CRAFTML

An Efficient Clustering-based Random Forest for Extreme Multi-label Learning

Motivation

- In normal classification, we have a model defined, which classifies or tags a data instance with only one class label.
- If there are multiple class labels, the classifier will choose only one(best) among those.
 - Questions that arise -
- What if there are multiple possible tags (labels) associated with the data?
- Can a data instance be classified/tagged with multiple possible class labels from the set?
- How the model should be designed and how can we calculate accuracy for that model?

Examples



Single Label Classification: Is there a house? Yes / No

Multi Label Classification:

House	Tree	Beach	Cloud	Mountain	Animal			
Yes	Yes	no	Yes	no	no			

Examples

Three Type of Classification Tasks



Binary Classification



- Spam
- Not spam

Multiclass Classification



- Dog
- Cat
- Horse
- Fish
- Bird
- ...

Multi-label Classification



- Dog
- Cat
- Horse
- Fish
- Bird
- ...

eXtreme Multi-label Learning (XML)

- eXtreme Multi-label Learning (XML) considers large sets of items described by a number of labels that can exceed one million.
- We can do this classification using many existing machine learning algorithms, but there are some disadvantages.

Problems with existing algorithms when large number of labels are present -

- Scalability issues
- Performance degradation.

How to overcome these problems?

3 common ways:

- Using optimization tricks like sparsification and parallelization.
- Reducing the data dimensionality for solving a smaller size problem.
- Tree based approach: Hierarchically partitioning the initial problem into small scale sub-problem.

Previous works

Optimization tricks and parallelization :
 PDSparse, PPDSparse, DISMEC

Dimensionality reduction:
 WSABIE, LEML, SLEEC, AnnexML

 Tree based approach : LPSR, FastXML, PFastReXML

CraftML

- CRAFTML is a random forest based algorithm with a very fast partitioning approach.
- The splitting conditions are based on all the features.
- CRAFTML randomly reduces both the feature and the label spaces to obtain diversity.
- It replaces random selections with random projections to preserve more information.

Node structure of the decision tree

```
struct Node{
   int number_of_children;
   int branch_value;
   int split_attribute;
   int leaf_value;
   struct Node *children[10];
};
```

number_of_children: number of children in each node branch_value: make branch decision based on this value split_attribute: splitting attribute (-1 for leaf node) leaf_value: class value at leaf node (-1 for decision node)

Random projection of the dataset:

We randomly project the label and feature vectors into lower dimensional spaces.

```
void chooseRandomFeatures(){
    vector<vector<double> > trainFileRandom( N , vector<double> (M, 0));
    int number_of_features = 50;
    for(int i=0; i<number_of_features; i++){
        int guess = rand() % (M-1);
        trainFileRandom[i]=train_file[guess];
    }
    train_file=trainFileRandom;
}</pre>
```

Note: In contrary to classical random forests which use bootstraps, each tree of CRAFTML is trained on the full initial dataset.

We only select (project) a subset of the feature and labels space.

K Means algorithm

We build a k-means based partitioning of the instances into k temporary subsets from their projected labels.

```
double** k means(){
     int minima[features]={INT MAX};
     int maxima[features]={INT MIN};
     int cluster[N];
     int t=20, k;
     double mean arr[K][features];
     for(int i=0; i<K; i++){
         for(int j=0; j<features; j++){</pre>
             int num = (rand() % (maxima[j] - minima[j] + 1)) + minima[j];
             mean arr[i][j]=num;
     for (int i = 0; i < t; i++) {
         for (int j = 0; j < N; j++) {
             double* dists = new double[k];
             for (int p = 0; p < k; p++) {
                 dists[p] = cosine distance( trainFile1[j], mean arr[p], M);
             cluster[j] = std::min element(dists, dists + k) - dists;
             delete[] dists;
```

When to stop?

- (i) the cardinality of the node's instance subset is lower than a given threshold.
- (ii) all the instances have the same features
- (iii) all the instances have the same labels

```
void decision(int *h_attr, int *h_data, node *root, int h_dataSize) {
   int threshold = 10;
   // stopping conditions

// checking whether the cardinality of the node's instance subset lower than a
   if(h_dataSize<=threshold)
        return;

// checking whether every instances have the same labels
   flag=1;
   for(int i=1;i<h_dataSize;i++){
        if(trainFile[h_data[i]][M-1]!=trainFile[h_data[i-1]][M-1]){
        flag=0;
        break;
    }
   }
   if(flag=1){
        root->val=trainFile[h_data[0]][M-1];
        return;
}
```

CraftML: Predictions

- For each tree, the input instance follows a root-to-leaf path.
- The path is determined by the successive decisions of the classifier.
- The prediction is the average label vector stored in the leaf reached.
- The forest aggregates the tree predictions with the average operator.

CraftML: Algorithm

```
Algorithm 1 trainTree
  Input: Training set with a feature matrix X and a label
   matrix Y.
   Initialize node v
   v.isLeaf \leftarrow testStopCondition(X, Y)
  if v.isLeaf = false then
      v.classif \leftarrow trainNodeClassifier(X, Y)
      (X_{child_i}, Y_{child_i})_{i=0,...,k-1} \leftarrow \text{split}(v.\text{classif}, X, Y)
      for i from 0 to k-1 do
         v.child_i \leftarrow trainTree(X_{child_i}, Y_{child_i})
      end for
   else
      v.\widehat{y} \leftarrow \text{computeMeanLabelVector}(Y)
   end if
   Output: node v
```

CraftML: Algorithm

Algorithm 2 trainNodeClassifier

Input: feature matrix (X_v) and label matrix (Y_v) of the instance set of the node v.

```
X_s, Y_s \leftarrow \operatorname{sampleRows}(X_v, Y_v, n_s)

X'_s \leftarrow X_s P_x # random feature projection

Y'_s \leftarrow Y_s P_y # random label projection

\mathbf{c} \leftarrow k\operatorname{-means}(Y'_s, k) # \mathbf{c} \in \{0, ..., k-1\}^{\min(n_v, n_s)}

for i from 0 to k-1 do

(classif)<sub>i,.</sub> \leftarrow computeCentroid(\{(X'_s)_{j,.}|c_j=i\})

end for

Output: Classifier classif (\in \mathbb{R}^{k \times d'_x}).
```

c is a vector where the j^{th} component c_j denotes the cluster index of the j^{th} instance associated to $(X'_s)_{j,..}$ and $(Y'_s)_{j,..}$.

CraftML: How to parallelize?

There are 2 parts where we can parallelize our code:

1. Buliding trees: While building the individual decision trees.

2. Predictions: While making the predictions from the different decision trees.

CraftML: Building Trees in parallel

- The trees are independant of one another.
- We keep the number of blocks equal to the number of trees.
- We build each of the trees parallely in a thread.
- The code snippet for creation of trees in 50 blocks is shown below :

```
#define NUMBER_OF_TREES 50
buildDecisionTree<<<<NUMBER_OF_TREES, 1>>>(device_data, number_of_features, number_of_samples)
```

```
global_ void buildDecisionTree(int trainData[BLOCK_SIZE][BLOCK_SIZE], int number_of_features,int number_of_samples){
   int bid = blockIdx.x;
   _shared_ int randomFeatures[50];
   // Choose Random Features
   for(int i=0; i<number_of_features; i++){
        randomFeatures[i] = rand() % number_of_features;
}
// Build Tree</pre>
```

CraftML: Predictions in parallel

- The trees built are independent of one another.
- We send the input instances to each of the threads which run in parallel.
- Each of the trees outputs the predicted set of labels.
- We take the majority of the individual outputs to make our predictions.

CraftML: Alternative approach

- The trees built are independant of one another
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- Each of the trees outputs the predicted set of labels.
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CraftML: Alternative approach

Advantages:

- When we have large amounts of data, like in eXtreme Multi Label Learning (XML), deciding the split attributes for each tree sequentially is tedious.
- Thus, we make the construction of each tree in parallel. Disadvantages:
- We incur a loss in computational time by building the different trees sequentially.

Implementation

Dataset:

- We take a popular multi label dataset: Yeast dataset
- It has 1500 samples, 103 features and 14 labels.
- A snapshot of a part of the dataset is shown below:

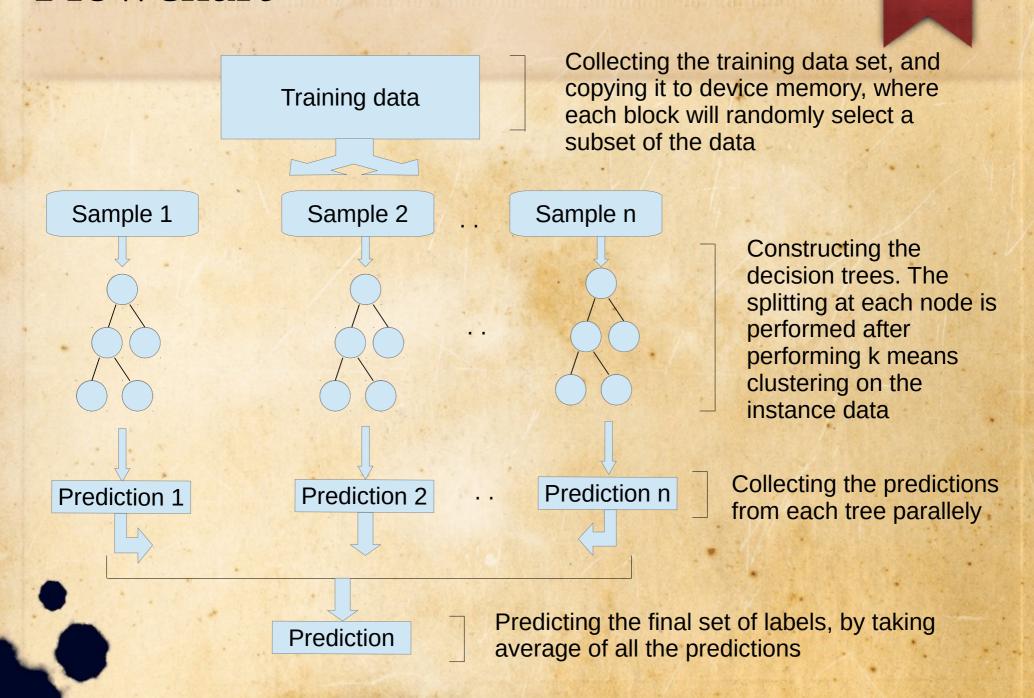
		Att1	Att2	Att3	Att4	Att5	Att6	Att7	Att8	Att9	Att10	 Class5	Class6	Class7	Class8	•
(0	0.093700	0.139771	0.062774	0.007698	0.083873	-0.119156	0.073305	0.005510	0.027523	0.043477	 0	0	0	0	1
1	1	-0.022711	-0.050504	-0.035691	-0.065434	-0.084316	-0.378560	0.038212	0.085770	0.182613	-0.055544	 0	0	1	1	1
2	2	-0.090407	0.021198	0.208712	0.102752	0.119315	0.041729	-0.021728	0.019603	-0.063853	-0.053756	 0	0	0	0	ı
:	3	-0.085235	0.009540	-0.013228	0.094063	-0.013592	-0.030719	-0.116062	-0.131674	-0.165448	-0.123053	 0	0	0	0	1
-	1	-0.088765	-0.026743	0.002075	-0.043819	-0.005465	0.004306	-0.055865	-0.071484	-0.159025	-0.111348	 0	0	0	0	1

Implementation

Steps:

- Collect the data in host memory.
- Copy data from host memory to device memory.
- Create number of blocks equal to the number of decision trees and call the kernel to build the trees.
- In each parallel block, create a projection into lower dimension space of the dataset.
- Build the trees in parallel.
- To make the predictions, take the test instance and make it traverse the root-to-leaf path of every tree.
- Take the average predictions from all the trees and output the final multi label prediction.
- Calculate the accuracy by comparing it to the ground truth labels.

Flowchart



Output

- We predict the labels for the different samples in the test data.
- The test data has 917 samples, with 103 features and also the ground truth values of the labels.
- We compare our predictions for each label (class) with it's ground truth value, and calculate the accuracy of each class.

```
Accuracy: 0.688113
          Accuracy: 0.571429
lass2
          Accuracy: 0.580153
lass3
          Accuracy: 0.640131
lass4
:lass5 :
          Accuracy: 0.693566
Class6 : Accuracy: 0.761178
          Accuracy: 0.817884
Class7 :
Class8
          Accuracy: 0.791712
          Accuracy: 0.912759
Class9
Class10 : Accuracy: 0.899673
           Accuracy: 0.900763
           Accuracy: 0.750273
Class12
           Accuracy: 0.74482
Class13
           Accuracy: 0.985823
lass14
```

Similar calculations of accuracy of each label is tested using the Random Forest module of sklearn and shown later.

Comparing with sklearn

- We run the sklearn Random Forest module on our dataset, using train validation split.
- Random forest classifier

```
[17] 1 from sklearn.ensemble import RandomForestClassifier
2 from sklearn.metrics import accuracy_score
3 clf = RandomForestClassifier(n_estimators=50, criterion='gini', max_depth=None, bootstrap=True)
```

Since number of samples is not very large, it gives us the output in comparable time.

Comparing with sklearn

- Output:
 - Accuracy of the classes

```
[18] 1 for category in y test:
     2 clf.fit(X train, y train[category])
        y pred = clf.predict(X test)
        print(category, " : ", accuracy score(y test[category],y pred))
 Class2 : 0.65
    Class3 : 0.74
    Class4 : 0.7166666666666667
    Class5 : 0.7366666666666667
    Class6 : 0.766666666666667
    Class7 : 0.85
    Class8 : 0.84
    Class9 : 0.946666666666667
    Class10 : 0.89333333333333333
    Class11 : 0.86333333333333333
    Class12 : 0.74333333333333333
    Class13 : 0.7133333333333334
    Class14 : 0.98
```



We see that we get comparable results.

Future scope

- Run CraftML algorithm on larger dimensional data, and more number of labels, and test the performance.
- Compute the multi label predictions parallely in each of the decision trees.
- Compare the performances and time gains with respect to sequential algorithm on multiple machines.

