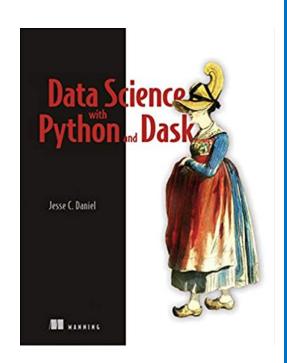
# UMD DATA605 - Big Data Systems Python Dask

GP Saggese gsaggese@umd.edu

with thanks to Alan Sussman, Amol Deshpande, and authors of www.mmds.org

### **Dask: Resources**

- Web resources:
  - Dask project
  - Dask examples
- Tutorial
  - Dask tutorial
  - Dask advanced tutorial
- Class project
- Mastery
  - Data science with Python and Dask, 2019
    - Amazon



### **Dataset Size Issues**

#### Small datasets

- < 1 GB
- Fits into RAM
- Manipulation doesn't require paging to disk

#### Medium dataset

- < 1TB</p>
- Doesn't fit into RAM
- Fits into local disk
  - Performance penalty imposed by using local disk
- Need multiple CPU cores
  - Difficult to take advantage of parallelism with Python / Pandas

#### Large dataset

- \_ > 1TB
- Doesn't fit into RAM
- Doesn't fit into local disk
- Need multiple servers
  - Python / Pandas were not built to operate on distributed datasets
  - Use frameworks for massive datasets
  - E.g., Hadoop, Spark, Dask, Ray

#### The thresholds are fuzzy and changing over time

- E.g., you can scale the computer 10x and get 10x bigger data sets
- Problem when scaling datasets
  - Long run times
  - Rewriting code in different language / API for datasets of different size
  - Need to think about what to do it and how to do it efficiently
  - Cumbersome framework



	outlook	temp	humidity	windy	play
0	sunny	hot	high	False	no
1	sunny	hot	high	True	no
2	overcast	hot	high	False	yes
3	rainy	mild	high	False	yes
4	rainy	cool	normal	False	yes
5	rainy	cool	normal	True	no
6	overcast	cool	normal	True	yes
7	sunny	mild	high	False	no
8	sunny	cool	normal	False	yes
9	rainy	mild	normal	False	yes
10	sunny	mild	normal	True	yes
11	overcast	mild	high	True	yes
12	overcast	hot	normal	False	yes
13	rainy	mild	high	True	no







### Dask

#### Dask is written in Python

- It scales natively Numpy, Pandas, sklearn
- Dask objects are wrappers (don't just mirror the interface) objects from the respective libraries
  - Dask DataFrame = composed of several Pandas DataFrame
  - Dask Array = composed of several Pandas numpy array
- Parallel parts are called "chunks" or "partitions"
  - Are queued to be worked on
  - Shipped between machines
  - Worked locally on a machine

#### Pros

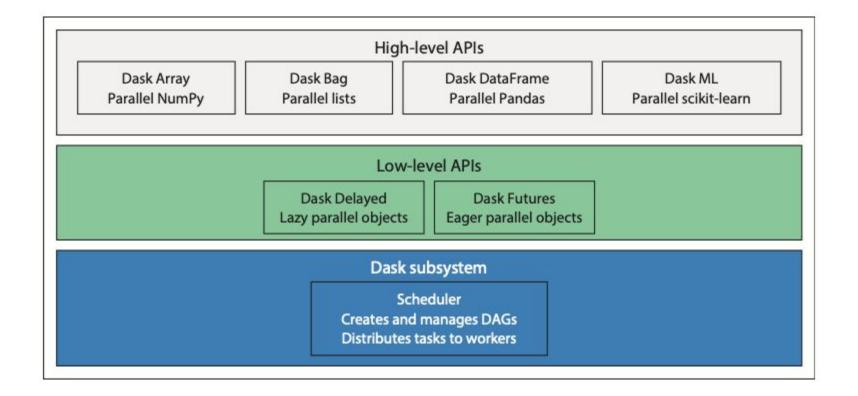
- Users don't need to learn a new language, but can use familiar interfaces
- Can focus on writing code that is optimized for parallelism
- Dask does the heavy lifting

#### Scaling Dask is easy

- Users can write a prototype task on a local machines and use a cluster when needed
- No need to refactor existing code
- No need to handle cluster-specific issues
  - E.g., resource management, data recovery, data movement
- Dask runs on multi-core and
- Dask can use cluster managers
  - E.g., Yarn, Mesos, Kubernetes, AWS ECS



# **Dask Layers**



# Scaling Up vs Scaling Out

#### Scaling up

- = replace equipment with larger, faster equipments
  - E.g., buy a larger pot
  - Replace knife with food processor

#### - Pros

You got better hardware, nothing else needs to change (e.g., code)

#### Cons

- There will be a time where you exceed the capacity of the current machines
- Cost: more powerful machines are expensive

#### Scaling out

- = divides the work between many workers in parallel
  - E.g., hire more cooks
  - Buy more knives

#### Pros

- Task scheduler organizes computation, assigning workers to each task
- More cost-effective solution since no specialized hardware is needed

#### Cons

- Need to write code to expose parallelism
- Costs of maintaining a cluster

## **Dask: Computation**

#### Lazy computations

- User defines the transformations on the data
- No need to wait for one computation to finish before defining the next
- Avoid loading the entire data in memory by operating in chunks
- E.g.,
  - Split a 2GB file into 32 64MB chunks
  - Operate on 8 chunks at a time
  - The max memory consumption doesn't exceed 512MB = (8 x 32)
- Each task tracks object dimensions and data types
  - No code is executed

#### compute()

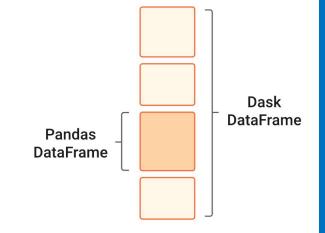
- Running a computation (aka materializing)
missing count pct = missing count.compute()

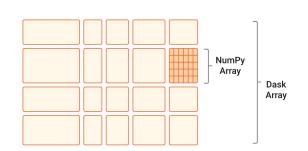
#### persist()

- As soon as a node in the graph emits results, its intermediate work is
- discarded to minimize memory usage
- If we need to do additional computation on intermediate nodes we need to re-run the graph
- persist() tells Dask to keep the intermediate result in memory
- This speeds up a large and complex DAG that needs to be reused many times

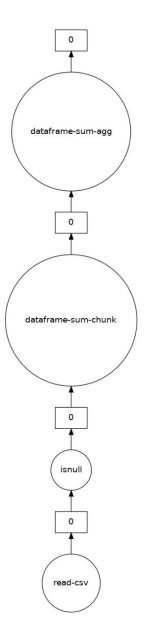
### **Dask: Data Structures**

- Dask DataFrame implements Pandas DataFrame
- Dask Array implements numpy ndarray
- Dag Bag coordinates Python lists of objects
  - Parallelize computations on unstructured or semi-structured data



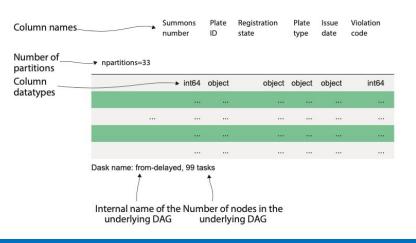


# **Dask Reading Data**



```
import dask.dataframe as dd
df = dd.read_csv('nyc-parking-tickets-2017.csv')
missing_values = df.isnull().sum()
missing_values
```

- dask.dataframe.read\_csv()
  - Doesn't load the data in memory with
  - Tries to infer the types of the columns
    - By randomly sampling some data
    - Best to explicitly set the data types
    - Even better is to use Parquet since it stores data and types together
- Partitions = chunks of data that can be worked independently
  - E.g., 33 partitions
  - Graph is composed of 99 tasks
  - Each partition reads data, splits data, initialize df object



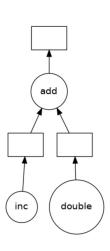
## Low Level APIs: Delayed

Handle computations that don't fit in native Dask data structures (e.g., Dask DataFrame)

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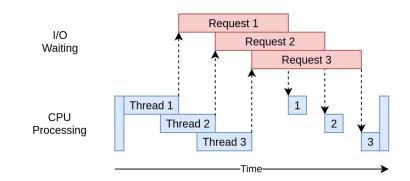
In the example there is parallelism that can be exploited

```
def inc(x):
    return x + 1
def double(x):
    return x * 2
def add(x, y):
    return x + y
data = [1, 2, 3, 4, 5]
output = []
for x in data:
    \# (x + 1) + (x * 2) = 3x + 1
    a = inc(x)
    b = double(x)
    c = add(a, b)
    # 1 -> 4
    # 2 -> 7
    # 3 -> 10
    # 4 -> 13
    # 5 -> 16
    output.append(c)
\# 4 + 7 + 10 + 13 + 16 = 20 + 20 + 10 = 50
total = sum(output)
print(total)
```



### Low Level APIs: Futures

- Futures is the most general way of specifying concurrency in Dask
  - Everything can be expressed in terms of futures
  - User can specify what's blocking and what's not blocking
- In parallel programming, a "future" encapsulates the asynchronous execution of a callable, representing the eventual result of the operation
- concurrent.futures
  - High-level interface for asynchronously executing callables
  - Thread pool or Process pool (same interface Executor)
- Dask extends concurrent.futures
  - Dask client can be used anywhere concurrent.futures can be used



```
def inc(x):
    return x + 1

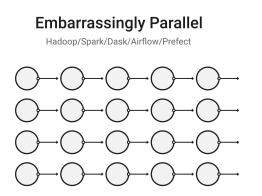
def add(x, y):
    return x + y

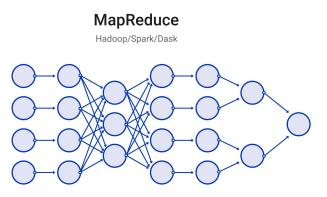
a = client.submit(inc, 10)
b = client.submit(inc, 20)
```

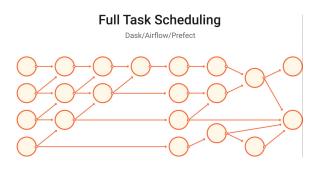
```
>>> a
<Future: status: pending, key: inc-b8aaf26b99466a7a1980efa1ade6701d>
>>> a
<Future: status: finished, type: int, key: inc-b8aaf26b99466a7a1980efa1ade6701d>
>>> a.result() # blocks until task completes and data arrives
11
```

# Different Types of Parallel Workload

- Break program in medium-size tasks of computation
  - E.g., a function call







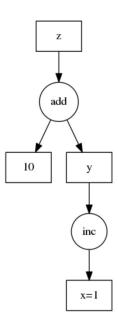
# **Encoding Task Graph**

Dask encodes tasks in terms of Python dicts and functions

```
def inc(i):
    return i + 1

def add(a, b):
    return a + b

x = 1
y = inc(x)
z = add(y, 10)
```



```
import dask.dataframe as dd

df = dd.read_csv('myfile.*.csv')

df = df + 100

df = df[df.name == 'Alice']
```

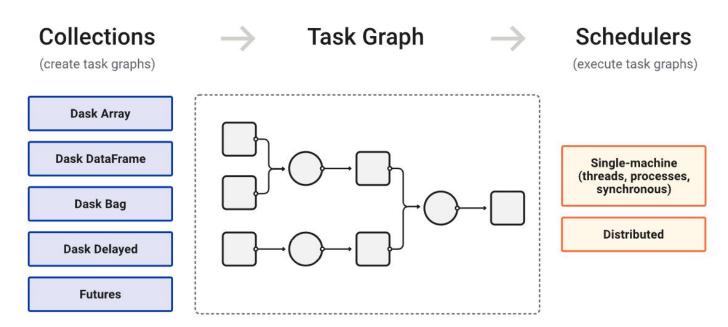
```
{
# From the dask.dataframe.read_csv call
('read-csv', 0): (pandas.read_csv, 'myfile.0.csv'),
('read-csv', 1): (pandas.read_csv, 'myfile.1.csv'),
('read-csv', 2): (pandas.read_csv, 'myfile.2.csv'),
('read-csv', 3): (pandas.read_csv, 'myfile.3.csv'),

# From the df + 100 call
('add', 0): (operator.add, ('read-csv', 0), 100),
('add', 1): (operator.add, ('read-csv', 1), 100),
('add', 2): (operator.add, ('read-csv', 2), 100),
('add', 3): (operator.add, ('read-csv', 3), 100),

# From the df[df.name == 'Alice'] call
('filter', 0): (lambda part: part[part.name == 'Alice'], ('add', 0)),
('filter', 1): (lambda part: part[part.name == 'Alice'], ('add', 1)),
('filter', 3): (lambda part: part[part.name == 'Alice'], ('add', 2)),
('filter', 3): (lambda part: part[part.name == 'Alice'], ('add', 3)),
}
```

## Task Scheduling

- Data collections (Bags, Arrays, DataFrame) and operations on them create task graphs
  - Nodes in the task graph are Python functions
  - Edges are dependencies (e.g., output from one task used as input in another task)
- Task graphs are then scheduled for execution on
  - a single machine or;
  - a cluster
- Single-machine scheduler
  - Use local process or thread pool
  - Simple but it can only run on a single machine
- Distributed scheduler
  - It can run locally or distributed across a cluster



# Task Scheduling

- Dask centralized task scheduler
  - Orchestrate the work dynamically
  - Not a static scheduling of operations like a relational DB
- · When the computation takes place, Dask dynamically assesses:
  - What work has been completed
  - What work is left to do
  - What resources are free
  - Where the data is located
- This dynamic approach handles:
  - Worker failure
  - Workers completing work at different speeds because of:
    - different computation
    - different hardware
    - different workloads on the servers
    - slower access to the data
  - Network unreliability

### Dask vs Spark

- Spark is a popular framework for analyzing large datasets
  - In-memory alternative to MapReduce / Hadoop
- Cons
  - Spark is a Java library, supporting Python through PySpark API
    - Python code is executed on JVM through `py4j`
    - Difficult to debug since execution occurs outside Python
  - Different DataFrame API than Pandas
    - Learn how to do things "the Spark way"
  - Optimized for MapReduce operations over a collection
  - Difficult to set-up and configure

### **Tutorial**

### **Tutorial**

- From the official documentation https://docs.dask.org/en/stable/10-minutes-to-dask.html