuters-vectors/tfidf-vectors/ \ 
-c reuters-initial-clusters \ 
-o reuters-kmeans-clusters \ 
-dm org.apache.mahout.common.distance.SquaredEuclideanDistanceMeasure \ 
-cd 1.0 -k 20 -x 20 -c1
The directory listing of the output folder looks something like this:

```
$ ls -l reuters-kmeans-clusters
drwxr-xr-x 4 user 5000 136 Feb 1 18:56 clusters-0
drwxr-xr-x 4 user 5000 136 Feb 1 18:56 clusters-1
drwxr-xr-x 4 user 5000 136 Feb 1 18:56 clusters-2
...
```

```
drwxr-xr-x 4 user 5000 136 Feb 1 18:59 clusteredPoints
```

```
$ bin/mahout clusteredump -dt sequencefile \
   -d reuters-vectors/dictionary.file-* \
   -s reuters-kmeans-clusters/clusters-19 -b 10 -n 10
```

Running ClusterDumper on the output folder corresponding to the last iteration produces output similar to the following:

**Id: 11736:**
Top Terms: debt, banks, brazil, bank, billion, he, payments, billion
dlrs, interest, foreign

**Id: 11235:**
Top Terms: amorphous, magnetic, metals, allied signal, 19.39, corrosion,
allied, molecular, mode, electronic components

... 

**Id: 20073:**
Top Terms: ibm, computers, computer, att, personal, pc, operating system,
in tel, machines, dos
Canopy clustering to estimate the number of clusters

Tell what size clusters to look for. The algorithm will find the number of clusters that have approximately that size. The algorithm uses two distance thresholds. This method prevents all points close to an already existing canopy from being the center of a new canopy.

Canopy clustering: if you start with a point (top left) and mark it as part of a canopy, all the points within distance $T_2$ (top right) are removed from the data set and prevented from becoming new canopies. The points within the outer circle (bottom-right) are also put in the same canopy, but they're allowed to be part of other canopies. This assignment process is done in a single pass on a mapper. The reducer computes the average of the centroid (bottom right) and merges close canopies.
Running canopy clustering

To run canopy generation over the Reuters data set, execute the canopy program using the Mahout launcher as follows:

$ bin/mahout canopy -i reuters-vectors/tfidf-vectors \
-o reuters-canopy-centroids \
-dm org.apache.mahout.common.distance.EuclideanDistanceMeasure \
-t1 1500 -t2 2000

Within a minute, CanopyDriver will generate the centroids in the output folder. Created less than 50 centroids.

$ bin/mahout kmeans -i reuters-vectors/tfidf-vectors \
-o reuters-kmeans-clusters \
-dm org.apache.mahout.common.distance.TanimotoDistanceMeasure \
-c reuters-canopy-centroids/clusters-0 -cd 0.1 -ow -x 20 -cl

After the clustering is done, use ClusterDumper to inspect the clusters. Some of them are listed here:

Id: 21523:name:
   Top Terms:
   tones, wheat, grain, said, usda, corn, us, sugar, export, agriculture
Id: 21409:name:
   Top Terms:
   stock, share, shares, shareholders, dividend, said, its, common, board,
   company
Id: 21155:name:
   Top Terms:
   oil, effective, crude, raises, prices, barrel, price, cts, said, dlr
Id: 19658:name:
   Top Terms:
   drug, said, aids, inc, company, its, patent, test, products, food
Id: 21323:name:
   Top Terms:
   7-apr-1987, 11, 10, 12, 07, 09, 15, 16, 02, 17
News clustering code

```java
public class NewsKMeansClustering {
    public static void main(String args[]) throws Exception {
        int minSupport = 2;
        int minDf = 5;
        int maxDFPercent = 95;
        int maxNgramSize = 2;
        int minLLRValue = 50;
        int reduceTasks = 1;
        int chunkSize = 200;
        int norm = 2;
        boolean sequentialAccessOutput = true;

        String inputDir = "inputDir";

        Configuration conf = new Configuration();
        FileSystem fs = FileSystem.get(conf);

        String outputDir = "newsClusters";
        HadoopUtil.delete(new Path(outputDir));
        Path tokenizedPath = new Path(outputDir,
                                  DocumentProcessor.TOKENIZED_DOCUMENT_OUTPUT_FOLDER);
        MyAnalyzer analyzer = new MyAnalyzer();
```
News clustering example —> finding related articles

Obama to Name 'Smart Grid' Projects
Wall Street Journal - Rebecca Smith - 1 hour ago
The Obama administration is expected Tuesday to name 100 utility projects that will share $3.4 billion in federal stimulus funding to speed deployment of advanced technology designed to cut energy use and make the electric-power grid...

Cobb firm wins "smart-grid" grant  Atlanta Journal Constitution
Obama putting $3.4B toward a 'smart' power grid  The Associate
Baltimore Sun - Bloomberg - New York Times - Reuters
all 594 news articles »  Email this story
News clustering code — II

DocumentProcessor.tokenizeDocuments(new Path(inputDir),
     analyzer.getClass().asSubclass(Analyzer.class),
     tokenizedPath, conf);

DictionaryVectorizer.createTermFrequencyVectors(tokenizedPath,
     new Path(outputDir), conf, minSupport, maxNGramSize, minLLRValue,
     2, true, reduceTasks,
     chunkSize, sequentialAccessOutput, false);

TFIDFConverter.processTfIdf(
     new Path(outputDir),
     DictionaryVectorizer.DOCUMENT_VECTOR_OUTPUT_FOLDER),
     new Path(outputDir), conf, chunkSize, minDf,
     maxDFPercent, norm, true, sequentialAccessOutput, false,
     reduceTasks);

Path vectorsFolder = new Path(outputDir, "tfidf-vectors");
Path canopyCentroids = new Path(outputDir, "canopy-centroids");
Path clusterOutput = new Path(outputDir, "clusters");

CanopyDriver.run(vectorsFolder, canopyCentroids,
     new EuclideanDistanceMeasure(), 250, 120,
     false, false);

KMeansDriver.run(conf, vectorsFolder,
     new Path(canopyCentroids, "clusters-0").
     clusterOutput, new TanimotoDistanceMeasure(), 0.01,
     20, true, false);
SequenceFile.Reader reader = new SequenceFile.Reader(fs,  
    new Path(clusterOutput  
        + Cluster.CLUSTERED_POINTS_DIR + "/part-00000"), conf);  

IntWritable key = new IntWritable();  
WeightedVectorWritable value = new WeightedVectorWritable();  
while (reader.next(key, value)) {  
    System.out.println(key.toString() + " belongs to cluster "  
        + value.toString());  
}  
reader.close();
Other clustering algorithms

Exclusive clustering

Overlapping clustering

Hierarchical clustering

Fiction

F

YA
Different clustering approaches

**FIXED NUMBER OF CENTERS**

**BOTTOM-UP APPROACH: FROM POINTS TO CLUSTERS VIA GROUPING**

Initial datapoints

Clusters formed after 1st iteration

**TOP-DOWN APPROACH: SPLITTING THE GIANT CLUSTER**
When to use Mahout for classification?

<table>
<thead>
<tr>
<th>System size in number of examples</th>
<th>Choice of classification approach</th>
</tr>
</thead>
<tbody>
<tr>
<td>&lt; 100,000</td>
<td>Traditional, non-Mahout approaches should work very well. Mahout may even be slower for training.</td>
</tr>
<tr>
<td>100,000 to 1 million</td>
<td>Mahout begins to be a good choice. The flexible API may make Mahout a preferred choice, even though there is no performance advantage.</td>
</tr>
<tr>
<td>1 million to 10 million</td>
<td>Mahout is an excellent choice in this range.</td>
</tr>
<tr>
<td>&gt; 10 million</td>
<td>Mahout excels where others fail.</td>
</tr>
</tbody>
</table>
The advantage of using Mahout for classification

- Scalable algorithm (Mahout wins!)
- Non-scalable algorithm
- Traditional data mining works here
- Scalable solutions required
**DEFINITION**  Computer classification systems are a form of machine learning that use learning algorithms to provide a way for computers to make decisions based on experience and, in the process, emulate certain forms of human decision making.
How does a classification system work?

Training examples with reference decisions

Predictors and target variables

Training algorithm

Model

Classification system

Copy

New examples

Predictors variables only

Model

Estimated target variable

Emulated decisions
### Key terminology for classification

<table>
<thead>
<tr>
<th>Key Idea</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>Model</td>
<td>A computer program that makes decisions; in classification, the output of the training algorithm is a model.</td>
</tr>
<tr>
<td>Training data</td>
<td>A subset of training examples labeled with the value of the target variable and used as input to the learning algorithm to produce the model.</td>
</tr>
<tr>
<td>Test data</td>
<td>A withheld portion of the training data with the value of the target variable hidden so that it can be used to evaluate the model.</td>
</tr>
<tr>
<td>Training</td>
<td>The learning process that uses training data to produce a model. That model can then compute estimates of the target variable given the predictor variables as inputs.</td>
</tr>
<tr>
<td>Training example</td>
<td>An entity with features that will be used as input for learning algorithm.</td>
</tr>
<tr>
<td>Feature</td>
<td>A known characteristic of a training or a new example; a feature is equivalent to a characteristic.</td>
</tr>
<tr>
<td>Variable</td>
<td>In this context, the value of a feature or a function of several features. This usage is somewhat different from the use of variable in a computer program.</td>
</tr>
<tr>
<td>Record</td>
<td>A container where an example is stored; such a record is composed of fields.</td>
</tr>
<tr>
<td>Field</td>
<td>Part of a record that contains the value of a feature (a variable).</td>
</tr>
<tr>
<td>Predictor variable</td>
<td>A feature selected for use as input to a classification model. Not all features need be used. Some features may be algorithmic combinations of other features.</td>
</tr>
<tr>
<td>Target variable</td>
<td>A feature that the classification model is attempting to estimate: the target variable is categorical, and its determination is the aim of the classification system.</td>
</tr>
</tbody>
</table>
Input and Output of a classification model

<table>
<thead>
<tr>
<th>T</th>
<th>x1 ... xn</th>
</tr>
</thead>
<tbody>
<tr>
<td>T</td>
<td>x1 ... xn</td>
</tr>
<tr>
<td>T</td>
<td>x1 ... xn</td>
</tr>
<tr>
<td>T</td>
<td>x1 ... xn</td>
</tr>
<tr>
<td>T</td>
<td>x1 ... xn</td>
</tr>
</tbody>
</table>

Learning algorithm

Model

<table>
<thead>
<tr>
<th>T</th>
<th>x1 ... xn</th>
</tr>
</thead>
<tbody>
<tr>
<td>?</td>
<td>x1 ... xn</td>
</tr>
<tr>
<td>?</td>
<td>x1 ... xn</td>
</tr>
<tr>
<td>?</td>
<td>x1 ... xn</td>
</tr>
<tr>
<td>?</td>
<td>x1 ... xn</td>
</tr>
</tbody>
</table>

Model

<table>
<thead>
<tr>
<th>T</th>
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</thead>
<tbody>
<tr>
<td>T</td>
<td>T</td>
</tr>
<tr>
<td>T</td>
<td>T</td>
</tr>
</tbody>
</table>

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## Four types of values for predictor variables

<table>
<thead>
<tr>
<th>Type of value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>Continuous</td>
<td>This is a floating-point value. This type of value might be a price, a weight, a time, or anything else that has a numerical magnitude and where this magnitude is the key property of the value.</td>
</tr>
<tr>
<td>Categorical</td>
<td>A categorical value can have one of a set of prespecified values. Typically the set of categorical values is relatively small and may be as small as two, although the set can be quite large. Boolean values are generally treated as categorical values. Another example might be a vendor ID.</td>
</tr>
<tr>
<td>Word-like</td>
<td>A word-like value is like a categorical value, but it has an open-ended set of possible values.</td>
</tr>
<tr>
<td>Text-like</td>
<td>A text-like value is a sequence of word-like values, all of the same kind. Text is the classic example of a text-like value, but a list of email addresses or URLs is also text-like.</td>
</tr>
</tbody>
</table>
Sample data that illustrates all four value types

<table>
<thead>
<tr>
<th>Name</th>
<th>Type</th>
<th>Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>from-address</td>
<td>Word-like</td>
<td>George <a href="mailto:george@fumble-tech.com">george@fumble-tech.com</a></td>
</tr>
<tr>
<td>in-address-book?</td>
<td>Categorical (TRUE, FALSE)</td>
<td>TRUE</td>
</tr>
<tr>
<td>non-spam-words</td>
<td>Text-like</td>
<td>Ted, Mahout, User, lunch</td>
</tr>
<tr>
<td>spam-words</td>
<td>Text-like</td>
<td>available</td>
</tr>
<tr>
<td>unknown-words</td>
<td>Continuous</td>
<td>0</td>
</tr>
<tr>
<td>message-length</td>
<td>Continuous</td>
<td>31</td>
</tr>
</tbody>
</table>
Supervised vs. Unsupervised Learning

Classification algorithms are related to, but still quite different from, clustering algorithms such as the k-means algorithm described in previous chapters. Classification algorithms are a form of supervised learning, as opposed to unsupervised learning, which happens with clustering algorithms. A supervised learning algorithm is one that’s given examples that contain the desired value of a target variable. Unsupervised algorithms aren’t given the desired answer, but instead must find something plausible on their own.

Supervised and unsupervised learning algorithms can often be usefully combined. A clustering algorithm can be used to create features that can then be used by a learning algorithm, or the output of several classifiers can be used as features by a clustering algorithm. Moreover, clustering systems often build a model that can be used to categorize new data. This clustering system model works much like the model produced by a classification system. The difference lies in what data was used to produce the model. For classification, the training data includes the target variables; for clustering, the training data doesn’t include target variables.
Work flow in a typical classification project

<table>
<thead>
<tr>
<th>Stage</th>
<th>Step</th>
</tr>
</thead>
<tbody>
<tr>
<td>1. Training the model</td>
<td>Define target variable.</td>
</tr>
<tr>
<td></td>
<td>Collect historical data.</td>
</tr>
<tr>
<td></td>
<td>Define predictor variables.</td>
</tr>
<tr>
<td></td>
<td>Select a learning algorithm.</td>
</tr>
<tr>
<td></td>
<td>Use the learning algorithm to train the model.</td>
</tr>
<tr>
<td>2. Evaluating the model</td>
<td>Run test data.</td>
</tr>
<tr>
<td></td>
<td>Adjust the input (use different predictor variables, different</td>
</tr>
<tr>
<td></td>
<td>algorithms, or both).</td>
</tr>
<tr>
<td>3. Using the model in production</td>
<td>Input new examples to estimate unknown target values.</td>
</tr>
<tr>
<td></td>
<td>Retrain the model as needed.</td>
</tr>
</tbody>
</table>
Classification Example 1 — Color-Fill

Position looks promising, especially the x-axis ==> predictor variable. Shape seems to be irrelevant. Target variable is “color-fill” label.
Target leak

• A target leak is a bug that involves unintentionally providing data about the target variable in the section of the predictor variables.
• Don’t confused with intentionally including the target variable in the record of a training example.
• Target leaks can seriously affect the accuracy of the classification system.
Classification Example 2 — Color-Fill (another feature)
Mahout classification algorithms include:

- Naive Bayesian
- Complementary Naive Bayesian
- Stochastic Gradient Descent (SDG)
- Random Forest
Comparing two types of Mahout Scalable algorithms

![Diagram comparing wall clock time vs. number of training examples for sequential and parallel algorithms.](image)
Step-by-step simple classification example

1. The data and the challenge
2. Training a model to find color-fill: preliminary thinking
3. Choosing a learning algorithm to train the model
4. Improving performance of the classifier
Classification Example 3
What may be a good predictor?
Homework #2

Recommendation:
1. Choose any two datasets from Yahoo Labs Ratings and Classification Data.
2. Try various recommendation algorithms provided by Mahout

Clustering:
Using datasets from:
1. Online news (e.g., New York Times article in September 2015, or other data sources)
2. Wikipedia articles
3. (optional) gather data from Twitter API, try clustering
Do clustering —> finding related documents

Classification:
1: Using the 20 newsgroups data (will be provided by TA), try various classification algorithms provided by Mahout, and discuss their performance
2: Do similar experiments on the Wikipedia data that you downloaded.
Questions?