

# [544] Spark MLlib

Tyler Caraza-Harter

# Outline

ML Review

Training/Predicting APIs

Demos

Decision Trees

Demos

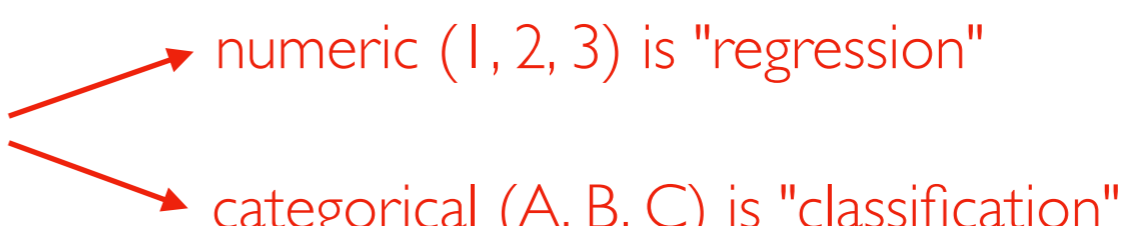
# Machine Learning, Major Ideas

Categories of Machine Learning:

- **Reinforcement learning:** agent makes series of actions to maximize reward
- **Unsupervised learning:** looking for generate patterns
- **Supervised learning:** train models to predict unknowns

**Models** are functions that return predictions:

```
def my_model(some_info):  
    ...  
    return some_prediction
```



numeric (1, 2, 3) is "regression"  
categorical (A, B, C) is "classification"

**Example:**

```
def weather_forecast(temp_today, temp_yesterday):  
    ...  
    return temp_tomorrow
```

# Learning from Data

	x1	x2	y
0	2	8	5
1	9	2	6
2	4	1	0
3	7	9	7
4	2	2	3
5	3	4	3
6	3	5	9
7	7	1	4
8	6	6	3
9	4	3	?
10	1	2	?
11	2	9	?

- feature columns: x1 and x2
- label column: y

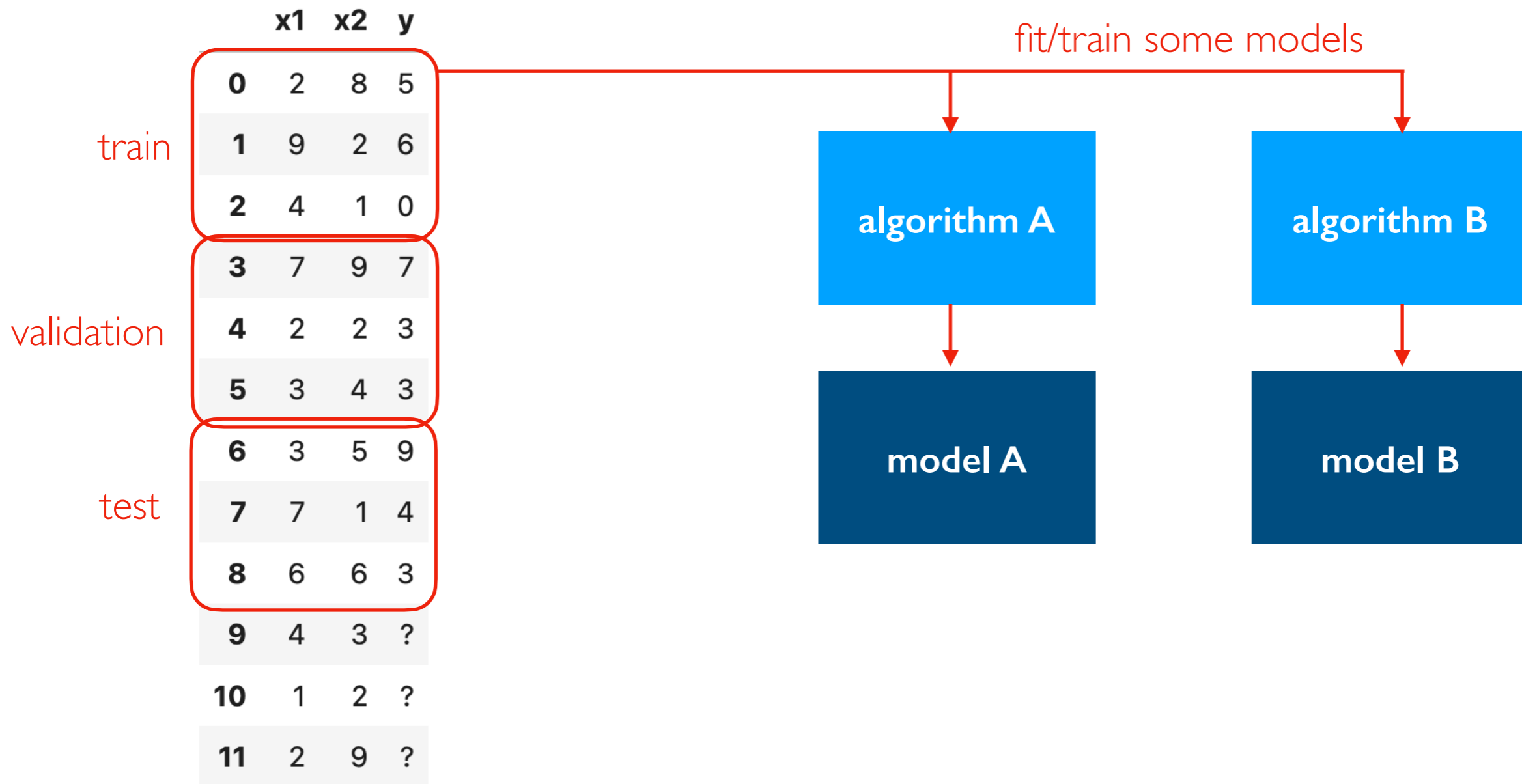
how can the cases where we DO know y help us predict the cases where we do not?

# Learning from Data

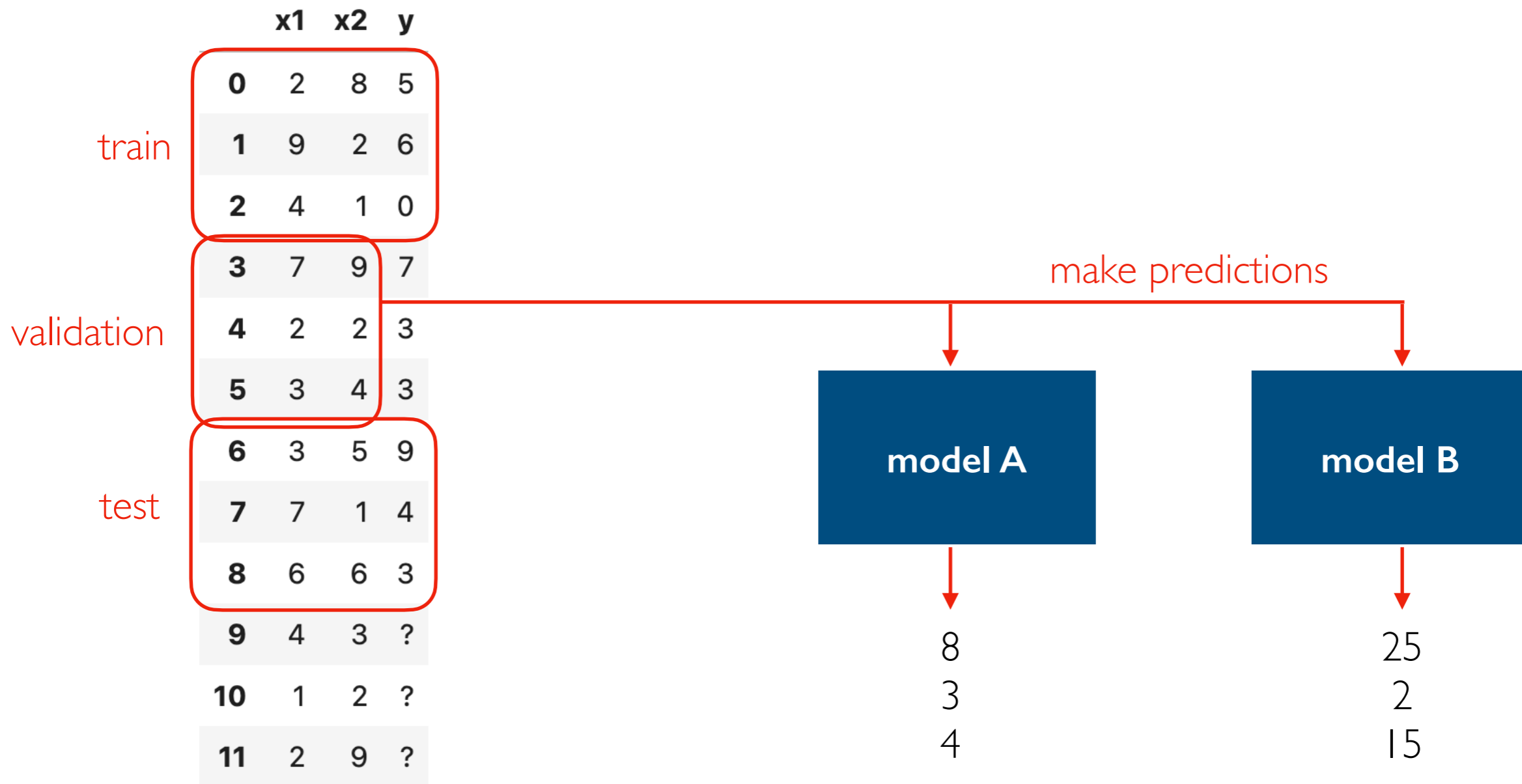
	x1	x2	y	
train	0	2	8	5
	1	9	2	6
	2	4	1	0
validation	3	7	9	7
	4	2	2	3
	5	3	4	3
test	6	3	5	9
	7	7	1	4
	8	6	6	3
	9	4	3	?
	10	1	2	?
	11	2	9	?

random split

# Learning from Data



# Learning from Data



# Learning from Data

	x1	x2	y	
train	0	2	8	5
	1	9	2	6
	2	4	1	0
validation	3	7	9	7
	4	2	2	3
	5	3	4	3
test	6	3	5	9
	7	7	1	4
	8	6	6	3
	9	4	3	?
	10	1	2	?
	11	2	9	?

which model predicts better?

winner!

model A

8  
3  
4

model B

25  
2  
15



# Learning from Data

	x1	x2	y	
train	0	2	8	5
	1	9	2	6
	2	4	1	0
validation	3	7	9	7
	4	2	2	3
	5	3	4	3
test	6	3	5	9
	7	7	1	4
	8	6	6	3
	9	4	3	?
	10	1	2	?
	11	2	9	?

why might the winning model do worse on the test data than the validation data?

winner!

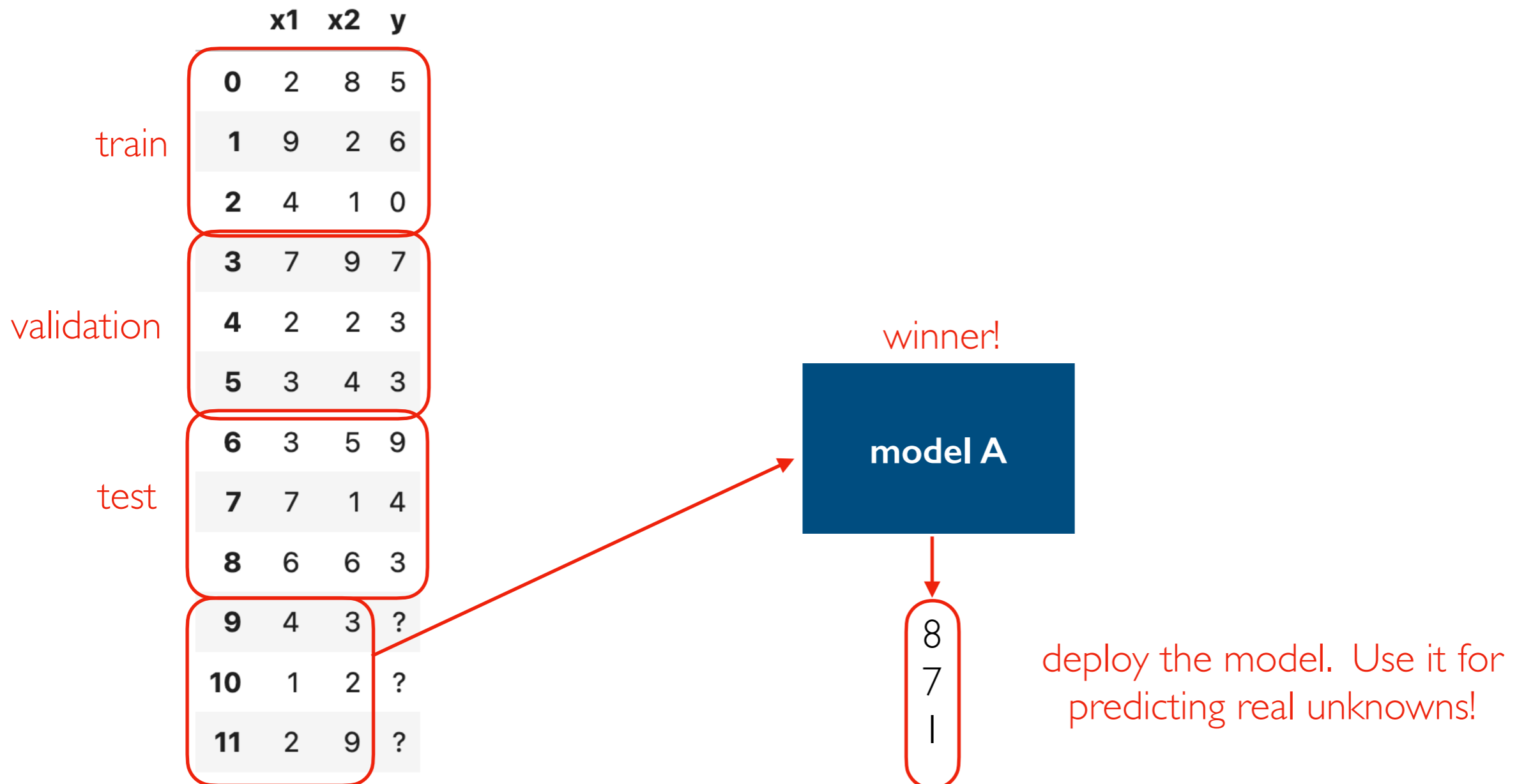
model A

10  
3  
3

how good does the chosen model do on the test data?

models that do good on train data but bad on validation/test data have "overfitted"

# Learning from Data



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- sklearn
- PyTorch
- Spark MLlib

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# Training

## scikit-learn

```
model = ????  
model.fit(X, y)  
# model parameters can relate X to y
```

## pytorch

```
model = ????  
# TODO: optimizer, loss function  
# training loop  
for epoch in range(????):  
    for X, y in ????:  
        ...  
# model parameters can relate X to y
```

- models are **mutable**
- fitting sets/improves parameters

## Spark MLlib

```
unfit_model = ????  
fit_model = unfit_model.fit(df)  
# fit_model params can relate x to y
```

- models are **immutable**
- fitting returns new model object

# Predicting

## scikit-learn

```
y = model.predict(X)
```

## pytorch

```
y = model(X)
```

## Spark MLlib

```
df2 = fit_model.transform(df)
```

# Data

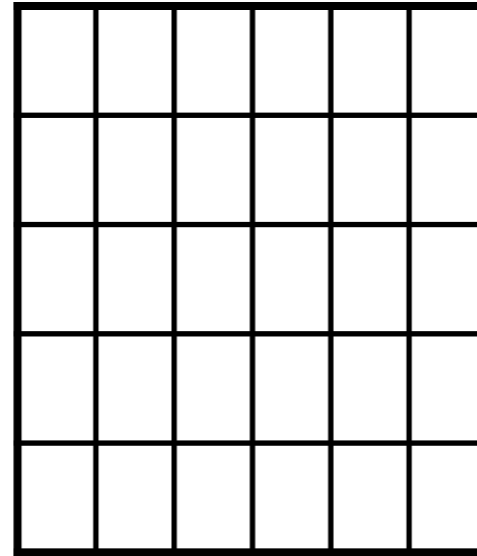
scikit-learn

```
y = model.predict(X)
```

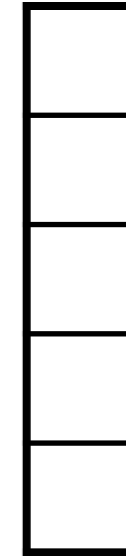
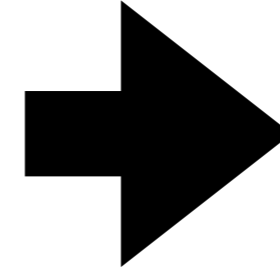
pytorch

```
y = model(X)
```

X (features)



y (label)

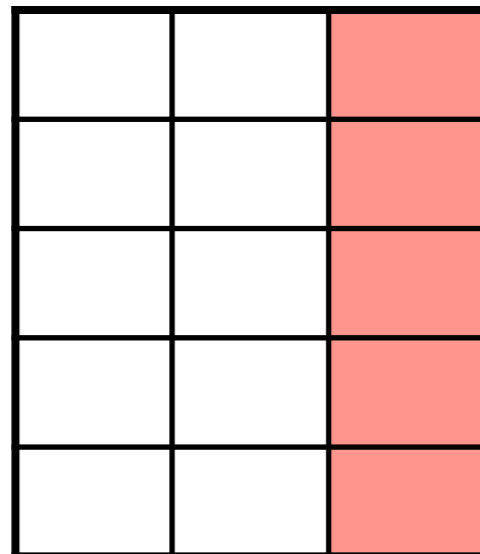


Spark MLlib

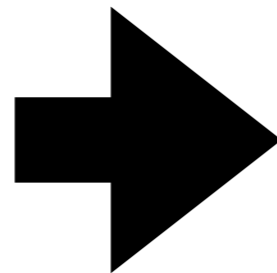
```
df2 = fit_model.transform(df)
```

X (features)

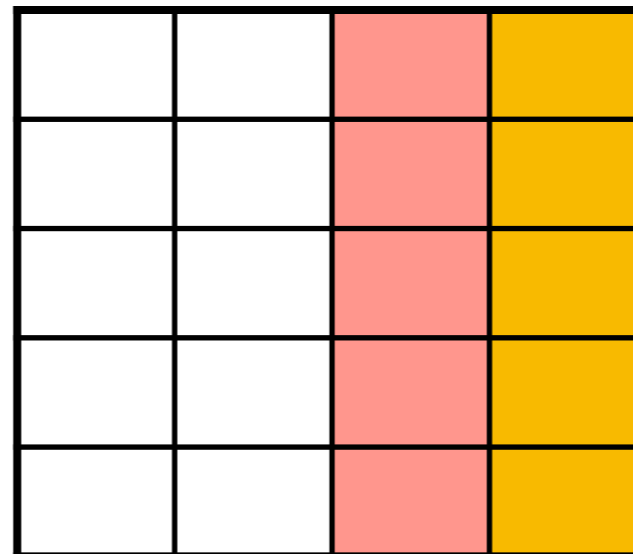
df



df2

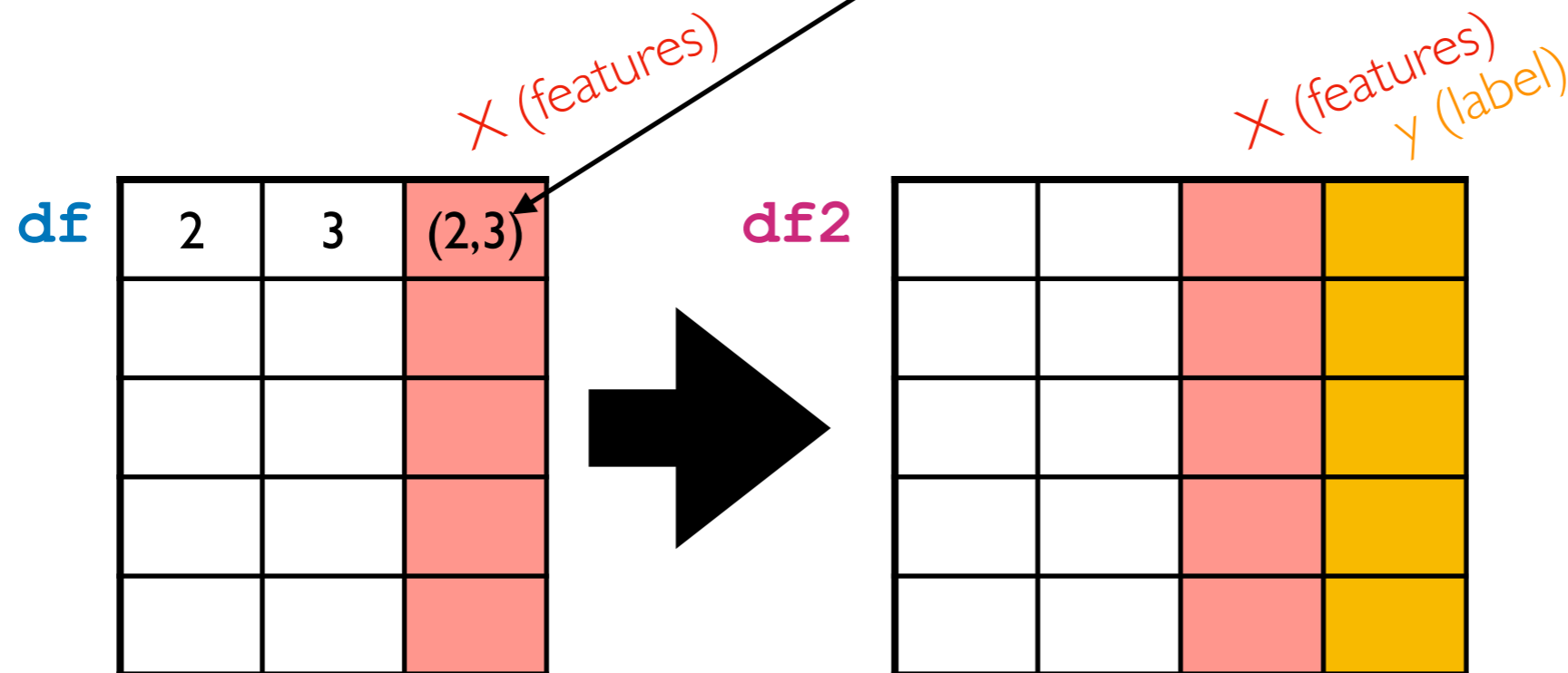


X (features)  
y (label)



# Features Column

- we only get one features column
- it usually contains vectors
- those vectors typically contain values from other columns
- example:  
(2, 3)



# Terminology

Spark and scikit-learn use many of the same terms, with very different meaning.

## Transformer (scikit-learn)

- has .transform method
- takes a DataFrame, returns a different DataFrame
- used as **preprocessing step** for a model

## Transformer (Spark)

- has .transform method
- takes a DataFrame, returns original **with 1 or more additional columns**
- **a fitted model is a transformer** that adds a prediction column

## Estimator (scikit-learn)

- has .fit and .predict methods
- .fit **modifies** the object
- makes predictions after learning params

## Estimator (Spark)

- has .fit method that **returns new object**
- an **unfitted model** is an estimator; calling .fit returns a **fitted model** (a transformer)



# Pipeline

Both scikit-learn and Spark: a pipeline is a series of stages (transformers/estimators). fit/transform/etc. are called as appropriate on each stage.

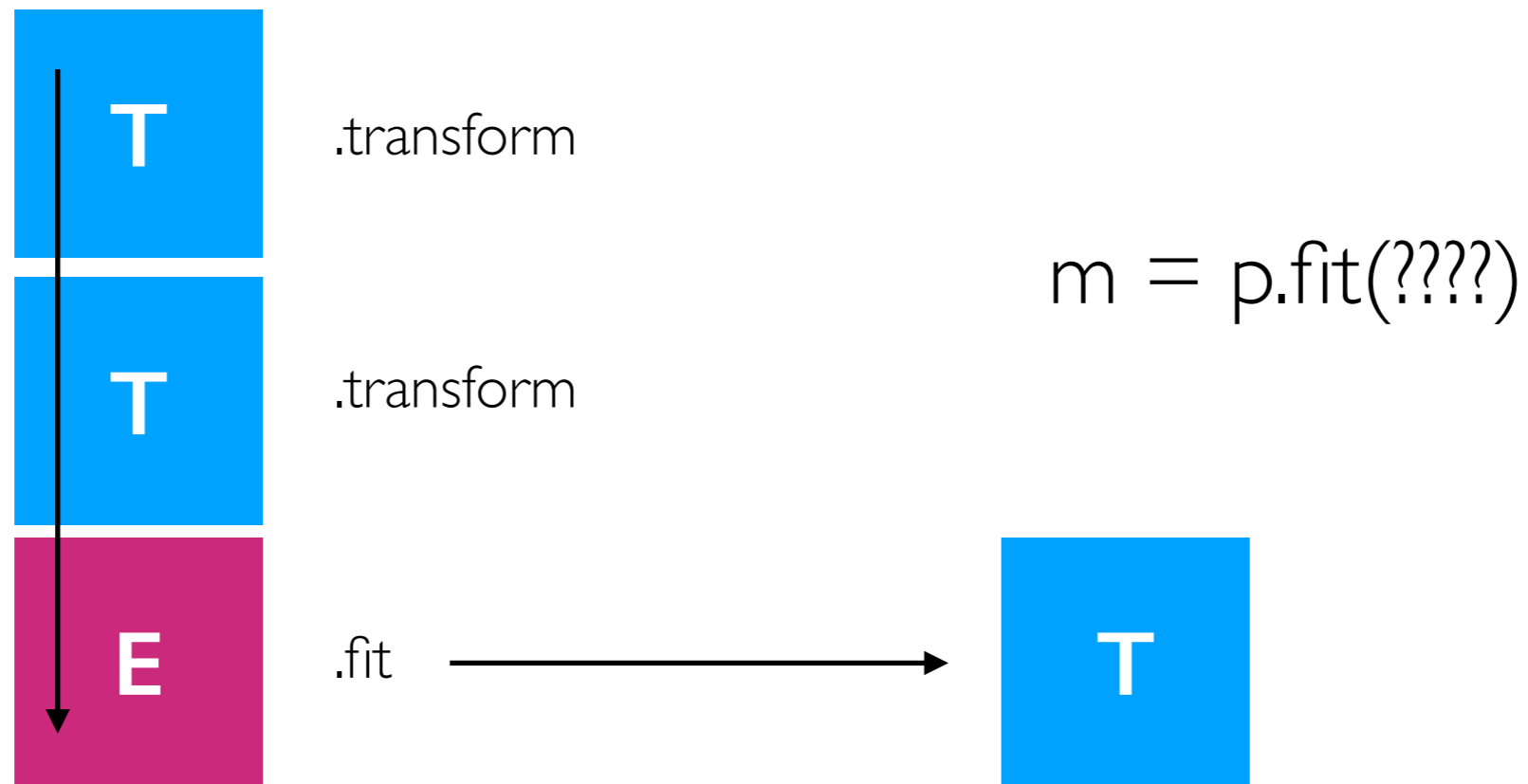
Pipeline (p)



# Pipeline (Spark Example)

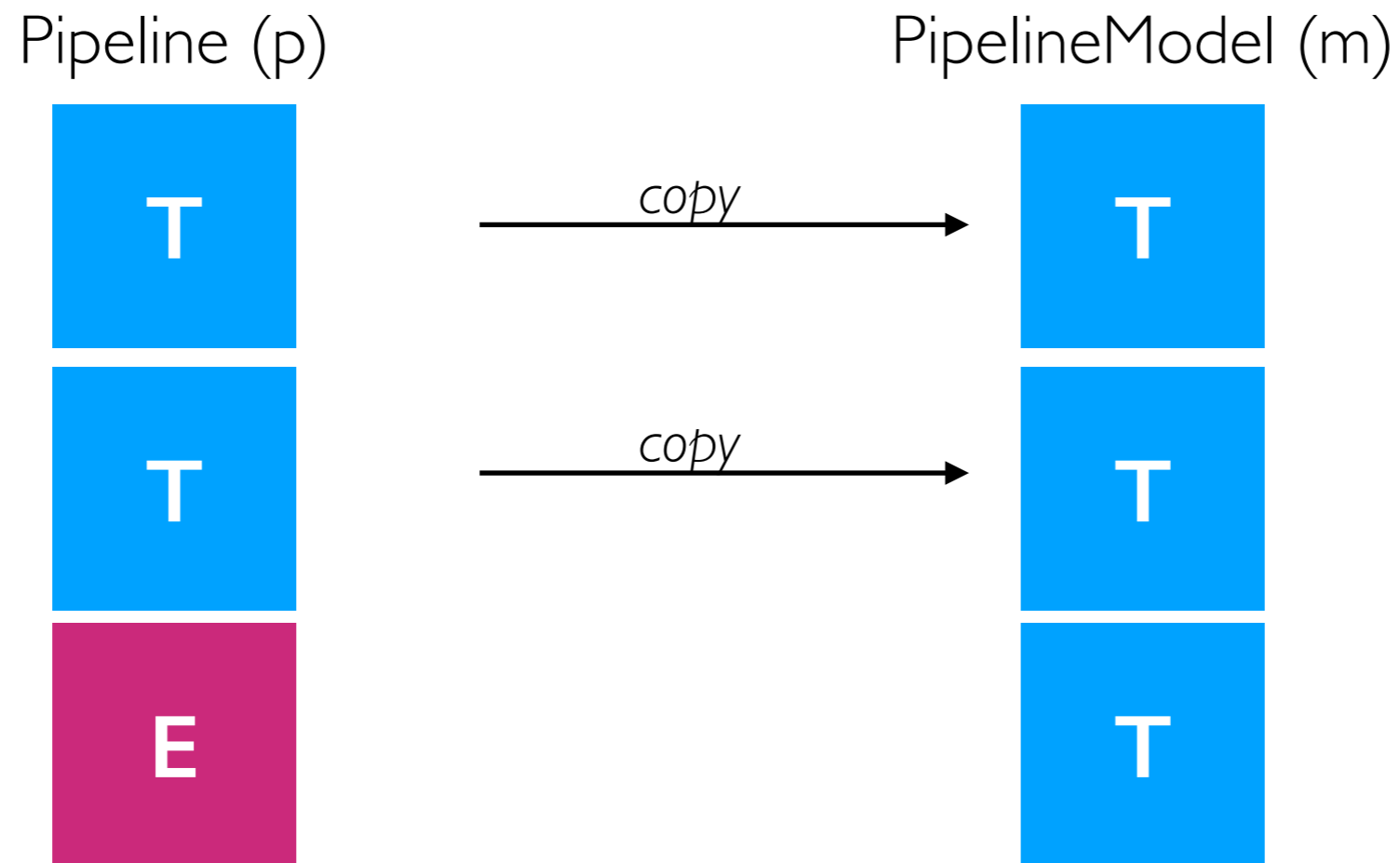
Both scikit-learn and Spark: a pipeline is a series of stages (transformers/estimators). fit/transform/etc. are called as appropriate on each stage.

Pipeline (p)



# Pipeline (Spark Example)

Both scikit-learn and Spark: a pipeline is a series of stages (transformers/estimators). fit/transform/etc. are called as appropriate on each stage.



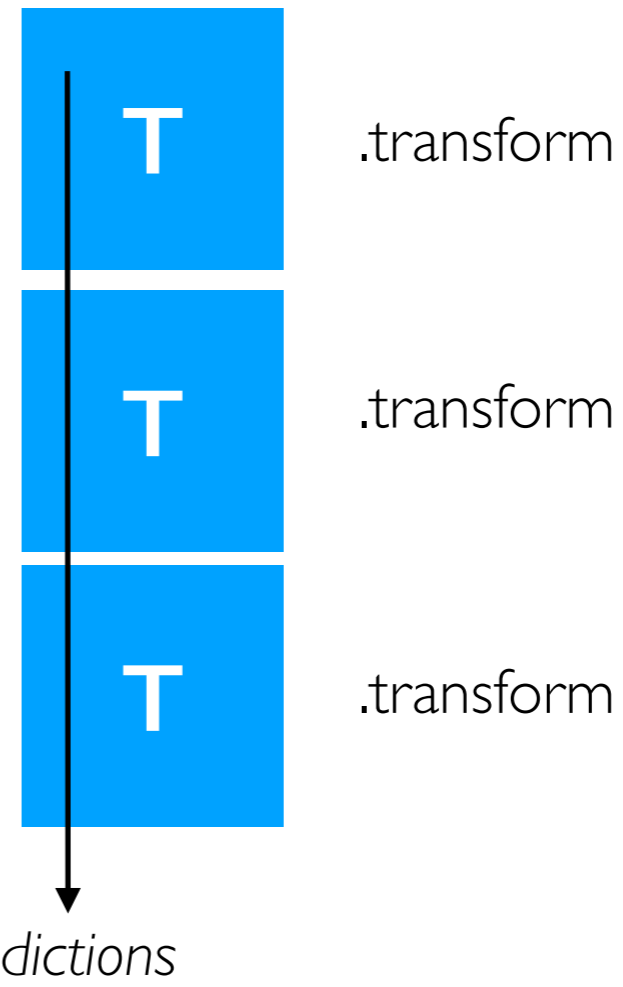
# Pipeline (Spark Example)

Both scikit-learn and Spark: a pipeline is a series of stages (transformers/estimators). fit/transform/etc. are called as appropriate on each stage.

Pipeline (p)



PipelineModel (m)



`m.transform(????)`

TopHat

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Demos

## Spark mllib packages

- pyspark.mllib -- based on RDDs
- [pyspark.ml](#) -- based on DataFrames

# Distributed ML Outline

ML Review

Training/Predicting APIs

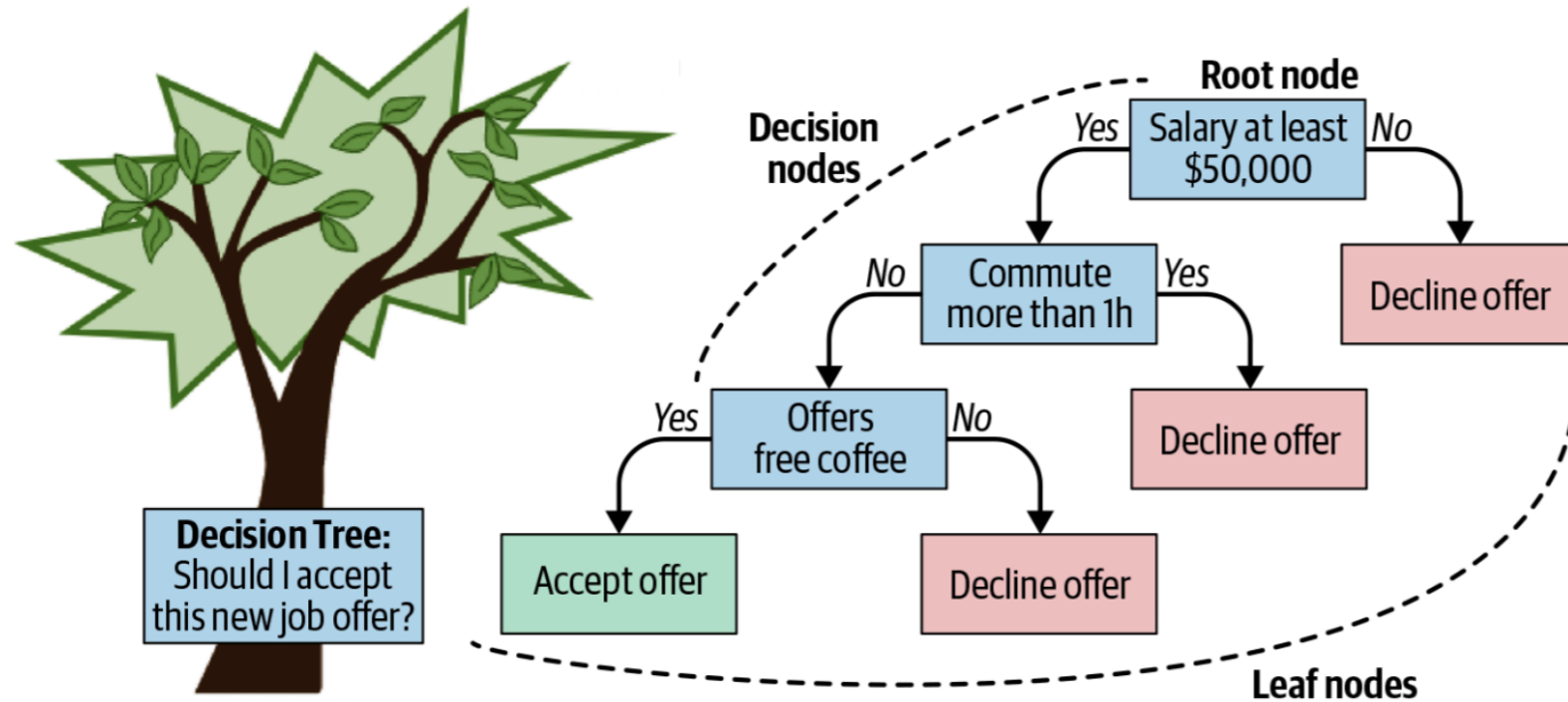
Demos

Decision Trees:

- Background
- Training in memory
- PLANET algorithm

Demos

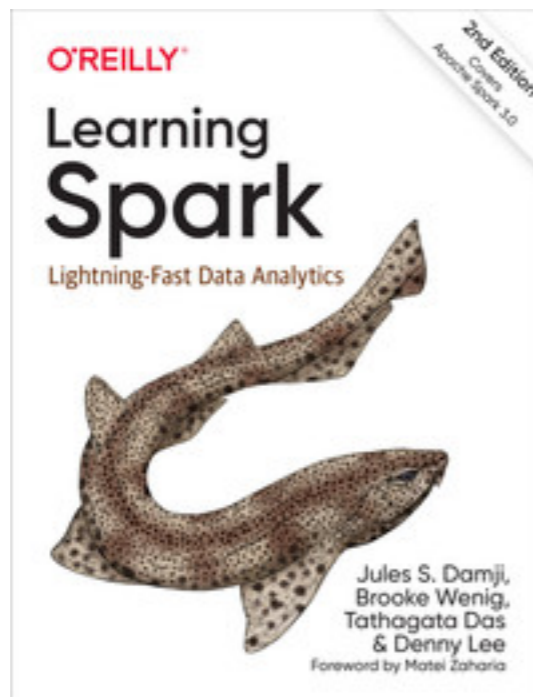
# Decision Trees



decision trees are like nested if/else statements

features and labels can be numeric or categorical

Figure 10-9. Decision tree example



```
def predict(row):  
    if row.salary < 50K:  
        return False  
    else:  
        if row.commute > 1h:  
            return False  
        else:  
            if row.coffee == "free":  
                return True  
            else:  
                return False
```



# Ensemble Methods

**Ensemble:** many simple models vote. Many simple decision trees (each trained on subset of rows/columns) together are often better than one big tree. Examples:

- random forest
- gradient-boosted trees

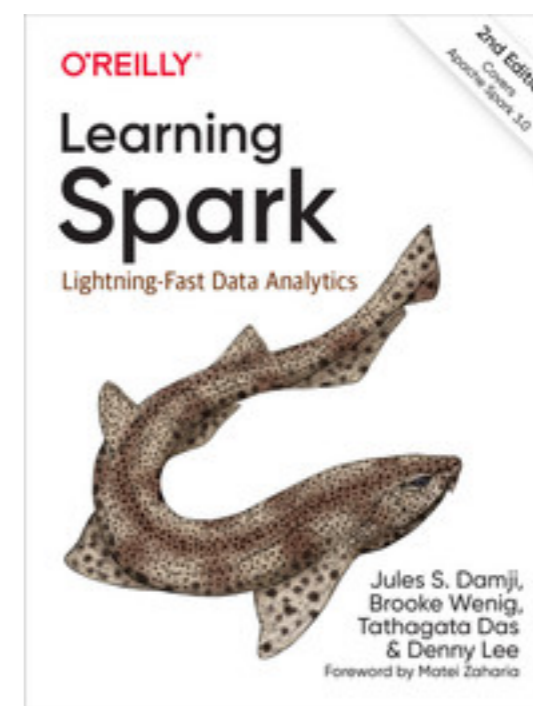
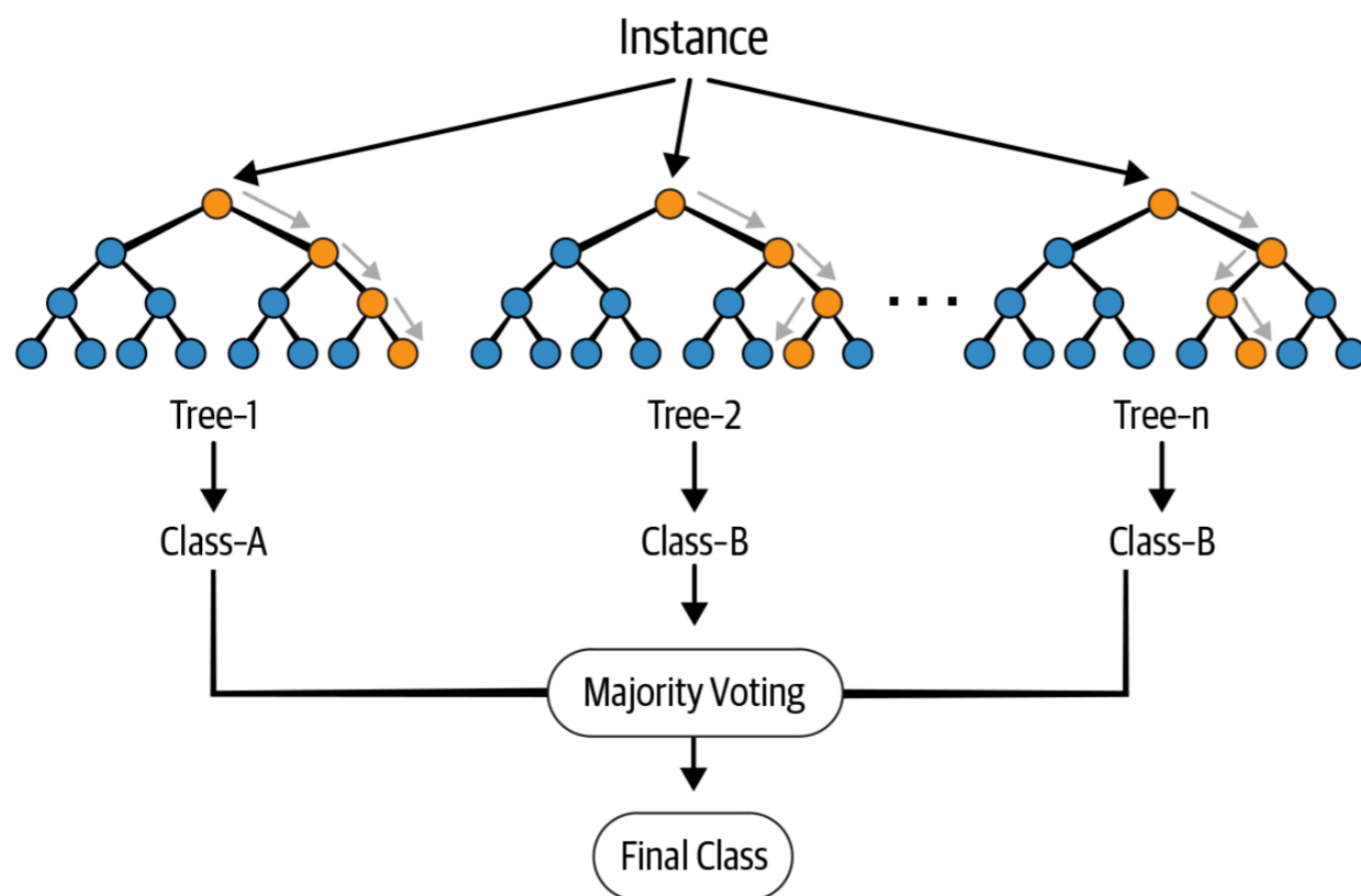


Figure 10-12. Random forest predictions

A Spark cluster can train many trees in a random forest simultaneously!

# Tree methods vs. Deep Learning

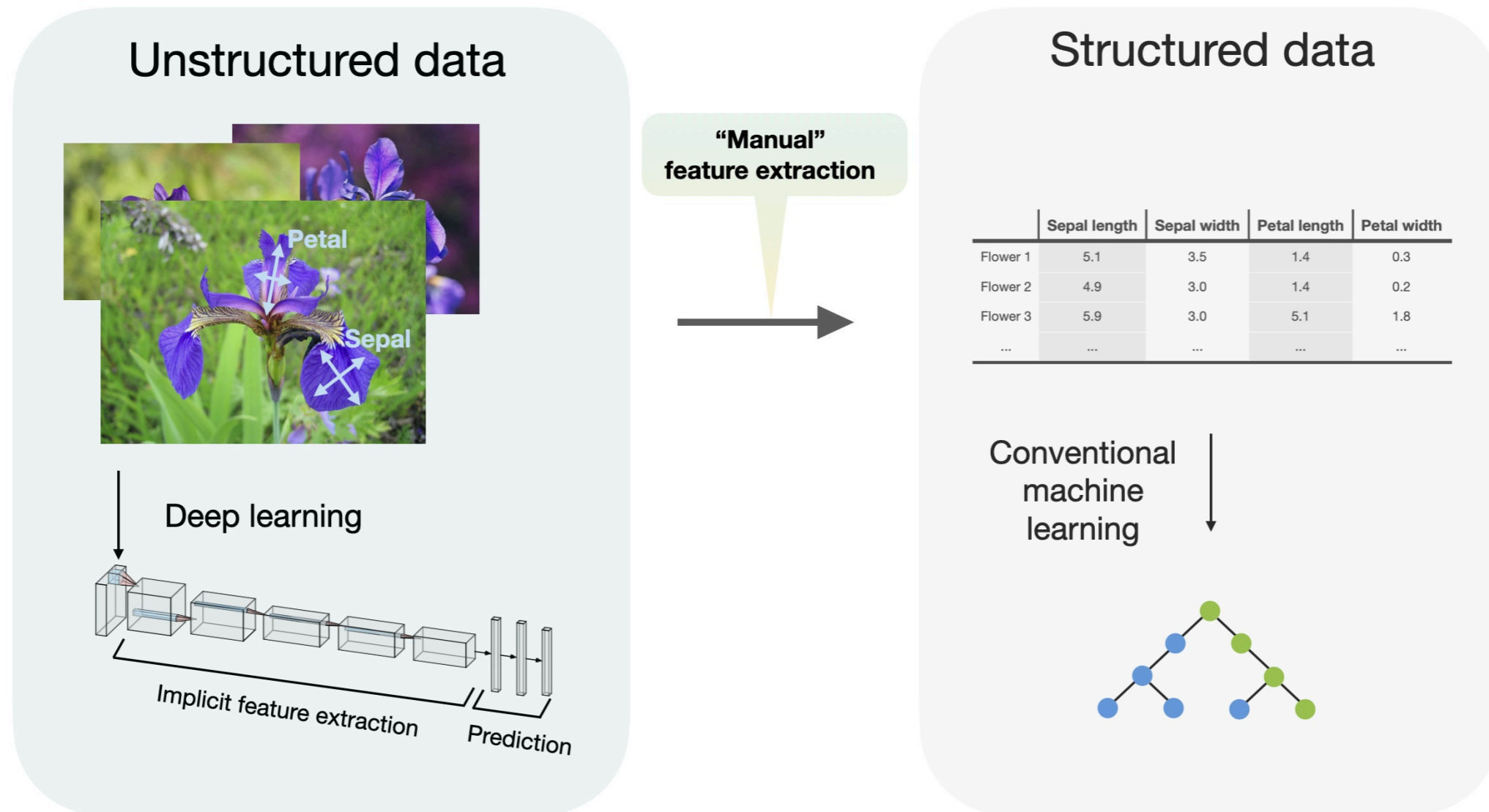


Image from Blog Post: A Short Chronology Of Deep Learning For Tabular Data  
Sebastian Raschka

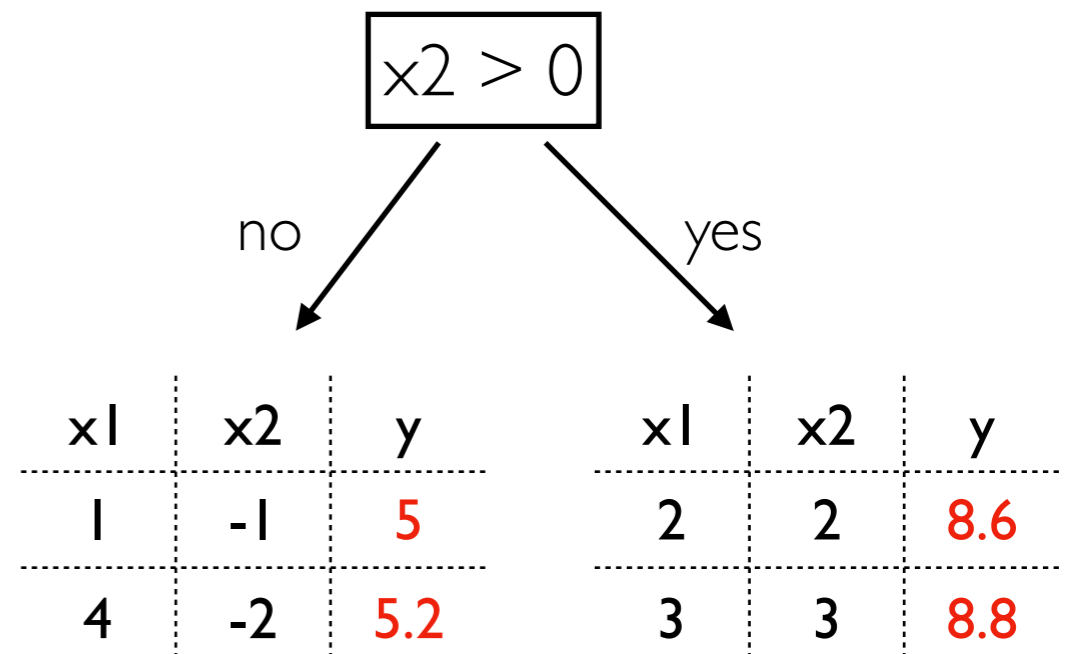
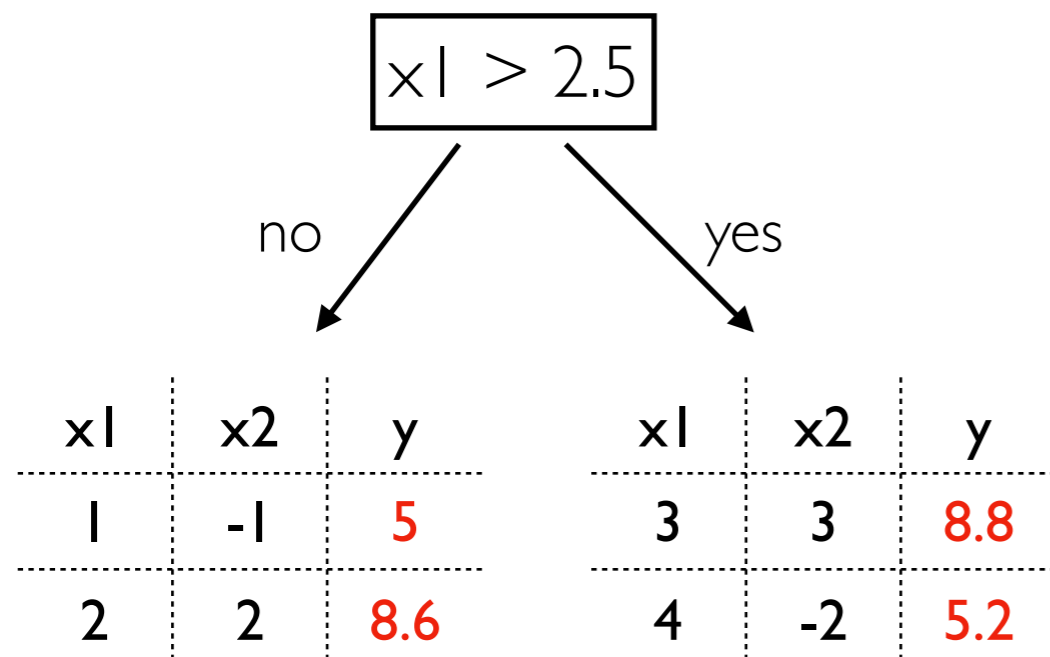
<https://sebastianraschka.com/blog/2022/deep-learning-for-tabular-data.html>

*Tree-based methods are still relevant in the age of deep learning because there are many important tabular datasets.*

# Is a Tree Good?

**data:**

x1	x2	y
1	-1	5
2	2	8.6
3	3	8.8
4	-2	5.2



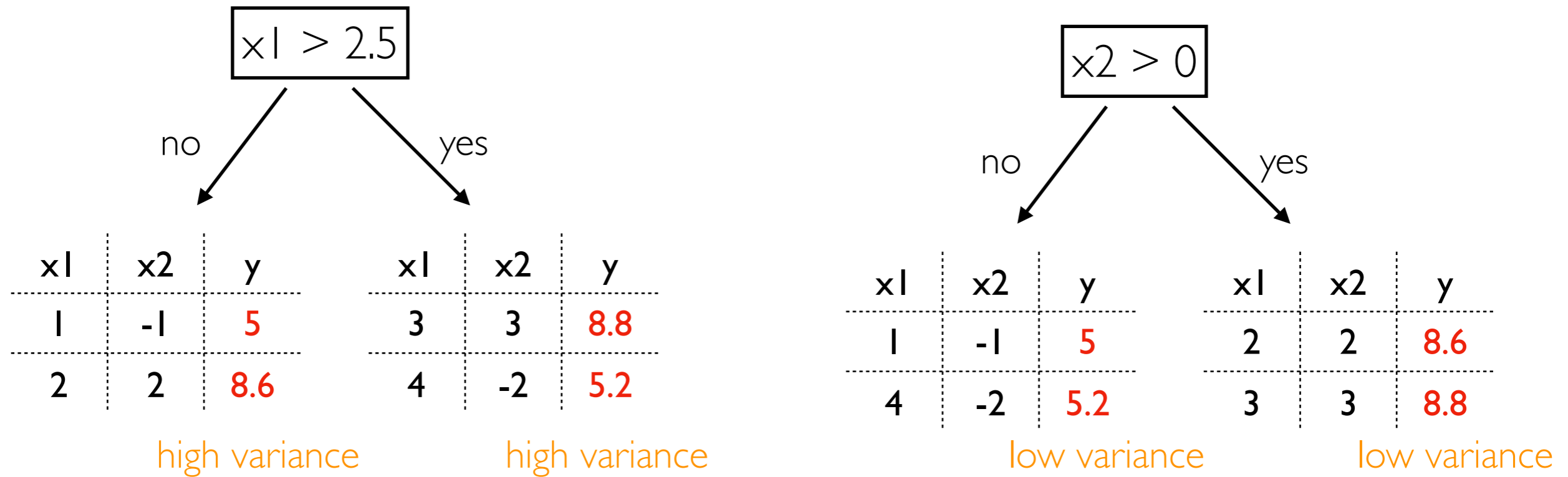
which tree asks better questions about x values if we want to predict y?

# Impurity

**data:**

x1	x2	y
1	-1	5
2	2	8.6
3	3	8.8
4	-2	5.2

*better tree*



impurity measures (like variance) measure how non-uniform label (y) values are in leaves

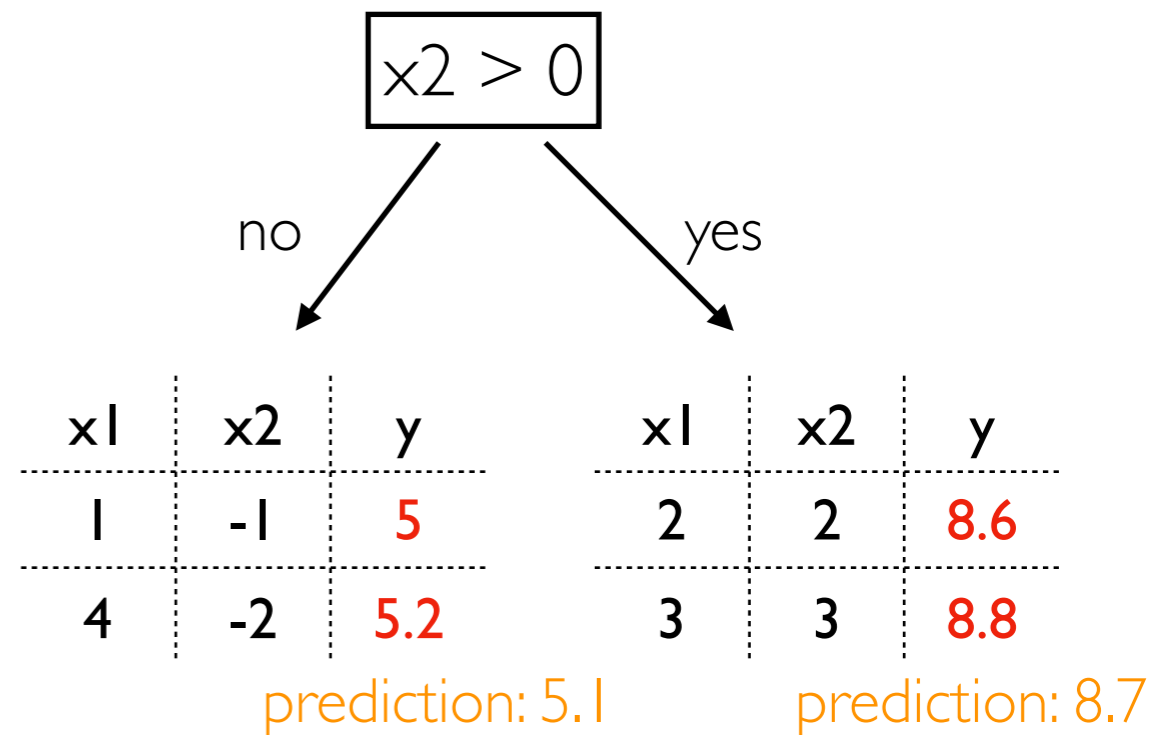
# Predictions

if a new data point lands in a leaf, assume it is similar to other rows in that leaf...

**data:**

x1	x2	y
1	-1	5
2	2	8.6
3	3	8.8
4	-2	5.2
3	50	????

*better tree*



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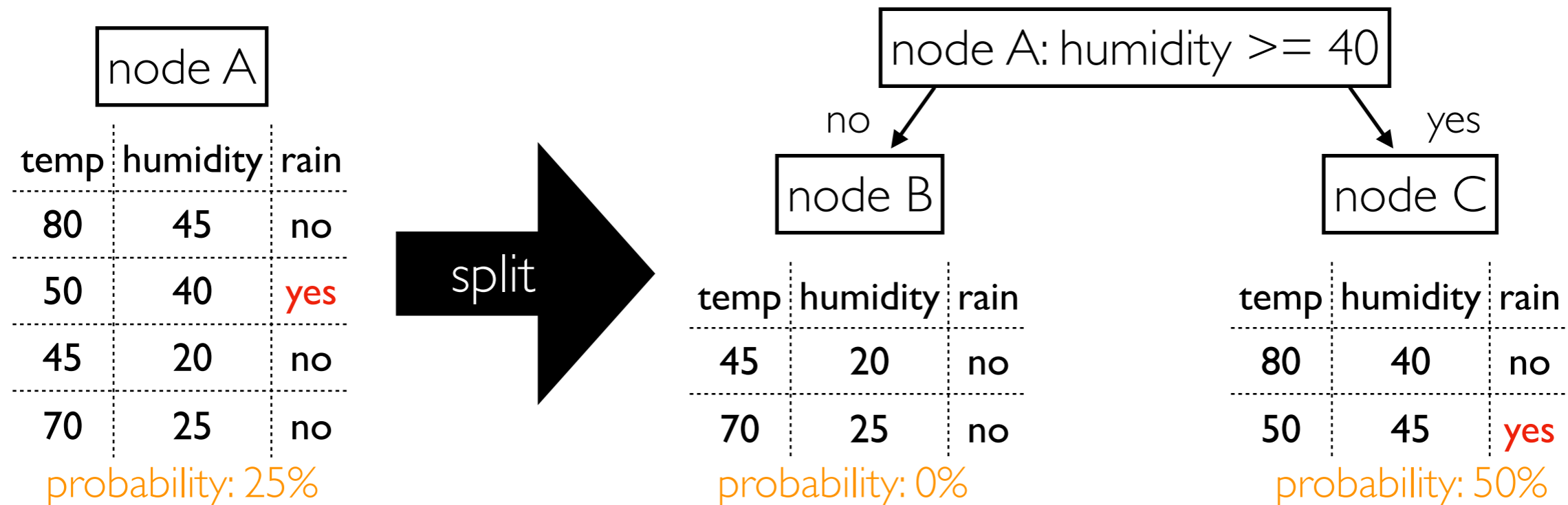
Demos

Decision Trees:

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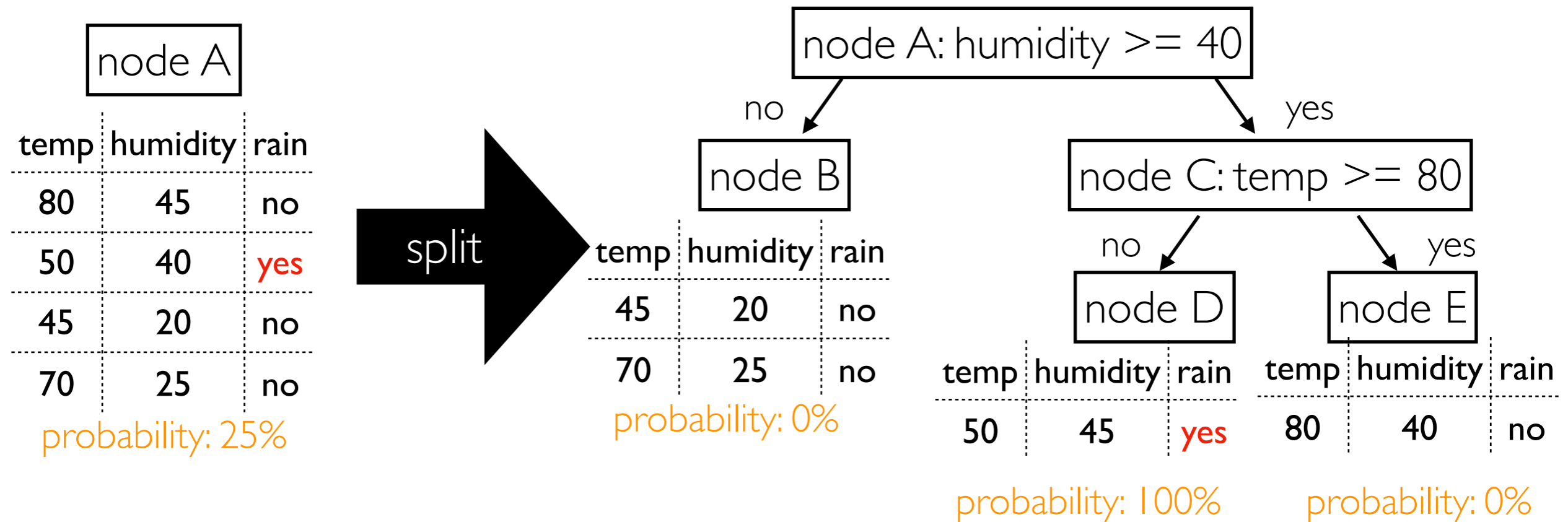
# Splitting Nodes



## Algorithm

- start with one node with all data
- find split point in some column to create two children
- identify another node, recursively split
- eventually stop

# Splitting Nodes

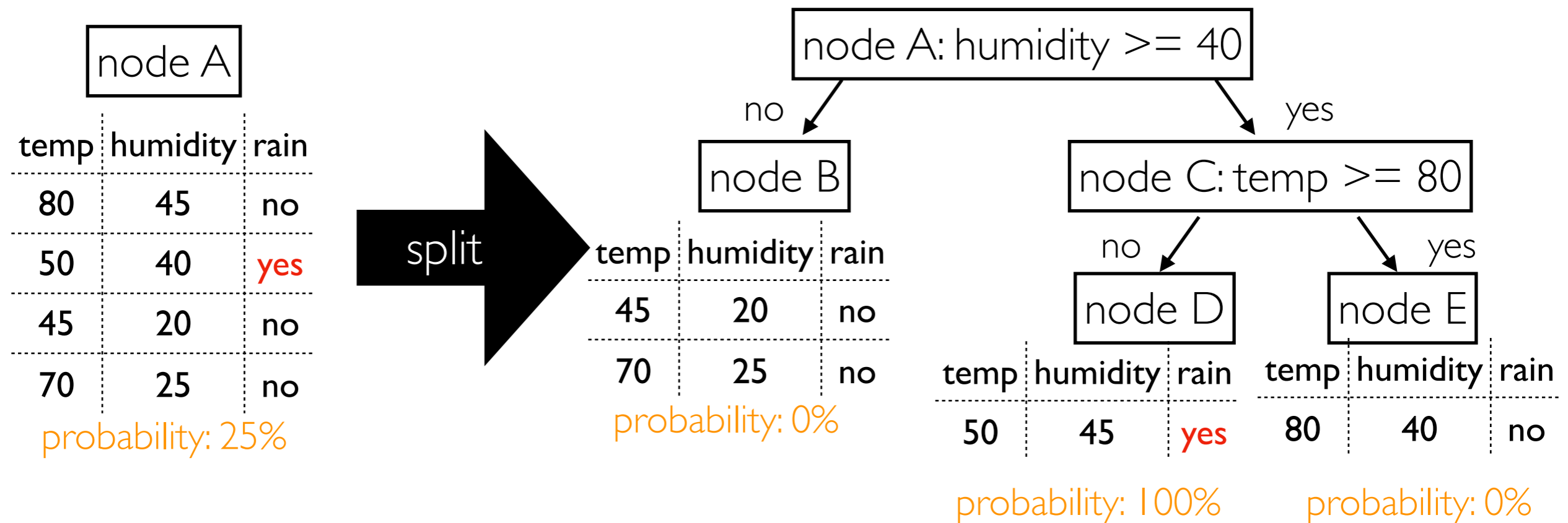


## Algorithm

- start with one node with all data
- find split point in some column to create two children
- identify another node, recursively split
- eventually stop



# When to Stop Splitting?



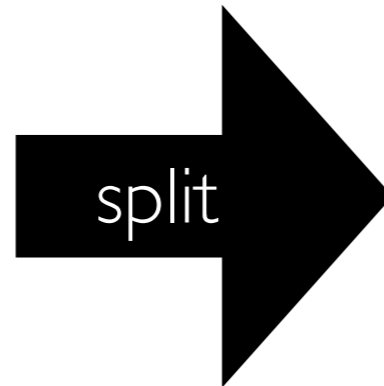
## Some Approaches

- set maximum tree height
- set minimum number of rows in node required for split
- prune tree later to get rid of unhelpful/excessive splitting

# Choosing Splits

node A		
temp	humidity	rain
80	45	no
50	40	yes
45	20	no
70	25	no

probability: 25%



## Which node to split?

- 2 feature columns
- 3 ways to divide 4 rows into big small
- $2*3 = 6$  choices
- try all, choose one that reduces impurity the most!
- how to do so efficiently?

# Choosing Splits

node A

temp	humidity	rain
80	45	no
50	40	yes
45	20	no
70	25	no

sort by each column

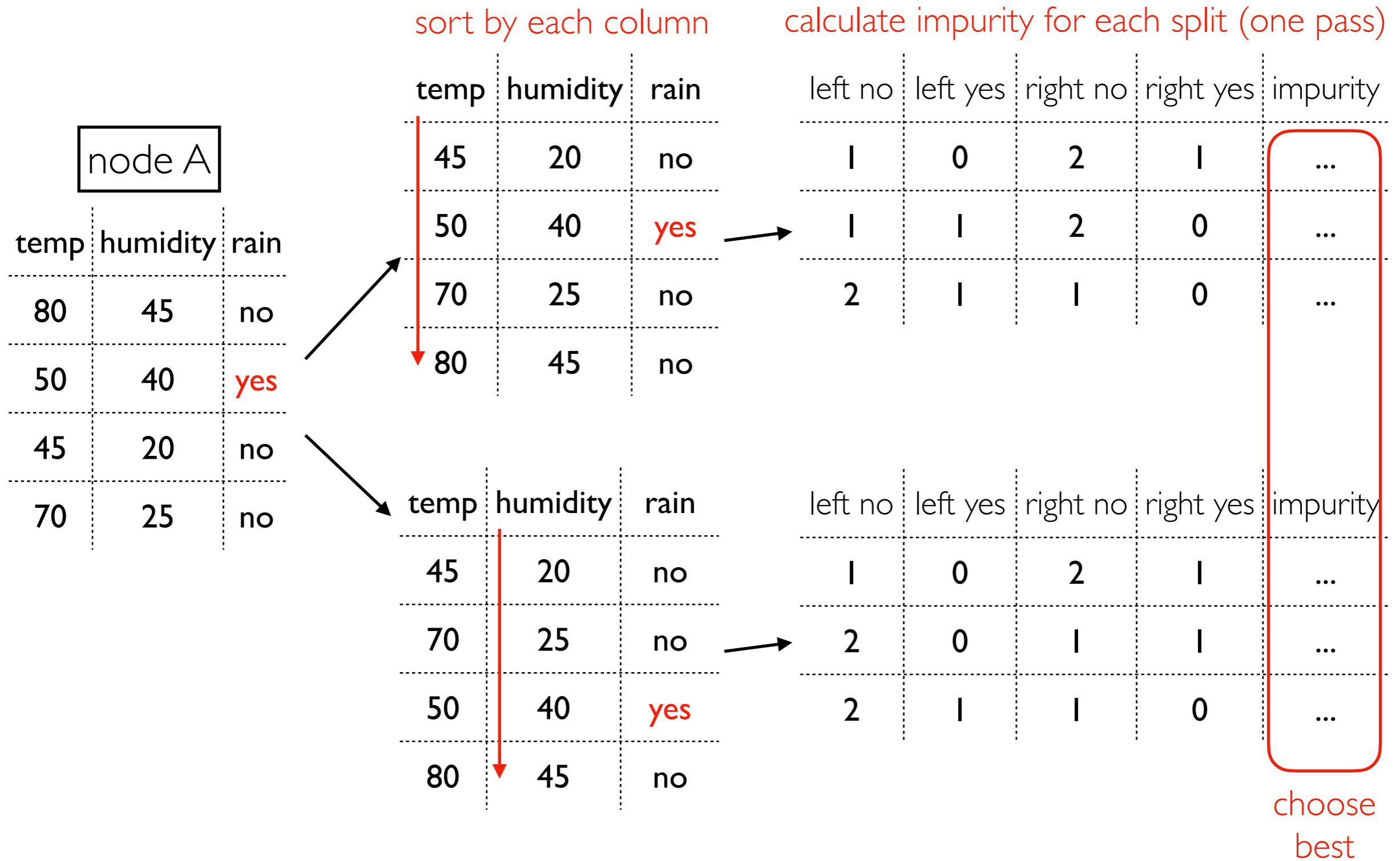
temp	humidity	rain
45	20	no
50	40	yes
70	25	no
80	45	no

calculate impurity for each split (one pass)

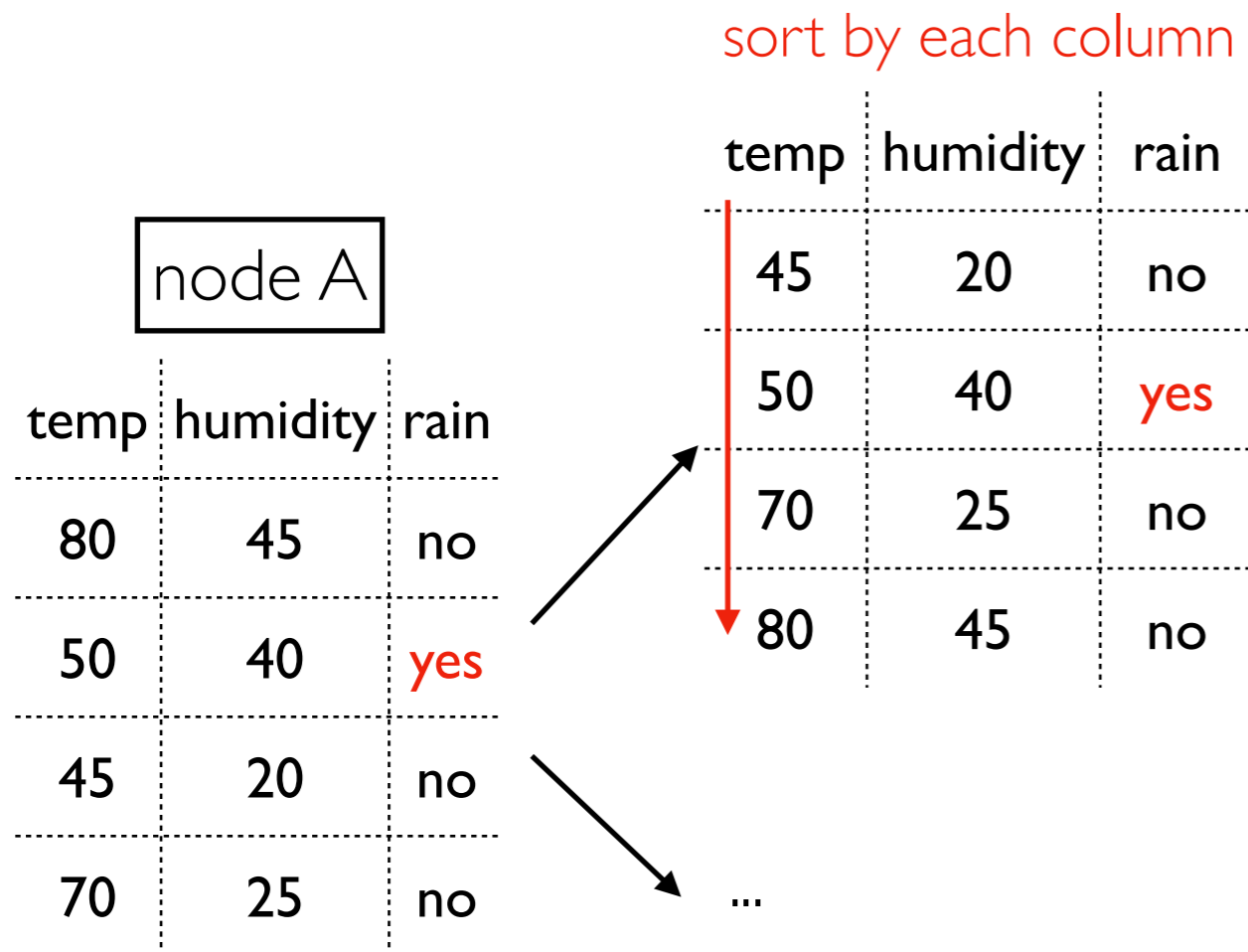
left no	left yes	right no	right yes	impurity
1	0	2	1	...
1	1	2	0	...
2	1	1	0	...

**Observation:** we can incrementally compute impurity for each split point by looking at just more row of data. Don't need to loop over all rows for every possible split point.

# Choosing Splits



# Challenge: Big Data



What if rows for a node are **too big to fit in RAM** on one worker?

- partitioned across many Spark workers
- maybe fits in cumulative RAM of many workers (maybe not)
- each sort would be expensive (network shuffle/exchange)
- as looping over every possible split point, we'll be computing on one worker at any given time (the one that has data around the split point). Not parallel!

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Demos

# PLANET Algorithm

## PLANET: Massively Parallel Learning of Tree Ensembles with MapReduce

Biswanath Panda, Joshua S. Herbach, Sugato Basu, Roberto J. Bayardo  
Google, Inc.

[bpanda, jsherbach, sugato]@google.com, bayardo@alum.mit.edu

### ABSTRACT

Classification and regression tree learning on massive datasets is a common data mining task at Google, yet many state of the art tree learning algorithms require training data to

plexities such as data partitioning, scheduling tasks across many machines, handling machine failures, and performing inter-machine communication. These properties have motivated many technology companies to run MapReduce

<https://static.googleusercontent.com/media/research.google.com/en//pubs/archive/36296.pdf>

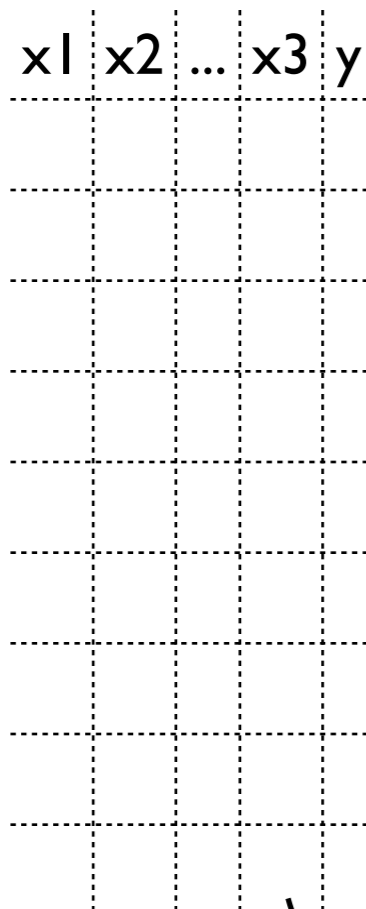
**PLANET**: Parallel Learner for Assembling Numerous Ensemble Trees

- originally implemented as MapReduce jobs
- Spark **DecisionTreeRegressor** and **DecisionTreeClassifier** use it too

Hybrid Approach

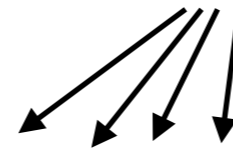
- in-memory splitting for nodes with few enough rows to fit in worker memory
- simplified distributed approach for nodes with lots of data

# Step 1: Compute Equi-Depth Histograms



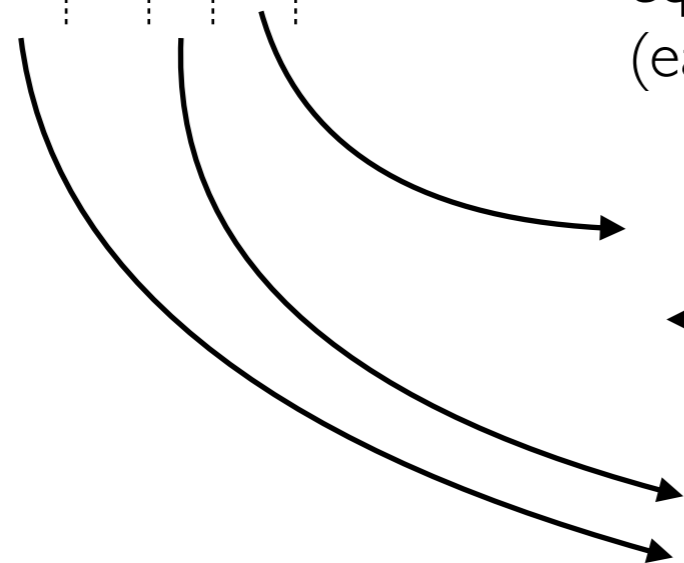
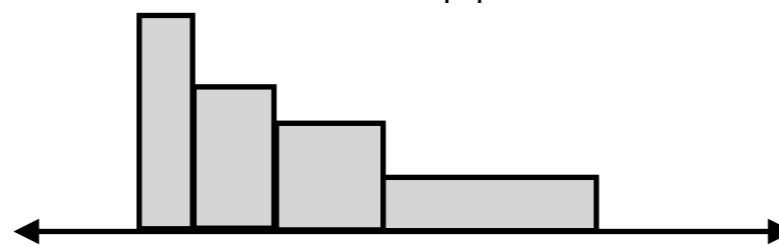
**Warning:** number of bins must be greater than biggest number of unique values in any categorical column

```
dt = DecisionTreeClassifier(labelCol="y")  
dt.setMaxBins(4)
```



**equi-depth histogram for x3 feature**

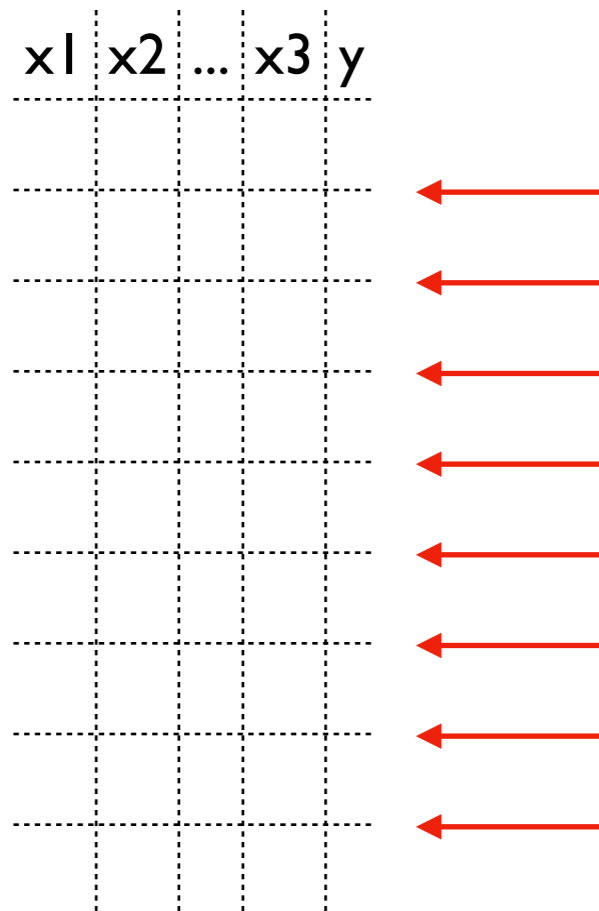
(each bucket has approx same number of samples)



*histograms for every feature...*

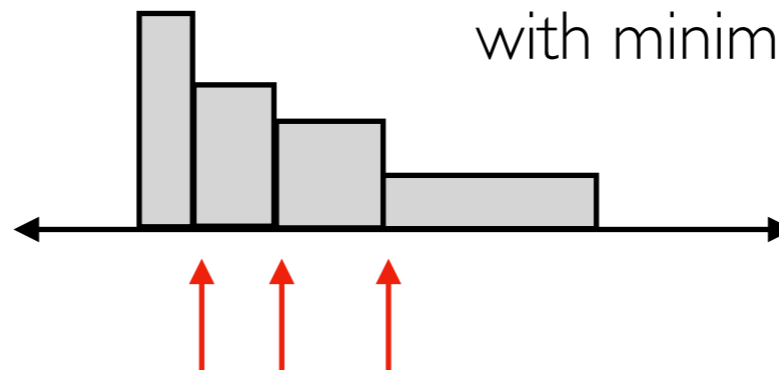


# Split Points: In-Mem vs. Distributed



split points for one-node,  
in-memory algorithm

- splits based on initial data
- every task has split info for every column
- small number of pre-defined split points make it easy for executors to collaborate with minimal coordination

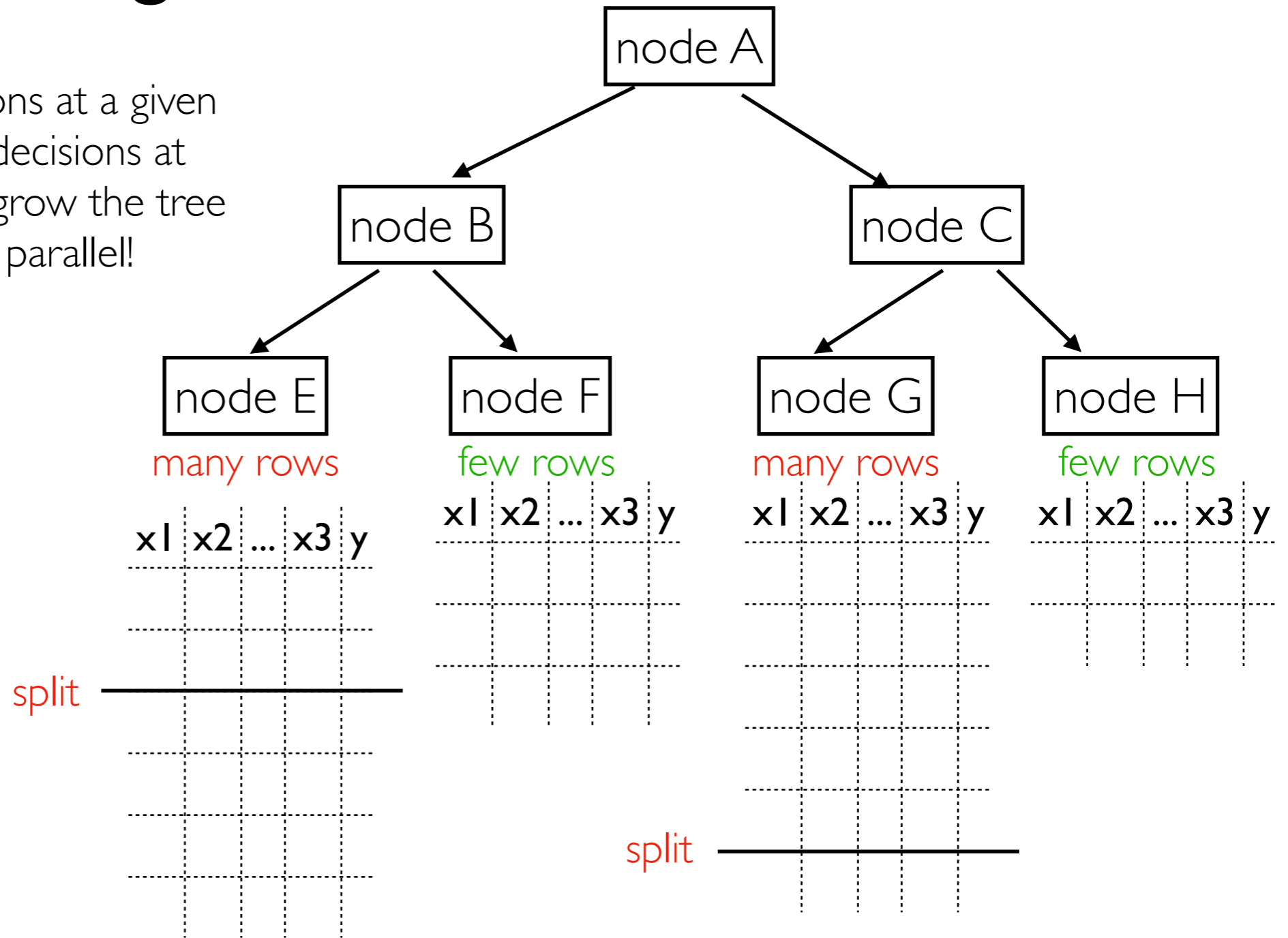


split points for distributed algorithm

# Parallel Splitting

**Observation:** split decisions at a given node are unrelated to decisions at other nodes. It's easy to grow the tree at different places in parallel!

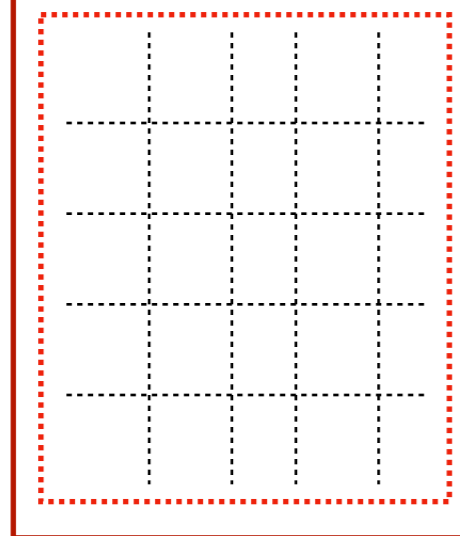
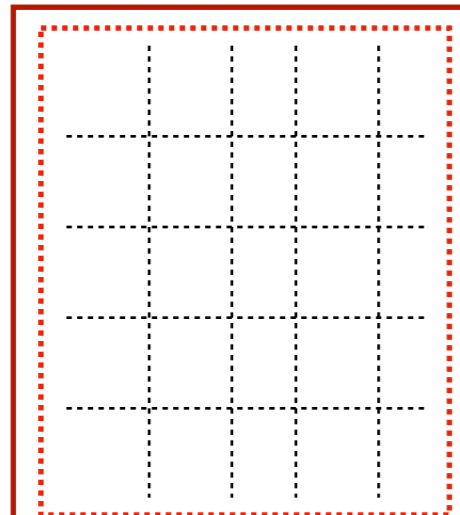
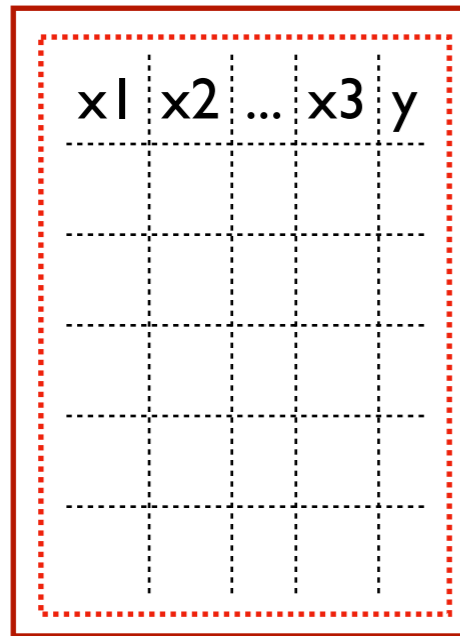
## Decision Tree



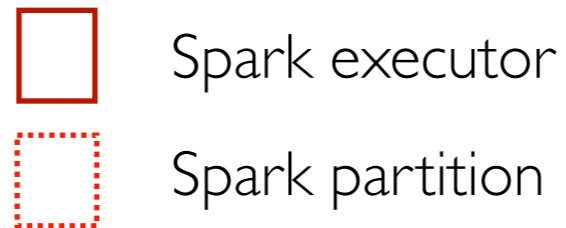
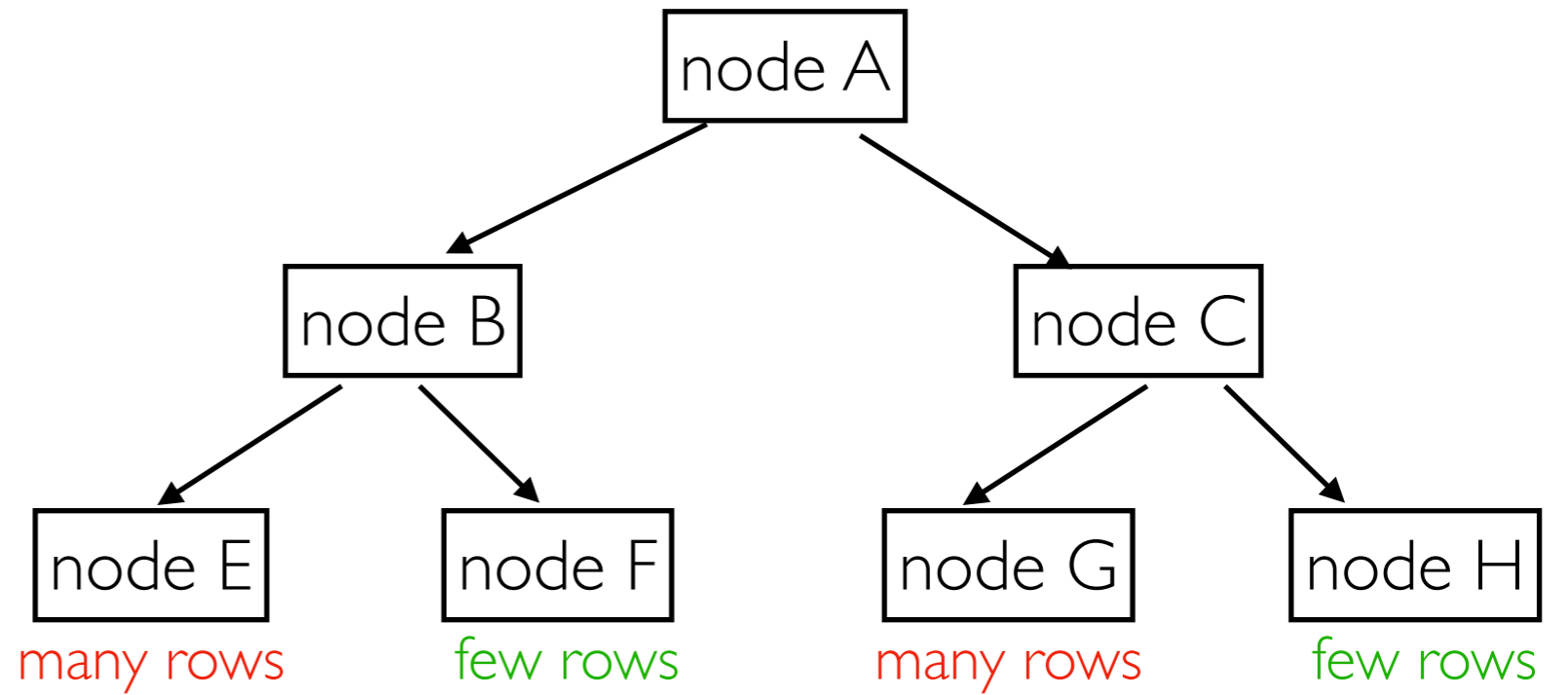
**Clarification:** nodes in the tree data structure DO NOT correspond to nodes in the Spark cluster.

## Logical View of Rows

# Physical Layout

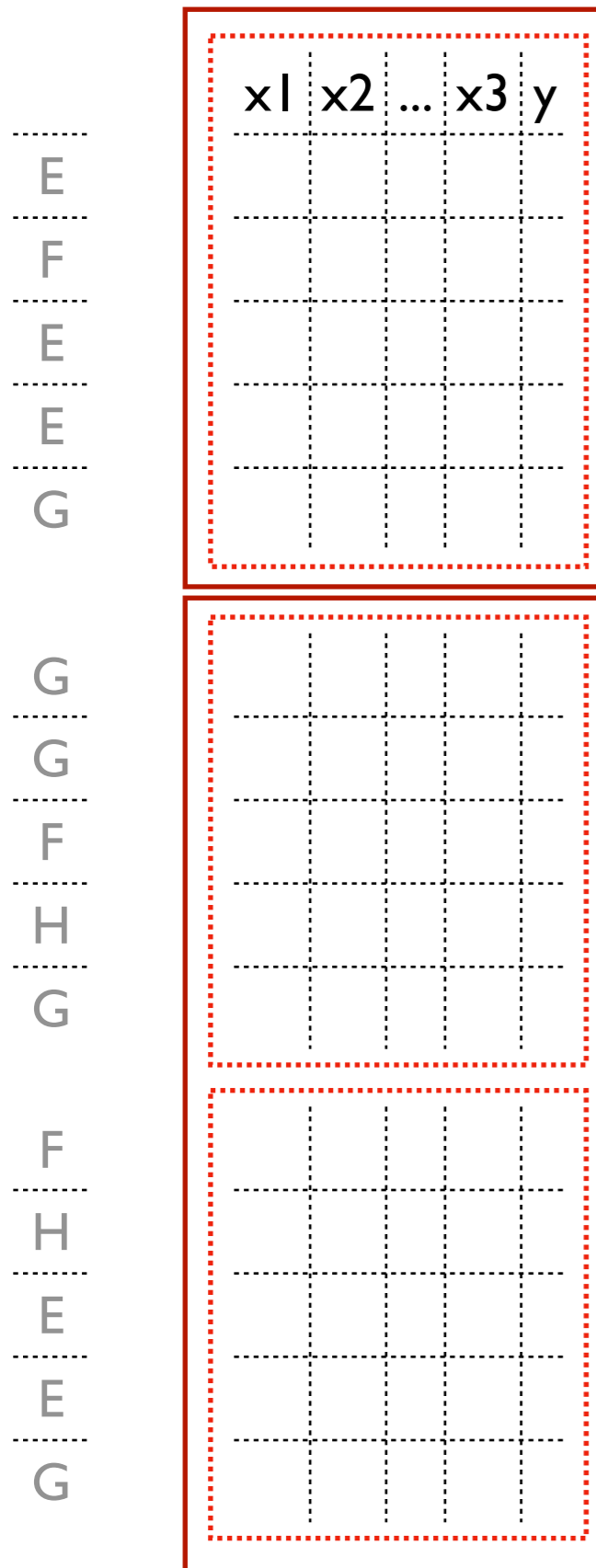


## Decision Tree

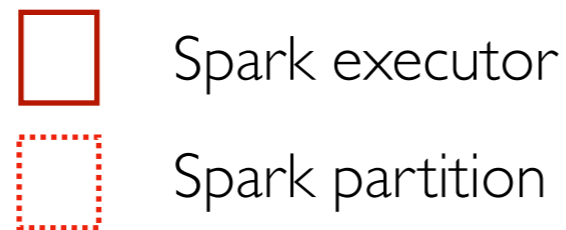
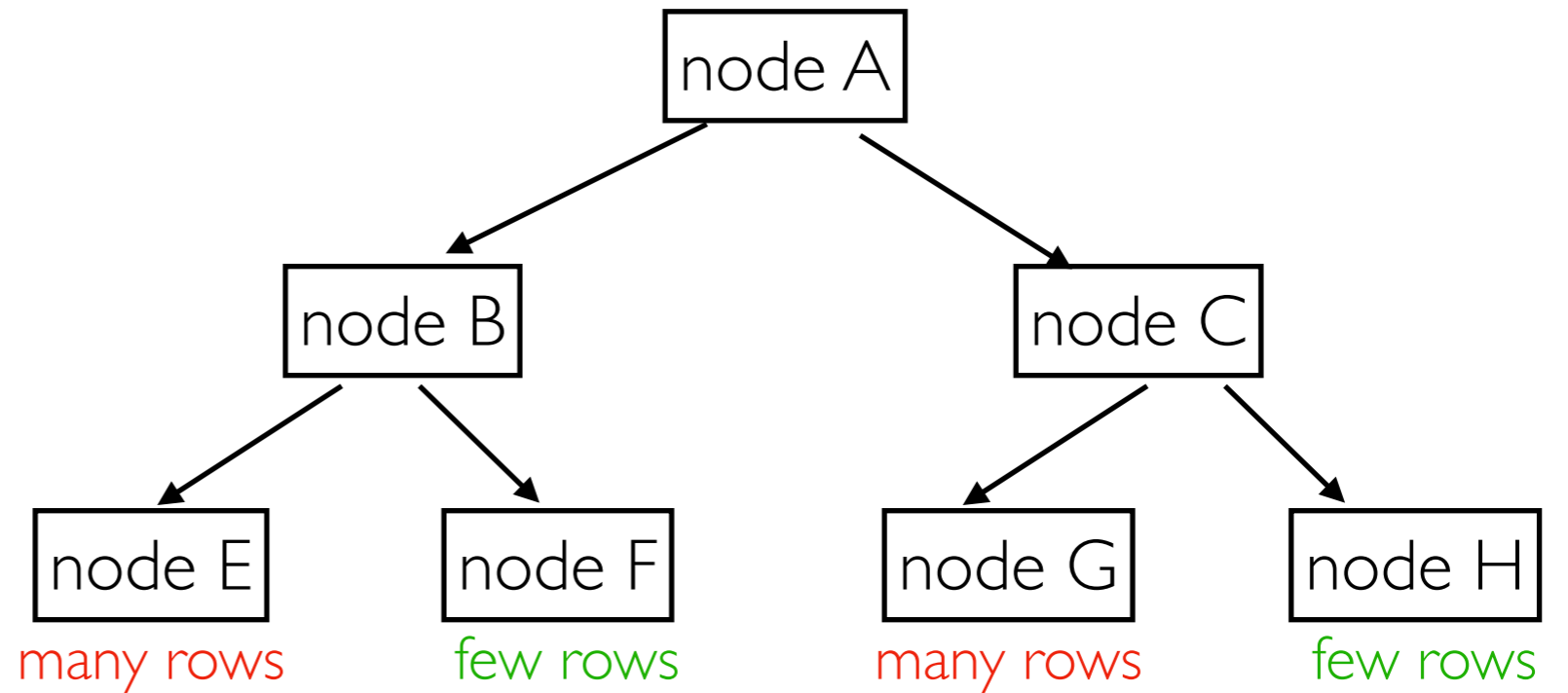


- all rows are in one big Spark DataFrame
- no particular order

# Physical Layout

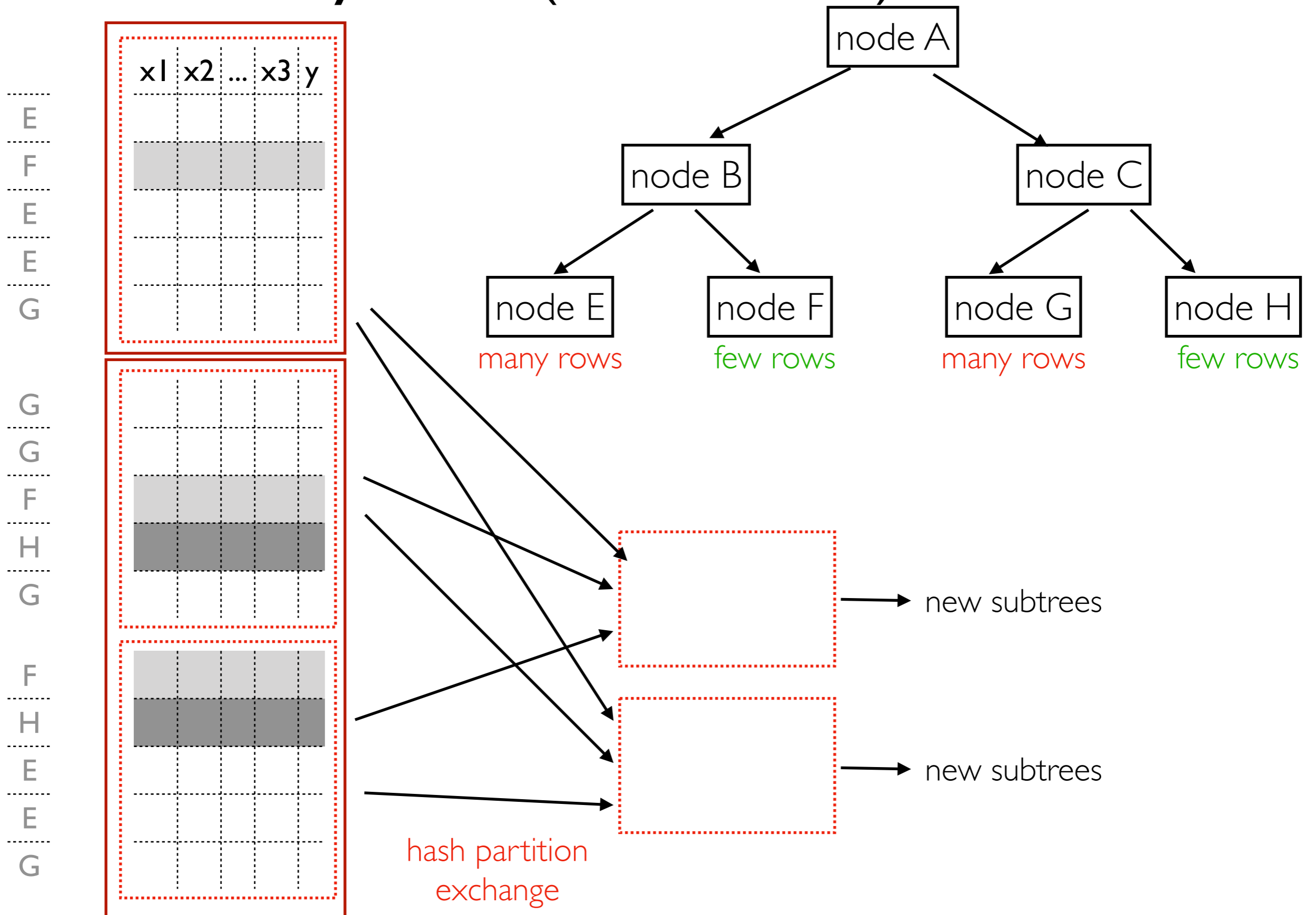


## Decision Tree

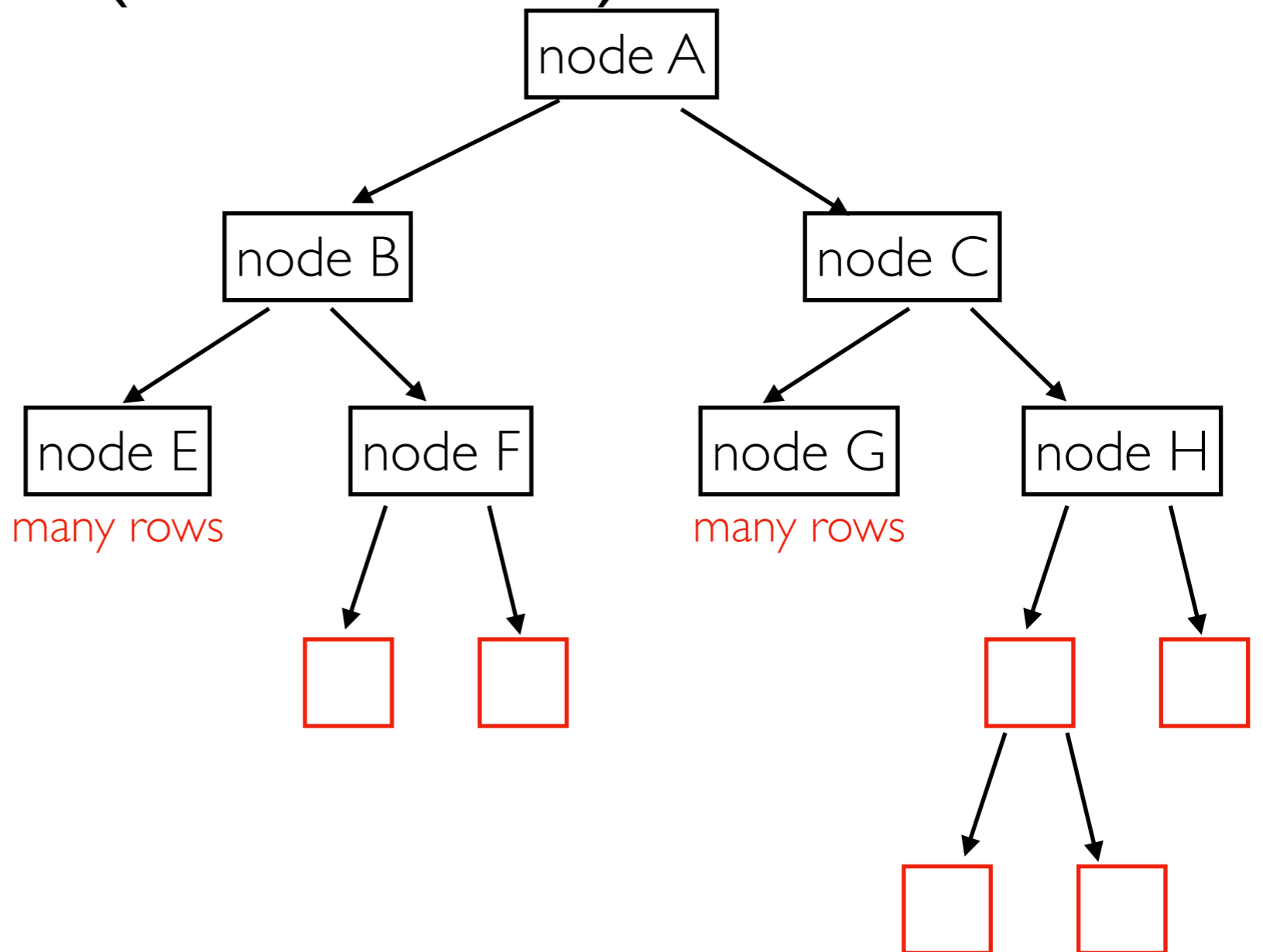
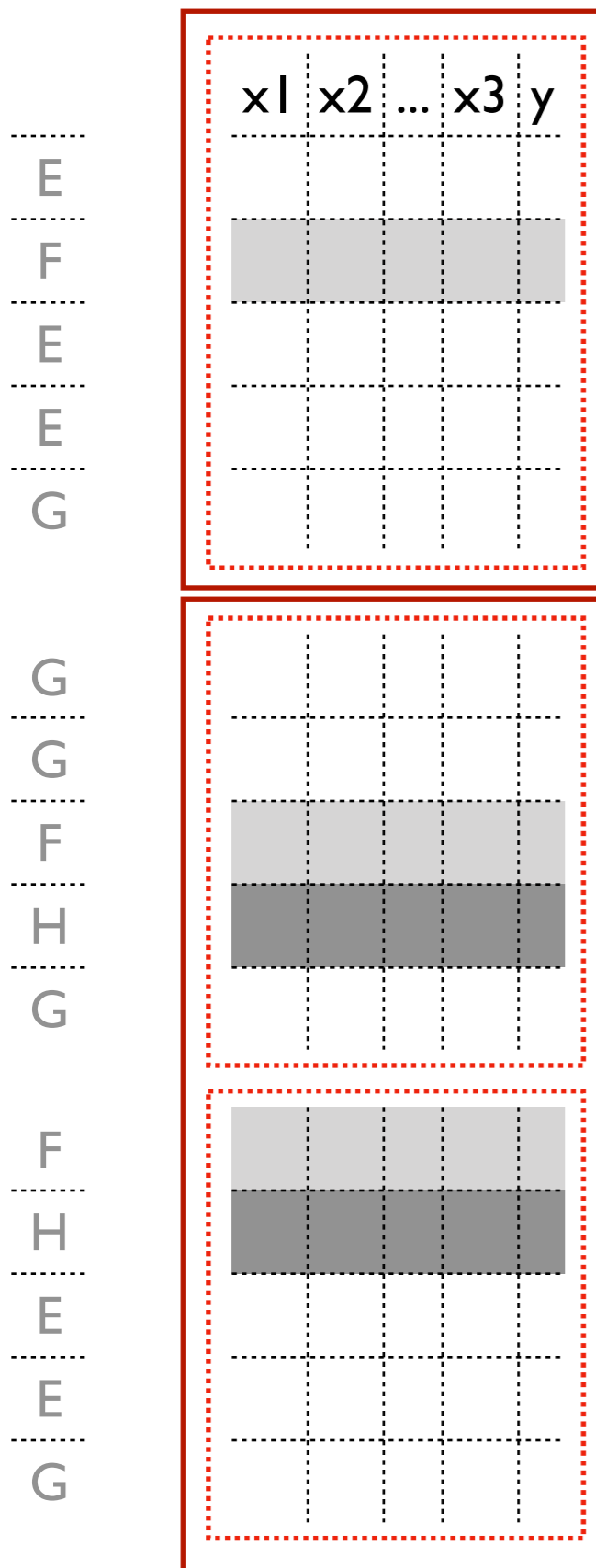


- all rows are in one big Spark DataFrame
- no particular order
- given current tree and  $x_1 \dots x_N$  values, we can infer what leaf node in the tree owns each row

# In Memory Build (small nodes)

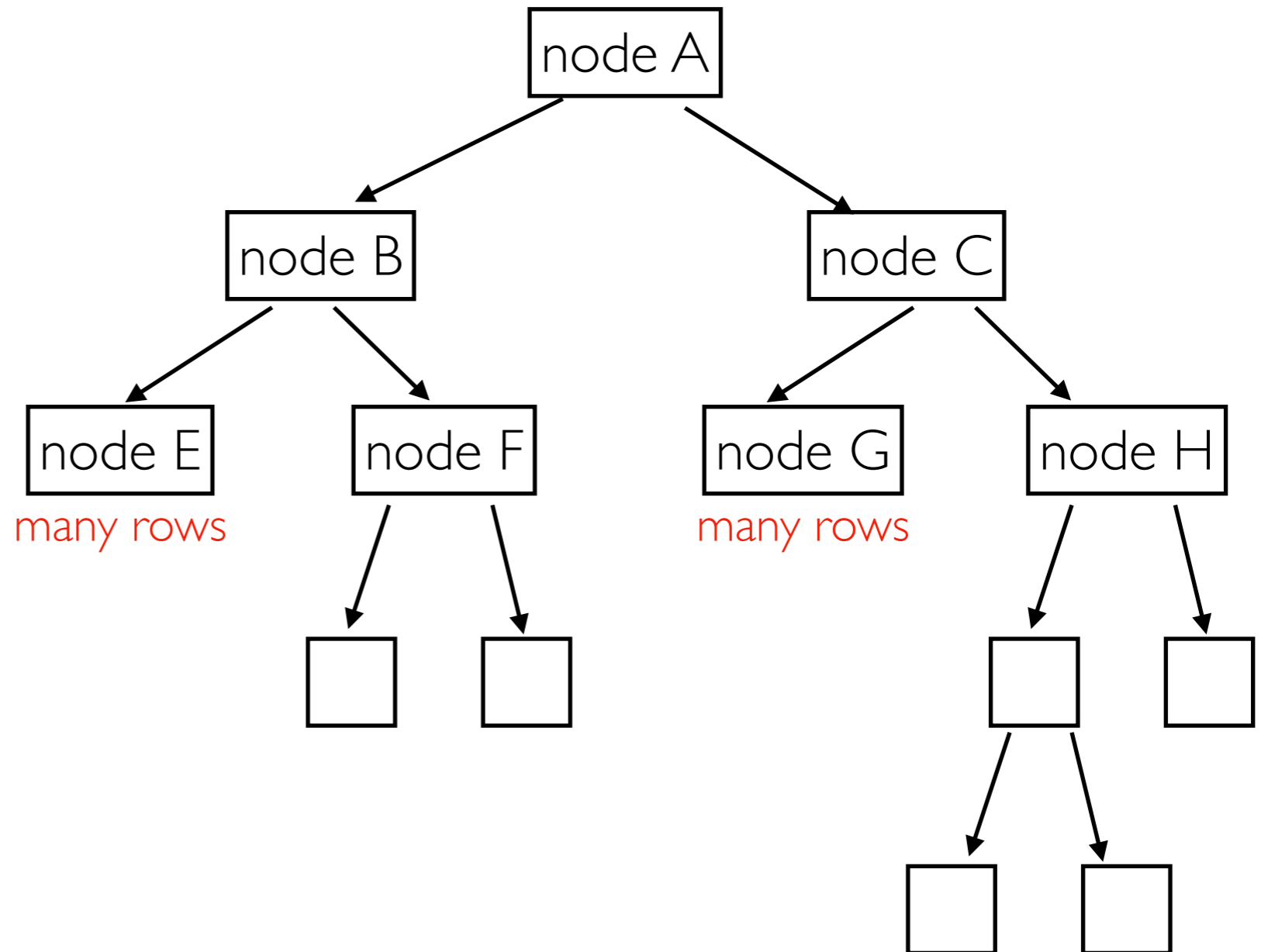
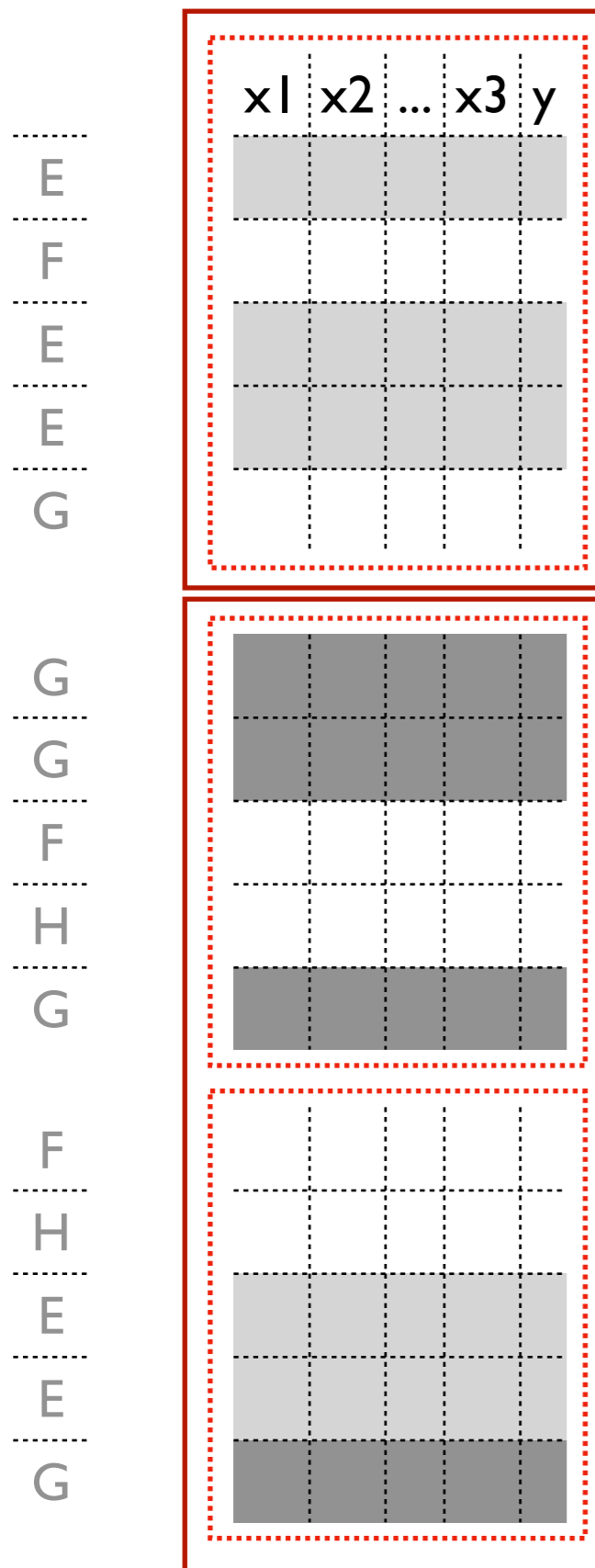


# In Memory Build (small nodes)



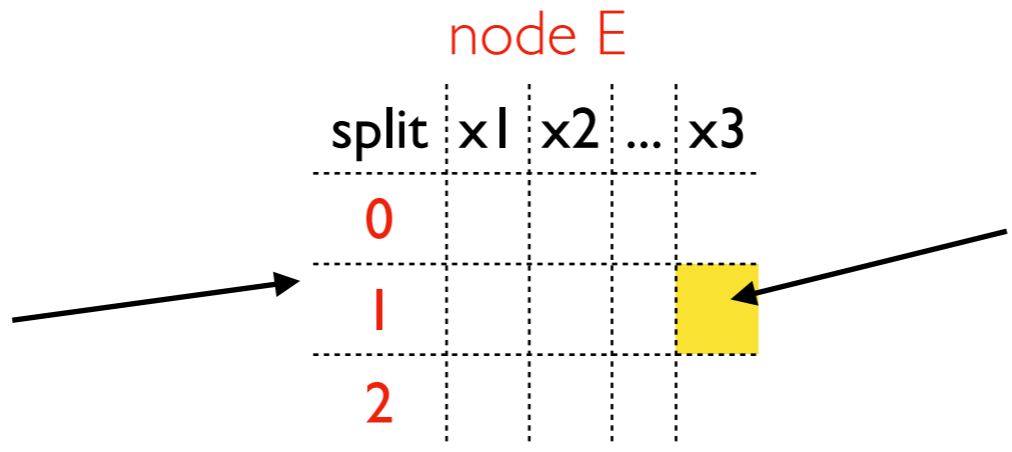
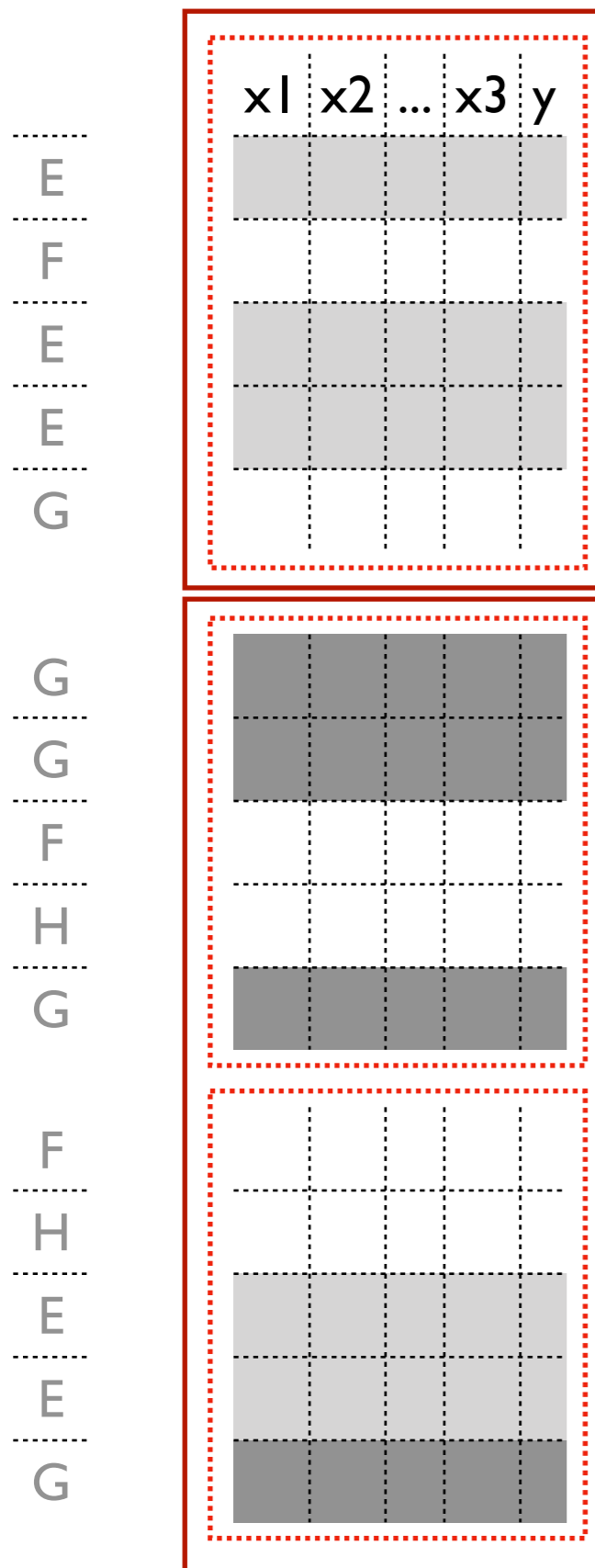
once in memory, splits keep happening recursively, so these nodes are done.

# Big Nodes



- don't move data!
- just output stats per partition for every split/feature option

# Big Nodes



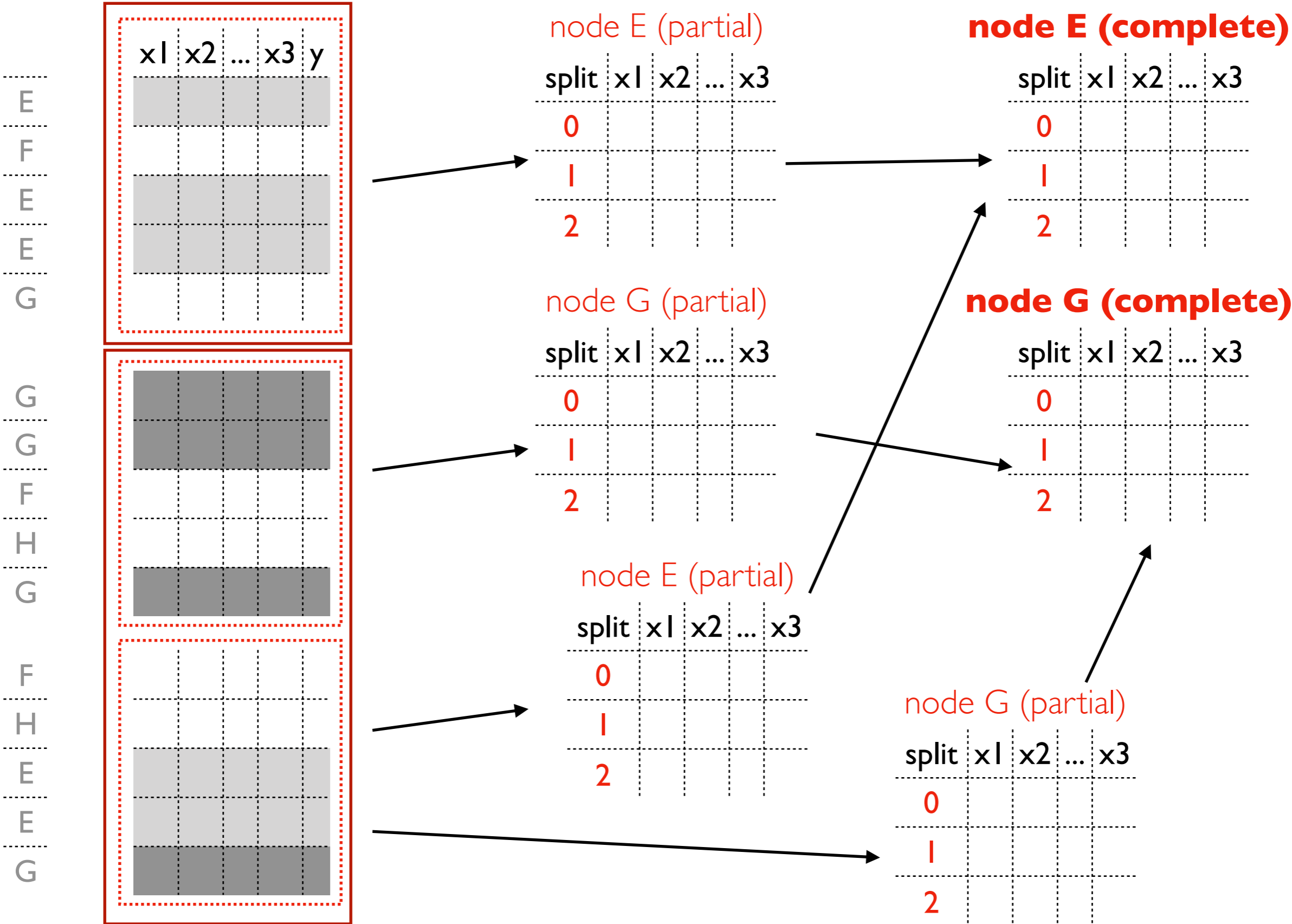
stats per feature/split combo

- left no: *number*
- left yes: *number*
- right no: *number*
- right yes: *number*

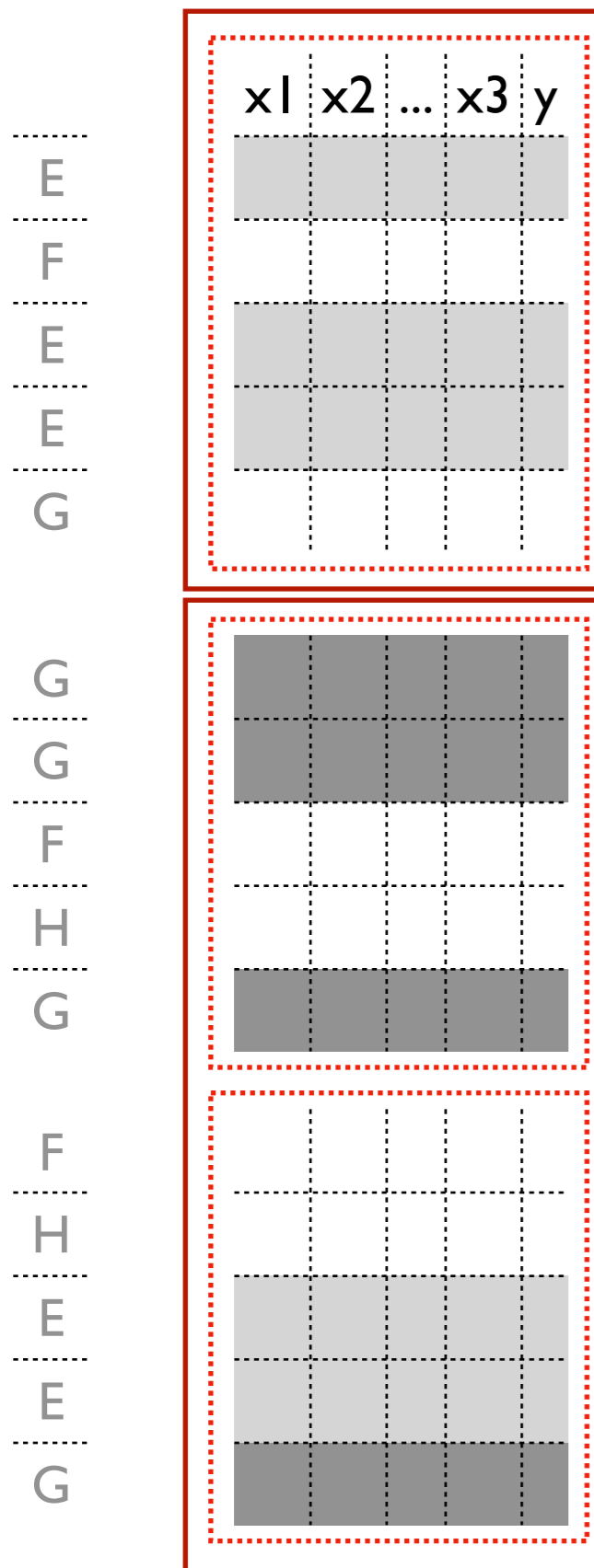
```
dt = DecisionTreeClassifier(labelCol="y")
dt.setMaxBins(3)
```



# Big Nodes



# Big Nodes



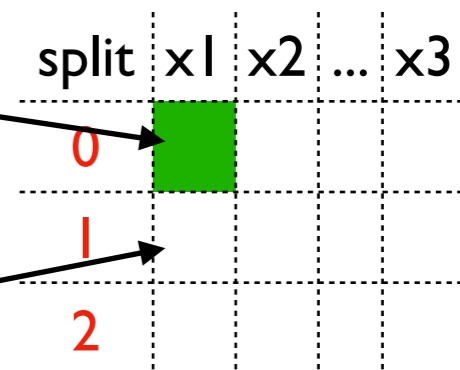
**best:**

left Y/N = 10/20  
right Y/N = 40/10

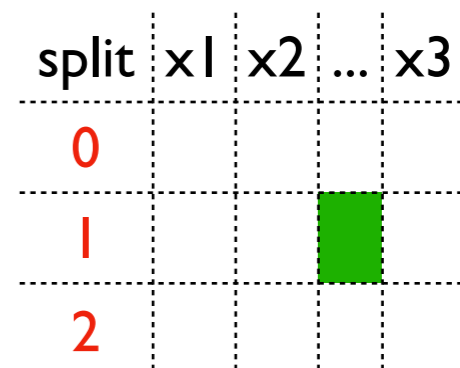
**not as good:**

left Y/N = 25/20  
right Y/N = 25/10

**node E (complete)**

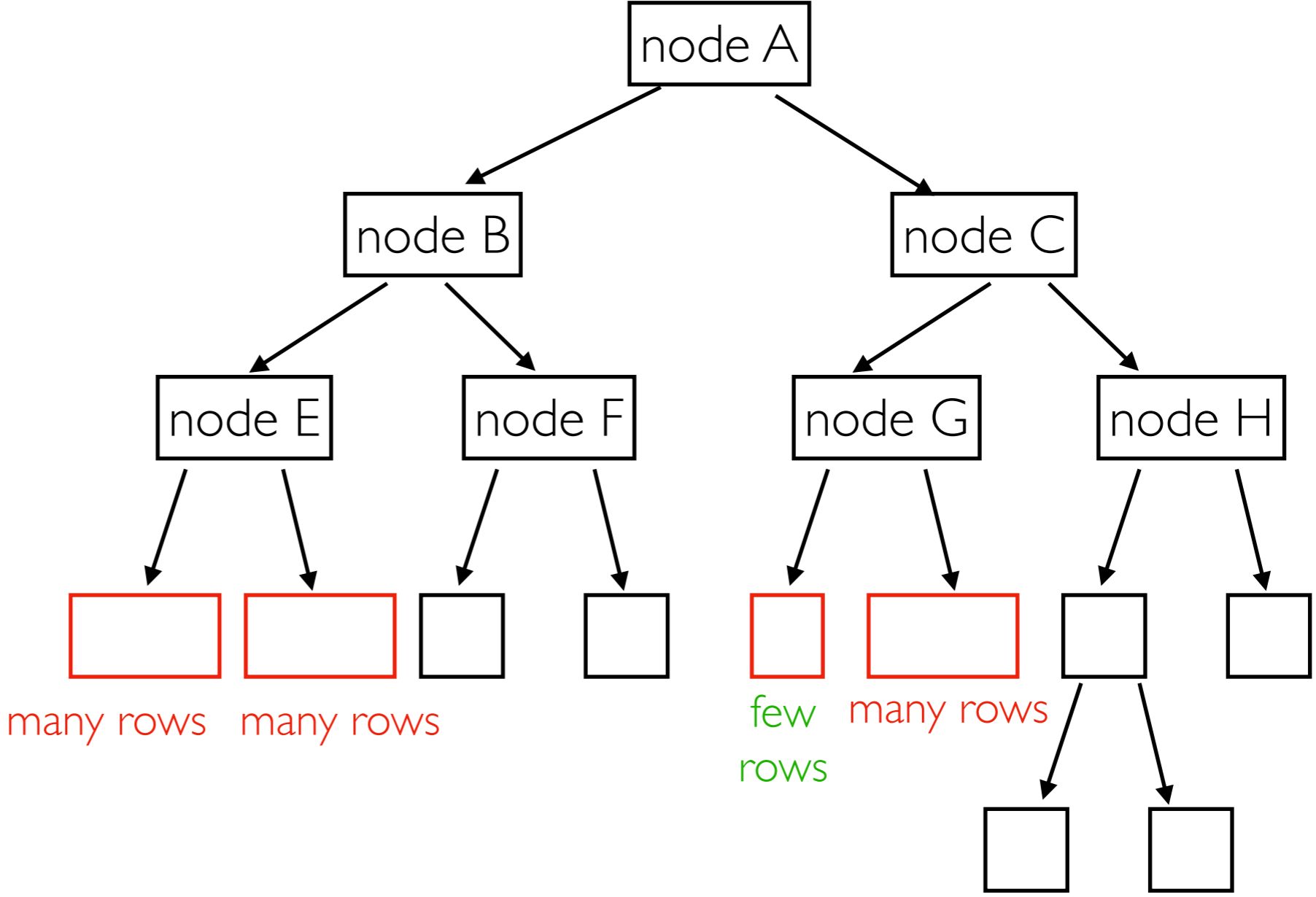
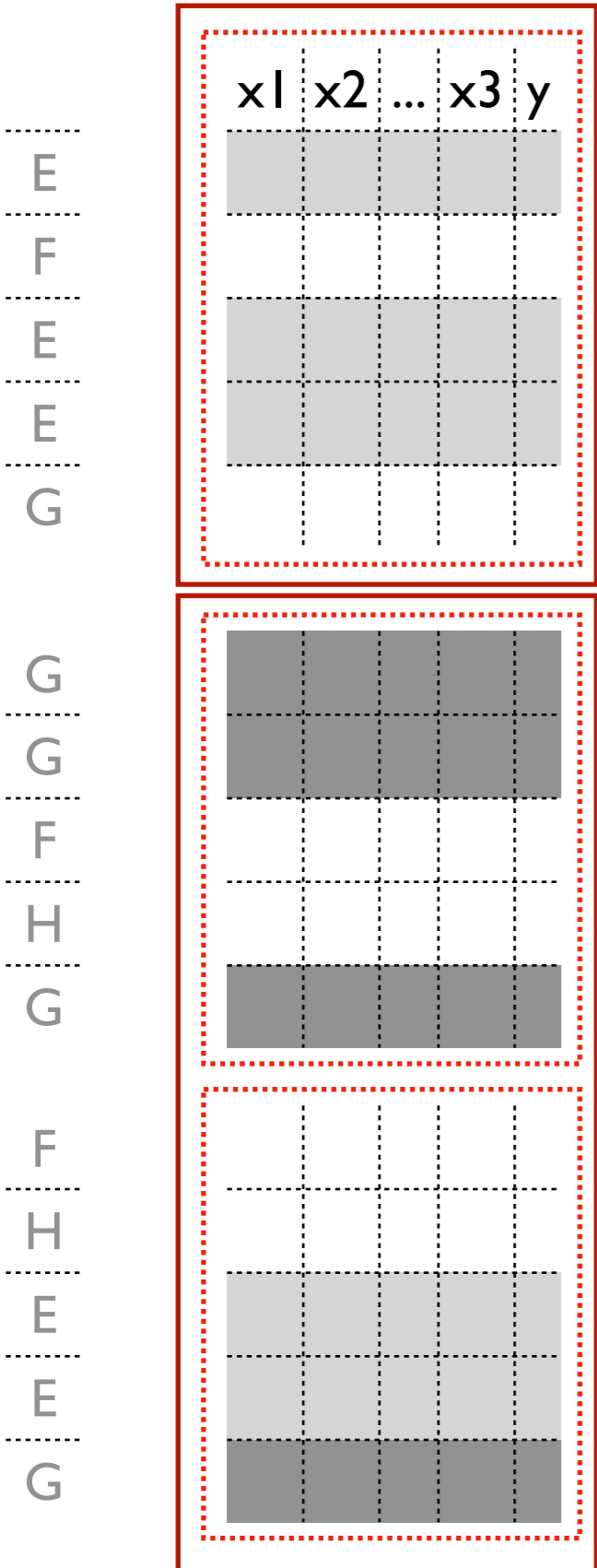


**node G (complete)**



- each **table** corresponds to a **node** we can split (we will choose best split for each node)
- each **column** represents a **feature** we could split on
- each **row** represents a **threshold** we could use for that split

# New Nodes



- we split E and G, creating 4 new nodes
- we DID NOT shuffle rows of data
- we DID shuffle statistics about split choices
- recursively keep splitting (either distributed or in-memory, depending on remaining size)

TopHat

# Distributed ML Outline

ML Review

Training/Predicting APIs

Demos

Decision Trees

Demos