Intro To Spark

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Spark Capabilities
(i.e. Hadoop shortcomings)

• Performance
  • First, use RAM
  • Also, be smarter

• Ease of Use
  • Python, Scala, Java first class citizens

• New Paradigms
  • SparkSQL
  • Streaming
  • MLlib
  • GraphX
  • ...more

But using Hadoop as the backing store is a common and sensible option.
Spark Formula

1. Create/Load RDD
   * Webpage visitor IP address log

2. Transform RDD
   * Filter out all non-U.S. IPs

3. But don’t do anything yet!
   * Wait until data is actually needed
   * Maybe apply more transforms (“Distinct IPs”)

4. Perform Actions that return data
   * Count “How many unique U.S. visitors?”
Simple Example

```
>>> lines_rdd = sc.textFile("nasa_19950801.tsv")
```

Spark Context

The first thing a Spark program requires is a context, which interfaces with some kind of cluster to use. Our pyspark shell provides us with a convenient `sc`, using the local filesystem, to start. Your standalone programs will have to specify one:

```python
from pyspark import SparkConf, SparkContext
conf = SparkConf().setMaster("local").setAppName("Test_App")
sc = SparkContext(conf = conf)
```

You would typically run these scripts like so:

```
spark-submit Test_App.py
```
Simple Example

Read into RDD

Transform

Actions

Lambdas

We’ll see a lot of these. A lambda is simply a function that is too simple to deserve its own subroutine. Anywhere we have a lambda we could also just name a real subroutine that could go off and do anything.

When all you want to do is see if “given an input variable line, is “stanford” in there?”, it isn’t worth the digression.

Most modern languages have adopted this nicety.
# Common Transformations

<table>
<thead>
<tr>
<th>Transformation</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td>map(func)</td>
<td>Return a new RDD by passing each element through <code>func</code>.</td>
</tr>
<tr>
<td>filter(func)</td>
<td>Return a new RDD by selecting the elements for which <code>func</code> returns true.</td>
</tr>
<tr>
<td>flatMap(func)</td>
<td><code>func</code> can return multiple items, and generate a sequence, allowing us to “flatten” nested entries (JSON) into a list.</td>
</tr>
<tr>
<td>distinct()</td>
<td>Return an RDD with only distinct entries.</td>
</tr>
<tr>
<td>sample(...)</td>
<td>Various options to create a subset of the RDD.</td>
</tr>
<tr>
<td>union(RDD)</td>
<td>Return a union of the RDDs.</td>
</tr>
<tr>
<td>intersection(RDD)</td>
<td>Return an intersection of the RDDs.</td>
</tr>
<tr>
<td>subtract(RDD)</td>
<td>Remove argument RDD from other.</td>
</tr>
<tr>
<td>cartesian(RDD)</td>
<td>Cartesian product of the RDDs.</td>
</tr>
<tr>
<td>parallelize(list)</td>
<td>Create an RDD from this (Python) list (using a spark context).</td>
</tr>
</tbody>
</table>

Full list at [http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD](http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD)
<table>
<thead>
<tr>
<th>Action</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td>collect()</td>
<td>Return all the elements from the RDD.</td>
</tr>
<tr>
<td>count()</td>
<td>Number of elements in RDD.</td>
</tr>
<tr>
<td>countByValue()</td>
<td>List of times each value occurs in the RDD.</td>
</tr>
<tr>
<td>reduce(func)</td>
<td>Aggregate the elements of the RDD by providing a function which combines any two into one (sum, min, max, ...).</td>
</tr>
<tr>
<td>first(), take(n)</td>
<td>Return the first, or first n elements.</td>
</tr>
<tr>
<td>top(n)</td>
<td>Return the n highest valued elements of the RDDs.</td>
</tr>
<tr>
<td>takeSample(...)</td>
<td>Various options to return a subset of the RDD.</td>
</tr>
<tr>
<td>saveAsTextFile(path)</td>
<td>Write the elements as a text file.</td>
</tr>
<tr>
<td>foreach(func)</td>
<td>Run the <code>func</code> on each element. Used for side-effects (updating accumulator variables) or interacting with external systems.</td>
</tr>
</tbody>
</table>
• Key/Value organization is a simple, but often very efficient schema, as we mentioned in our NoSQL discussion.

• Spark provides special operations on RDDs that contain key/value pairs. They are similar to the general ones that we have seen.

• On the language (Python, Scala, Java) side key/values are simply tuples. If you have an RDD whose elements happen to be tuples of two items, it is a Pair RDD and you can use the key/value operations that follow.
### Pair RDD Transformations

<table>
<thead>
<tr>
<th>Transformation</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td>reduceByKey(func)</td>
<td>Reduce values using <em>func</em>, but on a key by key basis. That is, combine values with the same key.</td>
</tr>
<tr>
<td>groupByKey()</td>
<td>Combine values with same key. Each key ends up with a list.</td>
</tr>
<tr>
<td>sortByKey()</td>
<td>Return an RDD sorted by key.</td>
</tr>
<tr>
<td>mapValues(func)</td>
<td>Use <em>func</em> to change values, but not key.</td>
</tr>
<tr>
<td>keys()</td>
<td>Return an RDD of only keys.</td>
</tr>
<tr>
<td>values()</td>
<td>Return an RDD of only values.</td>
</tr>
</tbody>
</table>

Note that all of the regular transformations are available as well.
Two Pair RDD Transformations

<table>
<thead>
<tr>
<th>Transformation</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td>subtractByKey(otherRDD)</td>
<td>Remove elements with a key present in other RDD.</td>
</tr>
<tr>
<td>join(otherRDD)</td>
<td>Inner join: Return an RDD containing all pairs of elements with matching keys in self and other. Each pair of elements will be returned as a ((k, (v_1, v_2))) tuple, where ((k, v_1)) is in self and ((k, v_2)) is in other.</td>
</tr>
<tr>
<td>leftOuterJoin(otherRDD)</td>
<td>For each element ((k, v)) in self, the resulting RDD will either contain all pairs ((k, (v, w))) for (w) in other, or the pair ((k, (v, None))) if no elements in other have key (k).</td>
</tr>
<tr>
<td>rightOuterJoin(otherRDD)</td>
<td>For each element ((k, w)) in other, the resulting RDD will either contain all pairs ((k, (v, w))) for (v) in this, or the pair ((k, (None, w))) if no elements in self have key (k).</td>
</tr>
<tr>
<td>cogroup(otherRDD)</td>
<td>Group data from both RDDs by key.</td>
</tr>
</tbody>
</table>

Full list at [http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD](http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD)
>>> x = sc.parallelize([('a', 1), ('b', 4)])
>>> y = sc.parallelize([('a', 2), ('a', 3)])
>>> z = x.join(y)
>>> z.collect()
[('a', (1, 2)), ('a', (1, 3))]

Who needs this? While the above isn't particularly motivating, we will shortly find ourselves in need of just this `join` operation. It pops up repeatedly in data manipulation whenever we want to combine data from two different sources which share keys (the first element).
As with transformations, all of the regular actions are available to Pair RDDs, and there are some additional ones that can take advantage of key/value structure.

<table>
<thead>
<tr>
<th>Action</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td>countByKey()</td>
<td>Count the number of elements for each key.</td>
</tr>
<tr>
<td>lookup(key)</td>
<td>Return all the values for this key.</td>
</tr>
</tbody>
</table>

Full list at [http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD](http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD)
We said one of the advantages of Spark is that we can control things for better performance. There are a multitude of optimization, performance, tuning and programmatic features to enable better control. We quickly look at a few of the most important.

- Persistence
- Partitioning
- Parallel Programming Capabilities
- Performance and Debugging Tools
Persistence

- Lazy evaluation implies by default that all the RDD dependencies will be computed when we call an action on that RDD.

- If we intend to use that data multiple times (say we are filtering some log, then dumping the results, but we will analyze it further) we can tell Spark to persist the data.

- We can specify different levels of persistence: MEMORY_ONLY, MEMORY_ONLY_SER, MEMORY_AND_DISK, MEMORY_AND_DISK_SER, DISK_ONLY

```python
>>> lines_rdd = sc.textFile("nasa_19950801.tsv")
>>> stanfordLines_rdd = lines.filter(lambda line: "stanford" in line)
>>> stanfordLines_rdd.persist(StorageLevel.MEMORY_AND_DISK)
>>> stanfordLines_rdd.count()
47

>>> stanfordLines_rdd.first(1)
[u.glim.stanford.edu\t-t\t807258894\tGET\t/shuttle/.../orbiters-logo.gif\tt200\tt1932\ttt']
...
... 
>>> stanfordLines.unpersist()
```

Do before first action.

Actions

Otherwise will just get evicted when out of memory (which is fine).
Partitions

Spark distributes the data of your RDDs across its resources. It tries to do some obvious things.

- With key/value pairs we can help keep that data grouped efficiently.
- We can create custom partitioners that beat the default (which is probably a hash or maybe range).
- Use persist() if you have partitioned your data in some smart way. Otherwise it will keep getting re-partitioned.
Spark has several parallel programming features that make it easier and more efficient to do operations in parallel in a more explicit way.

Accumulators are variables that allow many copies of a variable to exist on the separate worker nodes.

It is also possible to have replicated data that we would like all the workers to have access to. Perhaps a lookup table of IP addresses to country codes so that each worker can transform or filter on such information. Maybe we want to exclude all non-US IP entries in our logs. You might think of ways you could do this just by passing variables, but they would likely be expensive in actual operation (usually requiring multiple sends). The solution in Spark is to send an (immutable, read only) broadcast variable.

```
log = sc.textFile("logs")
blanks = sc.accumulator(0)
def tokenizeLog(line):
    global blanks  # write-only variable
    if (line =="")
        blanks += 1
    return line.split(" ")

entries = log.flatMap(tokenizeLog)
entries.saveAsTextFile("parsedlogs.txt")
print "Blank entries: %d" blanks.value
```

```
log = sc.textFile("log.txt")
IPtable = sc.broadcast(loadIPTable())
def countryFilter(IPentry, IPtable):
    return (IPentry.prefix() in IPTable)
USentries = log.filter(countryFilter)
```
Performance & Debugging

We will give unfortunately short shrift to performance and debugging, which are both important. Mostly, this is because they are very configuration and application dependent.

Here are a few things to at least be aware of:

• SparkConf() class. A lot of options can be tweaked here.

• Spark Web UI. A very friendly way to explore all of these issues.
Spark has an impressive, and growing, list of input/output formats it supports. Some important ones:

- Text
- CSV
- JSON
- Hadoop Interface
  - Sequence files (key/value)
  - Old and new Hadoop API
  - Compression (gzip...)
  - Database
  - HBase
  - MongoDB
- Protocol Buffers (Google thing)
Spark Streaming

Spark addresses the need for streaming processing of data with a API that divides the data into batches, which are then processed as RDDs.

There are features to enable:

- Fast recovery from failures or timeouts
- Load balancing
- Integration with static data and interactive queries
- Integration with other components (SQL, Machine Learning)

15% of the "global datasphere" (quantification of the amount of data created, captured, and replicated across the world) is currently real-time. That number is growing quickly both in absolute terms and as a percentage.
Shakespeare, a Data Analytics Favorite

Applying data analytics to the works of Shakespeare has become all the rage. Whether determining the legitimacy of his authorship (it wasn’t Marlowe) or if Othello is actually a comedy (perhaps), it is amazing how much publishable research has sprung from the recent analysis of 400 year old text.

We’re going to do some exercises here using a text file containing all of his works.
Who needs this Spark stuff?

As we do our first Spark exercises, you might think of several ways to accomplish these tasks that you already know. For example, Python Pandas is a fine way to do our following problem, and it will probably work on your laptop reasonably well.

However we are learning how to leverage scalable techniques that work on very big data. Shortly, we will encounter problems that are considerable in size, and you will leave this workshop knowing how to harness very large resources.

Searching the Complete Works of William Shakespeare for patterns is a lot different from searching the entire Web (perhaps as the 800TB Common Crawl dataset).

So everywhere you see an RDD, realize that it is a actually a parallel databank that could scale to PBs.
Some Simple Problems

We have an input file, Complete_Shakespeare.txt, that you can also find at http://www.gutenberg.org/ebooks/100.
You might find http://spark.apache.org/docs/latest/api/python/pyspark.html#pyspark.RDD useful to have in a browser window.

If you are starting from scratch on the login node:
1) interact  2) cd BigData/Shakespeare  3) module load spark  4) pyspark
...

>>> rdd = sc.textFile("Complete_Shakespeare.txt")

Let's try a few simple exercises.

1) Count the number of lines

2) Count the number of words (hint: Python "split" is a workhorse)

3) Count unique words

4) Count the occurrence of each word

5) Show the top 5 most frequent words

These last two are a bit more challenging. One approach is to think “key/value”. If you go that way, think about which data should be the key and don’t be afraid to swap it about with value. This is a very common manipulation when dealing with key/value organized data.
Some Simple Answers

```python
>> lines_rdd = sc.textFile("Complete_Shakespeare.txt")
>> lines_rdd.count()
124787
>> words_rdd = lines_rdd.flatMap(lambda x: x.split())
>> words_rdd.count()
904061
>> words_rdd.distinct().count()
67779
```

Next, I know I'd like to end up with a pair RDD of sorted word/count pairs:

```
(23407, 'the'), (19540,'I'), (15682, 'to'), (15649, 'of')...
```

Why not just `words_rdd.countByValue()`? We get back a massive Python unsorted dictionary of results:

```
... 1, u'precious-princely': 1, u'christenings?': 1, 'empire': 11, u'vaunts': 2, u'Lubber's': 1, u'poet.': 2, u'Toad!': 1, u'leaden': 15, u'captains': 1, u'leaf': 9, u'Barnes,': 1, u'lead': 101, u'"Hell": 1, u'wheat,': 3, u'lean': 28, u'Toad,' 1, u'trencher!': 2, u'1.F.2.': 1, u'leas': 2, u'leap': 17, ...
```

Where to go next? Sort this in Python or try to get back into an RDD? If this is truly BIG data, we want to remain as an RDD until we reach our final results.
Some Harder Answers

Turn these into k/v pairs

Reduce to get words counts

Flip keys and values so we can sort on wordcount instead of words.

$\text{results}_\text{rdd} = \text{lines}_\text{rdd}.\text{flatMap}(\lambda x: x.\text{split()}).\text{map}(\lambda x: (x,1)).\text{reduceByKey}(\lambda x, y: x+y).\text{map}(\lambda x: (x[1], x[0])).\text{sortByKey}(\text{False})$
To do research-level text analysis, we generally want to clean up our input. Here are some of the kinds of things you could do to get a more meaningful distinct word count.

1) Remove punctuation. Often punctuation is just noise, and it is here. Do a Map and/or Filter (some punctuation is attached to words, and some is not) to eliminate all punctuation from our Shakespeare data. Note that if you are familiar with regular expressions, Python has a ready method to use those.

2) Remove stop words. Stop words are common words that are also often uninteresting ("I", "the", "a"). You can remove many obvious stop words with a list of your own, and the MLlib that we are about to investigate has a convenient StopWordsRemover() method with default lists for various languages.

3) Stemming. Recognizing that various different words share the same root ("run", "running") is important, but not so easy to do simply. Once again, Spark brings powerful libraries into the mix to help. A popular one is the Natural Language Tool Kit. You should look at the docs, but you can give it a quick test quite easily:

```python
import nltk
from nltk.stem.porter import *
stemmer = PorterStemmer()
stems_rdd = words_rdd.map(lambda x: stemmer.stem(x))
```
MLib rolls in a lot of classic machine learning algorithms. We barely have time to touch upon this interesting topic today, but they include:

- Useful data types
- Basic Statistics
- Classification (including SVMs, Random Forests)
- Regression
- Algorithms (SGD,...)
- Clustering...
One of the reasons we use Spark is for easy access to powerful data analysis tools. The MLlib library gives us a machine learning library that is easy to use and utilizes the scalability of the Spark system.

It has supported APIs for Python (with NumPy), R, Java and Scala.

We will use the Python version in a generic manner that looks very similar to any of the above implementations.

There are good example documents for the clustering routine we are using here:

http://spark.apache.org/docs/latest/mllib-clustering.html

And an excellent API reference document here:

http://spark.apache.org/docs/latest/api/python/pyspark.mllib.html#pyspark.mllib.clustering.KMeans

I suggest you use these pages for all your Spark work.
Clustering is a very common operation for finding grouping in data and has countless applications. This is a very simple example, but you will find yourself reaching for a clustering algorithm frequently in pursuing many diverse machine learning objectives, sometimes as one part of a pipeline.
As intuitive as clustering is, it presents challenges to implement in an efficient and robust manner.

You might think this is trivial to implement in lower dimensional spaces.

But it can get tricky even there.

Sometimes you know how many clusters you have to start with. Often you don’t. How hard can it be to count clusters? How many are here?

We will start with 5000 2D points. We want to figure out how many clusters there are, and their centers. Let’s fire up pyspark and get to it...
Finding Clusters

*RDD map() takes a function to apply to the elements. We can certainly create our own separate function, but lambdas are a way many languages allow us to define trivial functions “in place”.

Using Python version 2.7.5
SparkContext available as sc
HiveContext available as sqlContext.

```python
>>> rdd1 = sc.textFile("5000_points.txt")
```

```python
>>> rdd2 = rdd1.map(lambda x: x.split())
```

```python
>>> rdd3 = rdd2.map(lambda x: [int(x[0]), int(x[1])])
```

```python
r288% module load spark
r288% pyspark
```
Finding Our Way

```python
>>> rdd1 = sc.textFile("5000_points.txt")
... rdd1.count()
5000
>>> rdd1.take(4)
[u' 664159  550946', u' 665845  557965', u' 597173  575538', u' 618600  551446']
>>> rdd2 = rdd1.map(lambda x:x.split())
>>> rdd2.take(4)
[[u'664159', u'550946'], [u'665845', u'557965'], [u'597173', u'575538'], [u'618600', u'551446']]
>>> rdd3 = rdd2.map(lambda x: [int(x[0]), int(x[1])])
>>> rdd3.take(4)
[[664159, 550946], [665845, 557965], [597173, 575538], [618600, 551446]]
```
Finding Clusters

Using Python version 2.7.5 (default, Nov 20 2015 02:00:19)
SparkContext available as sc, HiveContext available as sqlContext.

```python
>>> rdd1 = sc.textFile("5000_points.txt")
>>> rdd2 = rdd1.map(lambda x:x.split())
>>> rdd3 = rdd2.map(lambda x: [int(x[0]), int(x[1])])

>>> from pyspark.mllib.clustering import KMeans

class KMeans

Parameters:
  - rdd: Training points as an RDD of Vector or convertible sequence types.
  - k: Number of clusters to create.
  - maxIterations: Maximum number of iterations allowed. (default: 100)
  - runs: This parameter has no effect since Spark 2.0.0.
  - initializationMode: The initialization algorithm. This can be either "random" or "k-means||" (default: "k-means||")
  - seed: Random seed value for cluster initialization. Set as None to generate seed based on system time. (default: None)
  - initializationSteps: Number of steps for the k-means|| initialization mode. This is an advanced setting - the default of 5 is almost always enough. (default: 5)
  - epsilon: Distance threshold within which a center will be considered to have converged. If all centers move less than this Euclidian distance, iterations are stopped. (default: 1e-4)
  - initModel: Initial cluster centers can be provided as a KMeansModel object rather than using the random or k-means|| initializationModel. (default: None)
```
Finding Clusters

Let's see results for 1-30 cluster tries

<table>
<thead>
<tr>
<th>Clusters</th>
<th>Cost</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>5.76807041184e+14</td>
</tr>
<tr>
<td>2</td>
<td>3.43183673951e+14</td>
</tr>
<tr>
<td>3</td>
<td>2.23097486536e+14</td>
</tr>
<tr>
<td>4</td>
<td>1.64792608443e+14</td>
</tr>
<tr>
<td>5</td>
<td>1.19410028576e+14</td>
</tr>
<tr>
<td>6</td>
<td>7.97690150116e+13</td>
</tr>
<tr>
<td>7</td>
<td>7.16451594344e+13</td>
</tr>
<tr>
<td>8</td>
<td>4.81469246295e+13</td>
</tr>
<tr>
<td>9</td>
<td>4.23762700793e+13</td>
</tr>
<tr>
<td>10</td>
<td>3.65230706654e+13</td>
</tr>
<tr>
<td>11</td>
<td>3.16991867996e+13</td>
</tr>
<tr>
<td>12</td>
<td>2.94369408304e+13</td>
</tr>
<tr>
<td>13</td>
<td>2.04031903147e+13</td>
</tr>
<tr>
<td>14</td>
<td>1.37018893034e+13</td>
</tr>
<tr>
<td>15</td>
<td>5.91761561687e+12</td>
</tr>
<tr>
<td>16</td>
<td>1.3183365206e+13</td>
</tr>
<tr>
<td>17</td>
<td>1.39010717893e+13</td>
</tr>
<tr>
<td>18</td>
<td>8.22806178508e+12</td>
</tr>
<tr>
<td>19</td>
<td>8.22513516563e+12</td>
</tr>
<tr>
<td>20</td>
<td>7.79359299283e+12</td>
</tr>
<tr>
<td>21</td>
<td>7.79615059172e+12</td>
</tr>
<tr>
<td>22</td>
<td>7.70001662709e+12</td>
</tr>
<tr>
<td>23</td>
<td>7.24231610447e+12</td>
</tr>
<tr>
<td>24</td>
<td>7.21990743993e+12</td>
</tr>
<tr>
<td>25</td>
<td>7.0935133944e+12</td>
</tr>
<tr>
<td>26</td>
<td>6.92577879424e+12</td>
</tr>
<tr>
<td>27</td>
<td>6.53939015776e+12</td>
</tr>
<tr>
<td>28</td>
<td>6.37782690833e+12</td>
</tr>
<tr>
<td>29</td>
<td>6.37192522244e+12</td>
</tr>
</tbody>
</table>
for trials in range(10):
   ...
   print
   ...
   for clusters in range(12,18):
      ...
      model = KMeans.train(rdd3,clusters)
      ...
      print clusters, model.computeCost(rdd3)
for trials in range(10):
    for clusters in range(12, 16):
        model = KMeans.train(rdd3, clusters)
        cost = model.computeCost(rdd3)
        centers = model.clusterCenters
        if cost<1e+13:
            print clusters, cost
            for coords in centers:
                print int(coords[0]), int(coords[1])
            break

#Try ten times to find best result
#Only look in interesting range

#Let's grab cluster centers
#If result is good, print it out

Find the Centers

15 8.91761561687e+12
852058 157685
606574 574455
320602 161521
139395 558143
858947 546259
337264 562123
244654 847642
398870 404924
670929 862765
823421 731145
507818 175610
801616 321123
617926 399415
417799 787001
167856 347812
15 8.91765957989e+12
670929 862765
139395 558143
244654 847642
852058 157685
617601 399504
801616 321123
507818 175610
337264 562123
858947 546259
823421 731145
606574 574455
167856 347812
398555 404855
417799 787001
320602 161521
Fit?
16 Clusters

Series1
If you have another session window open on bridge's login node, you can edit this file, save it while you remain in the editor, and then run it again in the python shell window with execfile().

You do **not** need this second session to be on a compute node. Do not start another interactive session.
As mentioned earlier, an appreciation for having some defined structure to your data has come back into vogue. For one, because it simply makes sense and naturally emerges in many applications. Often even more important, it can greatly aid optimization, especially with the Java VM that Spark uses.

For both of these reasons, you will see that the newest set of APIs to Spark are DataFrame based. Sound leading-edge? This is simply SQL type columns. Very similar to Python pandas DataFrames (but based on RDDs, so not exactly).

We haven't prioritized them here because they aren't necessary, and some of the pieces aren't mature. But some of the latest features use them.
Creating DataFrames

It is very pretty intuitive to utilize DataFrames. Your elements just have labeled columns.

A row RDD is the basic way to go from RDD to DataFrame, and back, if necessary. A "row" is just a tuple.

```python
>>> row_rdd = sc.parallelize([ ("Joe","Pine St.","PA",12543), ("Sally","Fir Dr.","WA",78456),
                           ("Jose","Elm Pl.","ND",45698) ])  

>>> aDataFrameFromRDD = spark.createDataFrame( row_rdd, ["name", "street", "state", "zip"] )  
>>> aDataFrameFromRDD.show()
+--------+-------+--------+----+
| name   | street| state  | zip |
|--------+-------+--------+----|
| Joe    | Pine St.| PA     | 12543|
| Sally  | Fir Dr.| WA     | 78456|
| Jose   | Elm Pl.| ND     | 45698|
+--------+-------+--------+----+
```
Creating DataFrames

You will come across DataFrames created without a schema. They get default column names.

```python
>>> noSchemaDataFrame = spark.createDataFrame(row_rdd)
>>> noSchemaDataFrame.show()
+-----+--------+-----+-----+
| _1  | _2     | _3  | _4  |
+-----+--------+-----+-----+
| Joe | Pine St.| PA  |12543|
| Sally | Fir Dr.| WA |78456|
| Jose | Elm Pl.| ND |45698|
+-----+--------+-----+-----+
```

And you can create them inline as well.

```python
>>> directDataFrame = spark.createDataFrame([ ("Joe","Pine St." ,"PA",12543), ("Sally","Fir Dr." ,"WA",78456),
                                            ("Jose","Elm Pl." ,"ND",45698) ],
                                            ["name","street","state","zip"] )
```
Speaking of pandas, or SciPy, or...

Some of you may have experience with the many Python libraries that accomplish some of these tasks. Immediately relevant to today, pandas allows us to sort and query data, and SciPy provides some nice clustering algorithms. So why not just use them?

The answer is that Spark does these things in the context of having potentially huge, parallel resources at hand. We don't notice it as Spark is also convenient, but behind every Spark call:

- every RDD could be many TB in size
- every transform could use many thousands of cores and TB of memory
- every algorithm could also use those thousands of cores

So don't think of Spark as just a data analytics library because our exercises are modest. You are learning how to cope with Big Data.
How does all this fit together?