# mrjob Documentation

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**Steve Johnson** 

## Contents

1	Guid		
	1.1	Why mrjob?	3
	1.2	Fundamentals	
	1.3	Concepts	7
	1.4	Writing jobs	)
	1.5	Runners	3
	1.6	Spark	7
	1.7	Config file format and location	4
	1.8	Options available to all runners	)
	1.9	Hadoop-related options	5
	1.10	Spark runner options	9
	1.11	Configuration quick reference	)
	1.12	Cloud runner options	3
	1.13	Job Environment Setup Cookbook	3
	1.14	Hadoop Cookbook	1
	1.15	Testing jobs	1
	1.16	Cloud Dataproc	5
	1.17	Elastic MapReduce	0
	1.18	Python 2 vs. Python 3	5
	1.19	Contributing to mrjob	5
2	Refer		
	2.1	mrjob.ami - building custom AMIs	
	2.2	mrjob.cat - decompress files based on extension	)
	2.3	mrjob.cmd: The mrjob command-line utility	)
	2.4	mrjob.compat - Hadoop version compatibility	)
	2.5	mrjob.conf - parse and write config files	1
	2.6	mrjob.dataproc - run on Dataproc	3
	2.7	mrjob.emr - run on EMR	4
	2.8	mrjob.hadoop - run on your Hadoop cluster	5
	2.9	mrjob.inline - debugger-friendly local testing	5
	2.10	mrjob.job - defining your job	5
	2.11	mrjob.local - simulate Hadoop locally with subprocesses	3
	2.12	mrjob.parse - log parsing	3
	2.13	mrjob.protocol - input and output	9
	2.14	mrjob.spark.runner - run on any Spark cluster	
	2.15	mrjob.retry - retry on transient errors	
	2.16	mrjob.runner - base class for all runners	
		•	

	2.17	Jerosepp	127						
	2.18	J. T. T. J. T.	130						
	2.19	mrjob.util - general utility functions	132						
3	What's New								
	3.1	0.6.10	<b>137</b> 137						
	3.2	0.6.9	137						
	3.3	0.6.8	138						
	3.4	0.6.7	140						
	3.5	0.6.6	141						
	3.6	0.6.5	142						
	3.7	0.6.4	142						
	3.8	0.6.3	142						
	3.9	0.6.2	144						
	3.10	0.6.1	144						
	3.10		144						
		0.6.0	144						
	3.12	0.5.12							
	3.13	0.5.11	147						
	3.14	0.5.10	147						
		0.5.9	147						
		0.5.8	148						
		0.5.7	148						
		0.5.6	150						
		0.5.5	150						
		0.5.4	150						
	3.21	0.5.3	152						
		0.5.2	152						
		0.5.1	153						
		0.5.0	153						
		0.4.6	157						
		0.4.5	157						
		0.4.4	158						
	3.28	0.4.3	158						
		0.4.2	159						
		0.4.1	159						
		0.4.0	160						
	3.32	0.3.5	160						
		0.3.3	160						
		0.3.2	160						
	3.35	0.3	161						
4	Gloss	sary	165						
Python Module Index									

## mrjob lets you write MapReduce jobs in Python 2.7/3.4+ and run them on several platforms. You can:

- Write multi-step MapReduce jobs in pure Python
- Test on your local machine
- Run on a Hadoop cluster
- Run in the cloud using Amazon Elastic MapReduce (EMR)
- Run in the cloud using Google Cloud Dataproc (Dataproc)
- Easily run Spark jobs on EMR or your own Hadoop cluster

mrjob is licensed under the Apache License, Version 2.0.

To get started, install with pip:

pip install mrjob

and begin reading the tutorial below.

Contents 1

2 Contents

# **Guides**

# 1.1 Why mrjob?

#### 1.1.1 Overview

mrjob is the easiest route to writing Python programs that run on Hadoop. If you use mrjob, you'll be able to test your code locally without installing Hadoop or run it on a cluster of your choice.

Additionally, mrjob has extensive integration with Amazon Elastic MapReduce. Once you're set up, it's as easy to run your job in the cloud as it is to run it on your laptop.

Here are a number of features of mrjob that make writing MapReduce jobs easier:

- Keep all MapReduce code for one job in a single class
- Easily upload and install code and data dependencies at runtime
- Switch input and output formats with a single line of code
- · Automatically download and parse error logs for Python tracebacks
- Put command line filters before or after your Python code

If you don't want to be a Hadoop expert but need the computing power of MapReduce, mrjob might be just the thing for you.

# 1.1.2 Why use mrjob instead of X?

Where X is any other library that helps Hadoop and Python interface with each other.

- 1. mrjob has more documentation than any other framework or library we are aware of. If you're reading this, it's probably your first contact with the library, which means you are in a great position to provide valuable feedback about our documentation. Let us know if anything is unclear or hard to understand.
- 2. mrjob lets you run your code without Hadoop at all. Other frameworks require a Hadoop instance to function at all. If you use mrjob, you'll be able to write proper tests for your MapReduce code.
- 3. mrjob provides a consistent interface across every environment it supports. No matter whether you're running locally, in the cloud, or on your own cluster, your Python code doesn't change at all.
- 4. mrjob handles much of the machinery of getting code and data to and from the cluster your job runs on. You don't need a series of scripts to install dependencies or upload files.

- 5. mrjob makes debugging much easier. Locally, it can run a simple MapReduce implementation in-process, so you get a traceback in your console instead of in an obscure log file. On a cluster or on Elastic MapReduce, it parses error logs for Python tracebacks and other likely causes of failure.
- 6. mrjob automatically serializes and descrializes data going into and coming out of each task so you don't need to constantly json.loads() and json.dumps().

# 1.1.3 Why use X instead of mrjob?

The flip side to mrjob's ease of use is that it doesn't give you the same level of access to Hadoop APIs that Dumbo and Pydoop do. It's simplified a great deal. But that hasn't stopped several companies, including Yelp, from using it for day-to-day heavy lifting. For common (and many uncommon) cases, the abstractions help rather than hinder.

Other libraries can be faster if you use typedbytes. There have been several attempts at integrating it with mrjob, and it may land eventually, but it doesn't exist yet.

# 1.2 Fundamentals

#### 1.2.1 Installation

Install with pip:

```
pip install mrjob
```

or from a git clone of the source code:

```
python setup.py test && python setup.py install
```

# 1.2.2 Writing your first job

Open a file called mr\_word\_count.py and type this into it:

```
from mrjob.job import MRJob

class MRWordFrequencyCount(MRJob):

    def mapper(self, _, line):
        yield "chars", len(line)
        yield "words", len(line.split())
        yield "lines", 1

    def reducer(self, key, values):
        yield key, sum(values)

if __name__ == '__main__':
    MRWordFrequencyCount.run()
```

Now go back to the command line, find your favorite body of text (such mrjob's README.rst, or even your new file mr\_word\_count.py), and try this:

```
$ python mr_word_count.py my_file.txt
```

You should see something like this:

```
"chars" 3654
"lines" 123
"words" 417
```

Congratulations! You've just written and run your first program with mrjob.

#### What's happening

A job is defined by a class that inherits from MRJob. This class contains methods that define the steps of your job.

A "step" consists of a mapper, a combiner, and a reducer. All of those are optional, though you must have at least one. So you could have a step that's just a mapper, or just a combiner and a reducer.

When you only have one step, all you have to do is write methods called mapper(), combiner(), and reducer().

The mapper () method takes a key and a value as args (in this case, the key is ignored and a single line of text input is the value) and yields as many key-value pairs as it likes. The reduce () method takes a key and an iterator of values and also yields as many key-value pairs as it likes. (In this case, it sums the values for each key, which represent the numbers of characters, words, and lines in the input.)

**Warning:** Forgetting the following information will result in confusion.

The final required component of a job file is these two lines at the end of the file, **every time**:

```
if __name__ == '__main__':
    MRWordCounter.run() # where MRWordCounter is your job class
```

These lines pass control over the command line arguments and execution to mrjob. Without them, your job will not work. For more information, see *Hadoop Streaming and mrjob* and *Why can't I put the job class and run code in the same file?*.

# 1.2.3 Running your job different ways

The most basic way to run your job is on the command line:

```
$ python my_job.py input.txt
```

By default, output will be written to stdout.

You can pass input via stdin, but be aware that mrjob will just dump it to a file first:

```
$ python my_job.py < input.txt</pre>
```

You can pass multiple input files, mixed with stdin (using the – character):

```
$ python my_job.py input1.txt input2.txt - < input3.txt</pre>
```

By default, mrjob will run your job in a single Python process. This provides the friendliest debugging experience, but it's not exactly distributed computing!

You change the way the job is run with the -r/--runner option. You can use -r inline (the default), -r local, -r hadoop, or -r emr.

To run your job in multiple subprocesses with a few Hadoop features simulated, use -r local.

To run it on your Hadoop cluster, use -r hadoop.

1.2. Fundamentals 5

If you have Dataproc configured (see Dataproc Quickstart), you can run it there with -r dataproc.

Your input files can come from HDFS if you're using Hadoop, or GCS if you're using Dataproc:

```
$ python my_job.py -r dataproc gcs://my-inputs/input.txt
$ python my_job.py -r hadoop hdfs://my_home/input.txt
```

If you have Elastic MapReduce configured (see Elastic MapReduce Quickstart), you can run it there with -r emr.

Your input files can come from HDFS if you're using Hadoop, or S3 if you're using EMR:

```
$ python my_job.py -r emr s3://my-inputs/input.txt
$ python my_job.py -r hadoop hdfs://my_home/input.txt
```

If your code spans multiple files, see *Uploading your source tree*.

## 1.2.4 Writing your second job

Most of the time, you'll need more than one step in your job. To define multiple steps, override steps() to return a list of MRSteps.

Here's a job that finds the most commonly used word in the input:

```
from mrjob.job import MRJob
from mrjob.step import MRStep
import re
WORD_RE = re.compile(r"[\w']+")
class MRMostUsedWord (MRJob) :
    def steps(self):
        return [
            MRStep (mapper=self.mapper_get_words,
                   combiner=self.combiner_count_words,
                   reducer=self.reducer_count_words),
            MRStep(reducer=self.reducer_find_max_word)
        ]
    def mapper_get_words(self, _, line):
        # yield each word in the line
        for word in WORD_RE.findall(line):
            yield (word.lower(), 1)
    def combiner_count_words(self, word, counts):
        # optimization: sum the words we've seen so far
        yield (word, sum(counts))
    def reducer_count_words(self, word, counts):
        # send all (num_occurrences, word) pairs to the same reducer.
        # num_occurrences is so we can easily use Python's max() function.
        yield None, (sum(counts), word)
    # discard the key; it is just None
    def reducer_find_max_word(self, _, word_count_pairs):
        # each item of word_count_pairs is (count, word),
        # so yielding one results in key=counts, value=word
        yield max(word_count_pairs)
```

```
if __name__ == '__main__':
    MRMostUsedWord.run()
```

## 1.2.5 Configuration

mrjob has an overflowing cornucopia of configuration options. You'll want to specify some on the command line, some in a config file.

You can put a config file at /etc/mrjob.conf, ~/.mrjob.conf, or ./mrjob.conf for mrjob to find it without passing it via --conf-path.

Config files are interpreted as YAML if you have the yaml module installed. Otherwise, they are interpreted as JSON.

See Config file format and location for in-depth information. Here is an example file:

```
runners:
    emr:
    aws-region: us-west-2
    inline:
    local_tmp_dir: $HOME/.tmp
```

# 1.3 Concepts

# 1.3.1 MapReduce and Apache Hadoop

This section uses text from Apache's MapReduce Tutorial.

MapReduce is a way of writing programs designed for processing vast amounts of data, and a system for running those programs in a distributed and fault-tolerant way. Apache Hadoop is one such system designed primarily to run Java code.

A MapReduce job usually splits the input data-set into independent chunks which are processed by the map tasks in a completely parallel manner. The framework sorts the outputs of the maps, which are then input to the reduce tasks. Typically both the input and the output of the job are stored in a file system shared by all processing nodes. The framework takes care of scheduling tasks, monitoring them, and re-executing the failed tasks.

The MapReduce framework consists of a single master "job tracker" (Hadoop 1) or "resource manager" (Hadoop 2) and a number of worker nodes. The master is responsible for scheduling the jobs' component tasks on the worker nodes and re-executing the failed tasks. The worker nodes execute the tasks as directed by the master.

As the job author, you write map, combine, and reduce functions that are submitted to the job tracker for execution.

A *mapper* takes a single key and value as input, and returns zero or more (key, value) pairs. The pairs from all map outputs of a single step are grouped by key.

A *combiner* takes a key and a subset of the values for that key as input and returns zero or more (key, value) pairs. Combiners are optimizations that run immediately after each mapper and can be used to decrease total data transfer. Combiners should be idempotent (produce the same output if run multiple times in the job pipeline).

A *reducer* takes a key and the complete set of values for that key in the current step, and returns zero or more arbitrary (key, value) pairs as output.

After the reducer has run, if there are more steps, the individual results are arbitrarily assigned to mappers for further processing. If there are no more steps, the results are sorted and made available for reading.

1.3. Concepts 7

### An example

Consider a program that counts how many times words occur in a document. Here is some input:

```
The wheels on the bus go round and round, round and round and round and round The wheels on the bus go round and round, all through the town.
```

The inputs to the mapper will be (None, "one line of text"). (The key is None because the input is just raw text.)

The mapper converts the line to lowercase, removes punctuation, splits it on whitespace, and outputs (word, 1) for each item.

```
mapper input: (None, "The wheels on the bus go round and round,")
mapper output:
    "the", 1
    "wheels", 1
    "on", 1
    "the", 1
    "bus", 1
    "go", 1
    "round", 1
    "and", 1
    "round", 1
```

Each call to the combiner gets a word as the key and a list of 1s as the value. It sums the 1s and outputs the original key and the sum.

```
combiner input: ("the", [1, 1])
combiner output:
    "the", 2
```

The reducer is identical to the combiner; for each key, it simply outputs the original key and the sum of the values.

```
reducer input: ("round", [2, 4, 2])
reducer output:
    "round", 8
```

The final output is collected:

```
"all", 1
"and", 4
"bus", 2
"go", 2
"on", 2
"round", 8
"the", 5
"through", 1
"town", 1
"wheels", 2
```

Your algorithm may require several repetitions of this process.

# 1.3.2 Hadoop Streaming and mrjob

**Note:** If this is your first exposure to MapReduce or Hadoop, you may want to skip this section and come back later. Feel free to stick with it if you feel adventurous.

Although Hadoop is primarly designed to work with Java code, it supports other languages via *Hadoop Streaming*. This jar opens a subprocess to your code, sends it input via stdin, and gathers results via stdout.

In most cases, the input to a Hadoop Streaming job is a set of newline-delimited files. Each line of input is passed to your mapper, which outputs key-value pairs expressed as two strings separated by a tab and ending with a newline, like this:

```
key1\tvalue1\nkey2\tvalue2\n
```

Hadoop then sorts the output lines by key (the line up to the first tab character) and passes the sorted lines to the appropriate combiners or reducers.

mrjob is a framework that assists you in submitting your job to the Hadoop job tracker and in running each individual step under Hadoop Streaming.

## How your program is run

Depending on the way your script is invoked on the command line, it will behave in different ways. You'll only ever use one of these; the rest are for mrjob and Hadoop Streaming to use.

When you run with no arguments or with --runner, you invoke mrjob's machinery for running your job or submitting it to the cluster. Your mappers and reducers are not called in this process at all <sup>1</sup>.

This process creates a runner (see MRJobRunner), which then sends the job to Hadoop <sup>2</sup>.

It tells Hadoop something like this:

- Run a step with Hadoop Streaming.
- The command for the mapper is python my\_job.py --step-num=0 --mapper.
- The command for the combiner is python <code>my\_job.py --step-num=0 --combiner</code>.
- The command for the reducer is python my\_job.py --step-num=0 --reducer.

If you have a multi-step job, --step-num helps your script know which step is being run.

When Hadoop distributes tasks among the task nodes, Hadoop Streaming will use the appropriate command to process the data it is given.

**Note:** Prior to v0.6.7, your job would *also* run itself locally with the --steps switch, to get a JSON representation of the job's step. Jobs now pass that representation directly to the runner when they instantiate it. See mrjob.step represent Job Steps for more information.

# 1.4 Writing jobs

This guide covers everything you need to know to write your job. You'll probably need to flip between this guide and Runners to find all the information you need.

<sup>&</sup>lt;sup>1</sup> Unless you're using the inline runner, which is a special case for debugging.

<sup>&</sup>lt;sup>2</sup> Or when using the local runner, a simulation of Hadoop.

## 1.4.1 Defining steps

Your job will be defined in a file to be executed on your machine as a Python script, as well as on a Hadoop cluster as an individual map, combine, or reduce task. (See *How your program is run* for more on that.)

All dependencies must either be contained within the file, available on the task nodes, or uploaded to the cluster by mrjob when your job is submitted. (Runners explains how to do those things.)

The following two sections are more reference-oriented versions of Writing your first job and Writing your second job.

#### Single-step jobs

The simplest way to write a one-step job is to subclass MRJob and override a few methods:

```
from mrjob.job import MRJob
import re

WORD_RE = re.compile(r"[\w']+")

class MRWordFreqCount(MRJob):
    def mapper(self, _, line):
        for word in WORD_RE.findall(line):
            yield word.lower(), 1

    def combiner(self, word, counts):
        yield word, sum(counts)

    def reducer(self, word, counts):
        yield word, sum(counts)

if __name__ == '__main__':
    MRWordFreqCount.run()
```

(See Writing your first job for an explanation of this example.)

Here are all the methods you can override to write a one-step job. We'll explain them over the course of this document.

```
combiner()
reducer()
mapper_init()
combiner_init()
reducer_init()
mapper_final()
combiner_final()
```

• mapper()

- reducer\_final()
- mapper\_cmd()
- combiner\_cmd()

```
reducer_cmd()mapper_pre_filter()combiner_pre_filter()reducer_pre_filter()
```

#### Multi-step jobs

To define multiple steps, override steps() to return a list of MRSteps:

```
from mrjob.job import MRJob
from mrjob.step import MRStep
import re
WORD_RE = re.compile(r"[\w']+")
class MRMostUsedWord(MRJob):
    def mapper_get_words(self, _, line):
        # vield each word in the line
        for word in WORD_RE.findall(line):
            yield (word.lower(), 1)
    def combiner_count_words(self, word, counts):
        # sum the words we've seen so far
        yield (word, sum(counts))
    def reducer_count_words(self, word, counts):
        # send all (num_occurrences, word) pairs to the same reducer.
        # num_occurrences is so we can easily use Python's max() function.
        yield None, (sum(counts), word)
    # discard the key; it is just None
    def reducer_find_max_word(self, _, word_count_pairs):
        # each item of word_count_pairs is (count, word),
        # so yielding one results in key=counts, value=word
       yield max(word_count_pairs)
   def steps(self):
        return [
            MRStep (mapper=self.mapper_get_words,
                   combiner=self.combiner_count_words,
                   reducer=self.reducer_count_words),
           MRStep(reducer=self.reducer_find_max_word)
if __name__ == '__main__':
   MRMostUsedWord.run()
```

(This example is explained further in *Protocols*.)

The keyword arguments accepted by MRStep are the same as the method names listed in the previous section, plus a jobconf argument which takes a dictionary of jobconf arguments to pass to Hadoop.

**Note:** If this is your first time learning about mrjob, you should skip down to *Protocols* and finish this section later.

#### Setup and teardown of tasks

Remember from *How your program is run* that your script is invoked once per task by Hadoop Streaming. It starts your script, feeds it stdin, reads its stdout, and closes it. mrjob lets you write methods to run at the beginning and end of this process: the \*\_init() and \*\_final() methods:

```
mapper_init()
combiner_init()
reducer_init()
mapper_final()
combiner_final()
reducer_final()
```

(And the corresponding keyword arguments to MRStep.)

If you need to load some kind of support file, like a sqlite3 database, or perhaps create a temporary file, you can use these methods to do so. (See *File options* for an example.)

 $\star$ \_init() and  $\star$ \_final() methods can yield values just like normal tasks. Here is our word frequency count example rewritten to use these methods:

```
from mrjob.job import MRJob
from mrjob.step import MRStep
class MRWordFreqCount (MRJob):
    def init_get_words(self):
        self.words = {}
    def get_words(self, _, line):
        for word in WORD_RE.findall(line):
            word = word.lower()
            self.words.setdefault(word, 0)
            self.words[word] = self.words[word] + 1
    def final_get_words(self):
        for word, val in self.words.iteritems():
            yield word, val
    def sum_words(self, word, counts):
        yield word, sum(counts)
    def steps(self):
        return [MRStep (mapper_init=self.init_get_words,
                       mapper=self.get_words,
                       mapper_final=self.final_get_words,
                       combiner=self.sum_words,
                       reducer=self.sum_words)]
```

In this version, instead of yielding one line per word, the mapper keeps an internal count of word occurrences across all lines this mapper has seen so far. The mapper itself yields nothing. When Hadoop Streaming stops sending data to the map task, mrjob calls final\_get\_words(). That function emits the totals for this task, which is a much smaller set of output lines than the mapper would have output.

The optimization above is similar to using *combiners*, demonstrated in *Multi-step jobs*. It is usually clearer to use a combiner rather than a custom data structure, and Hadoop may run combiners in more places than just the ends of tasks.

Defining command line options has a partial example that shows how to load a sqlite3 database using mapper\_init().

#### Shell commands as steps

You can forego scripts entirely for a step by specifying it as a shell command. To do so, use mapper\_cmd, combiner\_cmd, or reducer\_cmd as arguments to MRStep, or override the methods of the same names on MRJob. (See mapper\_cmd(), combiner\_cmd(), and reducer\_cmd().)

```
Warning: The default inline runner does not support \star_cmd(). If you want to test locally, use the local runner (-r local).
```

You may mix command and script steps at will. This job will count the number of lines containing the string "kitty":

```
from mrjob.job import job

class KittyJob(MRJob):

   OUTPUT_PROTOCOL = JSONValueProtocol

   def mapper_cmd(self):
      return "grep kitty"

   def reducer(self, key, values):
      yield None, sum(1 for _ in values)

if __name__ == '__main__':
   KittyJob().run()
```

Step commands are run without a shell, so if you want to use pipes, etc, you'll need to run them in a subshell. For example:

```
class DemoJob(MRJob):
    def mapper_cmd(self):
        return 'sh -c "grep 'blah' | wc -l"'
```

**Note:** You may not use  $\star$ \_cmd() with any other options for a task such as  $\star$ \_filter(),  $\star$ \_init(),  $\star$ \_final(), or a regular mapper/combiner/reducer function.

**Note:** You might see an opportunity here to write your MapReduce code in whatever language you please. If that appeals to you, check out upload\_files for another piece of the puzzle.

#### Filtering task input with shell commands

You can specify a command to filter a task's input before it reaches your task using the mapper\_pre\_filter and reducer\_pre\_filter arguments to <code>MRStep</code>, or override the methods of the same names on <code>MRJob</code>. Doing so will cause mrjob to pipe input through that command before it reaches your mapper.

**Warning:** The default inline runner does not support  $*\_pre\_filter()$ . If you want to test locally, use the local runner (-r local).

Here's a job that tests filters using grep:

```
from mrjob.job import MRJob
from mrjob.protocol import JSONValueProtocol
from mrjob.step import MRStep
class KittiesJob (MRJob) :
   OUTPUT_PROTOCOL = JSONValueProtocol
   def test_for_kitty(self, _, value):
       yield None, 0 # make sure we have some output
        if 'kitty' not in value:
           yield None, 1
   def sum_missing_kitties(self, _, values):
        yield None, sum (values)
   def steps(self):
        return [
           MRStep(mapper_pre_filter='grep "kitty"',
                   mapper=self.test_for_kitty,
                   reducer=self.sum_missing_kitties)]
if name == ' main ':
   KittiesJob().run()
```

The output of the job should always be 0, since every line that gets to test\_for\_kitty() is filtered by **grep** to have "kitty" in it.

## 1.4.2 Protocols

Hadoop streaming assumes that all data is newline-delimited bytes. By default, mrjob assumes all output is in JSON format, but it can actually read and write lines in any format by using protocols.

(If you need to read non-line-based data, see *Passing entire files to the mapper*, below.)

Each job has an input protocol, an output protocol, and an internal protocol.

A protocol has a read() method and a write() method. The read() method converts bytes to pairs of Python objects representing the keys and values. The write() method converts a pair of Python objects back to bytes.

The *input protocol* is used to read the bytes sent to the first mapper (or reducer, if your first step doesn't use a mapper). The *output protocol* is used to write the output of the last step to bytes written to the output file. The *internal protocol* converts the output of one step to the input of the next if the job has more than one step.

You can specify which protocols your job uses like this:

```
class MyMRJob(mrjob.job.MRJob):
    # these are the defaults
    INPUT_PROTOCOL = mrjob.protocol.RawValueProtocol
```

```
INTERNAL_PROTOCOL = mrjob.protocol.JSONProtocol
OUTPUT_PROTOCOL = mrjob.protocol.JSONProtocol
```

The default input protocol is RawValueProtocol, which just reads in a line as a str. (The line won't have a trailing newline character because MRJob strips it.) So by default, the first step in your job sees (None, line) for each line of input <sup>3</sup>.

The default output and internal protocols are both JSONProtocol <sup>4</sup>, which reads and writes JSON strings separated by a tab character. (By default, Hadoop Streaming uses the tab character to separate keys and values within one line when it sorts your data.)

If your head hurts a bit, think of it this way: use <code>RawValueProtocol</code> when you want to read or write lines of raw text. Use <code>JSONProtocol</code> when you want to read or write key-value pairs where the key and value are <code>JSON-enoded</code> bytes.

**Note:** Hadoop Streaming does not understand JSON, or mrjob protocols. It simply groups lines by doing a string comparison on whatever comes before the first tab character.

See *mrjob.protocol* for the full list of protocols built-in to mrjob.

#### Data flow walkthrough by example

Let's revisit our example from *Multi-step jobs*. It has two steps and takes a plain text file as input.

The first step starts with mapper get words ():

```
def mapper_get_words(self, _, line):
    # yield each word in the line
    for word in WORD_RE.findall(line):
        yield (word.lower(), 1)
```

Since the input protocol is RawValueProtocol, the key will always be None and the value will be the text of the line.

The function discards the key and yields (word, 1) for each word in the line. Since the internal protocol is <code>JSONProtocol</code>, each component of the output is serialized to JSON. The serialized components are written to stdout separated by a tab character and ending in a newline character, like this:

```
"mrjob" 1
"is"    1
"a" 1
"python"    1
```

<sup>&</sup>lt;sup>3</sup> Experienced Pythonistas might notice that a str is a bytestring on Python 2, but Unicode on Python 3. That's right! RawValueProtocol is an alias for one of two different protocols depending on your Python version.

<sup>&</sup>lt;sup>4</sup> JSONProtocol is an alias for one of four different implementations; we try to use the (much faster) ujson library if it is available, and if not, rapidjson or simplejson before falling back to the built-in json implementation.

The next two parts of the step are the combiner and reducer:

```
def combiner_count_words(self, word, counts):
    # sum the words we've seen so far
    yield (word, sum(counts))

def reducer_count_words(self, word, counts):
    # send all (num_occurrences, word) pairs to the same reducer.
    # num_occurrences is so we can easily use Python's max() function.
    yield None, (sum(counts), word)
```

In both cases, bytes are describilized into (word, counts) by *JSONProtocol*, and the output is serialized as JSON in the same way (because both are followed by another step). It looks just like the first mapper output, but the results are summed:

```
"mrjob" 31
"is"    2
"a" 2
"Python"    1
```

The final step is just a reducer:

```
# discard the key; it is just None
def reducer_find_max_word(self, _, word_count_pairs):
    # each item of word_count_pairs is (count, word),
    # so yielding one results in key=counts, value=word
    yield max(word_count_pairs)
```

Since all input to this step has the same key (None), a single task will get all rows. Again, JSONProtocol will handle descrialization and produce the arguments to reducer\_find\_max\_word().

The output protocol is also JSONProtocol, so the final output will be:

```
31 "mrjob"
```

And we're done! But that's a bit ugly; there's no need to write the key out at all. Let's use <code>JSONValueProtocol</code> instead, so we only see the JSON-encoded value:

```
class MRMostUsedWord(MRJob):
    OUTPUT_PROTOCOL = JSONValueProtocol
```

Now we should have code that is identical to examples/mr\_most\_used\_word.py in mrjob's source code. Let's try running it (-q prevents debug logging):

```
$ python mr_most_used_word.py README.txt -q
"mrjob"
```

Hooray!

#### Specifying protocols for your job

Usually, you'll just want to set one or more of the class variables <code>INPUT\_PROTOCOL</code>, <code>INTERNAL\_PROTOCOL</code>, and <code>OUTPUT\_PROTOCOL</code>:

```
class BasicProtocolJob(MRJob):
    # get input as raw strings
    INPUT_PROTOCOL = RawValueProtocol
```

```
# pass data internally with pickle
INTERNAL_PROTOCOL = PickleProtocol
# write output as JSON
OUTPUT_PROTOCOL = JSONProtocol
```

If you need more complex behavior, you can override <code>input\_protocol()</code>, <code>internal\_protocol()</code>, or <code>output\_protocol()</code> and return a protocol object instance. Here's an example that sneaks a peek at <code>Defining command line options</code>:

Finally, if you need to use a completely different concept of protocol assignment, you can override pick\_protocols():

```
class WhatIsThisIDontEvenProtocolJob(MRJob):
    def pick_protocols(self, step_num, step_type):
        return random.choice([Protocololol, ROFLcol, Trolltocol, Locotorp])
```

#### Writing custom protocols

A protocol is an object with methods read(self, line) and write(self, key, value). The read() method takes a bytestring and returns a 2-tuple of decoded objects, and write() takes the key and value and returns bytes to be passed back to Hadoop Streaming or as output.

Protocols don't have to worry about adding or stripping newlines; this is handled automatically by MRJob.

Here is a simplified version of mrjob's JSON protocol:

```
import json

class JSONProtocol(object):

    def read(self, line):
        k_str, v_str = line.split('\t', 1)
        return json.loads(k_str), json.loads(v_str)

    def write(self, key, value):
        return '%s\t%s' % (json.dumps(key), json.dumps(value))
```

You can improve performance significantly by caching the serialization/deserialization results of keys. Look at the source code of <code>mrjob.protocol</code> for an example.

## 1.4.3 Passing entire files to the mapper

New in version 0.6.3.

Sometimes you need to read binary data (e.g. image files), or text-based data that has records longer than one line.

By using <code>mapper\_raw()</code>, you can pass entire files to your mapper, and read them however you want. Each mapper gets one file, and is passed both the path of a local copy of the file, and the URI where the original file is located on Hadoop's filesystem.

For example, if you want to read .wet files from Common Crawl data, you could handle them like this:

```
class MRCrawler(MRJob):
    def mapper_raw(self, wet_path, wet_uri):
        from warcio.archiveiterator import ArchiveIterator
    with open(wet_path, 'rb') as f:
        for record in ArchiveIterator(f):
        ...
```

To use a library like warcio, you'll need to ensure that it gets installed on your cluster. See *Using a virtualenv* for one way to do this.

Under the hood, mrjob is passes an input manifest (a list of URIs of input files) to Hadoop, and instructs Hadoop to send one line to each mapper. In most cases, this should be seamless, even to the point of telling you which file was being read when a task fails.

**Warning:** For all runners except EMR, mrjob uses **hadoop fs** to download files to the local filesystem, which means Hadoop has to invoke itself. If your cluster has tightly tuned memory requirements, this can sometimes cause an out-of-memory error.

# 1.4.4 Jar steps

You can run Java directly on Hadoop (bypassing Hadoop Streaming) by using <code>JarStep</code> instead of <code>MRStep()</code>.

For example, on EMR you can use a jar to run a script:

More interesting is combining MRStep and JarStep in the same job. Use mrjob.step.INPUT and mrjob.step.OUTPUT in args to stand for the input and output paths for that step. For example:

```
args=['naive-bayes', INPUT, OUTPUT]
)
]
```

Changed in version 0.6.6: mrjob no longer passes hadoop generic args (-D and -libjars) to JarSteps. If you want them, add mrjob.step.GENERIC\_ARGS to your JarStep's args, and mrjob will automatically interpolate them.

JarStep has no concept of *Protocols*. If your jar reads input from a *MRStep*, or writes input read by another *MRStep*, it is up to those steps to read and write data in the format your jar expects.

If you are writing the jar yourself, the easiest solution is to have it read and write mrjob's default protocol (lines containing two JSONs, separated by a tab).

If you are using a third-party jar, you can set custom protocols for the steps before and after it by overriding pick\_protocols().

**Warning:** If the first step of your job is a *JarStep* and you pass in multiple input paths, mrjob will replace *INPUT* with the input paths joined together with a comma. Not all jars can handle this!

Best practice in this case is to put all your input into a single directory and pass that as your input path.

# 1.4.5 Using other python modules and packages

New in version 0.6.4.

If you want to run Python code outside of the file containing your MRJob, you'll to make sure that code gets uploaded to Hadoop.

The easiest way to do this is with by setting the *DIRS* attribute in your job. Put the code you want to import in one or more packages (directories with an \_\_init\_\_.py file), and point *DIRS* at them:

```
class MRPackageUsingJob(MRJob):
   DIRS = ['mycode', '../someothercode']
   ...
```

And then import code from inside a mapper or reducer:

```
def mapper(self, key, value):
    from mycode.custom import important_business_logic
    from someotherlibrary import util_function
    ...
```

(If you want to import code from the top level of your script rather than inside a method, make sure it's in your PYTHONPATH, just like with any other code.)

DIRS is relative to the directory your script is in (not the current working directory). This works inside Hadoop because the current working directory is the same as the directory your script is in.

If you want to access individual Python modules or other support code, you can use FILES to upload them to your job's working directory inside Hadoop:

```
class MRFileUsingJob(MRJob):

FILES = ['mymodule.py', '../data/zipcodes.db']

def mapper(self, key, value):
    from mymodule import open_zipcode_db
```

```
with open_zipcode_db('zipcodes.db') as db:
...
```

For jobs with more complex dependencies (e.g. code that needs to be compiled), you may need to use the setup option. See Job Environment Setup Cookbook for more information.

## 1.4.6 Defining command line options

Recall from *How your program is run* that your script is executed in several contexts: once for the initial invocation, and once for each task. If you just add an option to your job's option parser, that option's value won't be propagated to other runs of your script. Instead, you can use mrjob's option API: <code>add\_passthru\_arg()</code> and <code>add\_file\_arg()</code>.

### Passthrough options

A *passthrough option* is an argparse option that mrjob is aware of. mrjob inspects the value of the option when you invoke your script and reproduces that value when it invokes your script in other contexts. The command line-switchable protocol example from before uses this feature:

When you run your script with —output-format=json, mrjob detects that you passed —output-format on the command line. When your script is run in any other context, such as on Hadoop, it adds —output-format=json to its command string.

 $add\_passthru\_arg \textit{()} takes the same arguments as \texttt{argparse.ArgumentParser.add\_argument().} For more information, see the argparse docs.$ 

#### Passing through existing options

Occasionally, it'll be useful for mappers, reducers, etc. to be able to see the value of other command-line options. For this, use <code>pass\_arg\_through()</code> with the corresponding command-line switch.

For example, you might wish to fetch supporting data for your job from different locations, depending on whether your job is running on EMR or locally:

```
class MRRunnerAwareJob(MRJob):
    def configure_args(self):
        super(MRRunnerAwareJob, self).configure_args()
        self.pass_arg_through('--runner')
```

```
def mapper_init(self):
    if self.options.runner == 'emr':
        self.data = ... # load from S3
    else:
        self.data = ... # load from local FS
```

**Note:** Keep in mind that self.options.runner (and the values of most options) will be None unless the user explicitly set them with a command-line switch.

### File options

A *file option* is like a passthrough option, but:

- 1. Its value must be a string or list of strings (action="store" or action="append"), where each string represents either a local path, or an HDFS or S3 path that will be accessible from the task nodes.
- 2. That file will be downloaded to each task's local directory and the value of the option will magically be changed to its path.

For example, if you had a map task that required a sqlite3 database, you could do this:

```
class SqliteJob(MRJob):

   def configure_args(self):
       super(SqliteJob, self).configure_args()
       self.add_file_arg('--database')

   def mapper_init(self):
       # make sqlite3 database available to mapper
       self.sqlite_conn = sqlite3.connect(self.options.database)
```

You could call it any of these ways, depending on where the file is:

In any of these cases, when your task runs, my\_db.sqlite3 will always be available in the task's working directory, and the value of self.options.database will always be set to its path.

See *Making files available to tasks* if you want to upload a file to your tasks' working directories without writing a custom command line option.

**Warning:** You **must** wait to read files until **after class initialization**. That means you should use the \*\_init() methods to read files. Trying to read files into class variables will not work.

## 1.4.7 Counters

Hadoop lets you track *counters* that are aggregated over a step. A counter has a group, a name, and an integer value. Hadoop itself tracks a few counters automatically. mrjob prints your job's counters to the command line when your job finishes, and they are available to the runner object if you invoke it programmatically.

To increment a counter from anywhere in your job, use the <code>increment\_counter()</code> method:

```
class MRCountingJob(MRJob):

   def mapper(self, _, value):
       self.increment_counter('group', 'counter_name', 1)
       yield _, value
```

At the end of your job, you'll get the counter's total value:

```
group:
    counter_name: 1
```

## 1.4.8 Input and output formats

Input and output formats are Java classes that determine how your job interfaces with data on Hadoop's filesystem(s).

Suppose we wanted to write a word frequency count job that wrote output into a separate directory based on the first letter of the word counted (a/part-\*, b/part-\*, etc.). We could accomplish this by using the MultipleValueOutputFormat class from the Open Source project nicknack.

First, we need to tell our job to use the custom output format by setting HADOOP\_OUTPUT\_FORMAT in our job class:

```
HADOOP_OUTPUT_FORMAT = 'nicknack.MultipleValueOutputFormat'
```

The output format class is part of a custom JAR, so we need to make sure that this JAR gets included in Hadoop's classpath. First download the jar to the same directory as your script, and add its name to LIBJARS:

```
LIBJARS = ['nicknack-1.0.0.jar']
```

(You can skip this step if you're using a format class that's built into Hadoop.)

Finally, output your data the way that your output format expects. MultipleValueOutputFormat expects the subdirectory name, followed by a tab, followed the actual line to write into the file.

First, we need to take direct control of how the job writes output by setting <code>OUTPUT\_PROTOCOL</code> to <code>RawValueProtocol</code>:

```
OUTPUT_PROTOCOL = RawValueProtocol
```

Then we need to format the line accordingly. In this case, let's continue output our final data in the standard format (two JSONs separated by a tab):

```
def reducer(self, word, counts):
   total = sum(counts)
   yield None, '\t'.join([word[0], json.dumps(word), json.dumps(total)])
```

Done! Here's the full, working job (this is mrjob.examples.mr\_nick\_nack):

```
import json
import re

from mrjob.job import MRJob
from mrjob.protocol import RawValueProtocol

WORD_RE = re.compile(r"[A-Za-z]+")

class MRNickNack(MRJob):

    HADOOP_OUTPUT_FORMAT = 'nicknack.MultipleValueOutputFormat'
```

```
LIBJARS = ['nicknack-1.0.0.jar']

OUTPUT_PROTOCOL = RawValueProtocol

def mapper(self, _, line):
    for word in WORD_RE.findall(line):
        yield (word.lower(), 1)

def reducer(self, word, counts):
    total = sum(counts)
    yield None, '\t'.join([word[0], json.dumps(word), json.dumps(total)])

if __name__ == '__main__':
    MRNickNack.run()
```

Input formats work the same way; just set <code>HADOOP\_INPUT\_FORMAT</code>. (You usually won't need to set <code>INPUT\_PROTOCOL</code> because it already defaults to <code>RawValueProtocol</code>.)

## 1.5 Runners

While the MRJob class is the part of the framework that handles the execution of your code in a MapReduce context, the **runner** is the part that packages and submits your job to be run, and reporting the results back to you.

In most cases, you will interact with runners via the command line and configuration files. When you invoke mrjob via the command line, it reads your command line options (the --runner parameter) to determine which type of runner to create. Then it creates the runner, which reads your configuration files and command line args and starts your job running in whatever context you chose.

Most of the time, you won't have any reason to construct a runner directly. Instead you'll invoke your Python script on the command line and it will make a runner automatically, you'll call mrjob run my\_script to have the mrjob command build a runner for your script (which may or may not be Python), or you'll write some sort of wrapper that calls my\_job.make\_runner().

Internally, the general order of operations is:

- Get a runner by calling make\_runner() on your job
- Call run () on your runner. This will:
  - Run your job with --steps to find out how many mappers/reducers to run
  - Copy your job and supporting files to Hadoop
  - Instruct Hadoop to run your job with the appropriate --mapper, --combiner, --reducer, and --step-num arguments

Each runner runs a single job once; if you want to run a job multiple times, make multiple runners.

Subclasses: DataprocJobRunner, EMRJobRunner, HadoopJobRunner, InlineMRJobRunner, LocalMRJobRunner

# 1.5.1 Testing locally

To test the job locally, just run:

1.5. Runners 23

```
python your_mr_job_sub_class.py < log_file_or_whatever > output
```

The script will automatically invoke itself to run the various steps, using InlineMRJobRunner (--runner=inline). If you want to simulate Hadoop more closely, you can use --runner=local, which doesn't add your working directory to the PYTHONPATH, sets a few Hadoop environment variables, and uses multiple subprocesses for tasks.

You can also run individual steps:

```
# test 1st step mapper:
python your_mr_job_sub_class.py --mapper
# test 2nd step reducer (step numbers are 0-indexed):
python your_mr_job_sub_class.py --reducer --step-num=1
```

By default, we read from stdin, but you can also specify one or more input files. It automatically decompresses .gz and .bz2 files:

```
python your_mr_job_sub_class.py log_01.gz log_02.bz2 log_03
```

See mrjob.examples for more examples.

# 1.5.2 Running on your own Hadoop cluster

- Set up a hadoop cluster (see http://hadoop.apache.org/docs/current/)
- Run your job with -r hadoop:

```
python your_mr_job_sub_class.py -r hadoop < input > output
```

**Note:** You don't need to install mrjob or any other libraries on the nodes of your Hadoop cluster, but they *do* at least need a version of Python that's compatible with your job.

# 1.5.3 Running on EMR

- Set up your Amazon account and credentials (see Configuring AWS credentials)
- Run your job with -r emr:

```
python your_mr_job_sub_class.py -r emr < input > output
```

## 1.5.4 Running on Dataproc

- Set up your Google account and credentials (see Getting started with Google Cloud)
- Run your job with -r dataproc:

```
python your_mr_job_sub_class.py -r dataproc < input > output
```

**Note:** Dataproc does not yet support Spark or libjars.

## 1.5.5 Configuration

Runners are configured by several methods:

- from mrjob.conf (see Config file format and location)
- from the command line
- by re-defining job\_runner\_kwargs() etc in your MRJob (see Job runner configuration)
- by instantiating the runner directly

In most cases, you should put all configuration in mrjob.conf and use the command line args or class variables to customize how individual jobs are run.

# 1.5.6 Running your job programmatically

It is fairly common to write an organization-specific wrapper around mrjob. Use <code>make\_runner()</code> to run an <code>MRJob</code> from another Python script. The context manager guarantees that all temporary files are cleaned up regardless of the success or failure of your job.

This pattern can also be used to write integration tests (see Testing jobs).

```
mr_job = MRWordCounter(args=['-r', 'emr'])
with mr_job.make_runner() as runner:
    runner.run()
    for key, value in mr_job.parse_output(runner.cat_output()):
        ... # do something with the parsed output
```

You instantiate the MRJob, use a context manager to create the runner, run the job, and cat its output, parsing that output with the job's output protocol.

parse\_output() and cat\_output() were introduced in version 0.6.0. In previous versions of mrjob, you'd iterate line by line instead, like this:

```
for line in runner.stream_output():
    key, value = mr_job.parse_output_line(line)
    ... # do something with the parsed output
```

#### Further reference:

- make runner()
- run()
- parse\_output()
- cat\_output()
- parse\_output\_line()
- stream\_output()

#### Limitations

**Note:** You should pay attention to the next sentence.

You cannot use the programmatic runner functionality in the same file as your job class. As an example of what not to do, here is some code that does not work.

1.5. Runners 25

Warning: The code below shows you what **not** to do.

```
from mrjob.job import MRJob

class MyJob(MRJob):
    # (your job)

# no, stop, what are you doing?!?!
mr_job = MyJob(args=[args])
with mr_job.make_runner() as runner:
    runner.run()
    # ... etc
```

If you try to do this, mrjob will give you an error message similar or identical to this one:

```
UsageError: make_runner() was called with --steps. This probably means you tried to use it from __main__, which doesn't work.
```

What you need to do instead is put your job in one file, and your run code in another. Here are two files that would correctly handle the above case.

```
# job.py
from mrjob.job import MRJob

class MyJob(MRJob):
    # (your job)

if __name__ == '__main__':
    MyJob.run()
```

```
# run.py
from job import MyJob
mr_job = MyJob(args=[args])
with mr_job.make_runner() as runner:
    runner.run()
# ... etc
```

#### Why can't I put the job class and run code in the same file?

The file with the job class is sent to Hadoop to be run. Therefore, the job file cannot attempt to start the Hadoop job, or you would be recursively creating Hadoop jobs!

The code that runs the job should only run *outside* of the Hadoop context.

The if \_\_name\_\_ == '\_\_main\_\_' block is only run if you invoke the job file as a script. It is not run when imported. That's why you can import the job class to be run, but it can still be invoked as an executable.

#### **Counters**

Counters may be read through the *counters* () method on the runner. The example below demonstrates the use of counters in a test case.

```
mr_counting_job.py
```

```
from mrjob.job import MRJob
from mrjob.step import MRStep
```

test\_counters.py

# 1.6 Spark

# 1.6.1 Why use mrjob with Spark?

mrjob augments Spark's native Python support with the following features familiar to users of mrjob:

- automatically upload input and other support files to HDFS, GCS, or S3 (see upload\_files, upload\_archives, and py\_files)
- run make and other command before running Spark tasks (see setup).
- passthrough and file arguments (see *Defining command line options*)
- automatically parse logs to explain errors and other Spark job failures
- easily pass through environment variables (see cmdenv)

1.6. Spark 27

- support for libjars
- automatic matching of Python version (see python\_bin)
- automatically set up Spark on EMR (see bootstrap\_spark)
- automatically making the mrjob library available to your job (see bootstrap\_mrjob)

## 1.6.2 mrjob spark-submit

If you already have a Spark script written, the easiest way to access mrjob's features is to run your job with mrjob spark-submit, just like you would normally run it with spark-submit. This can, for instance, make running a Spark job on EMR as easy as running it locally, or allow you to access features (e.g. setup) not natively supported by Spark.

For more details, see *mrjob spark-submit*.

## 1.6.3 Writing your first Spark MRJob

Another way to integrate mrjob with Spark is to add a *spark()* method to your *MRJob* class, and put your Spark code inside it. This will allow you to access features only available to MRJobs (e.g. *FILES*).

Here's how you'd implement a word frequency count job in Spark:

```
import re
from operator import add
from mrjob.job import MRJob
WORD_RE = re.compile(r"[\w']+")
class MRSparkWordcount (MRJob) :
    def spark(self, input_path, output_path):
        # Spark may not be available where script is launched
        from pyspark import SparkContext
        sc = SparkContext(appName='mrjob Spark wordcount script')
        lines = sc.textFile(input_path)
        counts = (
            lines.flatMap(self.get_words)
            .map(lambda word: (word, 1))
            .reduceByKey(add))
        counts.saveAsTextFile(output_path)
        sc.stop()
    def get_words(self, line):
        return WORD_RE.findall(line)
if __name__ == '__main__':
    MRSparkWordcount.run()
```

Since Spark already supports Python, mrjob takes care of setting up your cluster, passes in input and output paths, and otherwise gets out of the way. If you pass in multiple input paths, <code>input\_path</code> will be these paths joined by a comma (SparkContext.textFile() will accept this).

Note that pyspark is imported *inside* the *spark()* method. This allows your job to run whether pyspark is installed locally or not.

The *spark()* method can be used to execute arbitrary code, so there's nothing stopping you from using *SparkSession* instead of *SparkContext* in Spark 2, or writing a streaming-mode job rather than a batch one.

**Warning:** Prior to v0.6.8, to pass job methods into Spark (e.g. rdd.flatMap(self.get\_words)), you first had to call self.sandbox(); otherwise Spark would error because self was not serializable.

## 1.6.4 Running on your Spark cluster

By default, mrjob runs your job on the inline runner (see below). If you want to run your job on your own Spark cluster, run it with -r spark:

Use --spark-master (see spark\_master) to control where your job runs.

You can pass in spark options with -D (see jobconf) and set deploy mode (client or cluster) with --spark-deploy-mode. If you need to pass other arguments to **spark-submit**, use spark\_args.

The Spark runner can also run "classic" MRJobs (i.e. those made by defining <code>mapper()</code> etc. or with <code>MRSteps</code>) directly on Spark, allowing you to move off Hadoop without rewriting your jobs. See <code>below</code> for details.

**Warning:** If you don't set spark\_master, your job will run on Spark's default local[\*] master, which can't handle setup scripts or --files because it doesn't give tasks their own working directory.

**Note:** mrjob needs to know what master and deploy mode you're using, so it will override attempts to set spark master or deploy mode through jobconf (e.g. -D spark.master=...).

## 1.6.5 Using remote filesystems other than HDFS

By default, if you use a remote Spark master (i.e. not local or local-cluster), Spark will assume you want to use HDFS for your job's temp space, and that you will want to access it through **hadoop fs**.

Some Spark installations don't use HDFS at all. Fortunately, the Spark runner also supports S3 and GCS. Use spark\_tmp\_dir to specify a remote temp directory not on HDFS (e.g. --spark-tmp-dir s3a://bucket/path).

For more information on accessing S3 or GCS, see *Configuring AWS credentials* (S3) or *Configuring Google Cloud credentials* (GCS).

## 1.6.6 Other ways to run on Spark

#### Inline runner

Running your Spark job with -r inline (the default) will launch it directly through the pyspark library, effectively running it on the local[\*] master. This is convenient for debugging because exceptions will bubble up directly to your Python process.

1.6. Spark 29

The inline runner also builds a simulated working directory for your job, making it possible to test scripts that rely on certain files being in the working directory (it doesn't run setup scripts).

**Note:** If you don't have a local Spark installation, the pyspark library on PyPI is a pretty quick way to get one (pip install pyspark).

#### Local runner

Running your Spark job with -r local will launch it through **spark-submit** on a local-cluster master. local-cluster is designed to simulate a real Spark cluster, so setup will work as expected.

By default, the local runner launches Spark jobs with as many executors as your system has CPUs. Use --num-cores (see num\_cores to change this).

By default, the local runner gives each executor 1 GB of memory. If you need more, you can specify it through jobconf, e.g. -D spark.core.memory=4g.

#### **EMR** runner

Running your Spark job with -r emr will launch it in Amazon Elastic MapReduce (EMR), with the same seamless integration and features mrjob provides for Hadoop jobs on EMR.

The EMR runner will always run your job on the yarn Spark master in cluster deploy mode.

## **Hadoop runner**

Running your Spark job with -r hadoop will launch it on your own Hadoop cluster. This is not significantly different than the Spark runner. The main advantage of the Hadoop runner is that is has more knowledge about how to find logs and can be better at finding the relevant error if your job fails.

Unlike the Spark runner, the Hadoop runner's default spark master is yarn.

**Note:** mrjob does not yet support Spark on Google Cloud Dataproc.

## 1.6.7 Passing in libraries

Use --py-files to pass in .zip or .egg files full of Python code:

```
python your_mr_spark_job -r hadoop --py-files lib1.zip,lib2.egg
```

Or set py\_files in mrjob.conf.

## 1.6.8 Command-line options

Command-line options (passthrough options, etc) work exactly like they do with regular streaming jobs (even add\_file\_arg() on the local[\*] Spark master. See *Defining command line options*.

# 1.6.9 Uploading files to the working directory

upload\_files, FILES, and files uploaded via setup scripts all should work as expected (except on local masters because there is no working directory).

Note that you can give files a different name in the working directory (e.g. --files foo#bar) on all Spark masters, even though Spark treats that as a YARN-specific feature.

#### 1.6.10 Archives and directories

Spark treats ——archives as a YARN-specific feature. This means that upload\_archives, ARCHIVES, DIRS, etc. will be ignored on non-yarn Spark masters.

Future versions of mrjob may simulate archives on non-yarn masters using a setup script.

## 1.6.11 Multi-step jobs

There generally isn't a need to define multiple Spark steps (Spark lets you map/reduce as many times as you want). However, it may sometimes be useful to pre- or post-process Spark data using a *streaming* or *jar* step.

This is accomplished by overriding your job's steps () method and using the SparkStep class:

```
def steps():
    return [
        MRStep(mapper=self.preprocessing_mapper),
        SparkStep(spark=self.spark),
    ]
```

## 1.6.12 External Spark scripts

mrjob can also be used to launch external (non-mrjob) Spark scripts using the SparkScriptStep class, which specifies the path (or URI) of the script and its arguments.

As with <code>JarSteps</code>, you can interpolate input and output paths using <code>INPUT</code> and <code>OUTPUT</code> constants. For example, you could set your job's <code>steps()</code> method up like this:

# 1.6.13 Custom input and output formats

mrjob allows you to use input and output formats from custom JARs with Spark, just like you can *with streaming jobs*. First download your JAR to the same directory as your job, and add it to your job class with the LIBJARS attribute:

```
LIBJARS = ['nicknack-1.0.0.jar']
```

1.6. Spark 31

Then use Spark's own capabilities to reference your input or output format, keeping in mind the data types they expect.

For example, nicknack's MultipleValueOutputFormat expects <Text>, so if we wanted to integrate it with our wordcount example, we'd have to convert the count to a string:

# 1.6.14 Running "classic" MRJobs on Spark

The Spark runner provides near-total support for running "classic" MRJobs (the sort described in Writing your first job and Writing your second job) directly on any Spark installation, even though these jobs were originally designed to run on Hadoop Streaming. Support includes:

- \*\_init() and \*\_final() methods
- HADOOP\_INPUT\_FORMAT and HADOOP\_OUTPUT\_FORMAT
- SORT\_VALUES
- passthrough arguments
- increment\_counter()

Jobs will often run more quickly on Spark than Hadoop Streaming, so it's worth trying even if you don't plan to move off Hadoop in the forseeable future.

#### Multiple steps are run as a single job

If you have a job with multiple consecutive *MRSteps*, the Spark runner will run them all as a single Spark job. This is usually what you want (more efficient), but it can make debugging slightly more challenging (step failure exceptions give a range of steps, no way to access intermediate data).

To force the Spark runner to run steps separately, you can initialize each MRStep with a different jobconf dictionary.

#### No support for subprocesses

Pre-filters (e.g. mapper\_pre\_filter()) and command steps (e.g. reducer\_cmd()) are not supported because they require launching subprocesses.

It wouldn't be *impossible* to emulate this inside Spark, but then we'd essentially be turning Spark *into* Hadoop Streaming. (If you have a use case for this seemingly implausible feature, let us know through GitHub.)

### **Spark loves combiners**

Hadoop's "reduce" paradigm is a lot more heavyweight than Spark's; whereas a Spark reducer just wants to know how to combine two values into one, a Hadoop reducer expects to be able to see all the values for a given key, and to emit zero or more key-value pairs.

In fact, Spark reducers are a lot more like Hadoop combiners. The Spark runner knows how to translate something like:

```
def combiner(self, key, values):
    yield key, sum(values)
```

into Spark's reduce paradigm-basically it'll pass your combiner two values at a time, and hope it emits one. If your combiner does *not* behave like a Spark reducer function (emitting multiple or zero values), the Spark runner handles that gracefully as well.

### Counter emulation is almost perfect

Counters (see <code>increment\_counter()</code>) are a feature specific to Hadoop. mrjob emulates them on Spark anyway. If you have a multi-step job, mrjob will dutifully print out counters for each step and make them available through <code>counters()</code>.

The only drawback is that while Hadoop has the ability to "take back" counters produced by a failed task, there isn't a clean way to do this with Spark accumulators. Therefore, the counters produced by the Spark runner's Hadoop emulation may be overestimates.

#### Spark does not stream data

While Hadoop streaming (as its name implies) passes a stream of data to your job, Spark instead operates on *partitions*, which are loaded into memory.

A reducer like this can't run out of memory on Hadoop streaming, no matter how many values there are for key:

```
def reducer(self, key, values):
    yield key, sum(values)
```

However, on Spark, simply storing the partition that contains these values can cause Spark to run out of memory.

If this happens, you can let Spark use more memory (-D spark.executor.memory=10g) or add a combiner to your job.

#### **Compression emulation**

It's fairly common for people to request compressed output from Hadoop via configuration properties, for example:

```
python mr_your_job.py -D mapreduce.output.fileoutputformat.compress=true -D\
mapreduce.output.fileoutputformat.compress.codec=org.apache.hadoop.io.compress.BZip2Codec ...
```

This works with -r spark too; the Spark runner knows how to recognize these properties and pass the codec specified to Spark when it writes output.

1.6. Spark 33

## Spark won't split .gz files either

A common trick on Hadoop to ensure that segments of your data don't get split between mappers is to gzip each segment (since .gz is not a seekable compression format).

This works on Spark as well.

## Controlling number of output files

By default, Spark will write one output file per partition. This may give more output files than you expect, since Hadoop and Spark are tuned differently.

The Spark runner knows how to emulate the Hadoop configuration property that sets number of reducers on Hadoop (e.g. -D mapreduce.job.reduces=100), which will control the number of output files (assuming your last step has a reducer).

However, this is a somewhat heavyweight solution; once Spark runs a step's reducer, mrjob has to forbid Spark from re-partitioning until the end of the step.

A lighter weight solution is --max-output-files, allows you to limit the number of output files by running coalesce() just before writing output. Running your job with --max-output-files=100 would ensure it produces no more than 100 output files (but it could output less).

## Running classic MRJobs on Spark on EMR

It's often faster to run classic MRJobs on Spark than Hadoop Streaming. It's also convenient to be able to run on EMR rather than setting up your own Spark cluster (or SSH'ing in).

Can you do both? Yes! Run the job with the Spark runner, but tell it to use mrjob spark-submit to launch Spark jobs on EMR.

It looks something like this:

```
python mr_your_job.py -r spark \
   --spark-submit-bin 'mrjob spark-submit -r emr' \
   --spark-master yarn --spark-tmp-dir s3://your-bucket/tmp/ input1 input2
```

Note that because the Spark runner itself doesn't know the job is going to run on EMR, you have to give it a couple of hints so that it knows it's running on YARN (**--spark-master**) and that it needs to use S3 as its temp space (**--spark-tmp-dir**).

# 1.7 Config file format and location

We look for mrjob.conf in these locations:

- The location specified by MRJOB\_CONF
- ~/.mrjob.conf
- /etc/mrjob.conf

You can specify one or more configuration files with the --conf-path flag. See Options available to all runners for more information.

The point of mrjob.conf is to let you set up things you want every job to have access to so that you don't have to think about it. For example:

• libraries and source code you want to be available for your jobs

- where temp directories and logs should go
- · security credentials

mrjob.conf is just a YAML- or JSON-encoded dictionary containing default values to pass in to the constructors of the various runner classes. Here's a minimal mrjob.conf:

```
runners:
    emr:
    cmdenv:
    TZ: America/Los_Angeles
```

Now whenever you run mr\_your\_script.py -r emr, *EMRJobRunner* will automatically set TZ to America/Los\_Angeles in your job's environment when it runs on EMR.

If you don't have the yaml module installed, you can use JSON in your mrjob.conf instead (JSON is a subset of YAML, so it'll still work once you install yaml). Here's how you'd render the above example in JSON:

```
{
    "runners": {
        "emr": {
            "cmdenv": {
                "TZ": "America/Los_Angeles"
            }
        }
     }
}
```

# 1.7.1 Precedence and combining options

Options specified on the command-line take precedence over mrjob.conf. Usually this means simply overriding the option in mrjob.conf. However, we know that *cmdenv* contains environment variables, so we do the right thing. For example, if your mrjob.conf contained:

```
runners:
    emr:
    cmdenv:
        PATH: /usr/local/bin
        TZ: America/Los_Angeles
```

and you ran your job as:

```
mr_your_script.py -r emr --cmdenv TZ=Europe/Paris --cmdenv PATH=/usr/sbin
```

We'd automatically handle the PATH variables and your job's environment would be:

```
{'TZ': 'Europe/Paris', 'PATH': '/usr/sbin:/usr/local/bin'}
```

What's going on here is that *cmdenv* is associated with combine\_envs(). Each option is associated with an appropriate combiner function that that combines options in an appropriate way.

Combiner functions can also do useful things like expanding environment variables and globs in paths. For example, you could set:

```
runners:
   local:
    upload_files: &upload_files
     - $DATA_DIR/*.db
   hadoop:
    upload_files: *upload_files
```

```
emr:
    upload_files: *upload_files
```

and every time you ran a job, every job in your . db file in \$DATA\_DIR would automatically be loaded into your job's current working directory.

Also, if you specified additional files to upload with --file, those files would be uploaded in addition to the .db files, rather than instead of them.

See Configuration quick reference for the entire dizzying array of configurable options.

# 1.7.2 Option data types

The same option may be specified multiple times and be one of several data types. For example, the AWS region may be specified in mrjob.conf, in the arguments to EMRJobRunner, and on the command line. These are the rules used to determine what value to use at runtime.

Values specified "later" refer to an option being specified at a higher priority. For example, a value in mrjob.conf is specified "earlier" than a value passed on the command line.

When there are multiple values, they are "combined with" a *combiner function*. The combiner function for each data type is listed in its description.

### Simple data types

When these are specified more than once, the last non-None value is used.

**String** Simple, unchanged string. Combined with combine\_values().

**Command** String containing all ASCII characters to be parsed with shlex.split(), or list of command + arguments. Combined with combine\_cmds().

Path Local path with ~ and environment variables (e.g. \$TMPDIR) resolved. Combined with combine\_paths ().

### List data types

The values of these options are specified as lists. When specified more than once, the lists are concatenated together.

String list List of strings. Combined with combine\_lists().

**Path list** List of *paths*. Combined with *combine\_path\_lists()*.

Strings and non-sequence data types (e.g. numbers) are treated as single-item lists.

For example,

```
runners:
    emr:
    setup: /run/some/command with args
```

is equivalent to:

```
runners:
    emr:
    setup:
        - /run/some/command with args
```

### Dict data types

The values of these options are specified as dictionaries. When specified more than once, each has custom behavior described below.

Plain dict Values specified later override values specified earlier. Combined with combine\_dicts().

#### JobConf Dicts

New in version 0.6.6: Like plain dicts except that non-string values are converted into a format that Java understands. For example, the boolean value true here:

```
jobconf:
   mapreduce.output.fileoutputformat.compress: true
```

gets passed through to Hadoop in Java format (true), not Python format (True).

Keys whose values are None are not passed to Hadoop at all.

**Warning:** Prior to version 0.6.6, you should use "true" and "false", for boolean jobconf values in config files, not true and false.

Environment variable dict Values specified later override values specified earlier, except for those with keys ending in PATH, in which values are concatenated and separated by a colon (:) rather than overwritten. The later value comes first.

For example, this config:

```
runners:
    emr:
    cmdenv:
        PATH: /usr/bin
```

when run with this command:

```
python my_job.py --cmdenv PATH=/usr/local/bin
```

will result in the following value of cmdenv:

```
/usr/local/bin:/usr/bin
```

The function that handles this is combine\_envs().

The one exception to this behavior is in the local runner, which uses the local system separator (on Windows;, on everything else still:) instead of always using: In local mode, the function that combines config values is <code>combine\_local\_envs()</code>.

# 1.7.3 Using multiple config files

If you have several standard configurations, you may want to have several config files "inherit" from a base config file. For example, you may have one set of AWS credentials, but two code bases and default instance sizes. To accomplish this, use the include option:

```
~/mrjob.very-large.conf:
```

```
include: ~/.mrjob.base.conf
runners:
    emr:
        num_core_instances: 20
        core_instance_type: m1.xlarge
```

```
~/mrjob.very-small.conf:
```

```
include: $HOME/.mrjob.base.conf
runners:
    emr:
       num_core_instances: 2
       core_instance_type: m1.small
```

~/.mrjob.base.conf:

```
runners:
    emr:
    aws_access_key_id: HADOOPHADOOPBOBADOOP
    aws_secret_access_key: MEMIMOMADOOPBANANAFANAFOFADOOPHADOOP
    region: us-west-1
```

Options that are lists, commands, dictionaries, etc. combine the same way they do between the config files and the command line (with combiner functions).

You can use \$ENVIRONMENT\_VARIABLES and ~/file\_in\_your\_home\_dir inside include.

You can inherit from multiple config files by passing include a list instead of a string. Files on the right will have precedence over files on the left. To continue the above examples, this config:

~/.mrjob.everything.conf

will be equivalent to this one:

~/.mrjob.everything-2.conf

```
runners:
    emr:
    aws_access_key_id: HADOOPHADOOPBOBADOOP
    aws_secret_access_key: MEMIMOMADOOPBANANAFANAFOFADOOPHADOOP
    core_instance_type: m1.xlarge
    num_core_instances: 20
    region: us-west-1
```

In this case, ~/.mrjob.very-large.conf has taken precedence over ~/.mrjob.very-small.conf.

#### **Relative includes**

Relative include: paths are relative to the real (after resolving symlinks) path of the including conf file.

For example, you could do this:

~/.mrjob/base.conf:

```
runners:
```

~/.mrjob/default.conf:

```
include: base.conf
```

You could then load your configs via a symlink ~/.mrjob.conf to ~/.mrjob/default.conf and ~/.mrjob/base.conf would still be included (even though it's not in the same directory as the symlink).

## 1.7.4 Clearing configs

Sometimes, you just want to override a list-type config (e.g. setup) or a \*PATH environment variable, rather than having mrjob cleverly concatenate it with previous configs.

You can do this in YAML config files by tagging the values you want to take precedence with the !clear tag.

For example:

~/.mrjob.base.conf

```
runners:
    emr:
    aws_access_key_id: HADOOPHADOOPBOBADOOP
    aws_secret_access_key: MEMIMOMADOOPBANANAFANAFOFADOOPHADOOP
    cmdenv:
        PATH: /this/nice/path
        PYTHONPATH: /here/be/serpents
        USER: dave
    setup:
        - /run/this/command
```

~/.mrjob.conf

```
include: ~/mrjob.base.conf
runners:
    emr:
    cmdenv:
        PATH: !clear /this/even/better/path/yay
        PYTHONPATH: !clear
        setup: !clear
        - /run/this/other/command
```

is equivalent to:

If you specify multiple config files (e.g. -c ~/mrjob.base.conf -c ~/mrjob.conf), a !clear in a later file will override earlier files. include: is really just another way to prepend to the list of config files to load.

If you find it more readable, you may put the !clear tag before the key you want to clear. For example,

```
runners:
    emr:
    !clear setup:
        - /run/this/other/command
```

is equivalent to:

!clear tags in lists are ignored. You cannot currently clear an entire set of configs (e.g. runners: emr: !clear ... does not work).

# 1.8 Options available to all runners

The format of each item in this document is:

```
mrjob_conf_option_name (--command-line-option-name) [option_type] Default: default value
```

Description of option behavior

Options that take multiple values can be passed multiple times on the command line. All options can be passed as keyword arguments to the runner if initialized programmatically.

## 1.8.1 Making files available to tasks

Most jobs have dependencies of some sort - Python packages, Debian packages, data files, etc. This section covers options available to all runners that mrjob uses to upload files to your job's execution environments. See *File options* if you want to write your own command line options related to file uploading.

**Warning:** You **must** wait to read files until **after class initialization**. That means you should use the \*\_init() methods to read files. Trying to read files into class variables will not work.

```
bootstrap_mrjob (--bootstrap-mrjob, --no-bootstrap-mrjob) [boolean] Default: (automatic)
```

Should we automatically zip up the mrjob library and install it when we run job? By default, we do unless interpreter (deprecated) is set.

Set this to False if you've already installed mrjob on your Hadoop cluster or install it by some other method. mrjob used to be bootstrapped as a tarball.

```
py_files (--py-files) [path list] Default: []
```

List of .egg or .zip files to add to your job's PYTHONPATH.

This is based on a Spark feature, but it works just as well with streaming jobs.

New in version 0.5.7.

Changed in version 0.6.7: Deprecated --py-file in favor of --py-files

```
upload_archives (--archives) [path list] Default: []
```

A list of archives (e.g. tarballs) to unpack in the local directory of the mr\_job script when it runs. You can set the name in the job's working directory we unpack into by appending #nameinworkingdir to the path; otherwise we just use the name of the archive file (e.g. foo.tar.gz is unpacked to the directory foo.tar.gz/, and foo.tar.gz#stuff is unpacked to the directory stuff/).

Changed in version 0.5.7: This works with Spark on YARN as well.

Changed in version 0.6.7: Deprecated --archive in favor of --archives

```
upload_dirs (--dirs) [path list] Default: []
```

A list of directories to copy to the local directory of the mr\_job script when it runs (mrjob does this by tarballing the directory and submitting the tarball to Hadoop as an archive).

You can set the name in the job's working directory of the directory we copy by appending #nameinworkingdir to the path; otherwise we just use its name.

This works with Spark on YARN only.

New in version 0.5.8.

Changed in version 0.6.7: Deprecated --dir in favor of --dirs

```
upload_files (--files) [path list] Default: []
```

Files to copy to the local directory of the mr\_job script when it runs. You can set the name of the file in the job's working directory by appending #nameinworkingdir to the path; otherwise we just use the name of the file.

In the config file:

```
upload_files:
    - file_1.txt
    - file_2.sqlite
```

#### On the command line:

```
--files file_1.txt,file_2.sqlite
```

Changed in version 0.5.7: This works with Spark as well.

Changed in version 0.6.7: Deprecated --file in favor of --files

Changed in version 0.6.8: In Spark, can use #nameinworkingdir even when not on YARN.

## 1.8.2 Temp files and cleanup

```
cleanup (--cleanup) [string] Default: 'ALL'
```

List of which kinds of directories to delete when a job succeeds. Valid choices are:

- 'ALL': delete logs and local and remote temp files; stop cluster if on EMR and the job is not done when cleanup is run.
- 'CLUSTER': terminate EMR cluster if job not done when cleanup is run
- 'JOB': stop job if not done when cleanup runs (temporarily disabled)
- 'LOCAL\_TMP': delete local temp files only
- 'LOGS': delete logs only
- 'NONE': delete nothing
- 'REMOTE\_TMP': delete remote temp files only
- 'TMP': delete local and remote temp files, but not logs

In the config file:

```
cleanup: [LOGS, JOB]
```

#### On the command line:

```
--cleanup=LOGS, JOB
```

Changed in version 0.5.0: Options ending in TMP used to end in SCRATCH

```
cleanup_on_failure (--cleanup-on-failure) [string] Default: 'NONE'
```

Which kinds of directories to clean up when a job fails. Valid choices are the same as **cleanup**.

```
local_tmp_dir (--local-tmp-dir) [path] Default: value of tempfile.gettempdir()
```

Alternate local temp directory.

```
--local-tmp-dir '' tells mrjob to ignore the config file and use the default temp directory (tempfile.gettempdir())
```

Changed in version 0.6.6: Added *-local-tmp-dir* switch.

Changed in version 0.5.0: This option used to be named base\_tmp\_dir.

```
output_dir (--output-dir) [string] Default: (automatic)
```

An empty/non-existent directory where Hadoop streaming should put the final output from the job. If you don't specify an output directory, we'll output into a subdirectory of this job's temporary directory. You can control this from the command line with --output-dir. This option cannot be set from configuration files. If used with the hadoop runner, this path does not need to be fully qualified with hdfs:// URIs because it's understood that it has to be on HDFS.

```
cat_output (--cat-output, --no-cat-output) [boolean] Default: output if output_dir is not set
```

Should we stream job output to STDOUT after completion?

Changed in version 0.6.3: used to be --no-output.

```
step_output_dir (--step-output-dir) [string] Default: (automatic)
```

For a multi-step job, where to put output from job steps other than the last one. Each step's output will go into a numbered subdirectory of this one (0000/, 0001/, etc.)

This option can be useful for debugging. By default, intermediate output goes into HDFS, which is fastest but not easily accessible on EMR or Dataproc.

This option currently does nothing on local and inline runners.

## 1.8.3 Job execution context

```
cmdenv (--cmdenv) [environment variable dict] Default: { }
```

Dictionary of environment variables to pass to the job inside Hadoop streaming.

In the config file:

```
cmdenv:
PYTHONPATH: $HOME/stuff
TZ: America/Los_Angeles
```

On the command line:

```
--cmdenv PYTHONPATH=$HOME/stuff,TZ=America/Los_Angeles
```

Changed in version 0.5.7: This works with Spark too. In client mode (hadoop runner), these environment variables are passed directly to **spark-submit**.

```
interpreter (--interpreter) [string] Default: None
```

Deprecated since version 0.6.7: support for non-Python jobs is being phased out.

Non-Python command to launch your script with (e.g. 'ruby'). This will also be used to query the script about steps unless you set steps\_interpreter.

If you want to use an alternate Python command to run the job, use python\_bin.

This takes precedence over python\_bin and steps\_python\_bin.

### python\_bin (--python-bin) [command] Default: (automatic)

Name/path of alternate Python binary for wrapper scripts and mappers/reducers (e.g. 'python -v').

If you're on Python 3, this always defaults to 'python3'.

If you're on Python 2, this defaults to 'python' (except on EMR AMIs prior to 4.3.0, where it will be 'python2.7').

Likewise, if you're using PyPy, this defaults to 'pypy' or 'pypy3' depending on your version.

This option also affects which Python binary is used for file locking in setup scripts, so it might be useful to set even if you're using a non-Python interpreter (deprecated). It's also used by <code>EMRJobRunner</code> to compile mrjob after bootstrapping it (see bootstrap\_mrjob).

Changed in version 0.6.10: added 'pypy' and 'pypy3' as possible defaults

Note: mrjob does not auto-install PyPy for you on EMR; see *Installing PyPy* for how to do this

### setup (--setup) [string list] Default: []

A list of lines of shell script to run before each task (mapper/reducer).

This option is complex and powerful; the best way to get started is to read the Job Environment Setup Cookbook.

Using this option replaces your task with a shell "wrapper" script that executes the setup commands, and then executes the task as the last line of the script. This means that environment variables set by hadoop (e.g. \$mapred\_job\_id) are available to setup commands, and that you can pass environment variables to the task (e.g. \$PYTHONPATH) using export.

We use file locking around the setup commands (not the task) to ensure that multiple tasks running on the same node won't run them simultaneously (it's safe to run make). Before running the task, we cd back to the original working directory.

In addition, passing expressions like path#name will cause *path* to be automatically uploaded to the task's working directory with the filename *name*, marked as executable, and interpolated into the script by its absolute path on the machine running the script.

path may also be a URI, and  $\sim$  and environment variables within path will be resolved based on the local environment. name is optional.

You can indicate that an archive should be unarchived into a directory by putting a / after *name* (e.g. foo.tar.gz#foo/).

You can indicate that a directory should be copied into the job's working directory by putting a / after *path* (e.g. src-tree/#). You may optionally put a / after *name* as well (e.g. cd src-tree/#/subdir).

This works for Spark as well (except on the local[\*] master, where it doesn't make sense). The setup script is run before every executor, but only run before the driver in cluster mode.

**Note:** Uploading archives and directories (e.g. src-tree/#) to Spark's working directory still only works on YARN.

New in version 0.5.8: support for directories (above)

New in version 0.6.7: support for Spark on YARN only

New in version 0.6.8: full support for Spark

For more details of parsing, see parse\_setup\_cmd().

sh bin (--sh-bin) [command] Default: /bin/sh -ex

Name/path of alternate shell binary to use for setup and bootstrap. Needs to be backwards compatible with Bourne Shell (e.g. bash, zsh).

If you want to add an argument, use an absolute path (**/bin/bash -x**, not **bash -x**). Please do not pass multiple args to your shell binary (this plays poorly with Linux shebang syntax).

This is also used to wrap mappers, reducers, etc. that require piping one command into another (see e.g. mapper\_pre\_filter()).

Changed in version 0.5.9: Starting with EMR AMI 5.2.0, **/bin/sh** -**e** is broken, so we emulate the -**e** switch by using **/bin/sh** -**x** as our shell, and prepending **set** -**e** to any shell script generated by mrjob. **set** -**e** is not prepended if you set *sh\_bin* yourself; you could add it with setup.

Changed in version 0.6.7: Used to be **sh** -**ex** on local and Hadoop runners

Changed in version 0.6.8: Setting this to an empty value (--sh-bin '') means to use the default (used to cause an error).

## steps\_interpreter (--steps-interpreter) [command] Default: current Python interpreter

Deprecated since version 0.6.7: support for non-Python jobs is being phased out.

Alternate (non-Python) command to use to query the job about its steps. Usually it's good enough to set interpreter.

If you want to use an alternate Python command to get the job's steps, use steps\_python\_bin.

This takes precedence over steps\_python\_bin.

## steps\_python\_bin (--steps-python-bin) [command] Default: (current Python interpreter)

Deprecated since version 0.6.7: In most cases, runners no longer query jobs for steps, so this does nothing.

Name/path of alternate python binary to use to query the job about its steps. Rarely needed. If not set, we use sys.executable (the current Python interpreter).

## task\_python\_bin (--task-python-bin) [command] Default: same as python\_bin

Name/path of alternate python binary to run the job (invoking it with --mapper, --spark or anything other than --steps).

In most cases, you're better off setting python\_bin, which this defaults to. This option exists mostly to support running tasks inside Docker while using a normal Python binary in setup wrapper scripts.

#### 1.8.4 Other

```
conf_paths (-c, --conf-path, --no-conf) [path list] Default: see find_mrjob_conf()
```

List of paths to configuration files. This option cannot be used in configuration files, because that would cause a universe-ending causality paradox. Use *-no-conf* on the command line or *conf\_paths=[]* to force mrjob to load no configuration files at all. If no config path flags are given, mrjob will look for one in the locations specified in *Config file format and location*.

Config path flags can be used multiple times to combine config files, much like the **include** config file directive. Using --no-conf will cause mrjob to ignore all preceding config path flags.

For example, this line will cause mrjob to combine settings from left.conf and right .conf:

```
python my_job.py -c left.conf -c right.conf
```

This line will cause mrjob to read no config file at all:

```
python my_job.py --no-conf
```

This line will cause mrjob to read only right.conf, because --no-conf nullifies -c left.conf:

```
python my_job.py -c left.conf --no-conf -c right.conf
```

```
read_logs (--read-logs, --no-read-logs) [boolean] Default: True
```

New in version 0.6.5.

If set to false, don't list or read the contents of log files generated in the course of running your job.

The main impact of turning off *read\_logs* is that if your job fails, mrjob won't spend any time or effort determining why it failed. On EMR, this effectively disables counter fetching as well.

This option does not stop the Hadoop and Dataproc runners from reading the output of the job driver (i.e. hadoop jar ...), so you will continue to get counters and high-level Java errors on these platforms.

## 1.8.5 Options ignored by the local and inline runners

These options are ignored because they require a real instance of Hadoop:

- hadoop\_input\_format
- hadoop\_output\_format
- libjars
- partitioner

# 1.8.6 Options ignored by the inline runner

These options are ignored because the inline runner does not invoke the job as a subprocess:

- bootstrap\_mrjob
- py\_files
- python\_bin
- · read\_logs
- setup
- steps\_python\_bin

# 1.9 Hadoop-related options

Since mrjob is geared toward Hadoop, there are a few Hadoop-specific options. However, due to the difference between the different runners, the Hadoop platform, and Elastic MapReduce, they are not all available for all runners.

## 1.9.1 Options specific to the local and inline runners

```
hadoop_version (--hadoop-version) [string] Default: None
```

Set the version of Hadoop to simulate (this currently only matters for jobconf).

If you don't set this, the local and inline runners will run in a version-agnostic mode, where anytime the runner sets a simulated jobconf variable, it'll use *every* possible name for it (e.g. user.name *and* mapreduce.job.user.name).

```
num_cores (--num-cores) [integer] Default: None
```

Maximum number of tasks to handle at one time. If not set, defaults to the number of CPUs on your system.

This also affects the number of input file splits the runner makes (the only impact in inline mode).

New in version 0.6.2.

# 1.9.2 Options available to local, hadoop, and emr runners

These options are both used by Hadoop and simulated by the local and inline runners to some degree.

```
jobconf (-D, --jobconf) [jobconf dict] Default: { }
```

```
-D args to pass to hadoop streaming. This should be a map from property name to value. Equivalent to passing ['-D', 'KEY1=VALUE1', '-D', 'KEY2=VALUE2', ...] to hadoop_extra_args
```

Changed in version 0.6.6: added the -D switch on the command line, to match Hadoop.

Changed in version 0.6.6: boolean true and false values in config files are passed correctly to Hadoop (see *JobConf dicts*)

## 1.9.3 Options available to hadoop and emr runners

```
hadoop_extra_args (--hadoop-args) [string list] Default: []
```

Extra arguments to pass to hadoop streaming.

Changed in version 0.6.6: Deprecated --hadoop-arg in favor of --hadoop-args

```
hadoop_streaming_jar (--hadoop-streaming-jar) [string] Default: (automatic)
```

Path to a custom hadoop streaming jar.

On EMR, this can be either a local path or a URI (s3://...). If you want to use a jar at a path on the master node, use a file:// URI.

On Hadoop, mrjob tries its best to find your hadoop streaming jar, searching these directories (recursively) for a .jar file with hadoop followed by streaming in its name:

- \$HADOOP PREFIX
- \$HADOOP HOME
- \$HADOOP\_INSTALL
- \$HADOOP\_MAPRED\_HOME
- the parent of the directory containing the Hadoop binary (see hadoop\_bin), unless it's one of /, /usr or /usr/local
- \$HADOOP\_\*\_HOME (in alphabetical order by environment variable name)
- /home/hadoop/contrib
- /usr/lib/hadoop-mapreduce

(The last two paths allow the Hadoop runner to work out-of-the box inside EMR.)

#### libjars (--libjars) [string list] Default: []

List of paths of JARs to be passed to Hadoop with the -libjars switch.

~ and environment variables within paths will be resolved based on the local environment.

New in version 0.5.3.

Changed in version 0.6.7: Deprecated --libjar in favor of --libjars

Note: mrjob does not yet support *libjars* on Google Cloud Dataproc.

```
label (--label) [string] Default: script's module name, or no_script
```

Alternate label for the job

```
owner (--owner) [string] Default: getpass.getuser(), or no_user if that fails
```

Who is running this job (if different from the current user)

```
check_input_paths (--check-input-paths, --no-check-input-paths) [boolean] Default: True
```

Option to skip the input path check. With --no-check-input-paths, input paths to the runner will be passed straight through, without checking if they exist.

```
spark_args (--spark-args) [string list] Default: []
```

Extra arguments to pass to **spark-submit**.

**Warning:** Don't use this to set --master or --deploy-mode. On the Hadoop runner, you can change these with spark\_master and spark\_deploy\_mode. Other runners don't allow you to set these because they can only handle the defaults.

New in version 0.5.7.

Changed in version 0.6.6: Deprecated --spark-arg in favor of --spark-args

# 1.9.4 Options available to hadoop runner only

```
hadoop_bin (--hadoop-bin) [command] Default: (automatic)
```

Name/path of your **hadoop** binary (may include arguments).

mrjob tries its best to find hadoop, checking all of the following places for an executable file named hadoop:

- \$HADOOP\_PREFIX/bin
- \$HADOOP HOME/bin
- \$HADOOP\_INSTALL/bin
- \$HADOOP\_INSTALL/hadoop/bin
- \$PATH
- \$HADOOP\_\*\_HOME/bin (in alphabetical order by environment variable name)

If all else fails, we just use hadoop and hope for the best.

Changed in version 0.6.8: Setting this to an empty value (--hadoop-bin '') means to search for the Hadoop binary (used to effectively disable use of the hadoop command).

#### hadoop\_log\_dirs (--hadoop-log-dir) [path list] Default: (automatic)

Where to look for Hadoop logs (to find counters and probable cause of job failure). These can be (local) paths or URIs (hdfs:///...).

If this is *not* set, mrjob will try its best to find the logs, searching in:

- \$HADOOP\_LOG\_DIR
- \$YARN\_LOG\_DIR (on YARN only)
- hdfs:///tmp/hadoop-yarn/staging(on YARN only)
- <job output dir>/\_logs (usually this is on HDFS)
- \$HADOOP\_PREFIX/logs
- \$HADOOP\_HOME/logs
- \$HADOOP\_INSTALL/logs
- \$HADOOP\_MAPRED\_HOME/logs
- <dir containing hadoop bin>/logs (see hadoop\_bin), unless the hadoop binary is in /bin, /usr/bin, or /usr/local/bin
- \$HADOOP\_\*\_HOME/logs (in alphabetical order by environment variable name)
- /var/log/hadoop-yarn (on YARN only)
- /mnt/var/log/hadoop-yarn (on YARN only)
- /var/log/hadoop
- /mnt/var/log/hadoop

New in version 0.5.0.

Changed in version 0.5.3: Added paths in /var/log and /mnt/var/log/hadoop-yarn

#### hadoop\_tmp\_dir(--hadoop-tmp-dir) [path] Default: tmp/mrjob

Scratch space on HDFS. This path does not need to be fully qualified with hdfs:// URIs because it's understood that it has to be on HDFS.

Changed in version 0.5.0: This option used to be named hdfs scratch dir.

#### spark deploy mode (--spark-deploy-mode) [string] Default: 'client'

Deploy mode (client or cluster) to pass to the --deploy-mode argument of **spark-submit**.

New in version 0.6.6.

#### spark\_master (--spark-master) [string] Default: 'yarn'

Name or URL to pass to the --master argument of **spark-submit** (e.g. spark://host:port, yarn).

Note that archives (see upload\_archives) only work when this is set to yarn.

#### spark\_submit\_bin (--spark-submit-bin) [command] Default: (automatic)

Name/path of your **spark-submit** binary (may include arguments).

mrjob tries its best to find **spark-submit**, checking all of the following places for an executable file named spark-submit:

- \$SPARK\_HOME/bin
- \$PATH

- your pyspark installation's bin/directory
- /usr/lib/spark/bin
- /usr/local/spark/bin
- /usr/local/lib/spark/bin

If all else fails, we just use spark-submit and hope for the best.

New in version 0.5.7.

Changed in version 0.6.8: Searches for **spark-submit** in pyspark installation.

# 1.10 Spark runner options

All options from Options available to all runners and Hadoop-related options are available in the Spark runner.

In addition, the Spark runner has the following options in common with other runners:

- · aws access key id
- aws\_secret\_access\_key
- aws\_session\_token
- cloud\_fs\_sync\_secs
- cloud\_part\_size\_mb
- · gcs\_region
- project\_id
- s3\_endpoint
- s3\_region

Options unique to the Spark runner:

```
emulate_map_input_file (--emulate-map-input-file, --no-emulate-map-input-file) [boolean]
Default: False
```

Imitate Hadoop by setting \$mapreduce\_map\_input\_file to the path of the input file for the current partition. This helps support jobs that rely on <code>jobconf\_from\_env('mapreduce.map.input.file')</code>.

This feature only applies to the mapper of the job's first step, and is ignored by jobs that set <code>HADOOP\_INPUT\_FORMAT</code>.

New in version 0.6.9.

```
gcs_region (--gcs-region) [string] Default: None
```

The region to use when creating a temporary bucket on Google Cloud Storage.

Similar in meaning to region, but only used to configure GCS (not S3)

```
s3_region (--s3-region) [string] Default: None
```

The region to use when creating a temporary bucket on S3.

Similar in meaning to region, but only used to configure S3 (not GCS)

```
skip_internal_protocol (--skip-internal-protocol, --no-skip-internal-protocol) [boolean]
    Default: False
```

Don't emulate the job's internal protocol (used for communicating between job steps and tasks in the same step), instead relying on Spark to encode and decode data structures.

This should work for most but not all jobs, and make them run at least somewhat faster. Some things to keep in mind:

- data will no longer be "normalized" by being converted to and from string representation. For example, running a tuple through JSONProtocol (the default) implicitly converts it to a list because there are no tuples in JSON. With internal protocols skipped, it would remain a tuple.
- if your job uses SORT\_VALUES, keep in mind that your values will need to be comparable as Spark will be comparing them directly, rather than comparing their internal-protocol-encoded representation. This may also affect sorting order.

New in version 0.6.10.

```
spark_tmp_dir(--spark-tmp-dir) [string] Default: (automatic)
```

A place to put files where they are visible to Spark executors, similar to cloud\_tmp\_dir.

If running locally, defaults to a directory inside local\_tmp\_dir, and if running on a cluster, to tmp/mrjob on HDFS.

# 1.11 Configuration quick reference

# 1.11.1 Setting configuration options

You can set an option by:

- Passing it on the command line with the switch version (like --some-option)
- Passing it as a keyword argument to the runner constructor, if you are creating the runner programmatically
- Putting it in one of the included config files under a runner name, like this:

```
runners:
    local:
        python_bin: python3.6 # only used in local runner
    emr:
        python_bin: python3 # only used in Elastic MapReduce runner
```

See Config file format and location for information on where to put config files.

# 1.11.2 Options that can't be set from mrjob.conf (all runners)

There are some options that it makes no sense to set in the config file.

These options can be set via command-line switches:

Config	Command line	Default	Type
cat_output	-cat-output, -no-cat-output	output if output_dir is not set	boolean
conf_paths	-c, -conf-path, -no-conf	<pre>see find_mrjob_conf()</pre>	path list
output_dir	-output-dir	(automatic)	string
step_output_dir	-step-output-dir	(automatic)	string

These options can be set by overriding attributes or methods in your job class:

Option	Attribute	Method	Default
hadoop_input_format	HADOOP_INPUT_FORMAT	<pre>hadoop_input_format()</pre>	None
hadoop_output_format	HADOOP_OUTPUT_FORMAT	hadoop_output_format()	None
partitioner	PARTITIONER	partitioner()	None

These options can be set by overriding your job's <code>configure\_args()</code> to call the appropriate method:

Option	Method	Default
extra_args	<pre>add_passthru_arg()</pre>	[]
file_upload_args	add_file_arg()	[]

All of the above can be passed as keyword arguments to MRJobRunner.\_\_init\_\_() (this is what makes them runner options), but you usually don't want to instantiate runners directly.

# 1.11.3 Other options for all runners

These options can be passed to any runner without an error, though some runners may ignore some options. See the text after the table for specifics.

Config	Command line	Default	Туре
bootstrap_mrjob	-bootstrap-mrjob,	(automatic)	boolean
	–no-bootstrap-mrjob		
check_input_path	s –check-input-paths,	True	boolean
	-no-check-input-paths		
cleanup	-cleanup	'ALL'	string
cleanup_on_failur	e –cleanup-on-failure	'NONE'	string
cmdenv	-cmdenv	{}	environment
			variable dict
hadoop_extra_arg	s –hadoop-args	[]	string list
hadoop_streaming	_ <del>jar</del> adoop-streaming-jar	(automatic)	string
interpreter	-interpreter	None	string
jobconf	-D, –jobconf	{}	jobconf dict
label	-label	script's module name, or	string
		no_script	
libjars	–libjars	[]	string list
local_tmp_dir	-local-tmp-dir	value of	path
		tempfile.gettempdir()	
owner	-owner	getpass.getuser(),or	string
		no_user if that fails	
py_files	-py-files	[]	path list
python_bin	–python-bin	(automatic)	command
read_logs	-read-logs, -no-read-logs	True	boolean
setup	-setup	[]	string list
sh_bin	-sh-bin	/bin/sh -ex	command
spark_args	–spark-args	[]	string list
steps_interpreter	-steps-interpreter	current Python interpreter	command
steps_python_bin	-steps-python-bin	(current Python interpreter)	command
task_python_bin	–task-python-bin	same as python_bin	command
upload_archives	-archives	[]	path list
upload_dirs	-dirs	[]	path list
upload_files	-files	[]	path list

LocalMRJobRunner takes no additional options, but:

• bootstrap\_mrjob is False by default

- cmdenv uses the local system path separator instead of : all the time (so ; on Windows, no change elsewhere)
- python\_bin defaults to the current Python interpreter

In addition, it ignores hadoop\_input\_format, hadoop\_output\_format, hadoop\_streaming\_jar, and jobconf

InlineMRJobRunner works like LocalMRJobRunner, only it also ignores bootstrap\_mrjob, cmdenv, python\_bin, steps\_python\_bin, upload\_archives, and upload\_files.

# 1.11.4 Additional options for DataprocJobRunner

Config	Command line	Default	Туре
cluster_properties	-cluster-property	None	
core_instance_config	-core-instance-config	None	
gcloud_bin	-gcloud-bin	'gcloud'	command
master_instance_config	-master-instance-config	None	
network	-network	None	string
project_id	-project-id	read from credentials config file	string
service_account	-service-account	None	
service_account_scopes	-service-account-scopes	(automatic)	
task_instance_config	-task-instance-config	None	

# 1.11.5 Additional options for EMRJobRunner

Config	Command line	Default	Туре
additional_emr_info	-additional-emr-info	None	special
applications	-application, -applications	[]	string list
aws_access_key_id		None	string
aws_secret_access_key	-aws-secret-access-key	None	string
aws_session_token		None	string
bootstrap_actions	-bootstrap-actions	[]	string list
bootstrap_spark	-bootstrap-spark, -no-bootstrap-spark	(automatic)	boolean
cloud_log_dir	-cloud-log-dir	append logs to cloud_tmp_dir	string
core_instance_bid_price	-core-instance-bid-price	None	string
ebs_root_volume_gb	-ebs-root-volume-gb	None	integer
ec2_endpoint	-ec2-endpoint	(automatic)	string
ec2_key_pair	-ec2-key-pair	None	string
ec2_key_pair_file	-ec2-key-pair-file	None	path
emr_action_on_failure	-emr-action-on-failure	(automatic)	string
emr_configurations	-emr-configuration	[]	list of dicts
emr_endpoint	-emr-endpoint	infer from region	string
enable_emr_debugging	-enable-emr-debugging	False	boolean
hadoop_streaming_jar_on_emr	-hadoop-streaming-jar-on-emr	AWS default	string
iam_endpoint	-iam-endpoint	(automatic)	string
iam_instance_profile	-iam-instance-profile	(automatic)	string
iam_service_role	-iam-service-role	(automatic)	string
instance_fleets	-instance-fleet	None	
instance_groups	-instance-groups	None	
master_instance_bid_price	-master-instance-bid-price	None	string
mins_to_end_of_hour	-mins-to-end-of-hour	5.0	float
	•	Continued o	n next page

rable 1.1 – continued from previous page			
Config	Command line	Default	Type
pool_clusters	-pool-clusters	True	string
pool_name	-pool-name	'default'	string
pool_wait_minutes	-pool-wait-minutes	0	string
release_label	-release-label	None	string
s3_endpoint	-s3-endpoint	(automatic)	string
ssh_bin	-ssh-bin	'ssh'	command
tags	-tag	{}	dict
task_instance_bid_price	-task-instance-bid-price	None	string
visible_to_all_users	-visible-to-all-users, -no-visible-to-all-users	True	boolean

Table 1.1 – continued from previous page

## 1.11.6 Additional options for HadoopJobRunner

Config	Command line	Default	Туре
hadoop_bin	-hadoop-bin	(automatic)	command
hadoop_log_dirs	-hadoop-log-dir	(automatic)	path list
hadoop_tmp_dir	-hadoop-tmp-dir	tmp/mrjob	path
spark_deploy_mode	-spark-deploy-mode	'client'	string
spark_master	-spark-master	'yarn'	string
spark_submit_bin	-spark-submit-bin	(automatic)	command

# 1.12 Cloud runner options

These options are generally available whenever you run your job on a Hadoop cloud service (AWS Elastic MapReduce or Google Cloud Dataproc).

All options from Options available to all runners and Hadoop-related options are also available on cloud services.

# 1.12.1 Google credentials

See Getting started with Google Cloud for specific instructions about setting these options.

# 1.12.2 Choosing/creating a cluster to join

**cluster\_id** (--cluster-id) [string] **Default:** automatically create a cluster and use it

The ID of a persistent cluster to run jobs in (on Dataproc, this is the same thing as "cluster name").

It's fine for other jobs to be using the cluster; we give our job's steps a unique ID.

# 1.12.3 Job placement

region (--region) [string] Default: 'us-west-2' on EMR, 'us-west1' on Dataproc

Geographic region to run jobs in (e.g. us-central-1).

If mrjob create a temporary bucket, it will be created in this region as well.

If you set region, you do not need to set zone; a zone will be chosen for you automatically.

#### subnet (--subnet) [string] Default: None

Optional subnet(s) to launch your job in.

On Amazon EMR, this is the ID of a VPC subnet to launch cluster in (e.g. 'subnet-12345678'). This can also be a list of possible subnets if you are using instance\_fleets.

On Google Cloud Dataproc, this is the name of a subnetwork (e.g. 'default'). Specifying *subnet* rather than network will ensure that your cluster only has access to one specific geographic region, rather than the entire VPC.

Changed in version 0.6.8: --subnet '' un-sets the subnet on EMR (used to be ignored).

Changed in version 0.6.3: Works on Google Cloud Dataproc as well as AWS Elastic MapReduce.

New in version 0.5.3.

#### zone (--zone) [string] Default: None

Zone within a specific geographic region to run your job in.

If you set this, you do not neet to set region.

Changed in version 0.5.4: This option used to be named aws\_availability\_zone

## 1.12.4 Number and type of instances

instance\_type (--instance-type) [string] Default: m5.xlarge on EMR, n1-standard-1 on Dataproc

Type of instance that runs your Hadoop tasks.

Once you've tested a job and want to run it at scale, it's usually a good idea to use instances larger than the default. For EMR, see Amazon EC2 Instance Types and for Dataproc, see Machine Types.

**Note:** Many EC2 instance types can only run in a VPC (see subnet).

If you're running multiple nodes (see num\_core\_instances), this option *doesn't* apply to the master node because it's just coordinating tasks, not running them. Use master\_instance\_type instead.

Changed in version 0.6.10: Default on EMR is now m5.xlarge

Changed in version 0.6.6: Default on EMR was m4.large

Changed in version 0.5.6: On EMR, defaulted to ml.large if running Spark, ml.small if running on the (deprecated) 2.x AMIs, and ml.medium otherwise (ml.medium was previously the default in all cases).

Changed in version 0.5.4: This option used to be named *ec2\_instance\_type*.

#### core\_instance\_type (--core-instance-type) [string] Default: value of instance\_type

like instance\_type, but only for the core (worker) Hadoop nodes; these nodes run tasks and host HDFS. Usually you just want to use instance\_type.

Changed in version 0.5.4: This replaces the ec2\_core\_instance\_type and ec2\_slave\_instance\_type options.

# num\_core\_instances (--num-core-instances) [integer] Default: 0 on EMR, 2 on Dataproc

Number of core (worker) instances to start up. These run your job and host HDFS. This is in addition to the single master instance.

On Google Cloud Dataproc, this must be at least 2.

Changed in version 0.5.4: This option used to be named *num\_ec2\_core\_instances*.

#### master\_instance\_type (--master-instance-type) [string] Default: (automatic)

like instance\_type, but only for the master Hadoop node. This node hosts the task tracker/resource manager and HDFS, and runs tasks if there are no other nodes.

If you're running a single node (no num\_core\_instances or num\_task\_instances), this will default to the value of instance type.

Otherwise, on Dataproc, defaults to n1-standard-1, and on EMR defaults to m1.medium (exception: m1.small on the deprecated 2.x AMIs), which is usually adequate for all but the largest jobs.

Changed in version 0.5.4: This option used to be named *ec2\_master\_instance\_type*.

## task\_instance\_type (--task-instance-type) [string] Default: value of core\_instance\_type

like instance\_type, but only for the task (secondary worker) Hadoop nodes; these nodes run tasks but do not host HDFS. Usually you just want to use instance\_type.

Changed in version 0.5.4: This option used to be named *ec2\_task\_instance\_type*.

### num\_task\_instances (--num-task-instances) [integer] Default: 0

Number of task (secondary worker) instances to start up. These run your job but do not host HDFS.

You must have at least one core instance (see num\_core\_instances) to run task instances; otherwise there's nowhere to host HDFS.

Changed in version 0.5.4: This used to be called *num\_ec2\_task\_instances*.

## 1.12.5 Cluster software configuration

### image\_id (--image-id) [string] Default: None

ID of a custom machine image.

On EMR, this is complimentary with image\_version; you can install packages and libraries on your custom AMI, but it's up to EMR to install Hadoop, create the hadoop user, etc. image\_version may not be less than 5.7.0.

You can use <code>describe\_base\_emr\_images()</code> to identify Amazon Linux images that are compatible with EMR.

For more details about how to create a custom AMI that works with EMR, see Best Practices and Considerations.

**Note:** This is not yet implemented in the Dataproc runner.

New in version 0.6.5.

### image\_version (--image-version) [string] Default: '5.8.0' on EMR, '1.0' on Dataproc

Machine image version to use. This controls which Hadoop version(s) are available and which version of Python is installed, among other things.

See the AMI version docs (EMR) or the Dataproc version docs for more details.

You can use this instead of release\_label on EMR, even for 4.x+ AMIs; mrjob will just prepend emr- to form the release label.

Changed in version 0.5.7: Default on EMR used to be '3.11.0'.

Changed in version 0.5.4: This used to be called *ami\_version*.

**Warning:** The deprecated *ami\_version* alias for this option is completely ignored by mrjob 0.5.4 (it works in later 0.5.x versions).

Warning: The 2.x series of AMIs is deprecated by Amazon and not recommended.

**Warning:** The 1.x series of AMIs is no longer supported because they use Python 2.5.

## bootstrap (--bootstrap) [string list] Default: []

A list of lines of shell script to run once on each node in your cluster, at bootstrap time.

This option is complex and powerful. On EMR, the best way to get started is to read the EMR Bootstrapping Cookbook.

Passing expressions like path#name will cause *path* to be automatically uploaded to the task's working directory with the filename *name*, marked as executable, and interpolated into the script by their absolute path on the machine running the script.

path may also be a URI, and ~ and environment variables within path will be resolved based on the local environment. name is optional. For details of parsing, see parse\_setup\_cmd().

Unlike with setup, archives are not supported (unpack them yourself).

Remember to put sudo before commands requiring root privileges!

**bootstrap\_python** (--bootstrap-python, --no-bootstrap-python) [boolean] **Default:** True on Dataproc, as needed on EMR.

Attempt to install a compatible (major) version of Python at bootstrap time, including header files and pip (see *Installing Python packages with pip*).

The only reason to set this to False is if you want to customize Python/pip installation using bootstrap.

New in version 0.5.0.

Changed in version 0.5.4: no longer installs Python 3 on AMI version 4.6.0+ by default

```
extra cluster params (--extra-cluster-param) [dict] Default: {}
```

An escape hatch that allows you to pass extra parameters to the EMR/Dataproc API at cluster create time, to access API features that mrjob does not yet support.

For EMR, see the API documentation for RunJobFlow for the full list of options.

Option names are strings, and values are data structures. On the command line, --extra-cluster-param name=value:

```
--extra-cluster-param SupportedProducts='["mapr-m3"]'
--extra-cluster-param AutoScalingRole=HankPym
```

*value* can be either a JSON or a string (unless it starts with {, [, or ", so that we don't convert malformed JSON to strings). Parameters can be suppressed by setting them to null:

```
--extra-cluster-param LogUri=null
```

This also works with Google dataproc:

```
--extra-cluster-param labels='{"name": "wrench"}'
```

In the config file, extra\_cluster\_param is a dict:

```
runners:
    emr:
    extra_cluster_params:
        AutoScalingRole: HankPym
        LogUri: null # !clear works too
        SupportedProducts:
        - mapr-m3
```

Changed in version 0.6.8: You may use a *name* with dots in it to set (or unset) nested properties. For example: --extra-cluster-param Instances.EmrManagedMasterSecurityGroup=sg-foo.

New in version 0.6.0: This replaces the old *emr\_api\_params* option, which only worked with boto 2.

# 1.12.6 Monitoring your job

check\_cluster\_every (--check-cluster-every) [float] Default: 10 seconds on Dataproc, 30 seconds on EMR

How often to check on the status of your job, in seconds.

Changed in version 0.6.5: When the EMR client encounters a transient error, it will wait at least this many seconds before trying again.

Changed in version 0.5.4: This used to be called *check\_emr\_status\_every* 

```
ssh_tunnel (--ssh-tunnel, --no-ssh-tunnel) [boolean] Default: False
```

If True, create an ssh tunnel to the job tracker/resource manager and listen on a randomly chosen port.

On EMR, this requires you to set ec2\_key\_pair and ec2\_key\_pair\_file. See *Configuring SSH credentials* for detailed instructions.

On Dataproc, you don't need to set a key, but you do need to have the **gcloud** utility installed and set up (make sure you ran **gcloud auth login** and **gcloud config set project project\_id>). See Installing gcloud, gsutil, and other utilities.** 

Changed in version 0.6.3: Enabled on Google Cloud Dataproc

Changed in version 0.5.0: This option used to be named *ssh\_tunnel\_to\_job\_tracker*.

```
ssh_tunnel_is_open (--ssh-tunnel-is-open) [boolean] Default: False
```

if True, any host can connect to the job tracker through the SSH tunnel you open. Mostly useful if your browser is running on a different machine from your job runner.

Does nothing unless ssh\_tunnel is set.

```
ssh_bind_ports (--ssh-bind-ports) [list of integers] Default: range (40001, 40841)
```

A list of ports that are safe to listen on.

The main reason to set this is if your firewall blocks the default range of ports, or if you want to pick a single port for consistency.

On the command line, this looks like --ssh-bind-ports 2000[:2001][,2003,2005:2008,etc], where commas separate ranges and colons separate range endpoints.

# 1.12.7 Cloud Filesystem

```
cloud_fs_sync_secs (--cloud-fs-sync-secs) [float] Default: 5.0
```

How long to wait for cloud filesystem (e.g. S3, GCS) to reach eventual consistency? This is typically less than a second, but the default is 5 seconds to be safe.

Changed in version 0.5.4: This used to be called s3\_sync\_wait\_time

```
cloud_part_size_mb (--cloud-part-size-mb) [integer] Default: 100
```

Upload files to cloud filesystem in parts no bigger than this many megabytes (technically, mebibytes). Default is 100 MiB.

Set to 0 to disable multipart uploading entirely.

Currently, Amazon requires parts to be between 5 MiB and 5 GiB. mrjob does not enforce these limits.

Changed in version 0.6.3: Enabled on Google Cloud Storage. This used to be called *cloud\_upload\_part\_size*.

Changed in version 0.5.4: This used to be called s3\_upload\_part\_size.

```
cloud_tmp_dir (--cloud-tmp-dir) [string] Default: (automatic)
```

Directory on your cloud filesystem to use as temp space (e.g. s3://yourbucket/tmp/, gs://yourbucket/tmp/).

By default, mrjob looks for a bucket belong to you whose name starts with mrjob— and which matches region. If it can't find one, it creates one with a random name. This option is then set to *tmp*/ in this bucket (e.g. s3://mrjob-01234567890abcdef/tmp/).

Changed in version 0.5.4: This used to be called *s3\_tmp\_dir*.

Changed in version 0.5.0: This used to be called *s3\_scratch\_uri*.

#### 1.12.8 Auto-termination

```
max_mins_idle (--max-mins-idle) [float] Default: 10.0
```

Automatically terminate your cluster after it has been idle at least this many minutes. You cannot turn this off (clusters left idle rack up billing charges).

If your cluster is only running a single job, mrjob will attempt to terminate it as soon as your job finishes. This acts as an additional safeguard, as well as affecting *Cluster Pooling* on EMR.

Changed in version 0.6.5: EMR's idle termination script is more robust against sudo shutdown -h now being ignored, and logs the script's stdout and stderr /var/log/bootstrap-actions/mrjob-idle-termination.log.

Changed in version 0.6.3: Uses Dataproc's built-in cluster termination feature rather than a custom script. The API will not allow you to set an idle time of less than 10 minutes.

Changed in version 0.6.2: No matter how small a value you set this to, there is a grace period of 10 minutes between when the idle termination daemon launches and when it may first terminate the cluster, to allow Hadoop to accept your first job.

Changed in version 0.6.0: All clusters launched by mrjob now auto-terminate when idle. In previous versions, you needed to set max\_hours\_idle, set this option explicitly, or use *terminate-idle-clusters*.

```
max_hours_idle (--max-hours-idle) [float] Default: None
```

Deprecated since version 0.6.0: Use max\_mins\_idle instead.

# 1.13 Job Environment Setup Cookbook

Many jobs have significant external dependencies, both libraries and other source code.

Combining shell syntax with Hadoop's DistributedCache notation, mrjob's setup option provides a powerful, dynamic alternative to pre-installing your Hadoop dependencies on every node.

All our mrjob.conf examples below are for the hadoop runner, but these work equally well with the emr runner. Also, if you are using EMR, take a look at the EMR Bootstrapping Cookbook.

**Note:** Setup scripts don't work with Spark; try py\_files instead.

## 1.13.1 Uploading your source tree

**Note:** If you're using mrjob 0.6.4 or later, check out *Using other python modules and packages* first.

mrjob can automatically tarball your source directory and include it in your job's working directory. We can use setup scripts to upload the directory and then add it to PYTHONPATH.

Run your job with:

```
--setup 'export PYTHONPATH=$PYTHONPATH:your-src-code/#'
```

The / before the # tells mrjob that your-src-code is a directory. You may optionally include a / after the # as well (e.g. export PYTHONPATH=\$PYTHONPATH:your-source-code/#/your-lib).

If every job you run is going to want to use your-src-code, you can do this in your mr job. conf:

```
runners:
   hadoop:
    setup:
        - export PYTHONPATH=$PYTHONPATH:your-src-code/#
```

# 1.13.2 Uploading your source tree as an archive

Prior to mrjob 0.5.8, you had to archive directories yourself before uploading them.

```
tar -C your-src-code -f your-src-code.tar.gz -z -c .
```

Then, run your job with:

```
--setup 'export PYTHONPATH=$PYTHONPATH:your-src-code.tar.gz#/'
```

The / after the # (without one before it) is what tells mrjob that your-src-code.tar.gz is an archive that Hadoop should unpack.

To do the same thing in mrjob.conf:

```
runners:
   hadoop:
    setup:
        - export PYTHONPATH=$PYTHONPATH:your-src-code.tar.gz#/
```

## 1.13.3 Running a makefile inside your source dir

```
--setup 'cd your-src-dir.tar.gz#/' --setup 'make'
```

or, in mrjob.conf:

```
runners:
  hadoop:
  setup:
  - cd your-src-dir.tar.gz#
  - make
```

If Hadoop runs multiple tasks on the same node, your source dir will be shared between them. This is not a problem; mrjob automatically adds locking around setup commands to ensure that multiple copies of your setup script don't run simultaneously.

## 1.13.4 Making data files available to your job

Best practice for one or a few files is to use passthrough options; see add\_passthru\_arg().

You can also use upload\_files to upload file(s) into a task's working directory (or upload\_archives for tarballs and other archives).

If you're a setup purist, you can also do something like this:

```
--setup 'true your-file#desired-name'
```

since **true** has no effect and ignores its arguments.

## 1.13.5 Using a virtualenv

What if you can't install the libraries you need on your Hadoop cluster?

You could do something like this in your mrjob.conf:

```
runners:
   hadoop:
    setup:
        - virtualenv venv
        - . venv/bin/activate
        - pip install mr3po
```

However, now the locking feature that protects **make** becomes a liability; each task on the same node has its own virtualeny, but one task has to finish setting up before the next can start.

The solution is to share the virtualenv between all tasks on the same machine, something like this:

```
runners:
   hadoop:
    setup:
        - VENV=/tmp/$mapreduce_job_id
        - if [ ! -e $VENV ]; then virtualenv $VENV; fi
        - . $VENV/bin/activate
        - pip install mr3po
```

With Hadoop 1, you'd want to use \$mapred\_job\_id instead of \$mapreduce\_job\_id.

# 1.13.6 Other ways to use pip to install Python packages

If you have a lot of dependencies, best practice is to make a pip requirements file and use the -r switch:

```
--setup 'pip install -r path/to/requirements.txt#'
```

Note that **pip** can also install from tarballs (which is useful for custom-built packages):

```
--setup 'pip install $MY_PYTHON_PKGS/*.tar.gz#'
```

# 1.14 Hadoop Cookbook

## 1.14.1 Increasing the task timeout

**Warning:** Some EMR AMIs appear to not support setting parameters like timeout with jobconf at run time. Instead, you must use *Bootstrap-time configuration*.

If your mappers or reducers take a long time to process a single step, you may want to increase the amount of time Hadoop lets them run before failing them as timeouts.

You can do this with jobconf. For example, to set the timeout to one hour:

```
runners:
  hadoop: # also works for emr runner
  jobconf:
    mapreduce.task.timeout: 3600000
```

**Note:** If you're using Hadoop 1, which uses mapred.task.timeout, don't worry: this example still works because mrjob auto-converts your jobconf options between Hadoop versions.

# 1.14.2 Writing compressed output

To save space, you can have Hadoop automatically save your job's output as compressed files. Here's how you tell it to bzip them:

```
runners:
   hadoop: # also works for emr runner
   jobconf:
      # "true" must be a string argument, not a boolean! (Issue #323)
      mapreduce.output.fileoutputformat.compress: "true"
      mapreduce.output.fileoutputformat.compress.codec: org.apache.hadoop.io.compress.BZip2Codec
```

**Note:** You could also gzip your files with org.apache.hadoop.io.compress.GzipCodec. Usually bzip is a better option, as .bz2 files are splittable, and .gz files are not. For example, if you use .gz files as input, Hadoop has no choice but to create one mapper per .gz file.

# 1.15 Testing jobs

mrjob can run jobs without the help of Hadoop. This isn't very efficient, but it's a great way to test a job before submitting it to a cluster.

## 1.15.1 Inline runner

The inline runner (InlineMRJobRunner) is the default runner for mrjob (it's what's used when you run python mr\_your\_job.py <input> without any -r option). It runs your job in a single process so that you get faster feedback and simpler tracebacks.

## **Multiple splits**

The inline runner doesn't run mappers or reducers concurrently, but it does run at least two mappers and two reducers for each step. This can help catch bad assumptions about the MapReduce programming model.

For example, say we wanted to write a simple script that counted the number of lines of input:

```
from mrjob.job import MRJob

class MRCountLinesWrong(MRJob):

    def mapper_init(self):
        self.num_lines = 0

    def mapper(self, _, line):
        self.num_lines += 1

    def mapper_final(self):
        yield None, self.num_lines

if __name__ == '__main__':
    MRCountLinesWrong.run()
```

Looks good, but if we run it, we get more than one line count:

```
$ python -m mrjob.examples.mr_count_lines_wrong README.rst 2> /dev/null
null 77
null 60
```

Aha! Because there can be more than one mapper! It's fine to use mapper\_final() like this, but we need to reduce on a single key:

```
from mrjob.job import MRJob

class MRCountLinesRight (MRJob):

    def mapper_init(self):
        self.num_lines = 0

    def mapper(self, _, line):
        self.num_lines += 1

    def mapper_final(self):
        yield None, self.num_lines

    def reducer(self, key, values):
        yield key, sum(values)

if __name__ == '__main__':
    MRCountLinesRight.run()
```

```
$ python -m mrjob.examples.mr_count_lines_right README.rst 2> /dev/null
null 137
```

Thanks, inline runner!

### Isolated working directories

Just like Hadoop, the inline runner runs each mapper and reducer in its own (temporary) working directory. It *does* add the original working directory to \$PYTHONPATH so it can still access your local source tree.

## Simulating jobconf

The inline runner simulates jobconf variables/properties set by Hadoop (and their Hadoop 1 equivalents):

```
mapreduce.job.cache.archives (mapred.cache.archives)
mapreduce.job.cache.files (mapred.cache.files)
mapreduce.job.cache.local.archives (mapred.cache.localArchives)
mapreduce.job.cache.local.files (mapred.cache.localFiles)
mapreduce.job.id (mapred.job.id)
mapreduce.job.local.dir (job.local.dir)
mapreduce.map.input.file (map.input.file)
mapreduce.map.input.length (map.input.length)
mapreduce.map.input.start (map.input.start)
mapreduce.task.attempt.id (mapred.task.id)
mapreduce.task.id (mapred.tip.id)
```

- mapreduce.task.ismap(mapred.task.is.map)
- mapreduce.task.output.dir(mapred.work.output.dir)
- mapreduce.task.partition(mapred.task.partition)

You can use <code>jobconf\_from\_env()</code> to read these from your job's environment. For example:

```
from mrjob.compat import jobconf_from_env
from mrjob.job import MRJob

class MRCountLinesByFile(MRJob):
    def mapper(self, _, line):
        yield jobconf_from_env('mapreduce.map.input.file'), 1

    def reducer(self, path, ones):
        yield path, sum(ones)

if __name__ == '__main__':
    MRCountLinesByFile.run()
```

```
$ python -m mrjob.examples.mr_count_lines_by_file README.rst CHANGES.txt 2> /dev/null
"CHANGES.txt" 564
"README.rst" 137
```

1.15. Testing jobs 63

If you only want to simulate jobconf variables from a single version of Hadoop (for more stringent testing), you can set hadoop\_version.

### Setting number of mappers and reducers

Want more or less splits? You can tell the inline runner the same way you'd tell hadoop, with the mapreduce.job.maps and mapreduces.job.reduces jobconf options:

## 1.15.2 Local runnner

The local runner (LocalMRJobRunner; run using -r local) supports the above features, but, unlike the inline runner, it uses subprocesses.

This means it can be used to test options that don't make sense in a single-process context, including:

- python bin
- setup
- steps\_python\_bin

The local runner *does* run multiple subprocesses concurrently, but it's not really meant as a replacement for Hadoop; it's just for testing!

## 1.15.3 Anatomy of a test case

So, you've gotten a job working. Great! Here's how you write a regression test so that future developers won't break it

For this example we'll use a test of the \*\_init() methods from the mrjob test cases:

```
class MRInitJob(MRJob):

def __init__(self, *args, **kwargs):
    super(MRInitJob, self).__init__(*args, **kwargs)
    self.sum_amount = 0
    self.multiplier = 0
    self.combiner_multipler = 1

def mapper_init(self):
    self.sum_amount += 10

def mapper(self, key, value):
    yield(None, self.sum_amount)

def reducer_init(self):
    self.multiplier += 10
```

```
def reducer(self, key, values):
    yield(None, sum(values) * self.multiplier)

def combiner_init(self):
    self.combiner_multiplier = 2

def combiner(self, key, values):
    yield(None, sum(values) * self.combiner_multiplier)
```

Without using any mrjob features, we can write a simple test case to make sure our methods are behaving as expected:

```
from unittest import TestCase

class MRInitTestCase(TestCase):

    def test_mapper(self):
        j = MRInitJob()
        j.mapper_init()
        self.assertEqual(j.mapper(None, None).next(), (None, j.sum_amount))
```

To test the full job, you need to set up input, run the job, and check the collected output. The most straightforward way to provide input is to use the <code>sandbox()</code> method. Create a <code>BytesIO</code> object, populate it with data, initialize your job to read from stdin, and enable the sandbox with your <code>BytesIO</code> as stdin.

You'll probably also want to specify --no-conf so options from your local mrjob.conf don't pollute your testing environment.

This example reads from **stdin** (hence the – parameter):

```
from io import BytesIO

def test_init_funcs(self):
    num_inputs = 2
    stdin = BytesIO(b'x\n' * num_inputs)
    mr_job = MRInitJob(['--no-conf', '-'])
    mr_job.sandbox(stdin=stdin)
```

To run the job without leaving temp files on your system, use the <code>make\_runner()</code> context manager. <code>make\_runner()</code> creates the runner specified in the command line arguments and ensures that job cleanup is performed regardless of the success or failure of the job.

Run the job with run(). The job's output is available as a generator through cat\_output() and can be parsed with the job's output protocol using parse\_output():

```
results = []
with mr_job.make_runner() as runner:
    runner.run()
    for key, value in mrjob.parse_output(runner.cat_output()):
        results.append(value)

# these numbers should match if mapper_init, reducer_init, and
# combiner_init were called as expected
self.assertEqual(sorted(results)[0], num_inputs * 10 * 10 * 2)
```

**Warning:** Do not let your tests depend on the input lines being processed in a certain order. Both mrjob and Hadoop divide input non-deterministically.

Note: In mrjob versions prior to 0.6.0, you have to parse output line by line; see Running your job programmatically

1.15. Testing jobs 65

for an example.

# 1.16 Cloud Dataproc

## 1.16.1 Dataproc Quickstart

### **Getting started with Google Cloud**

Using mrjob with Google Cloud Dataproc is as simple creating an account, enabling Google Cloud Dataproc, and creating credentials.

#### Creating an account

- Go to cloud.google.com.
- Click the circle in the upper right, and select your Google account (if you don't have one sign up here. If you have multiple Google accounts, sign out first, and then sign into the account you want to use.
- Click Try it Free in the upper right
- Enter your name and payment information
- Wait a few minutes while your first project is created

#### **Enabling Google Cloud Dataproc**

- Go here (or search for "dataproc" under APIs & Services > Library in the upper left-hand menu)
- Click Enable

### **Configuring Google Cloud credentials**

- Go here (or pick **APIs & Services > Credentials** in the upper left-hand menu)
- Pick Create credentials > Service account key
- Select Compute engine default service account
- Click Create to download a JSON file.

Then you should either install and set up the optional **gcloud** utility (see below) or point **\$GOOGLE\_APPLICATION\_CREDENTIALS** at the file you downloaded (export GOOGLE\_APPLICATION\_CREDENTIALS="/path/to/Your Credentials.json").

### Installing gcloud, gsutil, and other utilities

mrjob does not require you to install the **gcloud** command in order to run jobs on Google Cloud Dataproc, unless you want to set up an SSH tunnel to the Hadoop resource manager (see ssh\_tunnel).

The **gcloud** command can be very useful for monitoring your job. The **gsutil** utility, packaged with it, is very helpful for dealing with Google Storage, the cloud filesystem that Google Cloud Dataproc uses.

To install gcloud and gsutil:

- Follow these three steps to install the utilities
- Log in with your Google credentials (these will launch a browser): \* gcloud auth login \* gcloud auth application-default init

On some versions of gcloud, you may have to manually configure project ID for ssh\_tunnel to work.

- run gcloud projects list to get your project ID
- gcloud config set project <project\_id>

It's also helpful to set **gcloud**'s region and zone to match mrjob's defaults:

- gcloud config set compute/region us-west1
- gcloud config set compute/zone us-west1-a
- gcloud config set dataproc/region us-west1

## **Running a Dataproc Job**

Running a job on Dataproc is just like running it locally or on your own Hadoop cluster, with the following changes:

- The job and related files are uploaded to GCS before being run
- The job is run on Dataproc (of course)
- Output is written to GCS before mrjob streams it to stdout locally
- The Hadoop version is specified by the Dataproc version

This the output of this command should be identical to the output shown in Fundamentals, but it should take much longer:

```
> python word_count.py -r dataproc README.txt
"chars" 3654
"lines" 123
"words" 417
```

#### Sending Output to a Specific Place

If you'd rather have your output go to somewhere deterministic on GCS, use --output-dir:

```
> python word_count.py -r dataproc README.rst \
> --output-dir=gs://my-bucket/wc_out/
```

## **Choosing Type and Number of GCE Instances**

When you create a cluster on Dataproc, you'll have the option of specifying a number and type of GCE instances, which are basically virtual machines. Each instance type has different memory, CPU, I/O and network characteristics, and costs a different amount of money. See Machine Types and Pricing for details.

Instances perform one of three roles:

- Master: There is always one master instance. It handles scheduling of tasks (i.e. mappers and reducers), but does not run them itself.
- Worker: You may have one or more worker instances. These run tasks and host HDFS.
- **Preemptible Worker**: You may have zero or more of these. These run tasks, but do *not* host HDFS. This is mostly useful because your cluster can lose task instances without killing your job (see Preemptible VMs).

By default, mrjob runs a single n1-standard-1, which is a cheap but not very powerful instance type. This can be quite adequate for testing your code on a small subset of your data, but otherwise give little advantage over running a job locally. To get more performance out of your job, you can either add more instances, use more powerful instances, or both.

Here are some things to consider when tuning your instance settings:

- Google Cloud bills you a 10-minute minimum even if your cluster only lasts for a few minutes (this is an artifact of the Google Cloud billing structure), so for many jobs that you run repeatedly, it is a good strategy to pick instance settings that make your job consistently run in a little less than 10 minutes.
- Your job will take much longer and may fail if any task (usually a reducer) runs out of memory and starts using swap. (You can verify this by using **vmstat**.) Restructuring your job is often the best solution, but if you can't, consider using a high-memory instance type.
- Larger instance types are usually a better deal if you have the workload to justify them. For example, a n1-highcpu-8 costs about 6 times as much as an n1-standard-1, but it has about 8 times as much processing power (and more memory).

The basic way to control type and number of instances is with the instance\_type and num\_core\_instances options, on the command line like this:

```
--instance-type n1-highcpu-8 --num-core-instances 4
```

or in mrjob.conf, like this:

```
runners:
   dataproc:
    instance_type: n1-highcpu-8
    num_core_instances: 4
```

In most cases, your master instance type doesn't need to be larger than n1-standard-1 to schedule tasks. *instance\_type* only applies to instances that actually run tasks. (In this example, there are 1 n1-standard-1 master instance, and 4 n1-highcpu-8 worker instances.) You *will* need a larger master instance if you have a very large number of input files; in this case, use the master\_instance\_type option.

If you want to run preemptible instances, use the task\_instance\_type and num\_task\_instances options.

# 1.16.2 Dataproc runner options

All options from Options available to all runners, Hadoop-related options, and Cloud runner options are available when running jobs on Google Cloud Dataproc.

## Google credentials

Basic credentials are not set in the config file; see Getting started with Google Cloud for details.

project\_id (--project-id) [string] Default: read from credentials config file

The ID of the Google Cloud Project to run under.

Changed in version 0.6.2: This used to be called *gcp\_project* 

```
service account (--service-account) [] Default: None
```

Optional service account to use when creating a cluster. For more information see Service Accounts.

New in version 0.6.3.

### service\_account\_scopes (--service-account-scopes) [] Default: (automatic)

Optional service account scopes to pass to the API when creating a cluster.

Generally it's suggested that you instead create a service\_account with the scopes you want.

New in version 0.6.3.

### Job placement

See also subnet, region, zone

```
network (--network) [string] Default: None
```

Name or URI of network to launch cluster in. Incompatible with with subnet.

New in version 0.6.3.

### **Cluster configuration**

```
cluster_properties (--cluster-property) [] Default: None
```

A dictionary of properties to set in the cluster's config files (e.g. mapred-site.xml). For details, see Cluster properties.

```
core_instance_config (--core-instance-config) [] Default: None
```

A dictionary of additional parameters to pass as config.worker\_config when creating the cluster. Follows the format of InstanceGroupConfig except that it uses <code>snake\_case</code> instead of <code>camel\_case</code>.

For example, to specify 100GB of disk space on core instances, add this to your config file:

```
runners:
    dataproc:
        core_instance_config:
        disk_config:
        boot_disk_size_gb: 100
```

To set this option on the command line, pass in JSON:

```
--core-instance-config '{"disk_config": {"boot_disk_size_gb": 100}}'
```

This option can be used to set number of core instances (num\_instances) or instance type (machine\_type\_uri), but usually you'll want to use num\_core\_instances and core\_instance\_type along with this option.

New in version 0.6.3.

### master\_instance\_config (--master-instance-config) [] Default: None

A dictionary of additional parameters to pass as config.master\_config when creating the cluster. See core\_instance\_config for more details.

New in version 0.6.3.

#### task\_instance\_config (--task-instance-config) [] Default: None

A dictionary of additional parameters to pass as config.secondary\_worker\_config when creating the cluster. See task\_instance\_config for more details.

To make task instances preemptible, add this to your config file:

```
runners:
   dataproc:
     task_instance_config:
     is_preemptible: true
```

Note that this config won't be applied unless you specify at least one task instance (either through num\_task\_instances or by passing num\_instances to this option).

New in version 0.6.3.

## 1.16.3 Other rarely used options

```
gcloud_bin (--gcloud-bin) [command] Default: 'gcloud'
```

Path to the gcloud binary; may include switches (e.g. 'gcloud -v' or ['gcloud', '-v']). Defaults to gcloud.

Used only as a way to create an SSH tunnel to the Resource Manager.

Changed in version 0.6.8: Setting this to an empty value (--gcloud-bin '') instructs mrjob to use the default (used to disable SSH).

# 1.17 Elastic MapReduce

## 1.17.1 Elastic MapReduce Quickstart

## **Configuring AWS credentials**

Configuring your AWS credentials allows mrjob to run your jobs on Elastic MapReduce and use S3.

- · Create an Amazon Web Services account
- Go to Security Credentials in the login menu (upper right, third from the right), say yes, you want to proceed, click on Access Keys, and then Create New Access Key. Make sure to copy the secret access key, as there is no way to recover it after creation.

Now you can either set the environment variables AWS\_ACCESS\_KEY\_ID and AWS\_SECRET\_ACCESS\_KEY, or set aws\_access\_key\_id and aws\_secret\_access\_key in your mrjob.conf file like this:

```
runners:
   emr:
   aws_access_key_id: <your key ID>
   aws_secret_access_key: <your secret>
```

## **Configuring SSH credentials**

Configuring your SSH credentials lets mrjob open an SSH tunnel to your jobs' master nodes to view live progress, see the job tracker in your browser, and fetch error logs quickly.

- Go to https://console.aws.amazon.com/ec2/home
- Make sure the **Region** dropdown (upper right, second from the right) matches the region you want to run jobs in (usually "Oregon").
- Click on **Key Pairs** (left sidebar, under **Network & Security**)

- Click on **Create Key Pair** (top left).
- Name your key pair EMR (any name will work but that's what we're using in this example)
- Save EMR.pem wherever you like (~/.ssh is a good place)
- Run chmod og-rwx /path/to/EMR.pem so that ssh will be happy
- Add the following entries to your mr job.conf:

```
runners:
    emr:
    ec2_key_pair: EMR
    ec2_key_pair_file: /path/to/EMR.pem # ~/ and $ENV_VARS allowed here
    ssh_tunnel: true
```

### **Running an EMR Job**

Running a job on EMR is just like running it locally or on your own Hadoop cluster, with the following changes:

- The job and related files are uploaded to S3 before being run
- The job is run on EMR (of course)
- Output is written to S3 before mrjob streams it to stdout locally
- The Hadoop version is specified by the EMR AMI version

This the output of this command should be identical to the output shown in Fundamentals, but it should take much longer:

> python word\_count.py -r emr README.txt "chars" 3654 "lines" 123 "words" 417

### Sending Output to a Specific Place

If you'd rather have your output go to somewhere deterministic on S3, use --output-dir:

```
> python word_count.py -r emr README.rst \
> --output-dir=s3://my-bucket/wc_out/
```

There are many other ins and outs of effectively using mrjob with EMR. See Advanced EMR usage for some of the ins, but the outs are left as an exercise for the reader. This is a strictly no-outs body of documentation!

### **Choosing Type and Number of EC2 Instances**

When you create a cluster on EMR, you'll have the option of specifying a number and type of EC2 instances, which are basically virtual machines. Each instance type has different memory, CPU, I/O and network characteristics, and costs a different amount of money. See Instance Types and Pricing for details.

Instances perform one of three roles:

- Master: There is always one master instance. It handles scheduling of tasks (i.e. mappers and reducers), but
  does not run them itself.
- Core: You may have one or more core instances. These run tasks and host HDFS.
- Task: You may have zero or more of these. These run tasks, but do *not* host HDFS. This is mostly useful because your cluster can lose task instances without killing your job (see *Spot Instances*).

There's a special case where your cluster *only* has a single master instance, in which case the master instance schedules tasks, runs them, and hosts HDFS.

By default, mrjob runs a single m1.medium, which is a cheap but not very powerful instance type. This can be quite adequate for testing your code on a small subset of your data, but otherwise give little advantage over running a job locally. To get more performance out of your job, you can either add more instances, use more powerful instances, or both.

Here are some things to consider when tuning your instance settings:

- Your job will take much longer and may fail if any task (usually a reducer) runs out of memory and starts using swap. (You can verify this by running mrjob boss j-CLUSTERID vmstat and then looking in j-CLUSTERID/\*/stdout.) Restructuring your job is often the best solution, but if you can't, consider using a high-memory instance type.
- Larger instance types are usually a better deal if you have the workload to justify them. For example, a cl.xlarge costs about 6 times as much as an ml.medium, but it has about 8 times as much processing power (and more memory).

The basic way to control type and number of instances is with the *instance\_type* and *num\_core\_instances* options, on the command line like this:

```
--instance-type c1.medium --num-core-instances 4
```

or in mrjob.conf, like this:

```
runners:
    emr:
    instance_type: c1.medium
    num_core_instances: 4
```

In most cases, your master instance type doesn't need to be larger than m1.medium to schedule tasks, so *instance\_type* only applies to the 4 instances that actually run tasks. You *will* need a larger master instance if you have a very large number of input files; in this case, use the *master\_instance\_type* option.

The *num\_task\_instances* option can be used to run 1 or more task instances (these run tasks but don't host HDFS). There are also *core\_instance\_type* and *task\_instance\_type* options if you want to set these directly.

# 1.17.2 Cluster Pooling

Clusters on EMR take several minutes to spin up, which can make development painfully slow.

To get around this, mrjob provides **cluster pooling.** If you set pool\_clusters to true, once your job completes, the cluster will stay open to accept additional jobs, and eventually shut itself down after it has been idle for a certain amount of time (by default, ten minutes; see max\_mins\_idle).

**Warning:** When using cluster pooling prior to v0.6.0, make sure to set max\_hours\_idle, or your cluster will never shut down.

**Note:** Pooling is a way to reduce latency, not to save money. Though pooling was originally created to optimize AWS's practice of billing by the full hour, this ended in October 2017.

Pooling is designed so that jobs run against the same mrjob.conf can share the same clusters. This means that the version of mrjob and bootstrap configuration. Other options that affect which cluster a job can join:

- image\_version/release\_label: must match
- image id (or lack thereof) must match

- applications: require at least these applications (extra ones okay)
- emr configurations: must match
- ec2\_key\_pair: if specified, only join clusters with the same key pair
- subnet: only join clusters with the same EC2 subnet ID (or lack thereof)

Pooled jobs will also only use clusters with the same **pool name**, so you can use the pool\_name option to partition your clusters into separate pools.

Pooling is flexible about instance type and number of instances; it will attempt to select the most powerful cluster available as long as the cluster's instances provide at least as much memory and at least as much CPU as your job requests.

Pooling is also somewhat flexible about EBS volumes (see instance\_groups). Each volume must have the same volume type, but larger volumes or volumes with more I/O ops per second are acceptable, as are additional volumes of any type.

Pooling cannot match configurations with explicitly set ebs\_root\_volume\_gb against clusters that use the default (or vice versa) because the EMR API does not report what the default value is.

If you are using instance\_fleets, your jobs will only join other clusters which use instance fleets. The rules are similar, but jobs will only join clusters whose fleets use the same set of instances or a subset; there is no concept of "better" instances.

mrjob's pooling won't add more than 1000 steps to a cluster, as the EMR API won't show more than this many steps. (For very old AMIs there is a stricter limit of 256 steps).

mrjob also uses an S3-based "locking" mechanism to prevent two jobs from simultaneously joining the same cluster. This is somewhat ugly but works in practice, and avoids mrjob depending on Amazon services other than EMR and S3.

**Warning:** If S3 eventual consistency takes longer than cloud\_fs\_sync\_secs, then you may encounter race conditions when using pooling, e.g. two jobs claiming the same cluster at the same time, or the idle cluster killer shutting down your job before it has started to run. Regions with read-after-write consistency (i.e. every region except US Standard) should not experience these issues.

You can allow jobs to wait for an available cluster instead of immediately starting a new one by specifying a value for *—pool-wait-minutes*. If none is found during that time, mrjob will start a new one.

## 1.17.3 EMR runner options

All options from Options available to all runners, Hadoop-related options, and Cloud runner options are available when running jobs on Amazon Elastic MapReduce.

### **Amazon credentials**

See Configuring AWS credentials and Configuring SSH credentials for specific instructions about setting these options.

aws\_access\_key\_id [string] Default: None

"Username" for Amazon web services.

There isn't a command-line switch for this option because credentials are supposed to be secret! Use the environment variable AWS\_ACCESS\_KEY\_ID instead.

### aws\_secret\_access\_key (--aws-secret-access-key) [string] Default: None

Your "password" on AWS.

There isn't a command-line switch for this option because credentials are supposed to be secret! Use the environment variable AWS\_SECRET\_ACCESS\_KEY instead.

#### aws\_session\_token [string] Default: None

Temporary AWS session token, used along with aws\_access\_key\_id and aws\_secret\_access\_key when using temporary credentials.

There isn't a command-line switch for this option because credentials are supposed to be secret! Use the environment variable AWS\_SESSION\_TOKEN instead.

Changed in version 0.5.10: this used to be called *aws\_security\_token*.

### ec2\_key\_pair (--ec2-key-pair) [string] Default: None

name of the SSH key you set up for EMR.

### ec2\_key\_pair\_file (--ec2-key-pair-file) [path] Default: None

path to file containing the SSH key for EMR

## iam\_instance\_profile (--iam-instance-profile) [string] Default: (automatic)

Name of an IAM instance profile to use for EC2 clusters created by EMR. See <a href="http://docs.aws.amazon.com/ElasticMapReduce/latest/DeveloperGuide/emr-iam-roles.html">http://docs.aws.amazon.com/ElasticMapReduce/latest/DeveloperGuide/emr-iam-roles.html</a> for more details on using IAM with EMR.

### iam\_service\_role (--iam-service-role) [string] Default: (automatic)

Name of an IAM role for the EMR service to use. See http://docs.aws.amazon.com/ElasticMapReduce/latest/DeveloperGuide/emriam-roles.html for more details on using IAM with EMR.

### Instance configuration

On EMR, there are three ways to configure instances:

- instance\_fleets
- instance\_groups
- individual instance options:
  - core\_instance\_bid\_price
  - core\_instance\_type
  - instance\_type
  - master\_instance\_bid\_price
  - master instance type
  - num\_core\_instances
  - num\_task\_instances
  - task\_instance\_bid\_price,
  - task\_instance\_type

If there is a conflict, whichever comes later in the config files takes precedence, and the command line beats config files. In the case of a tie, *instance\_fleets* beats *instance\_groups* beats other instance options.

You may set ebs\_root\_volume\_gb regardless of which style of instance configuration you use.

### instance\_fleets (--instance-fleet) [] Default: None

A list of instance fleet definitions to pass to the EMR API. Pass a JSON string on the command line or use data structures in the config file (which is itself basically JSON). For example:

```
runners:
  emr:
    instance fleets:
    - InstanceFleetType: MASTER
      InstanceTypeConfigs:
      - InstanceType: m1.medium
      TargetOnDemandCapacity: 1
    - InstanceFleetType: CORE
      TargetSpotCapacity: 2
      TargetOnDemandCapacity: 2
      LaunchSpecifications:
        SpotSpecification:
          TimeoutDurationMinutes: 20
          TimeoutAction: SWITCH_TO_ON_DEMAND
      InstanceTypeConfigs:
      - InstanceType: ml.medium
        BidPriceAsPercentageOfOnDemandPrice: 50
        WeightedCapacity: 1
      - InstanceType: m1.large
        BidPriceAsPercentageOfOnDemandPrice: 50
        WeightedCapacity: 2
```

#### instance groups (--instance-groups) [] Default: None

A list of instance group definitions to pass to the EMR API. Pass a JSON string on the command line or use data structures in the config file (which is itself basically JSON).

This allows for more fine-tuned EBS volume configuration than ebs\_root\_volume\_gb. For example:

```
runners:
    emr:
    instance_groups:
    - InstanceRole: MASTER
        InstanceType: m1.medium
- InstanceRole: CORE
        InstanceCount: 10
        InstanceType: c1.xlarge
        EbsConfiguration:
        EbsOptimized: true
        EbsBlockDeviceConfigs:
        - VolumeSpecification:
            SizeInGB: 100
            VolumeType: gp2
```

instance\_groups is incompatible with instance\_fleets and other instance options. See instance\_fleets for details.

### core\_instance\_bid\_price (--core-instance-bid-price) [string] Default: None

When specified and not "0", this creates the core Hadoop nodes as spot instances at this bid price. You usually only want to set bid price for task instances.

Changed in version 0.5.4: This option used to be named *ec2\_core\_instance\_bid\_price*.

```
master_instance_bid_price (--master-instance-bid-price) [string] Default: None
```

When specified and not "0", this creates the master Hadoop node as a spot instance at this bid price. You usually only want to set bid price for task instances unless the master instance is your only instance.

Changed in version 0.5.4: This option used to be named ec2\_master\_instance\_bid\_price.

```
task_instance_bid_price (--task-instance-bid-price) [string] Default: None
```

When specified and not "0", this creates the master Hadoop node as a spot instance at this bid price. (You usually only want to set bid price for task instances.)

Changed in version 0.5.4: This option used to be named ec2\_task\_instance\_bid\_price.

```
ebs_root_volume_gb (--ebs-root-volume-gb) [integer] Default: None
```

When specified (and not zero), sets the size of the root EBS volume, in GiB.

New in version 0.6.5.

### **Cluster software configuration**

See also bootstrap, image\_id, and image\_version.

```
applications (--application, --applications) [string list] Default: []
```

Additional applications to run on 4.x AMIs (e.g. 'Ganglia', 'Mahout', 'Spark').

You do not need to specify 'Hadoop'; mrjob will always include it automatically. In most cases it'll autodetect when to include 'Spark' as well.

See Applications in the EMR docs for more details.

New in version 0.5.2.

Changed in version 0.5.9: This used to be called *emr\_applications*.

Changed in version 0.6.7: Added --applications switch

```
bootstrap_actions (--bootstrap-actions) [string list] Default: []
```

A list of raw bootstrap actions (essentially scripts) to run prior to any of the other bootstrap steps. Any arguments should be separated from the command by spaces (we use <code>shlex.split()</code>). If the action is on the local filesystem, we'll automatically upload it to S3.

This has little advantage over bootstrap; it is included in order to give direct access to the EMR API.

#### bootstrap\_spark (--bootstrap-spark, --no-bootstrap-spark) [boolean] Default: (automatic)

Install Spark on the cluster. This works on AMI version 3.x and later.

By default, we automatically install Spark only if our job has Spark steps.

New in version 0.5.7.

In case you're curious, here's how mrjob determines you're using Spark:

- any SparkStep or SparkScriptStep in your job's steps (including implicitly through the spark method)
- "Spark" included in applications option
- any bootstrap action (see bootstrap\_actions) ending in /spark-install (this is how you install Spark
  on 3.x AMIs)

### emr\_configurations (--emr-configuration) [list of dicts] Default: []

Cluster configs for AMI version 4.x and later. For example:

```
runners:
    emr:
    emr_configurations:
    - Classification: core-site
    Properties:
        hadoop.security.groups.cache.secs: 250
```

On the command line, configurations should be JSON-encoded:

```
--emr-configuration '{"Classification": "core-site", ...}
```

See Configuring Applications in the EMR docs for more details.

New in version 0.5.3.

```
release_label (--release-label) [string] Default: None
```

EMR Release to use (e.g. emr-4.0.0). This overrides image\_version.

For more information about Release Labels, see Differences Introduced in 4.x.

New in version 0.5.0.

### Monitoring your job

See also check\_cluster\_every, ssh\_tunnel.

```
enable_emr_debugging (--enable-emr-debugging) [boolean] Default: False
```

store Hadoop logs in SimpleDB

### **Cluster pooling**

```
pool_clusters (--pool-clusters) [string] Default: True
```

Try to run the job on a WAITING pooled cluster with the same bootstrap configuration. Prefer the one with the most compute units. Use S3 to "lock" the cluster and ensure that the job is not scheduled behind another job. If no suitable cluster is *WAITING*, create a new pooled cluster.

**Warning:** If you use this in mrjob versions prior to 0.6.0, make sure to set max\_hours\_idle too, or your pooled clusters will run (costing you money) forever.

Changed in version 0.5.4: Pooling now gracefully recovers from joining a cluster that was in the process of shutting down (see max\_hours\_idle).

```
pool_name (--pool-name) [string] Default: 'default'
```

Specify a pool name to join. Does not imply pool\_clusters.

```
pool_wait_minutes (--pool-wait-minutes) [string] Default: 0
```

If pooling is enabled and no cluster is available, retry finding a cluster every 30 seconds until this many minutes have passed, then start a new cluster instead of joining one.

### S3 Filesystem

See also cloud\_tmp\_dir, cloud\_part\_size\_mb

cloud\_log\_dir (--cloud-log-dir) [string] Default: append logs to cloud\_tmp\_dir

Where on S3 to put logs, for example s3://yourbucket/logs/. Logs for your cluster will go into a subdirectory, e.g. s3://yourbucket/logs/j-CLUSTERID/.

Changed in version 0.5.4: This option used to be named s3\_log\_uri

### **API Endpoints**

**Note:** You usually don't want to set \*\_endpoint options unless you have a challenging network situation (e.g. you have to use a proxy to get around a firewall).

```
ec2_endpoint (--ec2-endpoint) [string] Default: (automatic)
```

New in version 0.6.5.

Force mrjob to connect to EC2 on this endpoint (e.g. ec2.us-qov-west-1.amazonaws.com).

```
emr_endpoint (--emr-endpoint) [string] Default: infer from region
```

Force mrjob to connect to EMR on this endpoint (e.g. us-west-1.elasticmapreduce.amazonaws.com).

```
iam_endpoint (--iam-endpoint) [string] Default: (automatic)
```

Force mrjob to connect to IAM on this endpoint (e.g. iam.us-qov.amazonaws.com).

```
s3_endpoint (--s3-endpoint) [string] Default: (automatic)
```

Force mrjob to connect to S3 on this endpoint, rather than letting it choose the appropriate endpoint for each S3 bucket.

**Warning:** If you set this to a region-specific endpoint (e.g. 's3-us-west-1.amazonaws.com') mrjob may not be able to access buckets located in other regions.

#### Other rarely used options

```
additional_emr_info (--additional-emr-info) [special] Default: None
```

Special parameters to select additional features, mostly to support beta EMR features. Pass a JSON string on the command line or use data structures in the config file (which is itself basically JSON).

```
emr_action_on_failure (--emr-action-on-failure) [string] Default: (automatic)
```

What happens if step of your job fails

- 'CANCEL\_AND\_WAIT' cancels all steps on the cluster
- 'CONTINUE' continues to the next step (useful when submitting several jobs to the same cluster)
- 'TERMINATE\_CLUSTER' shuts down the cluster entirely

The default is 'CANCEL\_AND\_WAIT' when using pooling (see pool\_clusters) or an existing cluster (see cluster id), and 'TERMINATE CLUSTER' otherwise.

hadoop\_streaming\_jar\_on\_emr (--hadoop-streaming-jar-on-emr) [string] Default: AWS default

Deprecated since version 0.5.4: Prepend file:// and pass that to hadoop\_streaming\_jar instead.

#### mins\_to\_end\_of\_hour (--mins-to-end-of-hour) [float] Default: 5.0

Deprecated since version 0.6.0: This option was created back when EMR billed by the full hour, and does nothing as of v0.6.0. If using versions prior to v0.6.0, it's recommended you set this to 60.0 to effectively disable this feature.

```
ssh bin (--ssh-bin) [command] Default: 'ssh'
```

Path to the ssh binary; may include switches (e.g. 'ssh -v' or ['ssh', '-v']). Defaults to ssh.

On EMR, mrjob uses SSH to tunnel to the job tracker (see ssh\_tunnel), as a fallback way of fetching job progress, and as a quicker way of accessing your job's logs.

Changed in version 0.6.8: Setting this to an empty value (--ssh-bin '') instructs mrjob to use the default value (used to effectively disable SSH).

```
tags (--tag) [dict] Default: {}
```

Metadata tags to apply to the EMR cluster after its creation. See Tagging Amazon EMR Clusters for more information on applying metadata tags to EMR clusters.

Tag names and values are strings. On the command line, to set a tag use --tag KEY=VALUE:

```
--tag team=development
```

In the config file, tags is a dict:

```
runners:
    emr:
    tags:
       team: development
    project: mrjob
```

Changed in version 0.5.4: This option used to be named *emr\_tags* 

If true (the default) EMR clusters will be visible to all IAM users. Otherwise, the cluster will only be visible to the IAM user that created it.

Deprecated since version 0.6.0: Hiding clusters from other users on the same account is not very useful. If you don't want to share pooled clusters, try pool\_name.

# 1.17.4 EMR Bootstrapping Cookbook

Bootstrapping allows you to run commands to customize EMR machines, at the time the cluster is created.

#### When to use bootstrap, and when to use setup

You can use bootstrap and setup together.

Generally, you want to use bootstrap for things that are part of your general production environment, and setup for things that are specific to your particular job. This makes things work as expected if you are using *Cluster Pooling*.

EMR will generally not allow you to use **sudo** in setup commands. See Job Environment Setup Cookbook for how to install libraries, etc. without using **sudo**.

### Installing Python packages with pip

The only tricky thing is making sure you install packages for the correct version of Python. Figure out which version of Python you'll be running on EMR (see python\_bin for defaults).

- If it's Python 2, use pip-2.7 (just plain pip also works on AMI 4.3.0 and later)
- If it's Python 3, use pip-3.4

For example, to install ujson on Python 2:

```
runners:
    emr:
    bootstrap:
    - sudo pip-2.7 install ujson
```

See PyPI for a the full list of available Python packages.

You can also install packages from a requirements file:

```
runners:
    emr:
    bootstrap:
    - sudo pip-2.7 install -r /local/path/of/requirements.txt#
```

Or a tarball:

```
runners:
    emr:
    bootstrap:
    - sudo pip-2.7 install /local/path/of/tarball.tar.gz#
```

**Warning:** If you're trying to run jobs on AMI version 3.0.0 (protip: don't do that) **pip** appears not to work due to out-of-date SSL certificate information.

### Installing PyPy

First, download the version of PyPy you want to use from Portable PyPy Distributions for Linux.

Then instruct EMR to un-tar it and link to the binary in /usr/bin. For example:

```
runners:
    emr:
    bootstrap:
        - sudo tar xvfj /local/path/to/pypy-7.1.1-linux_x86_64-portable.tar.bz2# -C /opt
        - sudo ln -s /opt/pypy-7.1.1-linux_x86_64-portable/bin/pypy /usr/bin/pypy
```

### Installing System Packages

EMR gives you access to a variety of different Amazon Machine Images, or AMIs for short (see image\_version).

### 3.x and 4.x AMIs

Starting with 3.0.0, EMR AMIs use Amazon Linux, which uses **yum** to install packages. For example, to install NumPy:

```
runners:
    emr:
    bootstrap:
        - sudo yum install -y python27-numpy
```

(Don't forget the -y!)

Amazon Linux currently has few packages for Python 3 libraries; if you're on Python 3, just use pip.

Here are the package lists for all the various versions of Amazon Linux used by EMR:

- 2015.09 (3.11.0 and 4.2.0-4.4.0)
- 2015.03 (3.7.0-3.10.0 and 4.0.0-4.1.0)
- 2014.09 (3.4.0-3.6.0)
- 2014.03 (3.1.0-3.3.2)
- 2013.09 (3.0.0-3.0.4)

**Note:** The package lists gloss over Python versions; wherever you see a package named python-<lib name>, you'll want to install python26-<lib name> or python27-<lib name> instead.

#### 2.x AMIs

Probably not worth the trouble. The 2.x AMIs are based on a version of Debian that is so old it has been "archived," which makes their package installer, **apt-get**, no longer work out-of-the-box. Moreover, Python system packages work for Python 2.6, not 2.7.

Instead, just use pip-2.7 to install Python libraries.

#### **Installing Python from source**

If you really must use a version of Python that's not available on EMR (e.g. Python 3.5 or a very specific patch version), you can download and compile Python from source.

**Note:** This adds an extra 5 to 10 minutes before the cluster can run your job.

Here's how you download and install a Python tarball:

```
runners:
    emr:
    bootstrap:
        - wget -S -T 10 -t 5 https://www.python.org/ftp/python/x.y.z/Python-x.y.z.tgz
        - tar xfz Python-x.y.z.tgz
        - cd Python-x.y.z; ./configure && make && sudo make install; cd ..
    bootstrap_python: false
    python_bin: /usr/local/bin/python
```

(Replace x.y.z with a specific version of Python.)

Python 3.4+ comes with pip by default, but earlier versions do not, so you'll want to tack on get-pip.py:

```
runners:
   emr:
   bootstrap:
   ...
```

```
- wget -S -T 10 -t 5 https://bootstrap.pypa.io/get-pip.py- sudo /usr/local/bin/python get-pip.py
```

Also, pip will be installed in /usr/local/bin, which is not in the path for sudo, so use its full path:

```
runners:
    emr:
    bootstrap:
        ...
        - sudo /usr/local/bin/pip install ...
```

## 1.17.5 Troubleshooting

Many things can go wrong in an EMR job, and the system's distributed nature can make it difficult to find the source of a problem. mrjob attempts to simplify the debugging process by automatically scanning logs for probable causes of failure.

In addition to looking at S3, mrjob can be configured to also use SSH to fetch error logs directly from the master and worker nodes. This can speed up debugging significantly (EMR only transfers logs to S3 every five minutes).

### Using persistent clusters

When troubleshooting a job, it can be convenient to use a persistent cluster to avoid having to wait for bootstrapping every run.

Warning: Make sure you either use the --max-hours-idle option or have mrjob terminate-idle-clusters in your crontab, or you will billed for unused CPU time on any clusters you forget to terminate.

First, use the **mrjob create-cluster** to create a persistent cluster:

```
$ mrjob create-cluster --max-hours-idle 1
Using configs in /Users/davidmarin/.mrjob.conf
Using s3://mrjob-35cdec11663cb1cb/tmp/ as our temp dir on S3
Creating persistent cluster to run several jobs in...
Creating temp directory /var/folders/zv/jmtt5bxs6xl3kzt38470hcxm0000gn/T/no_script.davidmarin.2016032
Copying local files to s3://mrjob-35cdec11663cb1cb/tmp/no_script.davidmarin.20160324.231018.720057/filesbyhp30KB81XE
```

Now you can use the cluster ID to start the troublesome job:

```
$ python mrjob/examples/mr_boom.py README.rst -r emr --cluster-id j-3BYHP30KB81XE
Using configs in /Users/davidmarin/.mrjob.conf
Using s3://mrjob-35cdec11663cb1cb/tmp/ as our temp dir on S3
Creating temp directory /var/folders/zv/jmtt5bxs6x13kzt38470hcxm0000gn/T/mr_boom.davidmarin.20160324
Copying local files to s3://mrjob-35cdec11663cb1cb/tmp/mr_boom.davidmarin.20160324.231045.501027/file
Adding our job to existing cluster j-3BYHP30KB81XE
Waiting for step 1 of 1 (s-SGVW9B5LEXF5) to complete...

PENDING (cluster is STARTING: Provisioning Amazon EC2 capacity)
PENDING (cluster is BOOTSTRAPPING: Running bootstrap actions)
PENDING (cluster is BOOTSTRAPPING: Running bootstrap actions)
PENDING (cluster is BOOTSTRAPPING: Running bootstrap actions)
```

```
PENDING (cluster is BOOTSTRAPPING: Running bootstrap actions)
 PENDING (cluster is BOOTSTRAPPING: Running bootstrap actions)
 Opening ssh tunnel to resource manager...
 Connect to resource manager at: http://localhost:40069/cluster
 RUNNING for 9.2s
 RUNNING for 42.3s
    0.0% complete
 RUNNING for 72.6s
    5.0% complete
 RUNNING for 102.9s
    5.0% complete
 RUNNING for 133.4s
  100.0% complete
 FAILED
Cluster j-3BYHP30KB81XE is WAITING: Cluster ready after last step failed.
Attempting to fetch counters from logs...
Looking for step log in /mnt/var/log/hadoop/steps/s-SGVW9B5LEXF5 on ec2-52-37-112-240.us-west-2.comp
 Parsing step log: ssh://ec2-52-37-112-240.us-west-2.compute.amazonaws.com/mnt/var/log/hadoop/steps
Counters: 9
    Job Counters
            Data-local map tasks=1
           Failed map tasks=4
           Launched map tasks=4
           Other local map tasks=3
           Total megabyte-seconds taken by all map tasks=58125312
            Total time spent by all map tasks (ms) = 75684
            Total time spent by all maps in occupied slots (ms) = 227052
            Total time spent by all reduces in occupied slots (ms)=0
            Total vcore-seconds taken by all map tasks=75684
Scanning logs for probable cause of failure...
Looking for task logs in /mnt/var/log/hadoop/userlogs/application_1458861299388_0001 on ec2-52-37-11
 Parsing task syslog: ssh://ec2-52-37-112-240.us-west-2.compute.amazonaws.com/mnt/var/log/hadoop/usc
 Parsing task stderr: ssh://ec2-52-37-112-240.us-west-2.compute.amazonaws.com/mnt/var/log/hadoop/usc
Probable cause of failure:
PipeMapRed failed!
java.lang.RuntimeException: PipeMapRed.waitOutputThreads(): subprocess failed with code 1
   at org.apache.hadoop.streaming.PipeMapRed.waitOutputThreads(PipeMapRed.java:330)
   at org.apache.hadoop.streaming.PipeMapRed.mapRedFinished(PipeMapRed.java:543)
   at org.apache.hadoop.streaming.PipeMapper.close(PipeMapper.java:130)
   at org.apache.hadoop.mapred.MapRunner.run(MapRunner.java:81)
   at org.apache.hadoop.streaming.PipeMapRunner.run(PipeMapRunner.java:34)
   at org.apache.hadoop.mapred.MapTask.runOldMapper(MapTask.java:432)
   at org.apache.hadoop.mapred.MapTask.run(MapTask.java:343)
   at org.apache.hadoop.mapred.YarnChild$2.run(YarnChild.java:175)
   at java.security.AccessController.doPrivileged(Native Method)
   at javax.security.auth.Subject.doAs(Subject.java:415)
   at org.apache.hadoop.security.UserGroupInformation.doAs(UserGroupInformation.java:1$48)
   at org.apache.hadoop.mapred.YarnChild.main(YarnChild.java:170)
(from lines 37-50 of ssh://ec2-52-37-112-240.us-west-2.compute.amazonaws.com/mnt/var/lody/hadoop/user
caused by:
Traceback (most recent call last):
 File "mr_boom.py", line 10, in <module>
   MRBoom.run()
 File "/usr/lib/python3.4/dist-packages/mrjob/job.py", line 430, in run
```

```
mr_job.execute()
File "/usr/lib/python3.4/dist-packages/mrjob/job.py", line 439, in execute
    self.run_mapper(self.options.step_num)
File "/usr/lib/python3.4/dist-packages/mrjob/job.py", line 499, in run_mapper
    for out_key, out_value in mapper_init() or ():
File "mr_boom.py", line 7, in mapper_init
    raise Exception('BOOM')
Exception: BOOM

(from lines 1-12 of ssh://ec2-52-37-112-240.us-west-2.compute.amazonaws.com/mnt/var/log/hadoop/userle
while reading input from s3://mrjob-35cdec11663cblcb/tmp/mr_boom.davidmarin.20160324.231045.501027/fg
Step 1 of 1 failed
Killing our SSH tunnel (pid 52847)
```

Now you can fix the bug and try again, without having to wait for a new cluster to bootstrap.

**Note:** mrjob *can* fetch logs from persistent jobs even without SSH set up, but it has to pause 10 minutes to wait for EMR to transfer logs to S3, which defeats the purpose of rapid iteration.

## 1.17.6 Advanced EMR usage

### **Spot Instances**

You can potentially save money purchasing EC2 instances for your EMR clusters from AWS's spot market. The catch is that if someone bids more for instances that you're using, they can be taken away from your cluster. If this happens, you aren't charged, but your job may fail.

You can specify spot market bid prices using the *core\_instance\_bid\_price*, *master\_instance\_bid\_price*, and *task\_instance\_bid\_price* options to specify a price in US dollars. For example, on the command line:

```
--task-instance-bid-price 0.42
```

or in mrjob.conf:

```
runners:
    emr:
    task_instance_bid_price: '0.42'
```

(Note the quotes; bid prices are strings, not floats!)

Amazon has a pretty thorough explanation of why and when you'd want to use spot instances here. The brief summary is that either you don't care if your job fails, in which case you want to purchase all your instances on the spot market, or you'd need your job to finish but you'd like to save time and money if you can, in which case you want to run task instances on the spot market and purchase master and core instances the regular way.

*Cluster Pooling* interacts with bid prices more or less how you'd expect; a job will join a pool with spot instances only if it requested spot instances at the same price or lower.

### **Custom Python packages**

See Installing Python packages with pip and Installing System Packages.

### **Bootstrap-time configuration**

Some Hadoop options, such as the maximum number of running map tasks per node, must be set at bootstrap time and will not work with *–jobconf*. You must use Amazon's *configure-hadoop* script for this. For example, this limits the number of mappers and reducers to one per node:

```
--bootstrap-action="s3://elasticmapreduce/bootstrap-actions/configure-hadoop \
-m mapred.tasktracker.map.tasks.maximum=1 \
-m mapred.tasktracker.reduce.tasks.maximum=1"
```

**Note:** This doesn't work on AMI 4.0.0 and later.

### **Manually Reusing Clusters**

In some cases, it may be useful to have more fine-grained control than *Cluster Pooling* provides; for example, to run several related jobs on the same cluster.

**Warning:** If you do this on mrjob versions prior to 0.6.0, make sure to set max\_hours\_idle, or your manually created clusters will run forever, costing you money.

mrjob includes a utility to create persistent clusters without running a job. For example, this command will create a cluster with 12 EC2 instances (1 master and 11 core), taking all other options from mrjob.conf:

```
$ mrjob create-cluster --num-core-instances=11 --max-hours-idle 1
...
j-CLUSTERID
```

You can then add jobs to the cluster with the --cluster-id switch or the cluster\_id option in *mrjob.conf* (see EMRJobRunner. init ()):

```
$ python mr_my_job.py -r emr --cluster-id=j-CLUSTERID input_file.txt > out
...
Adding our job to existing cluster j-CLUSTERID
...
```

Debugging will be difficult unless you complete SSH setup (see *Configuring SSH credentials*) since the logs will not be copied from the master node to S3 before either five minutes pass or the cluster terminates.

# 1.18 Python 2 vs. Python 3

## 1.18.1 Raw protocols

Both because we don't want to break mrjob for Python 2 users, and to make writing jobs simple, jobs read their input as strs by default (even though str means bytes in Python 2 and unicode in Python 3).

The way this works in mrjob is that RawValueProtocol is actually an alias for one of two classes, BytesValueProtocol if you're in Python 2, and TextValueProtocol if you're in Python 3.

If you care about this distinction, you may want to explicitly set <code>INPUT\_PROTOCOL</code> to one of these. If your input has a well-defined encoding, probably you want <code>BytesValueProtocol</code>, and if it's a bunch of text that's mostly ASCII, with like, some stuff that... might be UTF-8? (i.e. most log files), you probably want <code>TextValueProtocol</code>. But most of the time it'll just work.

## 1.18.2 Bytes vs. strings

The following things are bytes in any version of Python (which means you need to use the bytes type and/or  $b' \dots c'$  constant

- data read or written by Protocols
- lines yielded by cat\_output()/stream\_output()
- anything read from cat ()

The stdin, stdout, and stderr attributes of MRJobs are always bytestreams (so, for example, self.stderr defaults to sys.stderr.buffer in Python 3).

Everything else (including file paths, URIs, arguments to commands, and logging messages) are strings; that is, strs on Python 3, and either unicodes or ASCII strs on Python 2. Like with <code>RawValueProtocol</code>, most of the time it'll just work even if you don't think about it.

## 1.18.3 python\_bin

python\_bin defaults to **python3** in Python 3, and **python** in Python 2 (except on EMR AMIs prior to 4.3.0, where we use **python2.7**)

## 1.18.4 Your Hadoop cluster

Whatever version of Python you use, you'll have to have a compatible version of Python installed on your Hadoop cluster. mrjob does its best to make this work on Elastic MapReduce (see bootstrap\_python), but if you're running on your own Hadoop cluster, this is up to you.

# 1.19 Contributing to mrjob

## 1.19.1 Contribution guidelines

mrjob is developed using a standard Github pull request process. Almost all code is reviewed in pull requests.

The general process for working on mrjob is:

- Fork the project on Github
- · Clone your fork to your local machine
- Create a feature branch from master (e.g. git branch delete\_all\_the\_code)
- Write code, commit often
- · Write test cases for all changed functionality
- Submit a pull request against master on Github
- · Wait for code review!

It would also help to discuss your ideas on the mailing list so we can warn you of possible merge conflicts with ongoing work or offer suggestions for where to put code.

Things that will make your branch more likely to be pulled:

- · Comprehensive, fast test cases
- Detailed explanation of what the change is and how it works

- · Reference relevant issue numbers in the tracker
- API backward compatibility

If you add a new configuration option, please try to do all of these things:

- Add command line switches that allow full control over the option
- Document the option and its switches in the appropriate file under docs

## 1.19.2 A guick tour through the code

mrjob's modules can be put in four categories:

- · Reading command line arguments and config files, and invoking machinery accordingly
  - mrjob.conf: Read config files
  - mrjob.launch: Invoke runners based on command line and configs
  - mrjob.options: Define command line options
- · Interacting with Hadoop Streaming
  - mrjob. job: Python interface for writing jobs
  - mrjob.protocol: Defining data formats between Python steps
- Runners and support; submitting the job to various MapReduce environments
  - mrjob.runner: Common functionality across runners
  - mrjob.hadoop: Submit jobs to Hadoop
  - mr job. step: Define/implement interface between runners and script steps
  - Local
    - \* mrjob.inline: Run Python-only jobs in-process
    - \* mrjob.local: Run Hadoop Streaming-only jobs in subprocesses
  - Google Cloud Dataproc
    - \* mrjob.dataproc: Submit jobs to Dataproc
  - Amazon Elastic MapReduce
    - \* mrjob.emr: Submit jobs to EMR
    - \* mrjob.pool: Utilities for cluster pooling functionality
    - \* mrjob.retry: Wrapper for S3 and EMR connections to handle recoverable errors
- Interacting with different "filesystems"
  - mrjob.fs.base: Common functionality
  - mrjob.fs.composite: Support multiple filesystems; if one fails, "fall through" to another
  - mrjob.fs.gcs: Google Cloud Storage
  - mrjob.fs.hadoop: HDFS
  - mrjob.fs.local: Local filesystem
  - mrjob.fs.s3: S3
  - mrjob.fs.ssh: SSH

## • Utilities

- mrjob.compat: Transparently handle differences between Hadoop versions
- mrjob.logs: Log interpretation (counters, probable cause of job failure)
- mrjob.parse: Parsing utilities for URIs, command line options, etc.
- mrjob.util: Utilities for dealing with files, command line options, various other things

# Reference

# 2.1 mrjob.ami - building custom AMIs

Utilities for creating custom AMIs.

```
mrjob.ami.describe_base_emr_images(ec2_client)
```

Fetch a list of Amazon Linux AMI images that are usable with EMR, with the most recent first. This can take several seconds.

```
Parameters ec2_client - a boto3 EC2 client, which can be obtained from
mrjob.emr.EMRJobRunner.make_ec2_client() or boto3.client('ec2')
```

For the sake of consistency, we have somewhat stricter requirements than the AWS documentation. Specifically:

- •Amazon Linux (not Amazon Linux 2)
- •HVM virtualization
- •x86\_64 architecture
- •single EBS volume \* standard volume type (not GP2)
- •stable version (no "testing" or "rc", only numbers and dots)

This only returns images going back to September 2016 (prior to that, EC2 used a different naming convention).

This returns a dictionary for each image, in the same response format as ec2\_client.describe\_images(). The *ImageId* field contains the AMI ID, and *Description* contains a human-readable description.

# 2.2 mrjob.cat - decompress files based on extension

Emulating the way Hadoop handles input files, decompressing compressed files based on their file extension.

This module also functions as a **cat** substitute that can handle compressed files. It it used by <code>local</code> mode and can function without the rest of the mrjob library.

```
mrjob.cat.bunzip2_stream(fileobj, bufsize=1024)
Decompress gzipped data on the fly.
```

### **Parameters**

- fileobj object supporting read ()
- **bufsize** number of bytes to read from *fileobj* at a time.

**Warning:** This yields decompressed chunks; it does *not* split on lines. To get lines, wrap this in  $to\_lines()$ .

```
mrjob.cat.decompress (readable, path, bufsize=1024)
```

Take a *readable* which supports the .read() method correponding to the given path and returns an iterator that yields chunks of bytes, possibly decompressing based on *path*.

if readable appears to be a fileobj, pass it through as-is.

if readable does not have a read () method, assume that it's a generator that yields chunks of bytes

```
mrjob.cat.gunzip_stream(fileobj, bufsize=1024)
```

Decompress gzipped data on the fly.

#### **Parameters**

- fileobj object supporting read ()
- **bufsize** number of bytes to read from *fileobj* at a time. The default is the same as in gzip.

**Warning:** This yields decompressed chunks; it does *not* split on lines. To get lines, wrap this in  $to\_lines()$ .

```
mrjob.cat.to_chunks(readable, bufsize=1024)
```

Convert readable, which is any object supporting read () (e.g. fileobjs) to a stream of non-empty bytes.

If readable has an \_\_iter\_\_ method but not a read method, pass through as-is.

# 2.3 mrjob.cmd: The mrjob command-line utility

The mr job command provides a number of sub-commands that help you run and monitor jobs.

The **mrjob** command comes with Python-version-specific aliases (e.g. **mrjob-3**, **mrjob-3.4**), in case you choose to install mrjob for multiple versions of Python. You may also run it as **python -m mrjob.cmd** <subcommand>.

# 2.3.1 audit-emr-usage

Audit EMR usage over the past 2 weeks, sorted by cluster name and user.

Usage:

```
mrjob audit-emr-usage > report
```

```
-c CONF_PATHS, --conf-path CONF_PATHS
Path to alternate mrjob.conf file to read from
--no-conf Don't load mrjob.conf even if it's available
--emr-endpoint EMR_ENDPOINT
Force mrjob to connect to EMR on this endpoint (e.g. us-west-1.elasticmapreduce.amazonaws.com). Default is to infer this from region.
-h, --help show this help message and exit
--max-days-ago MAX_DAYS_AGO
Max number of days ago to look at jobs. By default, we
```

```
go back as far as EMR supports (currently about 2 months)

-q, --quiet Don't print anything to stderr

--region REGION GCE/AWS region to run Dataproc/EMR jobs in.

--s3-endpoint S3_ENDPOINT

Force mrjob to connect to S3 on this endpoint (e.g. s3 -us-west-1.amazonaws.com). You usually shouldn't set this; by default mrjob will choose the correct endpoint for each S3 bucket based on its location.

-v, --verbose print more messages to stderr
```

## 2.3.2 boss

Run a command on every node of a cluster. Store stdout and stderr for results in OUTPUT DIR.

Usage:

```
mrjob boss CLUSTER_ID [options] "command string"
```

### Options:

```
-c CONF_PATHS, --conf-path CONF_PATHS
                      Path to alternate mrjob.conf file to read from
--no-conf
                      Don't load mrjob.conf even if it's available
--ec2-key-pair-file EC2_KEY_PAIR_FILE
                      Path to file containing SSH key for EMR
--emr-endpoint EMR_ENDPOINT
                      Force mrjob to connect to EMR on this endpoint (e.g.
                      us-west-1.elasticmapreduce.amazonaws.com). Default is
                      to infer this from region.
-h, --help
                      show this help message and exit
-o OUTPUT_DIR, --output-dir OUTPUT_DIR
                      Specify an output directory (default: CLUSTER_ID)
-q, --quiet
                      Don't print anything to stderr
--region REGION
                      GCE/AWS region to run Dataproc/EMR jobs in.
--s3-endpoint S3_ENDPOINT
                      Force mrjob to connect to S3 on this endpoint (e.g. s3
                      -us-west-1.amazonaws.com). You usually shouldn't set
                      this; by default mrjob will choose the correct
                      endpoint for each S3 bucket based on its location.
                     Name/path of ssh binary. Arguments are allowed (e.g.
--ssh-bin SSH_BIN
                      --ssh-bin 'ssh -v')
-v, --verbose
                      print more messages to stderr
```

### 2.3.3 create-cluster

Create a persistent EMR cluster to run clusters in, and print its ID to stdout.

Usage:

```
mrjob create-cluster
```

```
--additional-emr-info ADDITIONAL_EMR_INFO
A JSON string for selecting additional features on EMR
--applications APPLICATIONS, --application APPLICATIONS
```

```
Additional applications to run on 4.x and 5.x AMIs,
                      separated by commas (e.g. "Ganglia, Spark")
--bootstrap BOOTSTRAP
                      A shell command to set up libraries etc. before any
                      steps (e.g. "sudo apt-get -qy install python3"). You
                      may interpolate files available via URL or locally
                      with Hadoop Distributed Cache syntax ("sudo yum
                      install -y foo.rpm#")
--bootstrap-action BOOTSTRAP_ACTIONS
                      Raw bootstrap action scripts to run before any of the
                      other bootstrap steps. You can use --bootstrap-action
                      more than once. Local scripts will be automatically
                      uploaded to S3. To add arguments, just use quotes:
                      "foo.sh arg1 arg2"
                      Automatically zip up the mrjob library and install it
--bootstrap-mrjob
                      when we run the mrjob. This is the default. Use --no-
                      bootstrap-mrjob if you've already installed mrjob on
                      your Hadoop cluster.
--no-bootstrap-mrjob Don't automatically zip up the mrjob library and
                      install it when we run this job. Use this if you've
                      already installed mrjob on your Hadoop cluster.
                      Attempt to install a compatible version of Python at
--bootstrap-python
                      bootstrap time. Currently this only does anything for
                      Python 3, for which it is enabled by default.
--no-bootstrap-python
                      Don't automatically try to install a compatible
                      version of Python at bootstrap time.
                      Auto-install Spark on the cluster (even if not
--bootstrap-spark
                      needed).
--no-bootstrap-spark Don't auto-install Spark on the cluster.
--cloud-fs-sync-secs CLOUD_FS_SYNC_SECS
                      How long to wait for remote FS to reach eventual
                      consistency. This is typically less than a second but
                      the default is 5.0 to be safe.
--cloud-log-dir CLOUD_LOG_DIR
                      URI on remote FS to write logs into
--cloud-part-size-mb CLOUD_PART_SIZE_MB
                      Upload files to cloud FS in parts no bigger than this
                      many megabytes. Default is 100 MiB. Set to 0 to
                      disable multipart uploading entirely.
--cloud-upload-part-size CLOUD_PART_SIZE_MB
                      Deprecated alias for --cloud-part-size-mb
--cloud-tmp-dir CLOUD_TMP_DIR
                      URI on remote FS to use as our temp directory.
-c CONF_PATHS, --conf-path CONF_PATHS
                      Path to alternate mrjob.conf file to read from
--no-conf
                      Don't load mrjob.conf even if it's available
--core-instance-bid-price CORE_INSTANCE_BID_PRICE
                      Bid price to specify for core nodes when setting them
                      up as EC2 spot instances (you probably only want to do
                      this for task instances).
--core-instance-type CORE_INSTANCE_TYPE
                      Type of GCE/EC2 core instance(s) to launch
--ebs-root-volume-gb EBS_ROOT_VOLUME_GB
                      Size of root EBS volume, in GiB. Must be an
                      integer. Set to 0 to use the default
--ec2-endpoint EC2_ENDPOINT
                      Force mrjob to connect to EC2 on this endpoint (e.g.
```

```
ec2.us-west-1.amazonaws.com). Default is to infer this
                      from region.
--ec2-key-pair EC2_KEY_PAIR
                      Name of the SSH key pair you set up for EMR
--emr-action-on-failure EMR_ACTION_ON_FAILURE
                      Action to take when a step fails (e.g.
                      TERMINATE_CLUSTER, CANCEL_AND_WAIT, CONTINUE)
--emr-api-param EMR_API_PARAMS
                      Does nothing. Use --extra-cluster-param instead
--no-emr-api-param EMR_API_PARAMS
                      Does nothing. Use --extra-cluster-param instead
--emr-configuration EMR_CONFIGURATIONS
                      Configuration to use on 4.x AMIs as a JSON-encoded
                      dict; see http://docs.aws.amazon.com/ElasticMapReduce/
                      latest/ReleaseGuide/emr-configure-apps.html for
                      examples
--emr-endpoint EMR_ENDPOINT
                      Force mrjob to connect to EMR on this endpoint (e.g.
                      us-west-1.elasticmapreduce.amazonaws.com). Default is
                      to infer this from region.
--enable-emr-debugging
                      Enable storage of Hadoop logs in SimpleDB
--disable-emr-debugging
                      Disable storage of Hadoop logs in SimpleDB (the
                      default)
--extra-cluster-param EXTRA_CLUSTER_PARAMS
                      extra parameter to pass to cloud API when creating a
                      cluster, to access features not currently supported by
                      mrjob. Takes the form <param>=<value>, where value is
                      JSON or a string. Use <param>=null to unset a
                      parameter
-h, --help
                      show this help message and exit
--iam-endpoint IAM_ENDPOINT
                      Force mrjob to connect to IAM on this endpoint (e.g.
                      iam.us-gov.amazonaws.com)
--iam-instance-profile IAM_INSTANCE_PROFILE
                      EC2 instance profile to use for the EMR cluster -- see
                      "Configure IAM Roles for Amazon EMR" in AWS docs
--iam-service-role IAM_SERVICE_ROLE
                      IAM service role to use for the EMR cluster -- see
                      "Configure IAM Roles for Amazon EMR" in AWS docs
--image-id IMAGE_ID ID of custom AWS machine image (AMI) to use
--image-version IMAGE_VERSION
                      version of EMR/Dataproc machine image to run
--instance-fleets INSTANCE_FLEETS
                      detailed JSON list of instance fleets, including EBS
                      configuration. See docs for --instance-fleets at
                      http://docs.aws.amazon.com/cli/latest/reference/emr
                      /create-cluster.html
--instance-groups INSTANCE_GROUPS
                      detailed JSON list of EMR instance configs, including
                      EBS configuration. See docs for --instance-groups at
                      http://docs.aws.amazon.com/cli/latest/reference/emr
                      /create-cluster.html
--instance-type INSTANCE_TYPE
                      Type of GCE/EC2 instance(s) to launch GCE - e.g.
                      n1-standard-1, n1-highcpu-4, n1-highmem-4 -- See
                      https://cloud.google.com/compute/docs/machine-types
```

```
EC2 - e.g. ml.medium, c3.xlarge, r3.xlarge -- See
                      http://aws.amazon.com/ec2/instance-types/
--label LABEL
                      Alternate label for the job, to help us identify it.
--master-instance-bid-price MASTER_INSTANCE_BID_PRICE
                      Bid price to specify for the master node when setting
                      it up as an EC2 spot instance (you probably only want
                      to do this for task instances).
--master-instance-type MASTER_INSTANCE_TYPE
                      Type of GCE/EC2 master instance to launch
--max-hours-idle MAX_HOURS_IDLE
                      Please use --max-mins-idle instead
--max-mins-idle MAX_MINS_IDLE
                      If we create a cluster, have it automatically
                      terminate itself after it's been idle this many
                      minutes
--mins-to-end-of-hour MINS_TO_END_OF_HOUR
                      If --max-mins-idle is set, control how close to the
                      end of an hour the cluster can automatically terminate
                      itself (default is 5 minutes)
--num-core-instances NUM_CORE_INSTANCES
                      Total number of core instances to launch
--num-task-instances NUM_TASK_INSTANCES
                      Total number of task instances to launch
--owner OWNER
                     User who ran the job (default is the current user)
--pool-clusters
                      Add to an existing cluster or create a new one that
                      does not terminate when the job completes.
                     Don't run job on a pooled cluster (the default)
--no-pool-clusters
--pool-name POOL_NAME
                      Specify a pool name to join. Default is "default"
-q, --quiet
                      Don't print anything to stderr
--region REGION
                      GCE/AWS region to run Dataproc/EMR jobs in.
--release-label RELEASE_LABEL
                      Release Label (e.g. "emr-4.0.0"). Overrides --image-
                      version
--s3-endpoint S3_ENDPOINT
                      Force mrjob to connect to S3 on this endpoint (e.g. s3
                      -us-west-1.amazonaws.com). You usually shouldn't set
                      this; by default mrjob will choose the correct
                      endpoint for each S3 bucket based on its location.
--subnet SUBNET
                      ID of Amazon VPC subnet/URI of Google Compute Engine
                      subnetwork to launch cluster in.
--subnets SUBNET
                      Like --subnet, but with a comma-separated list, to
                      specify multiple subnets in conjunction with
                      --instance-fleets (EMR only)
--tag TAGS
                      Metadata tags to apply to the EMR cluster; should take
                      the form KEY=VALUE. You can use --tag multiple times
--task-instance-bid-price TASK_INSTANCE_BID_PRICE
                      Bid price to specify for task nodes when setting them
                      up as EC2 spot instances
--task-instance-type TASK_INSTANCE_TYPE
                      Type of GCE/EC2 task instance(s) to launch
-v, --verbose
                      print more messages to stderr
--visible-to-all-users
                      Make your cluster is visible to all IAM users on the
                      same AWS account (the default)
--no-visible-to-all-users
                      Hide your cluster from other IAM users on the same AWS
                      account.
```

zone ZONE	GCE	zone/AWS	availability	zone	to	run	Dataproc/EMR	
	jobs	in.						

## 2.3.4 diagnose

Print probable cause of error for a failed step.

Currently this only works on EMR.

Usage:

```
mrjob diagnose [opts] j-CLUSTERID
```

### Options:

```
-c CONF_PATHS, --conf-path CONF_PATHS
                      Path to alternate mrjob.conf file to read from
--no-conf
                      Don't load mrjob.conf even if it's available
--emr-endpoint EMR_ENDPOINT
                      Force mrjob to connect to EMR on this endpoint (e.g.
                      us-west-1.elasticmapreduce.amazonaws.com). Default is
                      to infer this from region.
-h, --help
                      show this help message and exit
-q, --quiet
                      Don't print anything to stderr
--region REGION
                      GCE/AWS region to run Dataproc/EMR jobs in.
--s3-endpoint S3_ENDPOINT
                      Force mrjob to connect to S3 on this endpoint (e.g. s3
                      -us-west-1.amazonaws.com). You usually shouldn't set
                      this; by default mrjob will choose the correct
                      endpoint for each S3 bucket based on its location.
--step-id STEP_ID
                     ID of a particular failed step to diagnose
-v, --verbose
                      print more messages to stderr
```

New in version 0.6.1.

# 2.3.5 report-long-jobs

Report jobs running for more than a certain number of hours (by default, 24.0). This can help catch buggy jobs and Hadoop/EMR operational issues.

Suggested usage: run this as a daily cron job with the -q option:

```
0 0 * * * mrjob report-long-jobs
```

```
-c CONF_PATHS, --conf-path CONF_PATHS

Path to alternate mrjob.conf file to read from

--no-conf Don't load mrjob.conf even if it's available

--emr-endpoint EMR_ENDPOINT

Force mrjob to connect to EMR on this endpoint (e.g. us-west-1.elasticmapreduce.amazonaws.com). Default is to infer this from region.

-x EXCLUDE, --exclude EXCLUDE

Exclude clusters that match the specified tags.

Specifed in the form TAG_KEY, TAG_VALUE.

-h, --help show this help message and exit
```

```
--min-hours MIN_HOURS

Minimum number of hours a job can run before we report it. Default: 24.0

-q, --quiet Don't print anything to stderr
--region REGION GCE/AWS region to run Dataproc/EMR jobs in.
--s3-endpoint S3_ENDPOINT

Force mrjob to connect to S3 on this endpoint (e.g. s3 -us-west-1.amazonaws.com). You usually shouldn't set this; by default mrjob will choose the correct endpoint for each S3 bucket based on its location.
-v, --verbose print more messages to stderr
```

## 2.3.6 run

```
mrjob run (path to script or executable) [options]
```

Run a job. While you can use this command to invoke your jobs, you can just as easily call python my\_job.py [options].

# 2.3.7 s3-tmpwatch

Delete all files in a given URI that are older than a specified time. The time parameter defines the threshold for removing files. If the file has not been accessed for *time*, the file is removed. The time argument is a number with an optional single-character suffix specifying the units: m for minutes, h for hours, d for days. If no suffix is specified, time is in hours.

Suggested usage: run this as a cron job with the -q option:

```
0 0 * * * mrjob s3-tmpwatch -q 30d s3://your-bucket/tmp/
```

#### Usage:

```
mrjob s3-tmpwatch [options] <time-untouched> <URIs>
```

### Options:

```
-c CONF_PATHS, --conf-path CONF_PATHS
                     Path to alternate mrjob.conf file to read from
--no-conf
                     Don't load mrjob.conf even if it's available
-h, --help
                    show this help message and exit
-q, --quiet
                     Don't print anything to stderr
--region REGION
                     GCE/AWS region to run Dataproc/EMR jobs in.
--s3-endpoint S3_ENDPOINT
                      Force mrjob to connect to S3 on this endpoint (e.g. s3
                      -us-west-1.amazonaws.com). You usually shouldn't set
                      this; by default mrjob will choose the correct
                      endpoint for each S3 bucket based on its location.
-t, --test
                      Don't actually delete any files; just log that we
-v, --verbose
                      print more messages to stderr
```

# 2.3.8 spark-submit

A drop-in replacement for **spark-submit** that can use mrjob's runners. For example, you can submit your spark job to EMR just by adding -r emr.

This also adds a few mrjob features that are not standard with **spark-submit**, such as --cmdenv, --dirs, and --setup.

New in version 0.6.7.

Changed in version 0.6.8: added local, spark runners, made spark the default (was hadoop)

#### Usage

```
mrjob spark-submit [-r <runner>] [options] <python file | app jar>
[app arguments]
```

```
All runners:
  -r {emr,hadoop,local,spark}, --runner {emr,hadoop,local,spark}
                        Where to run the job (default: "spark")
  --class MAIN_CLASS
                        Your application's main class (for Java /
                        Scala apps).
  --name NAME
                        The name of your application.
  --jars LIBJARS
                        Comma-separated list of jars to include on the
                        driverand executor classpaths.
  --packages PACKAGES
                        Comma-separated list of maven coordinates of jars to
                        include on the driver and executor classpaths. Will
                        search the local maven repo, then maven central and
                        any additional remote repositories given by
                        --repositories. The format for the coordinates should
                        be groupId:artifactId:version.
  --exclude-packages EXCLUDE_PACKAGES
                        Comma-separated list of groupId:artifactId, to
                        exclude while resolving the dependencies provided in
                        --packages to avoid dependency conflicts.
  --repositories REPOSITORIES
                        Comma-separated list of additional remote
                        repositories to search for the maven coordinates
                        given with --packages.
  --py-files PY_FILES
                        Comma-separated list of .zip, .egg, or .py files to
                        placeon the PYTHONPATH for Python apps.
                        Comma-separated list of files to be placed in the
  --files UPLOAD_FILES
                        working directory of each executor. Ignored on
                        local[*] master.
                        Arbitrary environment variable to set inside Spark,
  --cmdenv CMDENV
                        in the format NAME=VALUE.
  --conf JOBCONF
                        Arbitrary Spark configuration property, in the format
                        PROP=VALUE.
                        A command to run before each Spark executor in the
  --setup SETUP
                        shell ("touch foo"). In cluster mode, runs before the
                        Spark driver as well. You may interpolate files
                        available via URL or on your local filesystem using
                        Hadoop Distributed Cache syntax (". setup.sh#"). To
                        interpolate archives (YARN only), use #/: "cd
                        foo.tar.gz#/; make.
  --properties-file PROPERTIES_FILE
                        Path to a file from which to load extra properties.
                        If not specified, this will look for conf/spark-
                        defaults.conf.
  --driver-memory DRIVER_MEMORY
                        Memory for driver (e.g. 1000M, 2G) (Default: 1024M).
  --driver-java-options DRIVER_JAVA_OPTIONS
                        Extra Java options to pass to the driver.
```

```
--driver-library-path DRIVER_LIBRARY_PATH
                        Extra library path entries to pass to the driver.
  --driver-class-path DRIVER_CLASS_PATH
                        Extra class path entries to pass to the driver. Note
                        that jars added with --jars are automatically
                        included in the classpath.
  --executor-memory EXECUTOR_MEMORY
                        Memory per executor (e.g. 1000M, 2G) (Default: 1G).
  --proxy-user PROXY_USER
                        User to impersonate when submitting the application.
                        This argument does not work with --principal /
                        --keytab.
 -c CONF_PATHS, --conf-path CONF_PATHS
                        Path to alternate mrjob.conf file to read from
 --no-conf
                        Don't load mrjob.conf even if it's available
 -q, --quiet
                        Don't print anything to stderr
  -v, --verbose
                       print more messages to stderr
  -h, --help
                        show this message and exit
Spark and Hadoop runners only:
  --master SPARK_MASTER
                        spark://host:port, mesos://host:port,
                        yarn,k8s://https://host:port, or local. Defaults to
                        local[*] on spark runner, yarn on hadoop runner.
  --deploy-mode SPARK_DEPLOY_MODE
                        Whether to launch the driver program locally
                        ("client") or on one of the worker machines inside
                        the cluster ("cluster") (Default: client).
Cluster deploy mode only:
  --driver-cores DRIVER_CORES
                        Number of cores used by the driver (Default: 1).
Spark standalone or Mesos with cluster deploy mode only:
                        If given, restarts the driver on failure.
 --supervise
Spark standalone and Mesos only:
 --total-executor-cores TOTAL_EXECUTOR_CORES
                        Total cores for all executors.
Spark standalone and YARN only:
  --executor-cores EXECUTOR_CORES
                        Number of cores per executor. (Default: 1 in YARN
                        mode, or all available cores on the worker in
                        standalone mode)
YARN-only:
  --archives UPLOAD_ARCHIVES
                        Comma-separated list of archives to be extracted into
                        the working directory of each executor.
  --dirs UPLOAD_DIRS
                        Comma-separated list of directors to be archived and
                        then extracted into the working directory of each
                        executor.
  --queue QUEUE_NAME
                        The YARN queue to submit to (Default: "default").
  --num-executors NUM_EXECUTORS
                        Number of executors to launch (Default: 2). If
                        dynamic allocation is enabled, the initial number of
                        executors will be at least NUM.
```

```
--principal PRINCIPAL

Principal to be used to login to KDC, while running onsecure HDFS.

--keytab KEYTAB

The full path to the file that contains the keytab for the principal specified above. This keytab will be copied to the node running the Application Master via the Secure Distributed Cache, for renewing the login tickets and the delegation tokens periodically.
```

This also supports the same runner-specific switches as MRJobs (e.g. --hadoop-bin, --region).

### 2.3.9 terminate-cluster

Terminate an existing EMR cluster.

Usage:

```
mrjob terminate-cluster [options] CLUSTER_ID
```

Terminate an existing EMR cluster.

**Options:** 

```
-c CONF_PATHS, --conf-path CONF_PATHS
                      Path to alternate mrjob.conf file to read from
--no-conf
                      Don't load mrjob.conf even if it's available
--emr-endpoint EMR_ENDPOINT
                      Force mrjob to connect to EMR on this endpoint (e.g.
                      us-west-1.elasticmapreduce.amazonaws.com). Default is
                      to infer this from region.
-h, --help
                      show this help message and exit
-q, --quiet
                      Don't print anything to stderr
--region REGION
                      GCE/AWS region to run Dataproc/EMR jobs in.
--s3-endpoint S3_ENDPOINT
                      Force mrjob to connect to S3 on this endpoint (e.g. s3
                      -us-west-1.amazonaws.com). You usually shouldn't set
                      this; by default mrjob will choose the correct
                      endpoint for each S3 bucket based on its location.
-t, --test
                      Don't actually delete any files; just log that we
                      would
-v, --verbose
                      print more messages to stderr
```

### 2.3.10 terminate-idle-clusters

Terminate idle EMR clusters that meet the criteria passed in on the command line (or, by default, clusters that have been idle for one hour).

Suggested usage: run this as a cron job with the -q option:

```
*/30 * * * mrjob terminate-idle-clusters -q
```

Changed in version 0.6.4: Skips termination-protected idle clusters, rather than crashing. (This was also backported to mrjob v0.5.12.)

```
-c CONF_PATHS, --conf-path CONF_PATHS
                      Path to alternate mrjob.conf file to read from
--no-conf
                      Don't load mrjob.conf even if it's available
--dry-run
                      Don't actually kill idle jobs; just log that we would
--emr-endpoint EMR_ENDPOINT
                      Force mrjob to connect to EMR on this endpoint (e.g.
                      us-west-1.elasticmapreduce.amazonaws.com). Default is
                      to infer this from region.
-h, --help
                      show this help message and exit
--max-hours-idle MAX_HOURS_IDLE
                      Please use --max-mins-idle instead.
--max-mins-idle MAX_MINS_IDLE
                      Max number of minutes a cluster can go without
                      bootstrapping, running a step, or having a new step
                      created. This will fire even if there are pending
                      steps which EMR has failed to start. Make sure you set
                      this higher than the amount of time your jobs can take
                      to start instances and bootstrap.
--max-mins-locked MAX_MINS_LOCKED
                      Max number of minutes a cluster can be locked while
                      idle.
--mins-to-end-of-hour MINS_TO_END_OF_HOUR
                      Deprecated, does nothing.
--pool-name POOL_NAME
                      Only terminate clusters in the given named pool.
--pooled-only
                     Only terminate pooled clusters
-q, --quiet
                     Don't print anything to stderr
--region REGION
                     GCE/AWS region to run Dataproc/EMR jobs in.
--s3-endpoint S3_ENDPOINT
                      Force mrjob to connect to S3 on this endpoint (e.g. s3
                      -us-west-1.amazonaws.com). You usually shouldn't set
                      this; by default mrjob will choose the correct
                      endpoint for each S3 bucket based on its location.
--unpooled-only
                      Only terminate un-pooled clusters
-v, --verbose
                      print more messages to stderr
```

# 2.4 mrjob.compat - Hadoop version compatibility

Utility functions for compatibility with different version of hadoop.

```
mrjob.compat.jobconf_from_dict (jobconf, name, default=None)
Get the value of a jobconf variable from the given dictionary.
```

#### **Parameters**

- jobconf (dict) jobconf dictionary
- name (string) name of the jobconf variable (e.g. 'user.name')
- default fallback value

If the name of the jobconf variable is different in different versions of Hadoop (e.g. in Hadoop 2, map.input.file is mapreduce.map.input.file), we'll automatically try all variants before giving up.

Return default if that jobconf variable isn't set

### mrjob.compat.jobconf\_from\_env(variable, default=None)

Get the value of a jobconf variable from the runtime environment.

For example, a MRJob could use jobconf\_from\_env('map.input.file') to get the name of the file a mapper is reading input from.

If the name of the jobconf variable is different in different versions of Hadoop (e.g. in Hadoop 2.0, map.input.file is mapreduce.map.input.file), we'll automatically try all variants before giving up.

Return default if that jobconf variable isn't set.

### mrjob.compat.map\_version(version, version\_map)

Allows you to look up something by version (e.g. which jobconf variable to use, specifying only the versions where that value changed.

version is a string

version\_map is a map from version (as a string) that a value changed to the new value.

For efficiency, *version\_map* can also be a list of tuples of (LooseVersion(version\_as\_string), value), with oldest versions first.

If version is less than any version in version\_map, use the value for the earliest version in version\_map.

### mrjob.compat.translate\_jobconf(variable, version)

Translate variable to Hadoop version version. If it's not a variable we recognize, leave as-is.

```
mrjob.compat.translate_jobconf_dict(jobconf, hadoop_version=None)
```

Translates the configuration property name to match those that are accepted in hadoop\_version. Prints a warning message if any configuration property name does not match the name in the hadoop version. Combines the original jobconf with the translated jobconf.

**Returns** a map consisting of the original and translated configuration property names and values.

```
mrjob.compat.translate_jobconf_for_all_versions(variable)
```

Get all known variants of the given jobconf variable. Unlike translate\_jobconf(), returns a list.

```
mrjob.compat.uses_yarn(version)
```

Basically, is this Hadoop 2? This also handles versions in the zero series (0.23+) where YARN originated.

```
mrjob.compat.version_gte(version, cmp_version_str)
```

Return True if version >= *cmp\_version\_str*.

# 2.5 mrjob.conf - parse and write config files

"mrjob.conf" is the name of both this module, and the global config file for mr job.

# 2.5.1 Reading and writing mrjob.conf

```
mrjob.conf.find_mrjob_conf()
```

Look for mrjob.conf, and return its path. Places we look:

- •The location specified by MRJOB\_CONF
- •~/.mrjob.conf
- •/etc/mrjob.conf

Return None if we can't find it.

```
mrjob.conf.load_opts_from_mrjob_conf(runner_alias, conf_path=None, al-
ready loaded=None)
```

Load a list of dictionaries representing the options in a given mrjob.conf for a specific runner, resolving includes. Returns [(path, values)]. If *conf\_path* is not found, return [(None, {})].

#### **Parameters**

- runner\_alias (str) String identifier of the runner type, e.g. emr, local, etc.
- conf path (str) location of the file to load
- already\_loaded (*list*) list of real (according to os.path.realpath()) conf paths that have already been loaded (used by load\_opts\_from\_mrjob\_confs()).

Relative include: paths are relative to the real (after resolving symlinks) path of the including conf file

This will only load each config file once, even if it's referenced from multiple paths due to symlinks.

```
mrjob.conf.load_opts_from_mrjob_confs (runner_alias, conf_paths=None)
```

Load a list of dictionaries representing the options in a given list of mrjob config files for a specific runner. Returns [(path, values), ...]. If a path is not found, use (None, {}) as its value.

If conf\_paths is None, look for a config file in the default locations (see find\_mrjob\_conf()).

#### **Parameters**

- runner\_alias (str) String identifier of the runner type, e.g. emr, local, etc.
- conf\_path locations of the files to load

This will only load each config file once, even if it's referenced from multiple paths due to symlinks.

# 2.5.2 Combining options

Combiner functions take a list of values to combine, with later options taking precedence over earlier ones. None values are always ignored.

```
mrjob.conf.combine_cmds(*cmds)
```

Take zero or more commands to run on the command line, and return the last one that is not None. Each command should either be a list containing the command plus switches, or a string, which will be parsed with shlex.split(). The string must either be a byte string or a unicode string containing no non-ASCII characters.

Returns either None or a list containing the command plus arguments.

```
mrjob.conf.combine dicts(*dicts)
```

Combine zero or more dictionaries. Values from dicts later in the list take precedence over values earlier in the list.

If you pass in None in place of a dictionary, it will be ignored.

```
mrjob.conf.combine_envs(*envs)
```

Combine zero or more dictionaries containing environment variables. Environment variable values may be wrapped in ClearedValue.

Environment variables later from dictionaries later in the list take priority over those earlier in the list.

For variables ending with PATH, we prepend (and add a colon) rather than overwriting. Wrapping a path value in ClearedValue disables this behavior.

Environment set to ClearedValue (None) will *delete* environment variables earlier in the list, rather than setting them to None.

If you pass in None in place of a dictionary in envs, it will be ignored.

```
mrjob.conf.combine_jobconfs(*jobconfs)
```

Like combine\_dicts(), but non-string values are converted to Java-readable string (e.g. True becomes 'true'). Keys whose value is None are blanked out.

```
mrjob.conf.combine_lists(*seqs)
```

Concatenate the given sequences into a list. Ignore None values.

Generally this is used for a list of commands we want to run; the "default" commands get run before any commands specific to your job.

Strings, bytes, and non-sequence objects (e.g. numbers) are treated as single-item lists.

Changed in version 0.5.3: Dicts are treated as single-item lists

```
mrjob.conf.combine_local_envs(*envs)
```

Same as combine\_envs(), except that paths are combined using the local path separator (e.g; on Windows rather than:).

```
mrjob.conf.combine_path_lists(*path_seqs)
```

Concatenate the given sequences into a list. Ignore None values. Resolve ~ (home dir) and environment variables, and expand globs that refer to the local filesystem.

Can take single strings as well as lists.

```
mrjob.conf.combine_paths(*paths)
```

Returns the last value in paths that is not None. Resolve ~ (home dir) and environment variables.

```
mrjob.conf.combine_values(*values)
```

Return the last value in *values* that is not None.

The default combiner; good for simple values (booleans, strings, numbers).

# 2.6 mrjob.dataproc - run on Dataproc

### 2.6.1 Job Runner

```
class mrjob.dataproc.DataprocJobRunner(**kwargs)
```

Runs an MRJob on Google Cloud Dataproc. Invoked when you run your job with -r dataproc.

DataprocJobRunner runs your job in an Dataproc cluster, which is basically a temporary Hadoop cluster.

Input, support, and jar files can be either local or on GCS; use gs://... URLs to refer to files on GCS.

This class has some useful utilities for talking directly to GCS and Dataproc, so you may find it useful to instantiate it without a script:

```
from mrjob.dataproc import DataprocJobRunner
...
```

## 2.6.2 GCS Utilities

```
 \begin{array}{c} \textbf{class} \, \texttt{mrjob.dataproc.GCSFilesystem} \, (\textit{local\_tmp\_dir} = \textit{None}, \, \, \textit{credentials} = \textit{None}, \, \, \textit{project\_id} = \textit{None}, \\ part\_size = \textit{None}, \, \textit{location} = \textit{None}, \, \textit{object\_ttl\_days} = \textit{None}) \\ \textbf{Filesystem for Google Cloud Storage} \, (\textbf{GCS}) \, \textbf{URIs} \\ \end{array}
```

### **Parameters**

• local\_tmp\_dir - deprecated, does nothing, do not use

- **credentials** an optional google.auth.credentials.Credentials, used to initialize the storage client
- project\_id an optional project ID, used to initialize the storage client
- part\_size Part size for multi-part uploading, in bytes, or None
- **location** Default location to use when creating a bucket
- object\_ttl\_days Default object expiry for newly created buckets

Changed in version 0.6.8: deprecated local\_tmp\_dir, added part\_size, location, object\_ttl\_days

# 2.7 mrjob.emr - run on EMR

### 2.7.1 Job Runner

```
class mrjob.emr.EMRJobRunner(**kwargs)
```

Runs an MRJob on Amazon Elastic MapReduce. Invoked when you run your job with -r emr.

EMRJobRunner runs your job in an EMR cluster, which is basically a temporary Hadoop cluster. Normally, it creates a cluster just for your job; it's also possible to run your job in a specific cluster by setting *cluster\_id* or to automatically choose a waiting cluster, creating one if none exists, by setting *pool\_clusters*.

Input, support, and jar files can be either local or on S3; use \$3://... URLs to refer to files on S3.

This class has some useful utilities for talking directly to S3 and EMR, so you may find it useful to instantiate it without a script:

```
from mrjob.emr import EMRJobRunner

emr_client = EMRJobRunner().make_emr_client()
clusters = emr_client.list_clusters()
...
```

### 2.7.2 EMR Utilities

```
EMRJobRunner.get_cluster_id()
Get the ID of the cluster our job is running on, or None.

EMRJobRunner.get_image_version()
Get the version of the AMI that our cluster is running, or None.

Changed in version 0.5.4: This used to be called get_ami_version()

EMRJobRunner.get_job_steps()
Fetch the steps submitted by this runner from the EMR API.

New in version 0.6.1.

EMRJobRunner.make_emr_client()
Create a boto3 EMR client.
```

Returns a botocore.client.EMR wrapped in a mrjob.retry.RetryWrapper

# 2.7.3 S3 Utilities

```
class mrjob.fs.s3.S3Filesystem(aws_access_key_id=None, aws_secret_access_key=None, aws_session_token=None, s3_endpoint=None, s3_region=None, part_size=None)
```

Filesystem for Amazon S3 URIs. Typically you will get one of these via EMRJobRunner().fs, composed with SSHFilesystem and LocalFilesystem.

### **Parameters**

- aws\_access\_key\_id Your AWS access key ID
- aws\_secret\_access\_key Your AWS secret access key
- aws\_session\_token session token for use with temporary AWS credentials
- s3\_endpoint If set, always use this endpoint
- **s3\_region** Default region for connections to the S3 API and newly created buckets.
- part\_size Part size for multi-part uploading, in bytes, or None

Changed in version 0.6.8: added part\_size

S3Filesystem.create\_bucket (bucket\_name, region=None)

Create a bucket on S3 with a location constraint matching the given region.

Changed in version 0.6.0: The *region* argument used to be called *location*.

S3Filesystem.get\_all\_bucket\_names()

Get a list of the names of all buckets owned by this user on S3.

New in version 0.6.0.

S3Filesystem.get\_bucket(bucket\_name)

Get the (boto3) bucket, connecting through the appropriate endpoint.

S3Filesystem.make\_s3\_client(region\_name=None)

Create a boto3 S3 client, wrapped in a mrjob.retry.RetryWrapper

**Parameters** region – region to use to choose S3 endpoint.

New in version 0.6.0.

S3Filesystem.make s3 resource(region name=None)

Create a boto3 S3 resource, with its client wrapped in a mrjob.retry.RetryWrapper

**Parameters region** – region to use to choose S3 endpoint

It's best to use  $get\_bucket$  () because it chooses the appropriate S3 endpoint automatically. If you are trying to get bucket metadata, use  $make\_s3\_client$  ().

New in version 0.6.0.

# 2.7.4 Other AWS clients

```
EMRJobRunner.make_ec2_client()
```

Create a boto3 EC2 client.

Returns a botocore.client.EC2 wrapped in a mrjob.retry.RetryWrapper

EMRJobRunner.make\_iam\_client()

Create a boto3 IAM client.

Returns a botocore.client.IAM wrapped in a mrjob.retry.RetryWrapper

# 2.8 mrjob.hadoop - run on your Hadoop cluster

```
class mrjob.hadoop.HadoopJobRunner(**kwargs)
```

Runs an MRJob on your Hadoop cluster. Invoked when you run your job with -r hadoop.

Input and support files can be either local or on HDFS; use hdfs://... URLs to refer to files on HDFS.

```
HadoopJobRunner.__init__(**kwargs)
```

HadoopJobRunner takes the same arguments as MRJobRunner, plus some additional options which can be defaulted in *mrjob.conf* .

# 2.8.1 Utilities

```
mrjob.hadoop.fully_qualify_hdfs_path (path)
If path isn't an hdfs:// URL, turn it into one.
```

# 2.9 mrjob.inline - debugger-friendly local testing

```
class mrjob.inline.InlineMRJobRunner(mrjob_cls=None, **kwargs)
```

Runs an MRJob in the same process, so it's easy to attach a debugger.

This is the default way to run jobs (we assume you'll spend some time debugging your job before you're ready to run it on EMR or Hadoop).

Unlike other runners, InlineMRJobRunner's run() method raises the actual exception that caused a step to fail (rather than StepFailedException).

To more accurately simulate your environment prior to running on Hadoop/EMR, use -r local (see LocalMRJobRunner).

New in version 0.6.8: can run *SparkSteps* via the pyspark library.

```
InlineMRJobRunner.__init__(mrjob_cls=None, **kwargs)
```

InlineMRJobRunner takes the same keyword args as MRJobRunner. However, please note that hadoop\_input\_format, hadoop\_output\_format, and partitioner are ignored because they require Java. If you need to test these, consider starting up a standalone Hadoop instance and running your job with -r hadoop.

# 2.10 mrjob.job - defining your job

# 2.10.1 One-step jobs

```
MRJob.mapper(key, value)
```

Re-define this to define the mapper for a one-step job.

Yields zero or more tuples of (out\_key, out\_value).

### **Parameters**

- **key** A value parsed from input.
- **value** A value parsed from input.

If you don't re-define this, your job will have a mapper that simply yields (key, value) as-is.

### By default (if you don't mess with *Protocols*):

- key will be None
- value will be the raw input line, with newline stripped.
- out\_key and out\_value must be JSON-encodable: numeric, unicode, boolean, None, list, or dict whose keys are unicodes.

# MRJob.reducer(key, values)

Re-define this to define the reducer for a one-step job.

Yields one or more tuples of (out\_key, out\_value)

#### **Parameters**

- key A key which was yielded by the mapper
- **value** A generator which yields all values yielded by the mapper which correspond to key.

## By default (if you don't mess with *Protocols*):

- out\_key and out\_value must be JSON-encodable.
- key and value will have been decoded from JSON (so tuples will become lists).

#### MRJob.combiner(key, values)

Re-define this to define the combiner for a one-step job.

Yields one or more tuples of (out\_key, out\_value)

## **Parameters**

- key A key which was yielded by the mapper
- value A generator which yields all values yielded by one mapper task/node which correspond to key.

### By default (if you don't mess with *Protocols*):

- out\_key and out\_value must be JSON-encodable.
- key and value will have been decoded from JSON (so tuples will become lists).

### MRJob.mapper\_init()

Re-define this to define an action to run before the mapper processes any input.

One use for this function is to initialize mapper-specific helper structures.

Yields one or more tuples of (out\_key, out\_value).

By default, out\_key and out\_value must be JSON-encodable; re-define INTERNAL\_PROTOCOL to change this.

#### MRJob.mapper\_final()

Re-define this to define an action to run after the mapper reaches the end of input.

One way to use this is to store a total in an instance variable, and output it after reading all input data. See mrjob.examples for an example.

Yields one or more tuples of (out key, out value).

By default, out\_key and out\_value must be JSON-encodable; re-define INTERNAL\_PROTOCOL to change this.

### MRJob.reducer\_init()

Re-define this to define an action to run before the reducer processes any input.

One use for this function is to initialize reducer-specific helper structures.

Yields one or more tuples of (out\_key, out\_value).

By default, out\_key and out\_value must be JSON-encodable; re-define INTERNAL\_PROTOCOL to change this.

### MRJob.reducer\_final()

Re-define this to define an action to run after the reducer reaches the end of input.

Yields one or more tuples of (out\_key, out\_value).

By default,  $out\_key$  and  $out\_value$  must be JSON-encodable; re-define  $INTERNAL\_PROTOCOL$  to change this.

### MRJob.combiner\_init()

Re-define this to define an action to run before the combiner processes any input.

One use for this function is to initialize combiner-specific helper structures.

Yields one or more tuples of (out\_key, out\_value).

By default, out\_key and out\_value must be JSON-encodable; re-define INTERNAL\_PROTOCOL to change this.

# MRJob.combiner\_final()

Re-define this to define an action to run after the combiner reaches the end of input.

Yields one or more tuples of (out\_key, out\_value).

By default, out\_key and out\_value must be JSON-encodable; re-define INTERNAL\_PROTOCOL to change this.

### MRJob.mapper\_cmd()

Re-define this to define the mapper for a one-step job **as a shell command.** If you define your mapper this way, the command will be passed unchanged to Hadoop Streaming, with some minor exceptions. For important specifics, see *Shell commands as steps*.

Basic example:

```
def mapper_cmd(self):
    return 'cat'
```

## MRJob.reducer\_cmd()

Re-define this to define the reducer for a one-step job **as a shell command.** If you define your mapper this way, the command will be passed unchanged to Hadoop Streaming, with some minor exceptions. For specifics, see *Shell commands as steps*.

Basic example:

```
def reducer_cmd(self):
    return 'cat'
```

## MRJob.combiner\_cmd()

Re-define this to define the combiner for a one-step job **as a shell command.** If you define your mapper this way, the command will be passed unchanged to Hadoop Streaming, with some minor exceptions. For specifics, see *Shell commands as steps*.

### Basic example:

```
def combiner_cmd(self):
    return 'cat'
```

#### MRJob.mapper\_pre\_filter()

Re-define this to specify a shell command to filter the mapper's input before it gets to your job's mapper in a one-step job. For important specifics, see *Filtering task input with shell commands*.

Basic example:

```
def mapper_pre_filter(self):
    return 'grep "ponies"'
```

#### MRJob.reducer\_pre\_filter()

Re-define this to specify a shell command to filter the reducer's input before it gets to your job's reducer in a one-step job. For important specifics, see *Filtering task input with shell commands*.

Basic example:

```
def reducer_pre_filter(self):
    return 'grep "ponies"'
```

### MRJob.combiner\_pre\_filter()

Re-define this to specify a shell command to filter the combiner's input before it gets to your job's combiner in a one-step job. For important specifics, see *Filtering task input with shell commands*.

Basic example:

```
def combiner_pre_filter(self):
    return 'grep "ponies"'
```

## MRJob.mapper\_raw (input\_path, input\_uri)

Re-define this to make Hadoop pass one input file to each mapper.

#### **Parameters**

- input path a local path that the input file has been copied to
- input\_uri the URI of the input file on HDFS, S3, etc

New in version 0.6.3.

# MRJob.spark (input\_path, output\_path)

Re-define this with Spark code to run. You can read input with input\_path and output with output\_path.

**Warning:** Prior to v0.6.8, to pass job methods into Spark (rdd.flatMap(self.some\_method)), you first had to call self.sandbox(); otherwise Spark would error because self was not serializable.

# 2.10.2 Multi-step jobs

# MRJob.steps()

Re-define this to make a multi-step job.

If you don't re-define this, we'll automatically create a one-step job using any of mapper(), mapper\_init(), mapper\_final(), reducer\_init(), reducer\_final(), and reducer() that you've re-defined. For example:

**Returns** a list of steps constructed with MRStep or other classes in mrjob.step.

# 2.10.3 Running the job

```
classmethod MRJob.run()
```

Entry point for running job from the command-line.

This is also the entry point when a mapper or reducer is run by Hadoop Streaming.

Does one of:

```
Print step information (--steps). See show_steps()
Run a mapper (--mapper). See run_mapper()
Run a combiner (--combiner). See run_combiner()
Run a reducer (--reducer). See run_reducer()
Run the entire job. See run_job()
MRJob.__init__ (args=None)
```

Entry point for running your job from other Python code.

You can pass in command-line arguments, and the job will act the same way it would if it were run from the command line. For example, to run your job on EMR:

```
mr_job = MRYourJob(args=['-r', 'emr'])
with mr_job.make_runner() as runner:
...
```

Passing in None is the same as passing in [] (if you want to parse args from sys.argv, call MRJob.run()).

For a full list of command-line arguments, run: python -m mrjob.job --help

```
MRJob.make_runner()
```

Make a runner based on command-line arguments, so we can launch this job on EMR, on Hadoop, or locally.

```
Return type mrjob.runner.MRJobRunner
```

# 2.10.4 Parsing output

```
MRJob.parse_output (chunks)
```

Parse the final output of this MRJob (as a stream of byte chunks) into a stream of (key, value).

```
MRJob.parse_output_line(line)
```

Parse a line from the final output of this MRJob into (key, value).

Deprecated since version 0.6.0: Use parse\_output () instead.

# 2.10.5 Counters and status messages

### MRJob.increment\_counter(group, counter, amount=1)

Increment a counter in Hadoop streaming by printing to stderr.

#### **Parameters**

- **group** (*str*) counter group
- counter (str) description of the counter
- amount (int) how much to increment the counter by

Commas in counter or group will be automatically replaced with semicolons (commas confuse Hadoop streaming).

## MRJob.set\_status(msg)

Set the job status in hadoop streaming by printing to stderr.

This is also a good way of doing a keepalive for a job that goes a long time between outputs; Hadoop streaming usually times out jobs that give no output for longer than 10 minutes.

# 2.10.6 Setting protocols

## MRJob. INPUT\_PROTOCOL = <class 'mrjob.protocol.BytesValueProtocol'>

Protocol for reading input to the first mapper in your job. Default: RawValueProtocol.

For example you know your input data were in JSON format, you could set:

```
INPUT_PROTOCOL = JSONValueProtocol
```

in your class, and your initial mapper would receive decoded JSONs rather than strings.

See mrjob.protocol for the full list of protocols.

# MRJob.INTERNAL\_PROTOCOL = <class 'mrjob.protocol.StandardJSONProtocol'>

Protocol for communication between steps and final output. Default: JSONProtocol.

For example if your step output weren't JSON-encodable, you could set:

```
INTERNAL_PROTOCOL = PickleProtocol
```

and step output would be encoded as string-escaped pickles.

See mrjob.protocol for the full list of protocols.

### MRJob.OUTPUT\_PROTOCOL = <class 'mrjob.protocol.StandardJSONProtocol'>

Protocol to use for writing output. Default: JSONProtocol.

For example, if you wanted the final output in repr, you could set:

```
OUTPUT_PROTOCOL = ReprProtocol
```

See *mrjob.protocol* for the full list of protocols.

# MRJob.input\_protocol()

Instance of the protocol to use to convert input lines to Python objects. Default behavior is to return an instance of INPUT PROTOCOL.

#### MRJob.internal protocol()

Instance of the protocol to use to communicate between steps. Default behavior is to return an instance of INTERNAL\_PROTOCOL.

#### MRJob.output protocol()

Instance of the protocol to use to convert Python objects to output lines. Default behavior is to return an instance of OUTPUT PROTOCOL.

## MRJob.pick\_protocols (step\_num, step\_type)

Pick the protocol classes to use for reading and writing for the given step.

#### **Parameters**

- **step\_num** (*int*) which step to run (e.g. 0 for the first step)
- step\_type (str) one of 'mapper', 'combiner', or 'reducer'

**Returns** (read\_function, write\_function)

By default, we use one protocol for reading input, one internal protocol for communication between steps, and one protocol for final output (which is usually the same as the internal protocol). Protocols can be controlled by setting <code>INPUT\_PROTOCOL</code>, <code>INTERNAL\_PROTOCOL</code>, and <code>OUTPUT\_PROTOCOL</code>.

Re-define this if you need fine control over which protocols are used by which steps.

# 2.10.7 Secondary sort

#### MRJob.SORT VALUES = None

Set this to True if you would like reducers to receive the values associated with any key in sorted order (sorted by their *encoded* value). Also known as secondary sort.

This can be useful if you expect more values than you can fit in memory to be associated with one key, but you want to apply information in a small subset of these values to information in the other values. For example, you may want to convert counts to percentages, and to do this you first need to know the total count.

Even though values are sorted by their encoded value, most encodings will sort strings in order. For example, you could have values like: ['A', <total>], ['B', <count\_name>, <count>], and the value containing the total should come first regardless of what protocol you're using.

See jobconf() and partitioner() for more about

# 2.10.8 Command-line options

See *Defining command line options* for information on adding command line options to your job. See Configuration quick reference for a complete list of all configuration options.

#### MRJob.configure args()

Define arguments for this script. Called from \_\_init\_\_ ().

Re-define to define custom command-line arguments or pass through existing ones:

```
def configure_args(self):
    super(MRYourJob, self).configure_args()

self.add_passthru_arg(...)
    self.add_file_arg(...)
    self.pass_arg_through(...)
    ...
```

## MRJob.add\_passthru\_arg(\*args, \*\*kwargs)

Function to create options which both the job runner and the job itself respect (we use this for protocols, for example).

Use it like you would use argparse. ArgumentParser.add\_argument():

```
def configure_args(self):
    super(MRYourJob, self).configure_args()
    self.add_passthru_arg(
         '--max-ngram-size', type=int, default=4, help='...')
```

If you want to pass files through to the mapper/reducer, use add\_file\_arg() instead.

If you want to pass through a built-in option (e.g. --runner, use pass\_arg\_through() instead.

```
MRJob.add_file_arg(*args, **kwargs)
```

Add a command-line option that sends an external file (e.g. a SQLite DB) to Hadoop:

```
def configure_args(self):
    super(MRYourJob, self).configure_args()
    self.add_file_arg('--scoring-db', help=...)
```

This does the right thing: the file will be uploaded to the working dir of the script on Hadoop, and the script will be passed the same option, but with the local name of the file in the script's working directory.

**Note:** If you pass a file to a job, best practice is to lazy-load its contents (e.g. make a method that opens the file the first time you call it) rather than loading it in your job's constructor or <code>load\_args()</code>. Not only is this more efficient, it's necessary if you want to run your job in a Spark executor (because the file may not be in the same place in a Spark driver).

**Note:** We suggest against sending Berkeley DBs to your job, as Berkeley DB is not forwards-compatible (so a Berkeley DB that you construct on your computer may not be readable from within Hadoop). Use SQLite databases instead. If all you need is an on-disk hash table, try out the sqlite3dbm module.

Changed in version 0.6.6: now accepts explicit type=str

Changed in version 0.6.8: fully supported on Spark, including local[\*] master

```
MRJob.pass_arg_through(opt_str)
```

Pass the given argument through to the job.

```
MRJob.load_args (args)
```

Load command-line options into self.options and self.\_script\_path.

```
Called from __init__ () after configure_args().
```

**Parameters**  $args(list\ of\ str) - a$  list of command line arguments. None will be treated the same as [].

Re-define if you want to post-process command-line arguments:

```
def load_args(self, args):
    super(MRYourJob, self).load_args(args)

self.stop_words = self.options.stop_words.split(',')
...
```

```
MRJob.is_task()
```

True if this is a mapper, combiner, reducer, or Spark script.

This is mostly useful inside <code>load\_args()</code>, to disable loading args when we aren't running inside Hadoop.

# 2.10.9 Uploading support files

#### MRJob.FILES = []

Optional list of files to upload to the job's working directory. These can be URIs or paths on the local filesystem.

Relative paths will be interpreted as relative to the directory containing the script (not the current working directory). Environment variables and ~ in paths will be expanded.

If you want a file to be uploaded to a filename other than it's own, append #<name> (e.g. data/foo.json#bar.json).

If you need to dynamically generate a list of files, override files () instead.

New in version 0.6.4.

### MRJob.**DIRS** = []

Optional list of directories to upload to the job's working directory. These can be URIs or paths on the local filesystem.

Relative paths will be interpreted as relative to the directory containing the script (not the current working directory). Environment variables and ~ in paths will be expanded.

If you want a directory to be copied with a name other than it's own, append #<name> (e.g. data/foo#bar).

If you need to dynamically generate a list of files, override dirs () instead.

New in version 0.6.4.

### MRJob.ARCHIVES = []

Optional list of archives to upload and unpack in the job's working directory. These can be URIs or paths on the local filesystem.

Relative paths will be interpreted as relative to the directory containing the script (not the current working directory). Environment variables and ~ in paths will be expanded.

By default, the directory will have the same name as the archive (e.g. foo.tar.gz/). To change the directory's name, append #<name>:

```
ARCHIVES = ['data/foo.tar.gz#foo']
```

If you need to dynamically generate a list of files, override archives () instead.

New in version 0.6.4.

# MRJob.files()

Like FILES, except that it can return a dynamically generated list of files to upload. Overriding this method disables FILES.

Paths returned by this method are relative to the working directory (not the script). Note that the job runner will *always* expand environment variables and ~ in paths returned by this method.

You do not have to worry about inadvertently disabling --files; this switch is handled separately.

New in version 0.6.4.

#### MRJob.dirs()

Like *DIRS*, except that it can return a dynamically generated list of directories to upload. Overriding this method disables *DIRS*.

Paths returned by this method are relative to the working directory (not the script). Note that the job runner will always expand environment variables and  $\sim$  in paths returned by this method.

You do not have to worry about inadvertently disabling --dirs; this switch is handled separately.

New in version 0.6.4.

#### MRJob.archives()

Like ARCHIVES, except that it can return a dynamically generated list of archives to upload and unpack. Overriding this method disables ARCHIVES.

Paths returned by this method are relative to the working directory (not the script). Note that the job runner will *always* expand environment variables and  $\sim$  in paths returned by this method.

You do not have to worry about inadvertently disabling --archives; this switch is handled separately.

New in version 0.6.4.

# 2.10.10 Job runner configuration

```
classmethod MRJob.mr_job_script()
```

Path of this script. This returns the file containing this class, or None if there isn't any (e.g. it was defined from the command line interface.)

# 2.10.11 Running specific parts of jobs

```
MRJob.run_job()
```

Run the all steps of the job, logging errors (and debugging output if --verbose is specified) to STDERR and streaming the output to STDOUT.

Called from run (). You'd probably only want to call this directly from automated tests.

```
MRJob.run_mapper(step_num=0)
```

Run the mapper and final mapper action for the given step.

```
Parameters step_num (int) – which step to run (0-indexed)
```

Called from run (). You'd probably only want to call this directly from automated tests.

```
MRJob.map_pairs (pairs, step_num=0)
```

Runs mapper\_init(), mapper()/mapper\_raw(), and mapper\_final() for one map task in one step.

Takes in a sequence of (key, value) pairs as input, and yields (key, value) pairs as output.

run\_mapper() essentially wraps this method with code to handle reading/decoding input and writing/encoding output.

New in version 0.6.7.

```
MRJob.run_reducer(step_num=0)
```

Run the reducer for the given step.

```
Parameters step_num (int) – which step to run (0-indexed)
```

Called from run (). You'd probably only want to call this directly from automated tests.

```
MRJob.reduce_pairs (pairs, step_num=0)
```

```
Runs reducer_init(), reducer(), and reducer_final() for one reduce task in one step.
```

Takes in a sequence of (key, value) pairs as input, and yields (key, value) pairs as output.

run\_reducer() essentially wraps this method with code to handle reading/decoding input and writing/encoding output.

New in version 0.6.7.

# MRJob.run\_combiner(step\_num=0)

Run the combiner for the given step.

### **Parameters** step\_num (*int*) – which step to run (0-indexed)

If we encounter a line that can't be decoded by our input protocol, or a tuple that can't be encoded by our output protocol, we'll increment a counter rather than raising an exception. If –strict-protocols is set, then an exception is raised

Called from run(). You'd probably only want to call this directly from automated tests.

### MRJob.combine\_pairs (pairs, step\_num=0)

Runs combiner init(), combiner(), and combiner final() for one reduce task in one step.

Takes in a sequence of (key, value) pairs as input, and yields (key, value) pairs as output.

run\_combiner() essentially wraps this method with code to handle reading/decoding input and writing/encoding output.

New in version 0.6.7.

### MRJob.show\_steps()

Print information about how many steps there are, and whether they contain a mapper or reducer. Job runners (see Runners) use this to determine how Hadoop should call this script.

Called from run(). You'd probably only want to call this directly from automated tests.

# 2.10.12 Hadoop configuration

## MRJob.HADOOP\_INPUT\_FORMAT = None

Optional name of an optional Hadoop InputFormat class, e.g. 'org.apache.hadoop.mapred.lib.NLineInputFormat'.

Passed to Hadoop with the *first* step of this job with the -input format option.

If you require more sophisticated behavior, try hadoop\_input\_format () or the hadoop\_input\_format argument to mrjob.runner.MRJobRunner.\_\_init\_\_().

### MRJob.hadoop\_input\_format()

Optional Hadoop InputFormat class to parse input for the first step of the job.

Normally, setting <code>HADOOP\_INPUT\_FORMAT</code> is sufficient; redefining this method is only for when you want to get fancy.

#### MRJob.HADOOP OUTPUT FORMAT = None

Optional name of an optional Hadoop OutputFormat class, e.g. 'org.apache.hadoop.mapred.FileOutputFormat'.

Passed to Hadoop with the *last* step of this job with the -output format option.

If you require more sophisticated behavior, try hadoop\_output\_format () or the hadoop\_output\_format argument to mrjob.runner.MRJobRunner.\_\_init\_\_ ().

#### MRJob.hadoop\_output\_format()

Optional Hadoop OutputFormat class to write output for the last step of the job.

Normally, setting <code>HADOOP\_OUTPUT\_FORMAT</code> is sufficient; redefining this method is only for when you want to get fancy.

### MRJob.JOBCONF = {}

Optional jobconf arguments we should always pass to Hadoop. This is a map from property name to value. e.g.:

```
{'stream.num.map.output.key.fields': '4'}
```

It's recommended that you only use this to hard-code things that affect the semantics of your job, and leave performance tweaks to the command line or whatever you use to launch your job.

#### MRJob.jobconf()

-D args to pass to hadoop streaming. This should be a map from property name to value. By default, returns *JOBCONF*.

Changed in version 0.6.6: re-defining longer clobbers command-line -- jobconf options.

### MRJob.LIBJARS = []

Optional list of paths of jar files to run our job with using Hadoop's -lib jars option.

~ and environment variables in paths be expanded, and relative paths will be interpreted as relative to the directory containing the script (not the current working directory).

If you require more sophisticated behavior, try overriding libjars().

New in version 0.5.3.

### MRJob.libjars()

Optional list of paths of jar files to run our job with using Hadoop's -libjars option. Normally setting *LIBJARS* is sufficient. Paths from *LIBJARS* are interpreted as relative to the directory containing the script (paths from the command-line are relative to the current working directory).

Note that ~ and environment variables in paths will always be expanded by the job runner (see libjars).

New in version 0.5.3.

Changed in version 0.6.6: re-defining this no longer clobbers the command-line --libjars option

### MRJob.PARTITIONER = None

Optional Hadoop partitioner class to use to determine how mapper output should be sorted and distributed to reducers. For example: 'org.apache.hadoop.mapred.lib.HashPartitioner'.

If you require more sophisticated behavior, try partitioner().

### MRJob.partitioner()

Optional Hadoop partitioner class to use to determine how mapper output should be sorted and distributed to reducers.

By default, returns PARTITIONER.

You probably don't need to re-define this; it's just here for completeness.

# 2.10.13 Hooks for testing

MRJob.sandbox (stdin=None, stdout=None, stderr=None)

Redirect stdin, stdout, and stderr for automated testing.

You can set stdin, stdout, and stderr to file objects. By default, they'll be set to empty BytesIO objects. You can then access the job's file handles through self.stdin, self.stdout, and self.stderr. See *Testing jobs* for more information about testing.

You may call sandbox multiple times (this will essentially clear the file handles).

stdin is empty by default. You can set it to anything that yields lines:

```
mr_job.sandbox(stdin=BytesIO(b'some_data\n'))
```

or, equivalently:

```
mr_job.sandbox(stdin=[b'some_data\n'])
```

For convenience, this sandbox() returns self, so you can do:

```
mr_job = MRJobClassToTest().sandbox()
```

### Simple testing example:

```
mr_job = MRYourJob.sandbox()
self.assertEqual(list(mr_job.reducer('foo', ['a', 'b'])), [...])
```

More complex testing example:

```
from BytesIO import BytesIO

from mrjob.parse import parse_mr_job_stderr
from mrjob.protocol import JSONProtocol

mr_job = MRYourJob(args=[...])

fake_input = '"foo"\t"bar"\n"foo"\t"baz"\n'
mr_job.sandbox(stdin=BytesIO(fake_input))

mr_job.run_reducer(link_num=0)

self.assertEqual(mrjob.stdout.getvalue(), ...)
self.assertEqual(parse_mr_job_stderr(mr_job.stderr), ...)
```

**Note:** If you are using Spark, it's recommended you only pass in io. BytesIO or other serializable alternatives to file objects. *stdin*, *stdout*, and *stderr* get stored as job attributes, which means if they aren't serializable, neither is the job instance or its methods.

# 2.11 mrjob.local - simulate Hadoop locally with subprocesses

```
class mrjob.local.LocalMRJobRunner(**kwargs)
```

Runs an MRJob locally, for testing purposes. Invoked when you run your job with -r local.

Unlike InlineMRJobRunner, this actually spawns multiple subprocesses for each task.

It's rare to need to instantiate this class directly (see \_\_\_init\_\_\_() for details).

New in version 0.6.8: can run Spark steps as well, on the local-cluster Spark master.

```
LocalMRJobRunner.__init__(**kwargs)
```

Arguments to this constructor may also appear in mrjob.conf under runners/local.

LocalMRJobRunner's constructor takes the same keyword args as MRJobRunner. However, please note:

- •cmdenv is combined with combine\_local\_envs()
- •python\_bin defaults to sys.executable (the current python interpreter)
- •hadoop\_input\_format, hadoop\_output\_format, and partitioner are ignored because they require Java. If you need to test these, consider starting up a standalone Hadoop instance and running your job with -r hadoop.

# 2.12 mrjob.parse - log parsing

Utilities for parsing errors, counters, and status messages.

```
mrjob.parse.is_s3_uri(uri)
```

Return True if *uri* can be parsed into an S3 URI, False otherwise.

```
mrjob.parse.is_uri(uri)
```

Return True if *uri* is a URI and contains: // (we only care about URIs that can describe files)

Changed in version 0.5.7: used to recognize anything containing a colon as a URI unless it was a Windows path  $(C: \setminus ...)$ .

```
mrjob.parse.parse_mr_job_stderr(stderr, counters=None)
```

Parse counters and status messages out of MRJob output.

#### **Parameters**

- **stderr** a filehandle, a list of lines (bytes), or bytes
- **counters** Counters so far, to update; a map from group (string to counter name (string) to count.

Returns a dictionary with the keys counters, statuses, other:

- •counters: counters so far; same format as above
- •statuses: a list of status messages encountered
- •other: lines (strings) that aren't either counters or status messages

```
mrjob.parse.parse_s3_uri(uri)
```

Parse an S3 URI into (bucket, key)

```
>>> parse_s3_uri('s3://walrus/tmp/')
('walrus', 'tmp/')
```

If uri is not an S3 URI, raise a ValueError

```
mrjob.parse.to_uri(path_or_uri)
```

If *path\_or\_uri* is not a URI already, convert it to a file:/// URI.

# 2.13 mrjob.protocol - input and output

Protocols translate raw bytes into key, value pairs.

Typically, protocols encode a key and value into bytes, and join them together with a tab character.

However, protocols with Value in their name ignore keys and simply read/write values (with key read in as None), allowing you to read and write data in arbitrary formats.

For more information, see *Protocols* and *Writing custom protocols*.

# **2.13.1 Strings**

```
class mrjob.protocol.RawValueProtocol
```

Just output value (a str), and discard key (key is read in as None).

This is the default protocol used by jobs to read input.

This is an alias for RawValueProtocol on Python 2 and TextValueProtocol on Python 3.

```
class mrjob.protocol.BytesValueProtocol
```

Read line (without trailing newline) directly into value (key is always None). Output value (bytes) directly, discarding key.

### This is the default protocol used by jobs to read input on Python 2.

**Warning:** Typical usage on Python 2 is to have your mapper parse (byte) strings out of your input files, and then include them in the output to the reducer. Since this output is then (by default) JSON-encoded, encoding will fail if the bytestrings are not UTF-8 decodable. If this is an issue, consider using <code>TextValueProtocol</code> instead.

### class mrjob.protocol.TextValueProtocol

Attempt to UTF-8 decode line (without trailing newline) into value, falling back to latin-1. (key is always None). Output value UTF-8 encoded, discarding key.

### This is the default protocol used by jobs to read input on Python 3.

This is a good solution for reading text files which are mostly ASCII but may have some other bytes of unknown encoding (e.g. logs).

If you wish to enforce a particular encoding, use <code>BytesValueProtocol</code> instead:

```
class MREncodingEnforcer(MRJob):
    INPUT_PROTOCOL = BytesValueProtocol

def mapper(self, _, value):
    value = value.decode('utf_8')
    ...
```

#### class mrjob.protocol.RawProtocol

Output key (str) and value (str), separated by a tab character.

This is an alias for <code>BytesProtocol</code> on Python 2 and <code>TextProtocol</code> on Python 3.

## class mrjob.protocol.BytesProtocol

Encode (key, value) (bytestrings) as key and value separated by a tab.

If key or value is None, don't include a tab. When decoding a line with no tab in it, value will be None.

When reading from a line with multiple tabs, we break on the first one.

Your key should probably not be None or have tab characters in it, but we don't check.

### class mrjob.protocol.TextProtocol

UTF-8 encode key and value (unicode strings) and join them with a tab character. When reading input, we fall back to latin-1 if we can't UTF-8 decode the line.

If key or value is None, don't include a tab. When decoding a line with no tab in it, value will be None.

When reading from a line with multiple tabs, we break on the first one.

Your key should probably not be None or have tab characters in it, but we don't check.

### 2.13.2 JSON

## class mrjob.protocol.JSONProtocol

Encode (key, value) as two JSONs separated by a tab.

### This is the default protocol used by jobs to write output and communicate between steps.

This is an alias for the first one of *UltraJSONProtocol*, *RapidJSONProtocol*, *SimpleJSONProtocol*, or *StandardJSONProtocol* for which the underlying library is available.

#### class mrjob.protocol.UltraJSONProtocol

Implements JSONProtocol using the ujson library.

**Warning:** ujson is about five times faster than the standard implementation, but is more willing to encode things that aren't strictly JSON-encodable, including sets, dictionaries with tuples as keys, UTF-8 encoded bytes, and objects (!). Relying on this behavior won't stop your job from working, but it can make your job dependent on ujson, rather than just using it as a speedup.

**Note:** ujson also differs from the standard implementation in that it doesn't add spaces to its JSONs ({"foo": "bar"} versus {"foo": "bar"}). This probably won't affect anything but test cases and readability.

### class mrjob.protocol.RapidJSONProtocol

Implements *JSONProtocol* using the rapidjson library.

New in version 0.5.10.

# class mrjob.protocol.SimpleJSONProtocol

Implements JSONProtocol using the simple json library.

#### class mrjob.protocol.StandardJSONProtocol

Implements JSONProtocol using Python's built-in JSON library.

**Note:** The built-in json library is (appropriately) strict about the JSON standard; it won't accept dictionaries with non-string keys, sets, or (on Python 3) bytestrings.

### class mrjob.protocol.JSONValueProtocol

Encode value as a JSON and discard key (key is read in as None).

This is an alias for the first one of UltraJSONValueProtocol, RapidJSONValueProtocol, SimpleJSONValueProtocol, or StandardJSONValueProtocol for which the underlying library is available.

### class mrjob.protocol.UltraJSONValueProtocol

Implements JSONValueProtocol using the ujson library.

### class mrjob.protocol.RapidJSONValueProtocol

Implements JSONValueProtocol using the rapidjson library.

New in version 0.5.10.

# ${\bf class} \; {\tt mrjob.protocol.SimpleJSONValueProtocol}$

Implements JSONValueProtocol using the simplejson library.

# ${\bf class} \ {\tt mrjob.protocol.StandardJSONValueProtocol}$

Implements JSONValueProtocol using Python's built-in JSON library.

# 2.13.3 Repr

## class mrjob.protocol.ReprProtocol

Encode (key, value) as two reprs separated by a tab.

This only works for basic types (we use mrjob.util.safeeval()).

**Warning:** The repr format changes between different versions of Python (for example, braces for sets in Python 2.7, and different string contants in Python 3). Plan accordingly.

#### class mrjob.protocol.ReprValueProtocol

Encode value as a repr and discard key (key is read in as None).

See ReprProtocol for details.

### 2.13.4 Pickle

### class mrjob.protocol.PickleProtocol

Encode (key, value) as two string-escaped pickles separated by a tab.

We string-escape the pickles to avoid having to deal with stray \t and \n characters, which would confuse Hadoop Streaming.

Ugly, but should work for any type.

**Warning:** Pickling is only *backwards*-compatible across Python versions. If your job uses this as an output protocol, you should use at least the same version of Python to parse the job's output. Vice versa for using this as an input protocol.

# class mrjob.protocol.PickleValueProtocol

Encode value as a string-escaped pickle and discard key (key is read in as None).

See PickleProtocol for details.

# 2.14 mrjob.spark.runner - run on any Spark cluster

# 2.14.1 Job Runner

# 2.15 mrjob.retry - retry on transient errors

**class** mrjob.retry.**RetryWrapper** (*wrapped*, *retry\_if*, *backoff*=15, *multiplier*=1.5, *max\_tries*=10) Handle transient errors, with configurable backoff.

This class can wrap any object. The wrapped object will behave like the original one, except that if you call a function and it raises a retriable exception, we'll back off for a certain number of seconds and call the function again, until it succeeds or we get a non-retriable exception.

RetryWrapper.\_\_init\_\_ (wrapped, retry\_if, backoff=15, multiplier=1.5, max\_tries=10)
Wrap the given object

#### **Parameters**

- wrapped the object to wrap
- retry\_if a method that takes an exception, and returns whether we should retry
- backoff (float) the number of seconds to wait the first time we get a retriable error
- multiplier (*float*) if we retry multiple times, the amount to multiply the backoff time by every time we get an error
- max\_tries (int) how many tries we get. 0 means to keep trying forever

# 2.16 mrjob.runner - base class for all runners

Abstract base class for all runners

```
MRJobRunner.__init__ (mr_job_script=None, conf_paths=None, extra_args=None, file_upload_args=None, hadoop_input_format=None, hadoop_output_format=None, input_paths=None, output_dir=None, partitioner=None, sort_values=None, stdin=None, steps=None, step_output_dir=None, **opts)

All runners take the following keyword arguments:
```

#### **Parameters**

- mr\_job\_script (*str*) the path of the .py file containing the MRJob. If this is None, you won't actually be able to run() the job, but other utilities (e.g. 1s()) will work.
- **conf\_paths** (*None or list*) List of config files to combine and use, or None to search for mrjob.conf in the default locations.
- extra\_args (*list of str*) a list of extra cmd-line arguments to pass to the mr\_job script. This is a hook to allow jobs to take additional arguments.
- **file\_upload\_args** a list of tuples of ('--ARGNAME', path). The file at the given path will be uploaded to the local directory of the mr\_job script when it runs, and then passed into the script with --ARGNAME. Useful for passing in SQLite DBs and other configuration files to your job.
- hadoop\_input\_format (str) name of an optional Hadoop InputFormat class. Passed to Hadoop along with your first step with the -inputformat option. Note that if you write your own class, you'll need to include it in your own custom streaming jar (see hadoop\_streaming\_jar).
- hadoop\_output\_format (str) name of an optional Hadoop OutputFormat class. Passed to Hadoop along with your first step with the –outputformat option. Note that if you write your own class, you'll need to include it in your own custom streaming jar (see hadoop\_streaming\_jar).
- input\_paths (list of str) Input files for your job. Supports globs and recursively walks directories (e.g. ['data/common/', 'data/training/\*.gz']). If this is left blank, we'll read from stdin
- output\_dir (str) An empty/non-existent directory where Hadoop should put the final output from the job. If you don't specify an output directory, we'll output into a subdirectory of this job's temporary directory. You can control this from the command line with --output-dir. This option cannot be set from configuration files. If used with the hadoop runner, this path does not need to be fully qualified with hdfs:// URIs because it's understood that it has to be on HDFS.
- partitioner (*str*) Optional name of a Hadoop partitioner class, e.g. 'org.apache.hadoop.mapred.lib.HashPartitioner'. Hadoop streaming will use this to determine how mapper output should be sorted and distributed to reducers.

- **sort\_values** (*bool*) if true, set partitioners and jobconf variables so that reducers to receive the values associated with any key in sorted order (sorted by their *encoded* value). Also known as secondary sort.
- **stdin** an iterable (can be a BytesIO or even a list) to use as stdin. This is a hook for testing; if you set stdin via sandbox(), it'll get passed through to the runner. If for some reason your lines are missing newlines, we'll add them; this makes it easier to write automated tests.
- **steps** a list of descriptions of steps to run (see mrjob.step represent Job Steps for description formats)
- **step\_output\_dir** (*str*) An empty/non-existent directory where Hadoop should put output from all steps other than the last one (this only matters for multi-step jobs). Currently ignored by local runners.

# 2.16.1 Running your job

```
MRJobRunner.run()
```

Run the job, and block until it finishes.

Raise StepFailedException if there are any problems (except on InlineMRJobRunner, where we raise the actual exception that caused the step to fail).

```
MRJobRunner.cat_output()
```

Stream the job's output, as a stream of bytes. If there are multiple output files, there will be an empty bytestring (b'') between them.

Like Hadoop input formats, we ignore files and subdirectories whose names start with "\_" or "." (e.g. \_SUCCESS, \_logs/, .part-00000.crc.

New in version 0.6.0: In previous versions, you'd use stream\_output().

Changed in version 0.6.8: Ignore file/dirnames starting with "." as well as "\_".

```
MRJobRunner.stream_output()
```

Like cat\_output() except that it groups bytes into lines. Equivalent to mrjob.util.to\_lines(runner.cat\_output()).

Deprecated since version 0.6.0.

```
MRJobRunner.cleanup (mode=None)
```

Clean up running jobs, temp files, and logs, subject to the *cleanup* option passed to the constructor.

If you create your runner in a with block, <code>cleanup()</code> will be called automatically:

**Parameters mode** – override *cleanup* passed into the constructor. Should be a list of strings from *CLEANUP\_CHOICES* 

```
mrjob.options.CLEANUP_CHOICES = ['ALL', 'CLOUD_TMP', 'CLUSTER', 'HADOOP_TMP', 'JOB', 'LOCAL_TMP', cleanup options:
```

• 'ALL': delete logs and local and remote temp files; stop cluster if on EMR and the job is not done when cleanup is run.

- •' CLOUD\_TMP': delete temp files on cloud storage (e.g. S3) only
- 'CLUSTER': terminate the cluster if on EMR and the job is not done on cleanup
- ' HADOOP\_TMP': delete temp files on HDFS only
- •' JOB': stop job if on EMR and the job is not done when cleanup runs
- 'LOCAL TMP': delete local temp files only
- •' LOGS': delete logs only
- ' NONE ': delete nothing
- •' TMP': delete local, HDFS, and cloud storage temp files, but not logs

Changed in version 0.5.0: LOCAL\_TMP used to be LOCAL\_SCRATCH HADOOP\_TMP is new (and used to be covered by LOCAL\_SCRATCH) CLOUD\_TMP used to be REMOTE\_SCRATCH

# 2.16.2 Run Information

MRJobRunner.counters()

Get counters associated with this run in this form:

```
[{'group name': {'counter1': 1, 'counter2': 2}},
{'group name': ...}]
```

The list contains an entry for every step of the current job.

```
MRJobRunner.get_hadoop_version()
```

Return the version number of the Hadoop environment as a string if Hadoop is being used or simulated. Return None if not applicable.

EMRJobRunner infers this from the cluster. HadoopJobRunner gets this from hadoop version. LocalMRJobRunner has an additional hadoop\_version option to specify which version it simulates. InlineMRJobRunner does not simulate Hadoop at all.

```
MRJobRunner.get_job_key()
```

Get the unique key for the job run by this runner. This has the format label.owner.date.time.microseconds

# 2.16.3 Configuration

```
MRJobRunner.get_opts()
```

Get options set for this runner, as a dict.

# 2.16.4 File management

```
MRJobRunner.fs
```

Filesystem object for the local filesystem.

```
class mrjob.fs.base.Filesystem
```

Some simple filesystem operations that are common across the local filesystem, S3, GCS, HDFS, and remote machines via SSH.

Different runners provide functionality for different filesystems via their fs attribute. Generally a runner will wrap one or more filesystems with mrjob.fs.composite.CompositeFilesystem.

Schemes supported:

```
•mrjob.fs.gcs.GCSFilesystem: qs://
    \bulletmrjob.fs.hadoop.HadoopFilesystem: hdfs:// and other URIs
    •mrjob.fs.local.LocalFilesystem: / (non-URIs)
    •mrjob.fs.s3.S3Filesystem: s3://, s3a://, s3n://,
    •mrjob.fs.ssh.SSHFilesystem: ssh://
can handle path (path)
    Can we handle this path at all?
cat (path_glob)
     cat all files matching path_glob, decompressing if necessary
     This yields bytes, which don't necessarily correspond to lines (see #1544). If multiple files are catted,
     yields b'' between each file.
du (path_glob)
     Get the total size of files matching path_glob
     Corresponds roughly to: hadoop fs -du path glob
exists(path glob)
     Does the given path/URI exist?
     Corresponds roughly to: hadoop fs -test -e path_glob
join (path, *paths)
     Join paths onto path (which may be a URI)
ls (path_glob)
     Recursively list all files in the given path.
     We don't return directories for compatibility with S3 (which has no concept of them)
     Corresponds roughly to: hadoop fs -ls -R path_glob
md5sum(path\_glob)
     Generate the md5 sum of the file at path
mkdir (path)
     Create the given dir and its subdirs (if they don't already exist). On cloud filesystems (e.g. S3), also create
     the corresponding bucket as needed
     Corresponds roughly to: hadoop fs -mkdir -p path
     New in version 0.6.8: creates buckets on cloud filesystems
put (src, path)
     Upload a file on the local filesystem (src) to path. Like with shutil.copyfile(), path should be the
     full path of the new file, not a directory which should contain it.
     Corresponds roughly to hadoop fs -put src path.
     New in version 0.6.8.
rm (path_glob)
     Recursively delete the given file/directory, if it exists
     Corresponds roughly to: hadoop fs -rm -R path_glob
touchz (path)
     Make an empty file in the given location. Raises an error if a non-zero length file already exists in that
```

location.

Correponds to: hadoop fs -touchz path

# 2.17 mrjob.step - represent Job Steps

Representations of job steps, to use in your MRJob's steps() method.

Because the runner just needs to know how to invoke your MRJob script, not how it works insternally, each step instance's description () method produces a simplified, JSON-able description of the step, to pass to the runner.

# 2.17.1 Steps

```
class mrjob.step.MRStep(**kwargs)
```

Represents steps handled by the script containing your job.

Used by MRJob. steps. See Multi-step jobs for sample usage.

Takes the following keyword arguments: combiner, combiner\_cmd, combiner\_final, combiner\_init, combiner\_pre\_filter, mapper, mapper\_cmd, mapper\_final, mapper\_init, mapper\_pre\_filter, mapper\_raw, reducer, reducer\_cmd, reducer\_final, reducer\_init, reducer\_pre\_filter. These should be set to None or a function with the same signature as the corresponding method in MRJob.

Also accepts *jobconf*, a dictionary with custom jobconf arguments to pass to hadoop.

A MRStep's description looks like:

```
{
    'type': 'streaming',
    'mapper': { ... },
    'combiner': { ... },
    'reducer': { ... },
    'jobconf': { ... }, # dict of Hadoop configuration properties
}
```

At least one of mapper, combiner and reducer need be included. jobconf is completely optional.

mapper, combiner, and reducer are either handled by the script containing your job definition, in which case they look like:

```
{
    'type': 'script',
    'pre_filter': 'grep -v bad', # optional cmd to filter input
}
```

or they simply run a command, which looks like:

```
{
   'type': 'command',
   'command': 'cut -f 1-2', # command to run, as a string
}
```

class mrjob.step.JarStep (jar, \*\*kwargs)

Represents a running a custom Jar as a step.

Accepts the following keyword arguments:

### **Parameters**

• jar – The local path to the Jar. On EMR, this can also be an s3:// URI, or file:// to reference a jar on the local filesystem of your EMR instance(s).

- args (optional) A list of arguments to the jar. Use mrjob.step.INPUT and OUTPUT to interpolate input and output paths.
- jobconf (optional) A dictionary of Hadoop properties
- main\_class (optional) The main class to run from the jar. If not specified, Hadoop will use the main class in the jar's manifest file.

jar can also be passed as a positional argument

See *Jar steps* for sample usage.

Sample description of a JarStep:

```
{
    'type': 'jar',
    'jar': 'binks.jar.jar',
    'main_class': 'MyMainMan', # optional
    'args': ['argh', 'argh'] # optional
    'jobconf': { ... } # optional
}
```

To give your jar access to input files, an empty output directory, configuration properties, and libjars managed by mrjob, you may include *INPUT*, *OUTPUT*, and *GENERIC\_ARGS* in *args*.

```
class mrjob.step.SparkStep(spark, **kwargs)
```

Represents running a Spark step defined in your job.

Accepts the following keyword arguments:

### **Parameters**

- spark function containing your Spark code with same function signature as spark ()
- jobconf (optional) A dictionary of Hadoop properties
- **spark\_args** (optional) an array of arguments to pass to spark-submit (e.g. ['--executor-memory', '2G']).

Sample description of a SparkStep:

```
{
    'type': 'spark',
    'jobconf': { ... }, # optional
    'spark_args': ['--executor-memory', '2G'], # optional
}
```

class mrjob.step.SparkJarStep (jar, main\_class, \*\*kwargs)

Represents a running a separate Jar through Spark

Accepts the following keyword arguments:

#### **Parameters**

- jar The local path to the Python script to run. On EMR, this can also be an s3:// URI, or file:// to reference a jar on the local filesystem of your EMR instance(s).
- main\_class Your application's main class (e.g. 'org.apache.spark.examples.SparkPi')
- args (optional) A list of arguments to the script. Use mrjob.step.INPUT and OUTPUT to interpolate input and output paths.
- jobconf (optional) A dictionary of Hadoop properties

• spark\_args - (optional) an array of arguments to pass to spark-submit (e.g. ['--executor-memory', '2G']).

*jar* and *main\_class* can also be passed as positional arguments

Sample description of a SparkJarStep:

```
{
  'type': 'spark_jar',
  'jar': 'binks.jar.jar',
  'main_class': 'MyMainMan', # optional
  'args': ['argh', 'argh'], # optional
  'jobconf': { ... }, # optional
  'spark_args': ['--executor-memory', '2G'], # optional
}
```

To give your Spark JAR access to input files and an empty output directory managed by mrjob, you may include *INPUT* and *OUTPUT* in *args*.

```
class mrjob.step.SparkScriptStep (script, **kwargs)
```

Represents a running a separate Python script through Spark

Accepts the following keyword arguments:

#### **Parameters**

- **script** The local path to the Python script to run. On EMR, this can also be an s3:// URI, or file:// to reference a jar on the local filesystem of your EMR instance(s).
- args (optional) A list of arguments to the script. Use mrjob.step.INPUT and OUTPUT to interpolate input and output paths.
- jobconf (optional) A dictionary of Hadoop properties
- **spark\_args** (optional) an array of arguments to pass to spark-submit (e.g. ['--executor-memory', '2G']).

script can also be passed as a positional argument

Sample description of a ScriptStep:

```
{
    'type': 'spark_script',
    'script': 'my_spark_script.py',
    'args': ['script_arg1', 'script_arg2'],
    'jobconf': { ... }, # optional
    'spark_args': ['--executor-memory', '2G'], # optional
}
```

To give your Spark script access to input files and an empty output directory managed by mrjob, you may include *INPUT* and *OUTPUT* in *args*.

# 2.17.2 Argument interpolation

Use these constants in your step's args and mrjob will automatically replace them before running your step.

```
mrjob.step.INPUT = '<input>'
```

If passed as an argument to <code>JarStep</code>, <code>SparkJarStep</code>, or <code>SparkScriptStep</code>, it'll be replaced with the step's input path(s). If there are multiple paths, they'll be joined with commas.

```
mrjob.step.OUTPUT = '<output>'
```

If this is passed as an argument to <code>JarStep</code>, <code>SparkJarStep</code>, or <code>SparkScriptStep</code>, it'll be replaced with the step's output path

```
mrjob.step.GENERIC_ARGS = '<generic args>'
```

If this is passed as an argument to JarStep, it'll be replaced with generic hadoop args (-D and -libjars)

# 2.18 mrjob.setup - job environment setup

Utilities for setting up the environment jobs run in by uploading files and running setup scripts.

The general idea is to use Hadoop DistributedCache-like syntax to find and parse expressions like /path/to/file#name\_in\_working\_dir into "path dictionaries" like {'type': 'file', 'path': 'path/to/file', 'name': 'name\_in\_working\_dir'}}.

You can then pass these into a <code>WorkingDirManager</code> to keep track of which files need to be uploaded, catch name collisions, and assign names to unnamed paths (e.g. <code>/path/to/file#</code>). Note that <code>WorkingDirManager.name()</code> can take a path dictionary as keyword arguments.

If you need to upload files from the local filesystem to a place where Hadoop can see them (HDFS or S3), we provide <code>UploadDirManager</code>.

Path dictionaries are meant to be immutable; all state is handled by manager classes.

### class mrjob.setup.UploadDirManager (prefix)

Represents a directory on HDFS or S3 where we want to upload local files for consumption by Hadoop.

*UploadDirManager* tries to give files the same name as their filename in the path (for ease of debugging), but handles collisions gracefully.

UploadDirManager assumes URIs to not need to be uploaded and thus does not store them. uri () maps URIs to themselves.

```
add(path)
```

Add a path. If path hasn't been added before, assign it a name. If path is a URI don't add it; just return the URI.

**Returns** the URI assigned to the path

```
path_to_uri()
```

Get a map from path to URI for all paths that were added, so we can figure out which files we need to upload.

```
uri (path)
```

Get the URI for the given path. If path is a URI, just return it.

# class mrjob.setup.WorkingDirManager

Represents the working directory of hadoop/Spark tasks (or bootstrap commands in the cloud).

To support Hadoop's distributed cache, paths can be for ordinary files, or for archives (which are automatically uncompressed into a directory by Hadoop).

When adding a file, you may optionally assign it a name; if you don't; we'll lazily assign it a name as needed. Name collisions are not allowed, so being lazy makes it easier to avoid unintended collisions.

If you wish, you may assign multiple names to the same file, or add a path as both a file and an archive (though not mapped to the same name).

```
add (type, path, name=None)
```

Add a path as either a file or an archive, optionally assigning it a name.

#### **Parameters**

- type either 'archive' or 'file'
- path path/URI to add
- name optional name that this path *must* be assigned, or None to assign this file a name later

### name (type, path, name=None)

Get the name for a path previously added to this WorkingDirManager, assigning one as needed.

This is primarily for getting the name of auto-named files. If the file was added with an assigned name, you must include it (and we'll just return *name*).

We won't ever give an auto-name that's the same an assigned name (even for the same path and type).

#### **Parameters**

- type either 'archive' or 'file'
- path path/URI
- name known name of the file

### name\_to\_path (type=None)

Get a map from name (in the setup directory) to path for all known files/archives, so we can build <code>-file</code> and <code>-archive</code> options to Hadoop (or fake them in a bootstrap script).

```
Parameters type - either 'archive' or 'file'
```

```
paths (type=None)
```

Get a set of all paths tracked by this WorkingDirManager.

```
mrjob.setup.name_uniquely (path, names_taken=(), proposed_name=None, unhide=False)
Come up with a unique name for path.
```

### **Parameters**

- names\_taken a dictionary or set of names not to use.
- **proposed\_name** name to use if it is not taken. If this is not set, we propose a name based on the filename.
- unhide make sure final name doesn't start with periods or underscores

If the proposed name is taken, we add a number to the end of the filename, keeping the extension the same. For example:

```
>>> name_uniquely('foo.tar.gz', set(['foo.tar.gz']))
'foo-1.tar.gz'
```

## mrjob.setup.parse\_legacy\_hash\_path(type, path, must\_name=None)

Parse hash paths from old setup/bootstrap options.

This is similar to parsing hash paths out of shell commands (see parse\_setup\_cmd()) except that we pass in path type explicitly, and we don't always require the # character.

#### **Parameters**

- type Type of the path ('archive' or 'file')
- path Path to parse, possibly with a #

• must\_name - If set, use *path*'s filename as its name if there is no '#' in *path*, and raise an exception if there is just a '#' with no name. Set *must\_name* to the name of the relevant option so we can print a useful error message. This is intended for options like upload\_files that merely upload a file without tracking it.

```
mrjob.setup.parse_setup_cmd(cmd)
```

Parse a setup/bootstrap command, finding and pulling out Hadoop Distributed Cache-style paths ("hash paths").

**Parameters** cmd (*string*) – shell command to parse

**Returns** a list containing dictionaries (parsed hash paths) and strings (parts of the original command, left unparsed)

Hash paths look like path#name, where *path* is either a local path or a URI pointing to something we want to upload to Hadoop/EMR, and *name* is the name we want it to have when we upload it; *name* is optional (no name means to pick a unique one).

If *name* is followed by a trailing slash, that indicates *path* is an archive (e.g. a tarball), and should be unarchived into a directory on the remote system. The trailing slash will *also* be kept as part of the original command.

If *path* is followed by a trailing slash, that indicates *path* is a directory and should be tarballed and later unarchived into a directory on the remote system. The trailing slash will also be kept as part of the original command. You may optionally include a slash after *name* as well (this will only result in a single slash in the final command).

New in version 0.5.8: support for directories (above)

Parsed hash paths are dicitionaries with the keys path, name, and type (either 'file', 'archive', or 'dir').

Most of the time, this function will just do what you expect. Rules for finding hash paths:

- •we only look for hash paths outside of quoted strings
- •path may not contain quotes or whitespace
- •path may not contain: or = unless it is a URI (starts with <scheme>://); this allows you to do stuff like export PYTHONPATH=\$PYTHONPATH:foo.egg#.
- •name may not contain whitespace or any of the following characters: '":; >< |=/#, so you can do stuff like sudo dpkg -i fooify.deb#; fooify bar

If you really want to include forbidden characters, you may use backslash escape sequences in *path* and *name*. (We can't guarantee Hadoop/EMR will accept them though!). Also, remember that shell syntax allows you to concatenate strings like""this.

Environment variables and  $\sim$  (home dir) in *path* will be resolved (use backslash escapes to stop this). We don't resolve *name* because it doesn't make sense. Environment variables and  $\sim$  elsewhere in the command are considered to be part of the script and will be resolved on the remote system.

# 2.19 mrjob.util - general utility functions

```
Utility functions for MRJob
```

```
mrjob.util.cmd_line(args)
```

build a command line that works in a shell.

```
mrjob.util.expand_path(path)
```

Resolve ~ (home dir) and environment variables in *path*.

If path is None, return None.

```
mrjob.util.file_ext(filename)
```

return the file extension, including the .

```
>>> file_ext('foo.tar.gz')
'.tar.gz'
```

```
mrjob.util.log_to_null(name=None)
```

Set up a null handler for the given stream, to suppress "no handlers could be found" warnings.

#### **Parameters**

- name (str) name of the logger, or None for the root logger
- **stream** (*file object*) stream to log to (default is sys.stderr)
- **format** (*str*) log message format (default is '%(message)s')
- level log level to use
- **debug** (bool) quick way of setting the log level: if true, use logging. DEBUG, otherwise use logging. INFO

```
mrjob.util.parse_and_save_options(option_parser, args)
```

Return a map from option name (dest) to a list of the arguments in args that correspond to that dest.

This won't modify option\_parser.

Deprecated since version 0.6.0.

```
mrjob.util.random identifier()
```

A random 16-digit hex string.

mrjob.util.read\_file(path, fileobj=None, yields\_lines=True, cleanup=None)

Yields lines from a file, possibly decompressing it based on file extension.

Currently we handle compressed files with the extensions .gz and .bz2.

#### **Parameters**

- path (*string*) file path. Need not be a path on the local filesystem (URIs are okay) as long as you specify *fileobj* too.
- **fileobj** file object to read from. Need not be seekable. If this is omitted, we open (path).
- **yields\_lines** Does iterating over *fileobj* yield lines (like file objects are supposed to)? If not, set this to False (useful for objects that correspond to objects on cluster filesystems)
- cleanup Optional callback to call with no arguments when EOF is reached or an exception is thrown.

Deprecated since version 0.6.0.

```
mrjob.util.read_input (path, stdin=None)
```

Stream input the way Hadoop would.

- •Resolve globs (foo\_\*.qz).
- •Decompress .gz and .bz2 files.
- •If path is '-', read from stdin
- •If path is a directory, recursively read its contents.

You can redefine *stdin* for ease of testing. *stdin* can actually be any iterable that yields lines (e.g. a list).

Deprecated since version 0.6.0.

mrjob.util.safeeval (expr, globals=None, locals=None)

Like eval, but with nearly everything in the environment blanked out, so that it's difficult to cause mischief.

globals and locals are optional dictionaries mapping names to values for those names (just like in eval ()).

mrjob.util.save\_current\_environment(\*args, \*\*kwds)

Context manager that saves os.environ and loads it back again after execution

mrjob.util.save\_cwd(\*args, \*\*kwds)

Context manager that saves the current working directory, and chdir's back to it after execution.

mrjob.util.save\_sys\_path(\*args, \*\*kwds)

Context manager that saves sys.path and restores it after execution.

mrjob.util.save\_sys\_std(\*args, \*\*kwds)

Context manager that saves the current values of *sys.stdin*, *sys.stdout*, and *sys.stderr*, and flushes these filehandles before and after switching them out.

mrjob.util.shlex\_split(s)

Wrapper around shlex.split(), but convert to str if Python version < 2.7.3 when unicode support was added.

mrjob.util.strip\_microseconds(delta)

Return the given datetime.timedelta, without microseconds.

Useful for printing datetime.timedelta objects.

mrjob.util.to lines(chunks)

Take in data as a sequence of bytes, and yield it, one line at a time.

Only breaks lines on  $\n$  (not  $\r$ ), and does not add a trailing newline.

For efficiency, passes through anything with a readline () attribute.

mrjob.util.unarchive(archive\_path, dest)

Extract the contents of a tar or zip file at archive\_path into the directory dest.

### **Parameters**

- archive\_path (str) path to archive file
- **dest** (*str*) path to directory where archive will be extracted

dest will be created if it doesn't already exist.

tar files can be gzip compressed, bzip2 compressed, or uncompressed. Files within zip files can be deflated or stored.

mrjob.util.unique(items)

Yield items from *item* in order, skipping duplicates.

mrjob.util.which (cmd, path=None)

Like the UNIX which command: search in *path* for the executable named *cmd. path* defaults to PATH. Returns None if no such executable found.

This is basically shutil.which() (which was introduced in Python 3.3) without the *mode* argument. Best practice is to always specify *path* as a keyword argument.

mrjob.util.zip\_dir(dir, out\_path, filter=None, prefix='')

Compress the given *dir* into a zip file at *out\_path*.

If we encounter symlinks, include the actual file, not the symlink.

# **Parameters**

- dir (str) dir to tar up
- out\_path (str) where to write the tarball too
- **filter** if defined, a function that takes paths (relative to *dir* and returns True if we should keep them
- **prefix** (*str*) subdirectory inside the tarball to put everything into (e.g. 'mrjob')

# What's New

For a complete list of changes, see CHANGES.txt

# 3.1 0.6.10

Adds official support for PyPy (that is any version of it compatible with Python 2.7/3.5+). If you launch a job in PyPy python\_bin will automatically default to pypy or pypy3 as appropriate.

Note that mrjob does not auto-install PyPy for you on EMR (Amazon Linux does not provide a PyPy package). Installing PyPy yourself at bootstrap time is fairly straightforward, see *Installing PyPy*.

The Spark harness can now be used on EMR, allowing you to run "classic" MRJobs in Spark, which is often faster. Essentially, you launch jobs in the Spark runner with --spark-submit-bin 'mrjob spark-submit -r emr'; see Running classic MRJobs on Spark on EMR for details.

The Spark runner can now optionally disable internal protocols when running "classic" MRJobs, eliminating the (usually) unnecessary effort of encoding data structures into JSON or other string representations and then decoding them. See skip\_internal\_protocol for details.

The EMR runner's default instance type is now m5.xlarge, which works with newer reasons and should make it easier to run Spark jobs. The EMR runner also now logs the DNS of the master node as soon as it is available, to make it easier to SSH in.

Finally, mrjob gives a much clearer error message if you attempt to read a YAML mrjob.conf file without PyYAML installed.

# 3.2 0.6.9

This fixes a bug introduced in 0.6.8 that could break archives or directories uploaded into Hadoop or Spark if the name of the unpacked archive didn't have an archive extension (e.g. .tar.gz).

The Spark runner can now optionally emulate Hadoop's mapreduce.map.input.file configuration property when running the mapper of the first step of a streaming job if you enable emulate\_map\_input\_file. This means that jobs that depend on jobconf\_from\_env('mapreduce.map.input.file') will still work.

The Spark runner also now uses the correct argument names when emulating <code>increment\_counter()</code>, and logs a warning if <code>spark\_tmp\_dir</code> doesn't match <code>spark\_master</code>.

mrjob spark-submit can now pass switches to the Spark script/JAR without explicitly separating them out with --.

The local and inline runners now more correctly emulate the *mapreduce.map.input.file* config property by making it a file://URL.

Deprecated methods add\_file\_option() and add\_passthrough\_option() can now take a type (e.g. int) as their type argument, to better emulate optparse.

# 3.3 0.6.8

# 3.3.1 Nearly full support for Spark

This release adds nearly full support for Spark, including mrjob-speific features like setup scripts and *passthrough options*. See *Why use mrjob with Spark?* for everything mrjob can do with Spark.

This release adds a SparkMRJobRunner (-r spark), which works with any Spark installation, does not require Hadoop, and can access any filesystem supported by both mrjob and Spark (HDFS, S3, GCS). The Spark runner is now the default for *mrjob spark-submit*.

What's *not* supported? mrjob does not yet support Spark on Google Cloud Dataproc. The Spark runner does not yet parse logs to determine probable cause of failure when your job fails (though it does give you the Spark driver output).

# 3.3.2 Spark Hadoop Streaming emulation

Not only does the Spark runner not need Hadoop to run Spark jobs, it doesn't need Hadoop to run most *Hadoop Streaming* jobs, as it knows how to run them directly on Spark. This means if you want to migrate to a non-Hadoop Spark cluster, you can take all your old MRJobs with you. See *Running "classic" MRJobs on Spark* for details.

The "experimental harness script" mentioned in 0.6.7 is now fully integrated into the Spark runner and is no longer supported as a separate feature.

# 3.3.3 Local runner support for Spark

The local and inline runner can now run Spark scripts locally for testing, analogous to the way they've supported Hadoop streaming scripts (except that they *do* require a local Spark installation). See *Other ways to run on Spark*.

# 3.3.4 Other Spark improvements

MRJobs are now Spark-serializable without calling sandbox() (there used to be a problematic reference to sys.stdin). This means you can always pass job methods to rdd.flatMap() etc.

setup scripts are no longer a YARN-specific feature, working on all Spark masters (except local[\*], which doesn't give executors a separate working directory).

Likewise, you can now specify a different name for files in the job's working directory (e.g. --file foo#bar) on all Spark masters.

**Note:** Uploading archives and directories still only works on YARN for now; Spark considers --archives a YARN-specific feature.

When running on a local Spark cluster, uses file://... rather than just the path of the file when necessary (e.g. with --py-files).

cat\_output() now ignores files and subdirectories starting with "." (used to only be "\_"). This allows mrjob to ignore Spark's checksum files (e.g. .part-00000.crc), and also brings mrjob in closer compliance to the way Hadoop input formats read directories.

spark.yarn.appMasterEnv.\* config properties are only set if you're actually running on YARN.

The values of spark\_master and spark\_deploy\_mode can no longer be overridden with configuration properties (-D spark.master=...). While not exactly a "feature," this means that mrjob always knows what Spark platform it's running on.

# 3.3.5 Filesystems

Every runner has an fs attribute that gives access to all the filesystems that runner supports.

Added a put () method to all filesystems, which allows uploading a single file (it used to be that each runner had custom logic for uploads).

It also used to be that if you wanted to create a bucket on S3 or GCS, you had to call create\_bucket(...) explicitly. Now mkdir() will automatically create buckets as needed.

If you still need to access methods specific to a filesystem, you should do so through fs.<name>, where <name> is the (lowercase) name of the storage service. For example the Spark runner's filesystem offers both runner.fs.s3.create\_bucket() and runner.fs.gcs.create\_bucket(). The old style of implicitly passing through FS-specific methods (runner.fs.create\_bucket(...)) is deprecated and going away in v0.7.0.

GCSFilesystem's constructor had a useless local\_tmp\_dir argument, which is now deprecated and going away in v0.7.0.

### 3.3.6 EMR

Fixed a bad bug introduced in 0.6.7 that could prevent mrjob from running on EMR with a non-default temp bucket.

You can now set sub-parameters with extra\_cluster\_params. For example, you can now do:

```
--extra-cluster-param Instances.EmrManagedMasterSecurityGroup=...
```

without clobbering the zone or instance group/fleet configs specified in Instances.

Running your job with --subnet '' now un-sets a subnet specified in your config file (used to be ignored).

If you are using cluster pooling with retries (pool\_wait\_minutes), mrjob now retains information about clusters that is immutable (e.g. AMI version), saving API calls.

# 3.3.7 Dependency upgrades

Bumped the required versions of several Google Cloud Python libraries to be more compatible with current versions of their sub-dependencies (Google libraries pin a fairly narrow range of dependencies). mrjob now requires:

- google-cloud-dataproc at least 0.3.0,
- google-cloud-logging at least 1.9.0, and
- google-cloud-storage at least 1.13.1.

Also dropped support for PyYAML 3.08; now we require at least PyYAML 3.10 (which came out in 2011).

**Note:** We are aware that the Google libraries' extensive dependencies can be a nuisance for mrjob users who don't use Google Cloud. Our tentative plan is to make dependencies specific to a third-party service (including google-cloud-\* and boto3) optional starting in v0.7.0.

3.3. 0.6.8

# 3.3.8 Other bugfixes

Fixed a long-standing bug that would cause the Hadoop runner to hang or raise cryptic errors if hadoop\_bin or spark\_submit\_bin is not executable.

Support files for mrjob.examples (e.g. stop\_words.txt for MRMostUsedWord) are now installed along with mrjob.

Setting a \*\_bin option to an empty value (e.g. --hadoop-bin) now always instructs mrjob to use the default, rather than disabling core features or creating cryptic errors. This affects gcloud\_bin, hadoop\_bin, sh\_bin, and ssh\_bin; the various \*python\_bin options already worked this way.

# 3.4 0.6.7

setup commands now work on Spark (at least on YARN).

Added the *mrjob spark-submit* subcommand, which works as a drop-in replacement for **spark-submit** but with mrjob runners (e.g EMR) and mrjob features (e.g. setup, cmdenv).

Fixed a bug that was causing idle timeout scripts to silently fail on 2.x EMR AMIs.

Fixed a bug that broke create\_bucket() on us-east-1, preventing new mrjob installations from launching on EMR in that region.

Fixed an ImportError from attempting to import os.SIGKILL on Windows.

The default instance type on EMR is now m4.large.

EMR's cluster pooling now knows the CPU and memory capacity of c5 and m5 instances, allowing it to join "better" clusters.

Added the plural form of several switches (separate multiple values with commas):

- --applications
- --archives
- --dirs
- --files
- --libjars
- --py-files

Except for —application, the singular version of these switches (—archive, —dir, —file, —libjar, —py-file) is deprecated for consistency with Hadoop and Spark

sh\_bin is now fully qualified by default (/bin/sh -ex, not sh -ex). sh\_bin may no longer be empty, and a warning is issued if it has more than one argument, to properly support shell script shebangs (e.g. #!/bin/sh -ex) on Linux.

Runners no longer call MRJobs with --steps; instead the job passes its step description to the runner on instantiation. --steps and steps python bin are now deprecated.

The Hadoop and EMR runner can now set SPARK\_PYTHON and SPARK\_DRIVER\_PYTHON to different values if need be (e.g. to match task\_python\_bin, or to support setup scripts in client mode).

The inline runner no longer attempts to run command substeps.

The inline and local runner no longer silently pretend to run non-streaming steps.

The Hadoop runner no longer has the bootstrap\_spark option, which did nothing.

interpreter and steps\_interpreter are deprecated, in anticipation in removing support for writing MRJobs in other programming languages.

Runners now issue a warning if they receive options that belong to other runners (e.g. passing image\_version to the Hadoop runner).

```
mrjob create-cluster now supports --emr-action-on-failure.
```

Updated deprecate escape sequences in mrjob code that would break on Python 3.8.

--help message for mrjob subcommands now correctly includes the subcommand in usage.

mrjob no longer raises AssertionError, instead raising ValueError.

Added an experimental harness script (in mrjob/spark) to run basic MRJobs on Spark, potentially without Hadoop:

```
spark-submit mrjob_spark_harness.py module.of.YourMRJob input_path output_dir
```

Added map\_pairs(), reduce\_pairs(), and combine\_pairs() methods to MRJob, to enable the Spark harness script.

### 3.5 0.6.6

Fixes a longstanding bug where boolean jobconf values were passed to Hadoop in Python format (True instead of true). You can now do safely do something like this:

```
runners:
    emr:
    jobconf:
        mapreduce.output.fileoutputformat.compress: true
```

whereas in prior versions of mrjob, you had to use "true" in quotes.

Added -D as a synonym for -- jobconf, to match Hadoop.

On EMR, if you have SSH set up (see *Configuring SSH credentials*) mrjob can fetch your history log directly from HDFS, allowing it to more quickly diagnose why your job failed.

Added a --local-tmp-dir switch. If you set local tmp dir to empty string, mrjob will use the system default.

You can now pass multiple arguments to Hadoop --hadoop-args (for example, --hadoop-args='-fs hdfs://namenode:port'), rather than having to use --hadoop-arg one argument at time. --hadoop-arg is now deprecated.

Similarly, you can use --spark-args to pass arguments to spark-submit in place of the now-deprecated --spark-arg.

mrjob no longer automatically passes generic arguments (-D and -libjars) to JarSteps, because this confuses some JARs. If you want mrjob to pass generic arguments to a JAR, add <code>GENERIC\_ARGS</code> to your <code>JarStep</code>'s <code>args</code> keyword argument, like you would with <code>INPUT</code> and <code>OUTPUT</code>.

The Hadoop runner now has a spark\_deploy\_mode option.

Fixed the usage: usage: typo in --help messages.

```
mrjob.job.MRJob.add_file_arg() can now take an explicit type=str (used to cause an error).
```

The deprecated optparse emulation methods add\_file\_option() and add\_passthrough\_option() now support type='str' (used to only accept type='string').

Fixed a permissions error that was breaking inline and local mode on some versions of Windows.

3.5. 0.6.6

#### 3.6 0.6.5

This release fixes an issue with self-termination of idle clusters on EMR (see max\_mins\_idle) where the master node sometimes simply ignored sudo shutdown -h now. The idle self termination script now logs to bootstrap-actions/mrjob-idle-termination.log.

**Note:** If you are using *Cluster Pooling*, it's highly recommended you upgrade to this version to fix the self-termination issue.

You can now turn off log parsing (on all runners) by setting read\_logs to false. This can speed up cases where you don't care why a job failed (e.g. integration tests) or where you'd rather use the *diagnose* tool after the fact.

You may specify custom AMIs with the image\_id option. To find Amazon Linux AMIs compatible with EMR that you can use as a base for your custom image, use <code>describe\_base\_emr\_images()</code>.

The default AMI on EMR is now 5.16.0.

New EMR clusters launched by mrjob will be automatically tagged with \_\_mrjob\_label (filename of your mrjob script) and \_\_mrjob\_owner (your username), to make it easier to understand your mrjob usage in CloudWatch etc. You can change the value of these tags with the label and owner options.

You may now set the root EBS volume size for EMR clusters directly with ebs\_root\_volume\_gb (you used to have to use instance\_groups or instance\_fleets).

API clients returned by EMRJobRunner now retry on SSL timeouts. EMR clients returned by  $mrjob.emr.EMRJobRunner.make\_emr\_client()$  won't retry faster than check\_cluster\_every, to prevent throttling.

Cluster pooling recovery (relaunching a job when your pooled cluster self-terminates) now works correctly on singlenode clusters.

### 3.7 0.6.4

This release makes it easy to attach static files to your MRJob with the FILES, DIRS, and ARCHIVES attributes.

In most cases, you no longer need setup scripts to access other python modules or packages from your job because you can use *DIRS* instead. For more details, see *Using other python modules and packages*.

For completeness, also added files (), dirs (), and archives () methods.

*terminate-idle-clusters* now skips termination-protected idle clusters, rather than crashing (this is fixed in 0.5.12, but not previous 0.6.x versions).

Python 3.3 is no longer supported.

mrjob now requires google-cloud-dataproc 0.2.0+ (this library used to be vendored).

### 3.8 0.6.3

### 3.8.1 Read arbitrary file formats

You can now pass entire files in any format to your mapper by defining mapper\_raw(). See Passing entire files to the mapper for an example.

### 3.8.2 Google Cloud Datatproc parity

mrjob now offers feature parity between Google Cloud Dataproc and Amazon Elastic MapReduce. Support for Spark and libjars will be added in a future release. (There is no plan to introduce *Cluster Pooling* with Dataproc.)

Specifically, DataprocJobRunner now supports:

- · fetching and parsing counters
- parsing logs for probable cause of failure
- job progress messages (% complete)
- · Jar steps
- these config options:
  - cloud part size mb (chunked uploading)
  - core\_instance\_config, master\_instance\_config, task\_instance\_config
  - hadoop\_streaming\_jar
  - network/subnet (running in a VPC)
  - service\_account (custom IAM account)
  - service\_account\_scopes (fine-grained permissions)
  - ssh\_tunnel/ssh\_tunnel\_is\_open (resource manager)

Improvements to existing Dataproc features:

- bootstrap scripts run in a temp dir, rather than /
- uses Dataproc's built-in auto-termination feature, rather than a script
- GCS filesystem:
  - cat () streams data rather than dumping to a temp file
  - exists () no longer swallows all exceptions

To get started, read Getting started with Google Cloud.

## 3.8.3 Other changes

mrjob no longer streams your job output to the command line if you specify output\_dir. You can control this with the --cat-output and --no-cat-output switches (--no-output is deprecated).

cloud\_upload\_part\_size has been renamed to cloud\_part\_size\_mb (the old name will work until v0.7.0).

mrjob can now recognize "not a valid JAR" errors from Hadoop and suggest them as probable cause of job failure.

mrjob no longer depends on <code>google-cloud</code> (which implies several other Google libraries). Its current Google library dependencies are <code>google-cloud-logging 1.5.0+</code> and <code>google-cloud-storage 1.9.0+</code>. Future versions of mrjob will depend on <code>google-cloud-dataproc 0.11.0+</code> (currently included with mrjob because it hasn't yet been released).

RetryWrapper now sets name when wrapping methods, making for easier debugging.

3.8. 0.6.3

### 3.9 0.6.2

mrjob is now orders of magnitude quicker at parsing logs, making it practical to diagnose rare errors from very large jobs. However, on some AMIs, it can no longer parse errors without waiting for logs to transfer to S3 (this may be fixed in a future version).

To run jobs on Google Cloud Dataproc, mrjob no longer requires you to install the **gcloud** util (though if you do have it installed, mrjob can read credentials from its configs). For details, see Dataproc Quickstart.

mrjob no longer requires you to select a Dataproc zone prior to running jobs. Auto zone placement (just set region and let Dataproc pick a zone) is now enabled, with the default being auto zone placement in us-west1. mrjob no longer reads zone and region from **gcloud**'s compute engine configs.

mrjob's Dataproc code has been ported from the <code>google-python-api-client</code> library (which is in maintenance mode) to <code>google-cloud-sdk</code>, resulting in some small changes to the GCS filesystem API. See CHANGES.txt for details.

Local mode now has a num\_cores option that allow you to control how tasks it handles simultaneously.

### 3.10 0.6.1

Added the *diagnose* tool (run mrjob diagnose j-CLUSTERID), which determines why a previously run job failed.

Fixed a serious bug that made mrjob unable to properly parse error logs in some cases.

Added the get\_job\_steps() method to EMRJobRunner.

#### 3.11 0.6.0

#### **3.11.1 Dropped Python 2.6**

mrjob now supports Python 2.7 and Python 3.3+. (Some versions of PyPy also work but are not officially supported.)

#### 3.11.2 boto3, not boto

mrjob now uses boto3 rather than boto to talk to AWS. This makes it much simpler to pass user-defined data structures directly to the API, enabling a number of features.

At least version 1.4.6 of boto3 is required to run jobs on EMR.

It is now possible to fully configure instances (including EBS volumes). See instance\_groups for an example.

mrjob also now supports Instance Fleets, which may be fully configured (including EBS volumes) through the instance\_fleets option.

Methods that took or returned boto objects (for example, make\_emr\_conn()) have been completely removed as there as no way to make a deprecated shim for them without keeping boto as a dependency. See *EMRJobRunner* and *S3Filesystem* for new method names.

Note that boto3 reads temporary credentials from \$AWS\_SESSION\_TOKEN, not \$AWS\_SECURITY\_TOKEN as in boto (see aws session token for details).

## 3.11.3 argparse, not optparse

mrjob now uses argparse to parse options, rather than optparse, which has been deprecated since Python 2.7. argparse has slightly different option-parsing logic. A couple of things you should be aware of:

- everything that starts with is assumed to be a switch. ——hadoop—arg=-verbose works, but ——hadoop—arg —verbose does not.
- positional arguments may not be split. mr\_wc.py CHANGES.txt LICENSE.txt -r local will work, but mr\_wc.py CHANGES.txt -r local LICENSE.txt will not.

Passthrough options, file options, etc. are now handled with add\_file\_arg(), add\_passthru\_arg(), configure\_args(), load\_args(), and pass\_arg\_through(). The old methods with "option" in their name are deprecated but still work.

As part of this refactor, OptionStore and its subclasses have been removed; options are now handled by runners directly.

### 3.11.4 Chunks, not lines

mrjob no longer assumes that job output will be line-based. If you *run your job programmatically*, you should read your job output with *cat\_output()*, which yields bytestrings which don't necessarily correspond to lines, and run these through *parse\_output()*, which will convert them into key/value pairs.

runner.fs.cat() also now yields arbitrary bytestrings, not lines. When it yields from multiple files, it will yield an empty bytestring (b'') between the chunks from each file.

read\_file() and read\_input() are now deprecated because they are line-based. Try decompress(),
to\_chunks(), and to\_lines().

#### 3.11.5 Better local/inline mode

The sim runners (inline and local mode) have been completely rewritten, making it possible to fix a number of outstanding issues.

Local mode now runs one mapper/reducer per CPU, using multiprocesssing, for faster results.

We only sort by reducer key (not the full line) unless SORT\_VALUES is set, exposing bad assumptions sooner.

The step\_output\_dir option is now supported, making it easier to debug issues in intermediate steps.

Files in tasks' (e.g. mappers') working directories are marked user-executable, to better imitate Hadoop Distributed Cache. When possible, we also symlink to a copy of each file/archive in the "cache," rather than copying them.

If os.symlink() raises an exception, we fall back to copying (this can be an issue in Python 3 on Windows).

Tasks are run more like they are in Hadoop; input is passed through stdin, rather than as script arguments. mrjob.cat is no longer executable because local mode no longer needs it.

### 3.11.6 Cloud runner improvements

Much of the common code for the "cloud" runners (Dataproc and EMR) has been merged, so that new features can be rolled out in parallel.

The bootstrap option (for both Dataproc and EMR) can now take archives and directories as well as files, like the setup option has since version 0.5.8.

The extra\_cluster\_params option allows you to pass arbitrary JSON to the API at cluster create time (in Dataproc and EMR). The old *emr\_api\_params* option is deprecated and disabled.

3.11. 0.6.0

max\_hours\_idle has been replaced with max\_mins\_idle (the old option is deprecated but still works). The default is 10 minutes. Due to a bug, smaller numbers of minutes might cause the cluster to terminate before the job runs.

It is no longer possible for mrjob to launch a cluster that sits idle indefinitely (except by setting max\_mins\_idle to an unreasonably high value). It is still a good idea to run *report-long-jobs* because mrjob can't tell if a running job is doing useful work or has stalled.

### 3.11.7 EMR now bills by the second, not the hour

Elastic MapReduce recently stopped billing by the full hour, and now bills by the second. This means that *Cluster Pooling* is no longer a cost-saving strategy, though developers might find it handy to reduce wait times when testing.

The mins\_to\_end\_of\_hour option no longer makes sense, and has been deprecated and disabled.

audit-emr-usage has been updated to use billing by the second when approximating time billed and waste.

**Note:** Pooling was enabled by default for some development versions of v0.6.0, prior to the billing change. This did not make it into the release; you must still explicitly turn on *cluster pooling*.

### 3.11.8 Other EMR changes

The default AMI is now 5.8.0. Note that this means you get Spark 2 by default.

Regions are now case-sensitive, and the EU alias for eu-west-1 no longer works.

Pooling no longer adds dummy arguments to the master bootstrap script, instead setting the \_\_mrjob\_pool\_hash and \_\_mrjob\_pool\_name tags on the cluster.

mrjob automatically adds the \_\_mrjob\_version tag to clusters it creates.

Jobs will not add tags to clusters they join rather than create.

enable\_emr\_debugging now works on AMI 4.x and later.

AMI 2.4.2 and earlier are no longer supported (no Python 2.7). There is no longer any special logic for the "latest" AMI alias (which the API no longer supports).

The SSH filesystem no longer dumps file contents to memory.

Pooling will only join a cluster with enough *running* instances to meet its specifications; *requested* instances no longer count.

Pooling is now aware of EBS (disk) setup.

Pooling won't join a cluster that has extra instance types that don't have enough memory or disk space to run your job.

Errors in bootstrapping scripts are no longer dumped as JSON.

visible\_to\_all\_users is deprecated.

### 3.11.9 Massive purge of deprecated code

About a hundred functions, methods, options, and more that were deprecated in v0.5.x have been removed. See CHANGES.txt for details.

## 3.12 0.5.12

This release came out after v0.6.3. It was mostly a backport from v0.6.x.

Python 2.6 and 3.3 are no longer supported.

mrjob.parse.parse\_s3\_uri() handles s3a:// URIs.

terminate-idle-clusters now skips termination-protected idle clusters, rather than crashing.

Since Amazon no longer bills by the full hour, the mins\_to\_end\_of\_hour option now defaults to 60, effectively disabling it.

When mrjob passes an environment dictionary to subprocesses, it ensures that the keys and values are always strs (this mostly affects Python 2 on Windows).

### 3.13 0.5.11

The report-long-jobs utility can now ignore certain clusters based on EMR tags.

This version deals more gracefully with clusters that use instance fleets, preventing crashes that may occur in some rare edge cases.

### 3.14 0.5.10

Fixed an issue where bootstrapping mrjob on Dataproc or EMR could stall if mrjob was already installed.

The *aws\_security\_token* option has been renamed to aws\_session\_token. If you want to set it via environment variable, you still have to use \$AWS\_SECURITY\_TOKEN because that's what boto uses.

Added protocol support for rapidjson; see RapidJSONProtocol and RapidJSONValueProtocol. If available, rapidjson will be used as the default JSON implementation if ujson is not installed.

The master bootstrap script on EMR and Dataproc now has the correct file extension (.sh, not .py).

#### 3.15 0.5.9

Fixed a bug that prevented setup scripts from working on EMR AMIs 5.2.0 and later. Our workaround should be completely transparent unless you use a custom shell binary; see sh\_bin for details.

The EMR runner now correctly re-starts the SSH tunnel to the job tracker/resource manager when a cluster it tries to run a job on auto-terminates. It also no longer requires a working SSH tunnel to fetch job progress (you still a working SSH; see ec2\_key\_pair\_file).

The *emr\_applications* option has been renamed to applications.

The *terminate-idle-clusters* utility is now slightly more robust in cases where your S3 temp directory is an different region from your clusters.

Finally, there a couple of changes that probably only matter if you're trying to wrap your Hadoop tasks (mappers, reducers, etc.) in **docker**:

• You can set *just* the python binary for tasks with task\_python\_bin. This allows you to use a wrapper script in place of Python without perturbing setup scripts.

3.12. 0.5.12

• Local mode now no longer relies on an absolute path to access the *mrjob.cat* utility it uses to handle compressed input files; copying the job's working directory into Docker is enough.

#### 3.16 0.5.8

You can now pass directories to jobs, either directly with the upload\_dirs option, or through setup commands. For example:

```
--setup 'export PYTHONPATH=$PYTHONPATH:your-src-code/#'
```

mrjob will automatically tarball these directories and pass them to Hadoop as archives.

For multi-step jobs, you can now specify where inter-step output goes with step\_output\_dir (--step-output-dir), which can be useful for debugging.

All job step types now take the jobconf keyword argument to set Hadoop properties for that step.

Jobs' --help printout is now better-organized and less verbose.

Made several fixes to pre-filters (commands that pipe into streaming steps):

- you can once again add pre-filters to a single step job by re-defining mapper\_pre\_filter(), combiner\_pre\_filter(), and/or reducer\_pre\_filter()
- local mode now ignores non-zero return codes from pre-filters (this matters for BSD grep)
- local mode can now run pre-filters on compressed input files

mrjob now respects sh\_bin when it needs to wrap a command in sh before passing it to Hadoop (e.g. to support pipes)

On EMR, mrjob now fetches logs from task nodes when determining probable cause of error, not just core nodes (the ones that run tasks and host HDFS).

Several unused functions in mr job.util are now deprecated:

- args\_for\_opt\_dest\_subset()
- bash\_wrap()
- populate\_option\_groups\_with\_options()
- scrape options and index by dest()
- tar\_and\_gzip()

bunzip2\_stream() and gunzip\_stream() have been moved from mrjob.util to mrjob.cat.

SSHFilesystem.ssh\_slave\_hosts() has been deprecated.

Option group attributes in MRJobs have been deprecated, as has the get\_all\_option\_groups () method.

#### 3.17 0.5.7

### 3.17.1 Spark and related changes

mrjob now supports running Spark jobs on your own Hadoop cluster or Elastic MapReduce. mrjob provides significant benefits over Spark's built-in Python support; see *Why use mrjob with Spark*? for details.

Added the py\_files option, to put .zip or .egg files in your job's PYTHONPATH. This is based on a Spark feature, but it works with streaming jobs as well. mrjob is now bootstrapped (see bootstrap\_mrjob) as a .zip file rather than a tarball. If for some reason, the bootstrapped mrjob library won't compile, you'll get much cleaner error messages.

The default AMI version on EMR (see image\_version) has been bumped from 3.11.0 to 4.8.2, as 3.11.0's Spark support is spotty.

On EMR, mrjob now defaults to the cheapest instance type that will work (see instance\_type). In most cases, this is m1.medium, but it needs to be m1.large for Spark worker nodes.

### 3.17.2 Cluster pooling

mrjob can now add up to 1,000 steps on *pooled clusters* on EMR (except on very old AMIs). mrjob now prints debug messages explaining why your job matched a particular pooled cluster when running in verbose mode (the -v option). Fixed a bug that caused pooling to fail when there was no need for a master bootstrap script (e.g. when running with --no-bootstrap-mrjob).

### 3.17.3 Other improvements

Log interpretation is much more efficient at determining a job's probable cause of failure (this works with Spark as well).

When running custom JARs (see JarStep) mrjob now repects libjars and jobconf.

The hadoop\_streaming\_jar option now supports environment variables and ~.

The *terminate-idle-clusters* tool now works with all step types, including Spark. (It's still recommended that you rely on the max\_hours\_idle option rather than this tool.)

mrjob now works in Anaconda3 Jupyter Notebook.

### 3.17.4 Bugfixes

Added several missing command-line switches, including --no-bootstrap-python on Dataproc. Made a major refactor that should prevent these kinds of issues in the future.

Fixed a bug that caused mrjob to crash when the ssh binary (see ssh bin) was missing or not executable.

Fixed a bug that erroneously reported failed or just-started jobs as 100% complete.

Fixed a bug where timestamps were erroneously recognized as URIs. mrjob now only recognizes strings containing :// as URIs (see is\_uri()).

# 3.17.5 Deprecation

The following are deprecated and will be removed in v0.6.0:

- JarStep. "INPUT"; use mrjob.step.INPUT instead
- JarStep. "OUTPUT"; use mrjob.step.OUTPUT instead
- non-strict protocols (see strict\_protocols)
- the *python archives* option (try *this* instead)
- is windows path()
- parse key value list()

3.17. 0.5.7

```
parse_port_range_list()scrape_options_into_new_groups()
```

### 3.18 0.5.6

Fixed a critical bug that caused Dataproc runner to always crash when determining Hadoop version.

Log interpretation now prioritizes task errors (e.g. a traceback from your Python script) as probable cause of failure, even if they aren't the most recent error. Log interpretation will now continue to download and parse task logs until it finds a non-empty stderr log.

Log interpretation also strips the "subprocess failed" Java stack trace that appears in task stderr logs from Hadoop 1.

## 3.19 0.5.5

Functionally equivalent to 0.5.4, except that it restores the deprecated *ami\_version* option as an alias for image\_version, making it easier to upgrade from earlier versions of mrjob.

Also slightly improves *Cluster Pooling* on EMR with updated information on memory and CPU power of various EC2 instance types, and by treating application names (e.g. "Spark") as case-insensitive.

### 3.20 0.5.4

## 3.20.1 Pooling and idle cluster self-termination

**Warning:** This release accidentally removed the *ami\_version* option instead of merely deprecating it. If you are upgrading from an earlier version of mrjob, use version 0.5.5 or later.

This release resolves a long-standing EMR API race condition that made it difficult to use *Cluster Pooling* and idle cluster self-termination (see max\_hours\_idle) together. Now if your pooled job unknowingly runs on a cluster that was in the process of shutting down, it will detect that and re-launch the job on a different cluster.

This means pretty much everyone running jobs on EMR should now enable pooling, with a configuration like this:

```
runners:
    emr:
    max_hours_idle: 1
    pool_clusters: true
```

You may also run the terminate-idle-clusters script periodically, but (barring any bugs) this shouldn't be necessary.

### 3.20.2 Generic EMR option names

Many options to the EMR runner have been made more generic, to make it easier to share code with the Dataproc runner (in most cases, the new names are also shorter and easier to remember):

old option name	new option name
ami_version	image_version
aws_availablity_zone	zone
aws_region	region
check_emr_status_every	check_cluster_every
ec2_core_instance_bid_price	core_instance_bid_price
ec2_core_instance_type	core_instance_type
ec2_instance_type	instance_type
ec2_master_instance_bid_price	master_instance_bid_price
ec2_master_instance_type	master_instance_type
ec2_slave_instance_type	core_instance_type
ec2_task_instance_bid_price	task_instance_bid_price
ec2_task_instance_type	task_instance_type
emr_tags	tags
num_ec2_core_instances	num_core_instances
num_ec2_task_instances	num_task_instances
s3_log_uri	cloud_log_dir
s3_sync_wait_time	cloud_fs_sync_secs
s3_tmp_dir	cloud_tmp_dir
s3_upload_part_size	cloud_upload_part_size

The old option names and command-line switches are now deprecated but will continue to work until v0.6.0. (Exception: *ami\_version* was accidentally removed; if you need it, use 0.5.5 or later.)

num\_ec2\_instances has simply been deprecated (it's just num\_core\_instances plus one).

hadoop\_streaming\_jar\_on\_emr has also been deprecated; in its place, you can now pass a file:// URI to hadoop\_streaming\_jar to reference a path on the master node.

## 3.20.3 Log interpretation

Log interpretation (counters and probable cause of job failure) on Hadoop is more robust, handing a wider variety of log4j formats and recovering more gracefully from permissions errors. This includes fixing a crash that could happen on Python 3 when attempting to read data from HDFS.

Log interpretation used to be partially broken on EMR AMI 4.3.0 and later due to a permissions issue; this is now fixed.

## 3.20.4 pass\_through\_option()

You can now pass through *existing* command-line switches to your job; for example, you can tell a job which runner launched it. See pass\_through\_option() for details.

If you don't do this, self.options.runner will now always be None in your job (it used to confusingly default to 'inline').

## 3.20.5 Stop logging credentials

When mrjob is run in verbose mode (the -v option), the values of all runner options are debug-logged to stderr. This has been the case since the very early days of mrjob.

Unfortunately, this means that if you set your AWS credentials in mrjob.conf, they get logged as well, creating a surprising potential security vulnerability. (This doesn't happen for AWS credentials set through environment variables.)

3.20. 0.5.4

Starting in this version, the values of aws\_secret\_access\_key and aws\_security\_token are shown as '...' if they are set, and all but the last four characters of aws access key id are blanked out as well (e.g. '...YNDR').

### 3.20.6 Other improvements and bugfixes

The ssh tunnel to the resource manager on EMR (see ssh\_tunnel) now connects to its correct *internal* IP; this resolves a firewall issue that existed on some VPC setups.

Uploaded files will no longer be given names starting with \_ or ., since Hadoop's input processing treats these files as "hidden".

The EMR idle cluster self-termination script (see max\_hours\_idle) now only runs on the master node.

The audit-emr-usage command-line tool should no longer constantly trigger throttling warnings.

bootstrap\_python no longer bothers trying to install Python 3 on EMR AMI 4.6.0 and later, since it is already installed.

The --ssh-bind-ports command-line switch was broken (starting in 0.4.5!), and is now fixed.

### 3.21 0.5.3

This release adds support for custom libjars (such as nicknack), allowing easy access to custom input and output formats. This works on Hadoop and EMR (including on a cluster that's already running).

In addition, jobs can specify needed libjars by setting the LIBJARS attribute or overriding the libjars () method. For examples, see *Input and output formats*.

The Hadoop runner now tries *even harder* to find your log files without needing additional configuration (see hadoop\_log\_dirs).

The EMR runner now supports Amazon VPC subnets (see subnet), and, on 4.x AMIs, Application Configurations (see emr\_configurations).

If your EMR cluster fails during bootstrapping, mrjob can now determine the probable cause of failure.

There are also some minor improvements to SSH tunneling and a handful of small bugfixes; see CHANGES.txt for details.

#### 3.22 0.5.2

This release adds basic support for Google Cloud Dataproc which is Google's Hadoop service, roughly analogous to EMR. See Dataproc Quickstart. Some features are not yet implemented:

- fetching counters
- finding probable cause of errors
- running Java JARs as steps

Added the *emr\_applications* option, which helps you configure 4.x AMIs.

Fixed an EMR bug (introduced in v0.5.0) where we were waiting for steps to complete in the wrong order (in a multi-step job, we wouldn't register that the first step had finished until the last one had).

Fixed a bug in SSH tunneling (introduced in v0.5.0) that made connections to the job tracker/resource manager on EMR time out when running on a 2.x AMI inside a VPC (Virtual Private Cluster).

Fixed a bug (introduced in v0.4.6) that kept mrjob from interpreting ~ (home directory) in includes in mrjob.conf.

It is now again possible to run tool modules deprecated in v0.5.0 directly (e.g. **python** -m mrjob.tools.emr.create\_job\_flow). This is still a deprecated feature; it's recommended that you use the appropriate mrjob subcommand instead (e.g. mrjob create-cluster).

### 3.23 0.5.1

Fixes a bug in the previous relase that broke *SORT\_VALUES* and any other attempt by the job to set the partitioner. The --partitioner switch is now deprecated (the choice of partitioner is part of your job semantics).

Fixes a bug in the previous release that caused *strict\_protocols* and check\_input\_paths to be ignored in mrjob.conf. (We would much prefer you fixed jobs that are using "loose protocols" rather than setting strict\_protocols: false in your config file, but we didn't break this on purpose, we promise!)

mrjob terminate-idle-clusters now correctly handles EMR debugging steps (see enable\_emr\_debugging) set up by boto 2.40.0.

Fixed a bug that could result in showing a blank probable cause of error for pre-YARN (Hadoop 1) jobs.

ssh\_bind\_ports now defaults to a range object (xrange on Python 2), so that when you run on emr in verbose mode (-r emr -v), debug logging devotes one line to the value of ssh\_bind\_ports rather than 840.

### 3.24 0.5.0

### 3.24.1 Python versions

mrjob now fully supports Python 3.3+ in a way that should be transparent to existing Python 2 users (you don't have to suddenly start handling unicode instead of str). For more information, see Python 2 vs. Python 3.

If you run a job with Python 3, mrjob will automatically install Python 3 on ElasticMapreduce AMIs (see boot-strap\_python).

When you run jobs on EMR in Python 2, mrjob attempts to match your minor version of Python as well (either python2.6 or python2.7); see python\_bin for details.

**Note:** If you're currently running Python 2.7, and *using yum to install python libraries*, you'll want to use the Python 2.7 version of the package (e.g. python27-numpy rather than python-numpy).

The mrjob command is now installed with Python-version-specific aliases (e.g. mrjob-3, mrjob-3.4), in case you install mrjob for multiple versions of Python.

### 3.24.2 Hadoop

mrjob should now work out-of-the box on almost any Hadoop setup. If **hadoop** is in your path, or you set any commonly-used \$HADOOP\_\* environment variable, mrjob will find the Hadoop binary, the streaming jar, and your logs, without any help on your part (see hadoop\_bin, hadoop\_log\_dirs, hadoop\_streaming\_jar).

mrjob has been updated to fully support Hadoop 2 (YARN), including many updates to  $\tt HadoopFilesystem$ . Hadoop 1 is still supported, though anything prior to  $\tt Hadoop~0.20.203$  is not (mrjob is actually a few months older than  $\tt Hadoop~0.20.203$ , so this used to matter).

3.23. 0.5.1

#### 3.24.3 3.x and 4.x AMIs

mrjob now fully supports the 3.x and 4.x Elastic MapReduce AMIs, including SSH tunneling to the resource mananager, fetching counters and finding probable cause of job failure.

The default *ami\_version* (see image\_version) is now 3.11.0. Our plan is to continue updating this to the lastest (non-broken) 3.x AMI for each 0.5.x release of mrjob.

The default instance\_type is now m1.medium (m1.small is too small for the 3.x and 4.x AMIs)

You can specify 4.x AMIs with either the new release\_label option, or continue using ami\_version; both work.

mrjob continues to support 2.x AMIs. However:

**Warning:** 2.x AMIs are deprecated by AWS, and based on a very old version of Debian (squeeze), which breaks **apt-get** and exposes you to security holes.

Please, please switch if you haven't already.

### 3.24.4 AWS Regions

The new default *aws\_region* (see region) is us-west-2 (Oregon). This both matches the default in the EMR console and, according to Amazon, is carbon neutral.

An edge case that might affect you: EC2 key pairs (i.e. SSH credentials) are region-specific, so if you've set up SSH but not explicitly specified a region, you may get an error saying your key pair is invalid. The fix is simply to *create* new SSH keys for the us-west-2 (Oregon) region.

#### 3.24.5 S3

#### mrjob is much smarter about the way it interacts with S3:

- automatically creates temp bucket in the same region as jobs
- connects to S3 buckets on the endpoint matching their region (no more 307 errors)
  - EMRJobRunner and S3Filesystem methods no longer take s3\_conn args (passing around a single S3 connection no longer makes sense)
- no longer uses the temp bucket's location to choose where you run your job
- rm() no longer has special logic for \*\_\$folder\$ keys
- 1s () recurses "subdirectories" even if you pass it a URI without a trailing slash

### 3.24.6 Log interpretation

The part of mrjob that fetches counters and tells you what probably caused your job to fail was basically unmaintainable and has been totally rewritten. Not only do we now have solid support across Hadoop and EMR AMI versions, but if we missed anything, it should be straightforward to add it.

Once casualty of this change was the **mrjob fetch-logs** command, which means mrjob no longer offers a way to fetch or interpret logs from a *past* job. We do plan to re-introduce this functionality.

### 3.24.7 Protocols

Protocols are now strict by default (they simply raise an exception on unencodable data). "Loose" protocols can be re-enabled with the --no-strict-protocols switch; see *strict\_protocols* for why this is a bad idea.

Protocols will now use the much faster ujson library, if installed, to encode and decode JSON. This is especially recommended for simple jobs that spend a significant fraction of their time encoding and data.

**Note:** If you're using EMR, try out this bootstrap recipe to install u json.

mrjob will fall back to the simplejson library if ujson is not installed, and use the built-in json module if neither is installed.

You can now explicitly specify which JSON implementation you wish to use (e.g. StandardJSONProtocol, SimpleJSONProtocol, UltraJSONProtocol).

### 3.24.8 Status messages

We've tried to cut the logging messages that your job prints as it runs down to the basics (either useful info, like where a temp directory is, or something that tells you why you're waiting). If there are any messages you miss, try running your job with -v.

When a step in your job fails, mrjob no longer prints a useless stacktrace telling you where in the code the runner raised an exception about your step failing. This is thanks to StepFailedException, which you can also catch and interpret if you're *running jobs programmatically*.

## 3.24.9 Deprecation

Many things that were deprecated in 0.4.6 have been removed:

- options:
  - IF\_SUCCESSFUL cleanup option (use ALL)
  - iam\_job\_flow\_role (use iam\_instance\_profile)
- · functions and methods:
  - positional arguments to mrjob.job.MRJob.mr() (don't even use mr(); use
     mrjob.step.MRStep)
  - mrjob.job.MRJob.jar() (use mrjob.step.JarStep)
  - step\_args and name arguments to mrjob.step.JarStep (use args instead of step\_args, and don't use name at all)
  - mrjob.step.MRJobStep (use mrjob.step.MRStep)
  - mrjob.compat.get\_jobconf\_value() (use to jobconf\_from\_env())
  - mrjob.job.MRJob.parse\_counters()
  - mrjob.job.MRJob.parse\_output()
  - mrjob.conf.combine\_cmd\_lists()
  - mrjob.fs.s3.S3Filesystem.get\_s3\_folder\_keys()

3.24. 0.5.0

mrjob.compat functions supports\_combiners\_in\_hadoop\_streaming(), supports\_new\_distributed\_cache\_options(), and uses\_generic\_jobconf(), which only existed to support very old versions of Hadoop, were removed without deprecation warnings (sorry!).

To avoid a similar wave of deprecation warnings in the future, the name of every part of mrjob that isn't meant to be a stable interface provided by the library now starts with an underscore. You can still use these things (or copy them; it's Open Source), but there's no guarantee they'll exist in the next release.

If you want to get ahead of the game, here is a list of things that are deprecated starting in mrjob 0.5.0 (do these *after* upgrading mrjob):

- options:
  - base\_tmp\_dir is now local\_tmp\_dir
  - cleanup options LOCAL\_SCRATCH and REMOTE\_SCRATCH are now LOCAL\_TMP and REMOTE\_TMP
  - emr\_job\_flow\_id is now cluster\_id
  - emr\_job\_flow\_pool\_name is now pool\_name
  - hdfs\_scratch\_dir is now hadoop\_tmp\_dir
  - pool\_emr\_job\_flows is now pool\_clusters
  - s3\_scratch\_uri is now cloud\_tmp\_dir
  - ssh\_tunnel\_to\_job\_tracker is now simply ssh\_tunnel
- · functions and methods:
  - mrjob.job.MRJob.is\_mapper\_or\_reducer() is now is\_task()
  - Filesystem method path\_exists() is now simply exists()
  - Filesystem method path\_join() is now simply join()
  - Use runner.fs explicitly when accessing filesystem methods (e.g. runner.fs.ls(), not runner.ls())
- mrjob subcommands mrjob create-job-flow is now mrjob create-cluster mrjob terminate-idle-job-flows is now mrjob terminate-idle-clusters mrjob terminate-job-flow is now mrjob terminate-cluster

### 3.24.10 Other changes

- mrjob now requires boto 2.35.0 or newer (chances are you're already doing this). Later 0.5.x releases of mrjob may require newer versions of boto.
- visible to all users now defaults to True
- HadoopFilesystem.rm() uses -skipTrash
- new iam\_endpoint option
- custom hadoop\_streaming\_jars are properly uploaded
- JOB cleanup on EMR is temporarily disabled
- mrjob now follows symlinks when 1s () ing the local filesystem (beware recursive symlinks!)
- The interpreter option disables bootstrap\_mrjob by default (interpreter is meant for non-Python jobs)
- Cluster Pooling now respects ec2\_key\_pair

- cluster self-termination (see max\_hours\_idle) now respects non-streaming jobs
- LocalFilesystem now rejects URIs rather than interpreting them as local paths
- local and inline runners no longer have a default hadoop\_version, instead handling jobconf in a version-agnostic way
- steps\_python\_bin now defaults to the current Python interpreter.
- minor changes to mrjob.util:
  - file\_ext() takes filename, not path
  - gunzip\_stream() now yields chunks of bytes, not lines
  - moved random\_identifier() method here from mrjob.aws
  - buffer\_iterator\_to\_line\_iterator() is now named to\_lines(), and no longer appends
     a trailing newline to data.

### 3.25 0.4.6

include: in conf files can now use relative paths in a meaningful way. See Relative includes.

List and environment variable options loaded from included config files can be totally overridden using the !clear tag. See *Clearing configs*.

Options that take lists (e.g. setup) now treat scalar values as single-item lists. See this example.

Fixed a bug that kept the pool\_wait\_minutes option from being loaded from config files.

### 3.26 0.4.5

This release moves mrjob off the deprecated DescribeJobFlows EMR API call.

**Warning:** AWS *again* broke older versions mrjob for at least some new accounts, by returning 400s for the deprecated DescribeJobFlows API call. If you have a newer AWS account (circa July 2015), you must use at least this version of mrjob.

The new API does not provide a way to tell when a job flow (now called a "cluster") stopped provisioning instances and started bootstrapping, so the clock for our estimates of when we are close to the end of a billing hour now start at cluster creation time, and are thus more conservative.

Related to this change, terminate\_idle\_job\_flows no longer considers job flows in the STARTING state idle; use report\_long\_jobs to catch jobs stuck in this state.

terminate\_idle\_job\_flows performs much better on large numbers of job flows. Formerly, it collected all job flow information first, but now it terminates idle job flows as soon as it identifies them.

collect\_emr\_stats and job\_flow\_pool have not been ported to the new API and will be removed in v0.5.0.

Added an aws\_security\_token option to allow you to run mrjob on EMR using temporary AWS credentials.

Added an emr\_tags (see tags) option to allow you to tag EMR job flows at creation time.

EMRJobRunner now has a get\_ami\_version() method.

The hadoop\_version option no longer has any effect in EMR. This option only every did anything on the 1.x AMIs, which mrjob no longer supports.

3.25. 0.4.6

Added many missing switches to the EMR tools (accessible from the **mrjob** command). Formerly, you had to use a config file to get at these options.

You can now access the mrboss tool from the command line: mrjob boss <args>.

Previous 0.4.x releases have worked with boto as old as 2.2.0, but this one requires at least boto 2.6.0 (which is still more than two years old). In any case, it's recommended that you just use the latest version of boto.

This branch has a number of additional deprecation warnings, to help prepare you for mrjob v0.5.0. Please heed them; a lot of deprecated things really are going to be completely removed.

### 3.27 0.4.4

mrjob now automatically creates and uses IAM objects as necessary to comply with new requirements from Amazon Web Services.

(You do not need to install the AWS CLI or run aws emr create-default-roles as the link above describes; mrjob takes care of this for you.)

**Warning:** The change that AWS made essentially broke all older versions of mrjob for all new accounts. If the first time your AWS account created an Elastic MapReduce cluster was on or after April 6, 2015, you should use at least this version of mrjob.

If you must use an old version of mrjob with a new AWS account, see this thread for a possible workaround.

```
--iam-job-flow-role has been renamed to --iam-instance-profile.
```

New --iam-service-role option.

### 3.28 0.4.3

This release also contains many, many bugfixes, one of which probably affects you! See CHANGES.txt for details.

Added a new subcommand, mrjob collect-emr-active-stats, to collect stats about active jobflows and instance counts.

```
--iam-job-flow-role option allows setting of a specific IAM role to run this job flow.
```

You can now use --check-input-paths and --no-check-input-paths on EMR as well as Hadoop.

Files larger than 100MB will be uploaded to S3 using multipart upload if you have the *filechunkio* module installed. You can change the limit/part size with the --s3-upload-part-size option, or disable multipart upload by setting this option to 0. You can now require protocols to be strict from mrjob.conf; this means unencodable input/output will result in an exception rather than the job quietly incrementing a counter. It is recommended you set this for all runners:

```
runners:
    emr:
        strict_protocols: true
    hadoop:
        strict_protocols: true
    inline:
        strict_protocols: true
    local:
        strict_protocols: true
```

You can use --no-strict-protocols to turn off strict protocols for a particular job.

Tests now support pytest and tox.

Support for Python 2.5 has been dropped.

### 3.29 0.4.2

JarSteps, previously experimental, are now fully integrated into multi-step jobs, and work with both the Hadoop and EMR runners. You can now use powerful Java libraries such as Mahout in your MRJobs. For more information, see *Jar steps*.

Many options for setting up your task's environment (--python-archive, --setup-cmd and --setup-script) have been replaced by a powerful --setup option. See the Job Environment Setup Cookbook for examples.

Similarly, many options for bootstrapping nodes on EMR (--bootstrap-cmd, --bootstrap-file, --bootstrap-python-package and --bootstrap-script) have been replaced by a single --bootstrap option. See the EMR Bootstrapping Cookbook.

This release also contains many bugfixes, including problems with boto 2.10.0+, bz2 decompression, and Python 2.5.

### 3.30 0.4.1

The SORT\_VALUES option enables secondary sort, ensuring that your reducer(s) receive values in sorted order. This allows you to do things with reducers that would otherwise involve storing all the values in memory, such as:

- Receiving a grand total before any subtotals, so you can calculate percentages on the fly. See mr\_next\_word\_stats.py for an example.
- Running a window of fixed length over an arbitrary amount of sorted values (e.g. a 24-hour window over timestamped log data).

The max\_hours\_idle option allows you to spin up EMR job flows that will terminate themselves after being idle for a certain amount of time, in a way that optimizes EMR/EC2's full-hour billing model.

For development (not production), we now recommend always using *job flow pooling*, with max\_hours\_idle enabled. Update your *mrjob.conf* like this:

```
runners:
    emr:
    max_hours_idle: 0.25
    pool_emr_job_flows: true
```

**Warning:** If you enable pooling *without* max\_hours\_idle (or cronning terminate\_idle\_job\_flows), pooled job flows will stay active forever, costing you money!

You can now use --no-check-input-paths with the Hadoop runner to allow jobs to run even if hadoop fs -ls can't see their input files (see check\_input\_paths).

Two bits of straggling deprecated functionality were removed:

- Built-in *protocols* must be instantiated to be used (formerly they had class methods).
- Old locations for *mrjob.conf* are no longer supported.

This version also contains numerous bugfixes and natural extensions of existing functionality; many more things will now Just Work (see CHANGES.txt).

3.29. 0.4.2

### 3.31 0.4.0

The default runner is now *inline* instead of *local*. This change will speed up debugging for many users. Use *local* if you need to simulate more features of Hadoop.

The EMR tools can now be accessed more easily via the *mrjob* command. Learn more here.

Job steps are much richer now:

- You can now use mrjob to run jar steps other than Hadoop Streaming. More info
- You can filter step input with UNIX commands. More info
- In fact, you can use arbitrary UNIX commands as your whole step (mapper/reducer/combiner). More info

If you Ctrl+C from the command line, your job will be terminated if you give it time. If you're running on EMR, that should prevent most accidental runaway jobs. *More info* 

mrjob v0.4 requires boto 2.2.

We removed all deprecated functionality from v0.2:

- -hadoop-\*-format
- -\*-protocol switches
- MRJob.DEFAULT\_\*\_PROTOCOL
- MRJob.get\_default\_opts()
- MRJob.protocols()
- PROTOCOL\_DICT
- IF SUCCESSFUL
- DEFAULT\_CLEANUP
- S3Filesystem.get\_s3\_folder\_keys()

We love contributions, so we wrote some guidelines to help you help us. See you on Github!

### 3.32 0.3.5

The *pool\_wait\_minutes* (--pool-wait-minutes) option lets your job delay itself in case a job flow becomes available. Reference: Configuration quick reference

The JOB and JOB\_FLOW cleanup options tell mrjob to clean up the job and/or the job flow on failure (including Ctrl+C). See <code>CLEANUP\_CHOICES</code> for more information.

#### 3.33 0.3.3

You can now include one config file from another.

### 3.34 0.3.2

The EMR instance type/number options have changed to support spot instances:

- core\_instance\_bid\_price
- core\_instance\_type
- master\_instance\_bid\_price
- master\_instance\_type
- slave instance type (alias for core instance type)
- task instance bid price
- task\_instance\_type

There is also a new ami\_version option to change the AMI your job flow uses for its nodes.

For more information, see mrjob.emr.EMRJobRunner.\_\_init\_\_().

The new report\_long\_jobs tool alerts on jobs that have run for more than X hours.

### 3.35 0.3

#### **3.35.1 Features**

### **Support for Combiners**

You can now use combiners in your job. Like <code>mapper()</code> and <code>reducer()</code>, you can redefine <code>combiner()</code> in your subclass to add a single combiner step to run after your mapper but before your reducer. (MRWordFreqCount does this to improve performance.) <code>combiner\_init()</code> and <code>combiner\_final()</code> are similar to their mapper and reducer equivalents.

You can also add combiners to custom steps by adding keyword argumens to your call to steps().

More info: One-step jobs, Multi-step jobs

#### \* init(), \* final() for mappers, reducers, combiners

Mappers, reducers, and combiners have \*\_init() and \*\_final() methods that are run before and after the input is run through the main function (e.g. mapper\_init() and mapper\_final()).

More info: One-step jobs, Multi-step jobs

#### **Custom Option Parsers**

It is now possible to define your own option types and actions using a custom OptionParser subclass.

#### **Job Flow Pooling**

EMR jobs can pull job flows out of a "pool" of similarly configured job flows. This can make it easier to use a small set of job flows across multiple automated jobs, save time and money while debugging, and generally make your life simpler.

More info: Cluster Pooling

#### **SSH Log Fetching**

mrjob attempts to fetch counters and error logs for EMR jobs via SSH before trying to use S3. This method is faster, more reliable, and works with persistent job flows.

More info: Configuring SSH credentials

#### New EMR Tool: fetch\_logs

3.35. 0.3

If you want to fetch the counters or error logs for a job after the fact, you can use the new fetch\_logs tool.

More info: mrjob.tools.emr.fetch\_logs

#### New EMR Tool: mrboss

If you want to run a command on all nodes and inspect the output, perhaps to see what processes are running, you can use the new mrboss tool.

More info: mrjob.tools.emr.mrboss

## 3.35.2 Changes and Deprecations

#### Configuration

The search path order for mrjob.conf has changed. The new order is:

- The location specified by MRJOB\_CONF
- ~/.mrjob.conf
- ~/.mrjob (deprecated)
- mrjob.conf in any directory in PYTHONPATH (deprecated)
- /etc/mrjob.conf

If your mrjob.conf path is deprecated, use this table to fix it:

Old Location	New Location
~/.mrjob	~/.mrjob.conf
somewhere in PYTHONPATH	Specify in MRJOB_CONF

More info: mrjob.conf

#### **Defining Jobs (MRJob)**

Mapper, combiner, and reducer methods no longer need to contain a yield statement if they emit no data.

The --hadoop-\*-format switches are deprecated. Instead, set your job's Hadoop formats with  $HADOOP\_INPUT\_FORMAT/HADOOP\_OUTPUT\_FORMAT$  or  $hadoop\_input\_format()/hadoop\_output\_format()$ . Hadoop formats can no longer be set from mrjob.conf.

In addition to <code>--jobconf</code>, you can now set jobconf values with the <code>JOBCONF</code> attribute or the <code>jobconf()</code> method. To read jobconf values back, use <code>mrjob.compat.jobconf\_from\_env()</code>, which ensures that the correct name is used depending on which version of Hadoop is active.

You can now set the Hadoop partioner class with --partitioner, the PARTITIONER attribute, or the partitioner() method.

More info: Hadoop configuration

#### **Protocols**

Protocols can now be anything with a read() and write() method. Unlike previous versions of mrjob, they can be **instance methods** rather than class methods. You should use instance methods when defining your own protocols.

The --\*protocol switches and DEFAULT\_\*PROTOCOL are deprecated. Instead, use the \*\_PROTOCOL attributes or redefine the \*\_protocol() methods.

Protocols now cache the decoded values of keys. Informal testing shows up to 30% speed improvements.

More info: Protocols

#### **Running Jobs**

#### All Modes

All runners are Hadoop-version aware and use the correct jobconf and combiner invocation styles. This change should decrease the number of warnings in Hadoop 0.20 environments.

All \*\_bin configuration options (hadoop\_bin, python\_bin, and ssh\_bin) take lists instead of strings so you can add arguments (like ['python', '-v']). More info: Configuration quick reference

Cleanup options have been split into cleanup and cleanup\_on\_failure. There are more granular values for both of these options.

Most limitations have been lifted from passthrough options, including the former inability to use custom types and actions.

The job\_name\_prefix option is gone (was deprecated).

All URIs are passed through to Hadoop where possible. This should relax some requirements about what URIs you can use.

Steps with no mapper use **cat** instead of going through a no-op mapper.

Compressed files can be streamed with the cat () method.

#### **EMR Mode**

The default Hadoop version on EMR is now 0.20 (was 0.18).

The instance\_type option only sets the instance type for slave nodes when there are multiple EC2 instance. This is because the master node can usually remain small without affecting the performance of the job.

#### **Inline Mode**

Inline mode now supports the cmdenv option.

#### **Local Mode**

Local mode now runs 2 mappers and 2 reducers in parallel by default.

There is preliminary support for simulating some jobconf variables. The current list of supported variables is:

```
• mapreduce.job.cache.archives
```

- mapreduce.job.cache.files
- mapreduce.job.cache.local.archives
- mapreduce.job.cache.local.files
- mapreduce.job.id
- mapreduce.job.local.dir
- mapreduce.map.input.file
- mapreduce.map.input.length
- mapreduce.map.input.start

3.35. 0.3

- mapreduce.task.attempt.id
- mapreduce.task.id
- mapreduce.task.ismap
- mapreduce.task.output.dir
- mapreduce.task.partition

#### **Other Stuff**

boto 2.0+ is now required.

The Debian packaging has been removed from the repostory.

# Glossary

- **combiner** A function that converts one key and a list of values that share that key (not necessarily all values for the key) to zero or more key-value pairs based on some function. See Concepts for details.
- **Hadoop Streaming** A special jar that lets you run code written in any language on Hadoop. It launches a subprocess, passes it input on stdin, and receives output on stdout. Read more here.
- **input protocol** The *protocol* that converts the input file to the key-value pairs seen by the first step. See *Protocols* for details.
- **internal protocol** The *protocol* that converts the output of one step to the intput of the next. See *Protocols* for details.
- **mapper** A function that converts one key-value pair to zero or more key-value pairs based on some function. See Concepts for details.
- **output protocol** The *protocol* that converts the output of the last step to the bytes written to the output file. See *Protocols* for details.
- **protocol** An object that converts a stream of bytes to and from Python objects. See *Protocols* for details.
- **reducer** A function that converts one key and all values that share that key to zero or more key-value pairs based on some function. See Concepts for details.
- **step** One *mapper*, *combiner*, and *reducer*. Any of these may be omitted from a mrjob step as long as at least one is included.

#### **Appendices**

genindex

modindex

search

### m

```
mrjob.ami,89
mrjob.cat,89
mrjob.compat, 100
mrjob.conf, 101
mrjob.dataproc, 103
mrjob.emr, 104
mrjob.fs.base, 125
mrjob.hadoop, 106
mrjob.inline, 106
mrjob.job, 106
mrjob.local, 118
mrjob.parse, 118
mrjob.protocol, 119
mrjob.retry, 122
mrjob.runner, 123
mrjob.setup, 130
mrjob.spark.runner, 122
mrjob.step, 127
mrjob.tools.diagnose, 95
mrjob.tools.emr.audit_usage, 90
mrjob.tools.emr.create_cluster,91
mrjob.tools.emr.mrboss, 91
mrjob.tools.emr.report_long_jobs,95
mrjob.tools.emr.s3\_tmpwatch, 96
mrjob.tools.emr.terminate_cluster,99
mrjob.tools.emr.terminate_idle_clusters,
mrjob.tools.spark_submit,96
mrjob.util, 132
```

168 Python Module Index

Symbols	combine_dicts() (in module mrjob.conf), 102
\$AWS_SECURITY_TOKEN, 144, 147	combine_envs() (in module mrjob.conf), 102
\$AWS_SESSION_TOKEN, 144	combine_jobconfs() (in module mrjob.conf), 102
\$HADOOP_*, 153	combine_lists() (in module mrjob.conf), 103
\$PYTHONPATH, 63	combine_local_envs() (in module mrjob.conf), 103
\$mapreduce_map_input_file, 49	combine_pairs() (mrjob.job.MRJob method), 116
init() (mrjob.hadoop.HadoopJobRunner method),	combine_path_lists() (in module mrjob.conf), 103
106	combine_paths() (in module mrjob.conf), 103
init() (mrjob.inline.InlineMRJobRunner method),	combine_values() (in module mrjob.conf), 103
106	combiner, 165
init() (mrjob.job.MRJob method), 110	combiner() (mrjob.job.MRJob method), 107
init() (mrjob.local.LocalMRJobRunner method), 118	combiner_cmd() (mrjob.job.MRJob method), 108
init() (mrjob.retry.RetryWrapper method), 122	combiner_final() (mrjob.job.MRJob method), 108
init() (mrjob.runner.MRJobRunner method), 123	combiner_init() (mrjob.job.MRJob method), 108
	combiner_pre_filter() (mrjob.job.MRJob method), 109
A	configure_args() (mrjob.job.MRJob method), 112
add() (mrjob.setup.UploadDirManager method), 130	counters() (mrjob.runner.MRJobRunner method), 125
add() (mrjob.setup.WorkingDirManager method), 130	create_bucket() (mrjob.emr.S3Filesystem method), 105
add_file_arg() (mrjob.job.MRJob method), 113	D
add_passthru_arg() (mrjob.job.MRJob method), 112	D
ARCHIVES (mrjob.job.MRJob attribute), 114	DataprocJobRunner (class in mrjob.dataproc), 103
archives() (mrjob.job.MRJob method), 114	decompress() (in module mrjob.cat), 90
AWS_ACCESS_KEY_ID, 70, 73	describe_base_emr_images() (in module mrjob.ami), 89
AWS_SECRET_ACCESS_KEY, 70, 74	DIRS (mrjob.job.MRJob attribute), 114
AWS_SESSION_TOKEN, 74	dirs() (mrjob.job.MRJob method), 114
TWO_DEBBIOTY_TOTALTY, / T	du() (mrjob.fs.base.Filesystem method), 126
В	_
bunzip2_stream() (in module mrjob.cat), 89	E
BytesProtocol (class in mrjob.protocol), 120	EMRJobRunner (class in mrjob.emr), 104
Bytes Value Protocol (class in mrjob.protocol), 119	environment variable
Bytes variation (class in mijos.protocor), 119	\$AWS_SECURITY_TOKEN, 144, 147
C	\$AWS_SESSION_TOKEN, 144
	\$HADOOP_*, 153
can_handle_path() (mrjob.fs.base.Filesystem method),	\$PYTHONPATH, 63
	\$mapreduce_map_input_file, 49
cat() (mrjob.fs.base.Filesystem method), 126 cat_output() (mrjob.runner.MRJobRunner method), 124	AWS_ACCESS_KEY_ID, 70, 73
	AWS_SECRET_ACCESS_KEY, 70, 74
cleanup() (mrjob.runner.MRJobRunner method), 124	AWS_SESSION_TOKEN, 74
CLEANUP_CHOICES (in module mrjob.options), 124	MRJOB_CONF, 34, 101, 162
cmd_line() (in module mrjob.util), 132	PATH, 35, 134
combine_cmds() (in module mrjob.conf), 102	•

PYTHONPATH, 24, 59, 162 TZ, 35 exists() (mrjob.fs.base.Filesystem method), 126	is_s3_uri() (in module mrjob.parse), 118 is_task() (mrjob.job.MRJob method), 113 is_uri() (in module mrjob.parse), 119
expand_path() (in module mrjob.util), 132	J
File_ext() (in module mrjob.util), 132 FILES (mrjob.job.MRJob attribute), 114 files() (mrjob.job.MRJob method), 114 Filesystem (class in mrjob.fs.base), 125 find_mrjob_conf() (in module mrjob.conf), 101 fs (mrjob.runner.MRJobRunner attribute), 125 fully_qualify_hdfs_path() (in module mrjob.hadoop), 106	JarStep (class in mrjob.step), 127  JOBCONF (mrjob.job.MRJob attribute), 116  jobconf() (mrjob.job.MRJob method), 116  jobconf_from_dict() (in module mrjob.compat), 100  jobconf_from_env() (in module mrjob.compat), 100  join() (mrjob.fs.base.Filesystem method), 126  JSONProtocol (class in mrjob.protocol), 120  JSONValueProtocol (class in mrjob.protocol), 121
G	L
GCSFilesystem (class in mrjob.dataproc), 103 GENERIC_ARGS (in module mrjob.step), 130 get_all_bucket_names() (mrjob.emr.S3Filesystem method), 105 get_bucket() (mrjob.emr.S3Filesystem method), 105 get_cluster_id() (mrjob.emr.EMRJobRunner method),	LIBJARS (mrjob.job.MRJob attribute), 117 libjars() (mrjob.job.MRJob method), 117 load_args() (mrjob.job.MRJob method), 113 load_opts_from_mrjob_conf() (in module mrjob.conf),
104	102 LocalMRJobRunner (class in mrjob.local), 118
get_hadoop_version() (mrjob.runner.MRJobRunner method), 125	log_to_null() (in module mrjob.util), 133
get_image_version() (mrjob.emr.EMRJobRunner method), 104	log_to_stream() (in module mrjob.util), 133 ls() (mrjob.fs.base.Filesystem method), 126
get_job_key() (mrjob.runner.MRJobRunner method), 125 get_job_steps() (mrjob.emr.EMRJobRunner method), 104	M make_ec2_client() (mrjob.emr.EMRJobRunner method),
get_opts() (mrjob.runner.MRJobRunner method), 125 gunzip_stream() (in module mrjob.cat), 90	105 make_emr_client() (mrjob.emr.EMRJobRunner method),
Н	104 make_iam_client() (mrjob.emr.EMRJobRunner method),
Hadoop Streaming, <b>165</b> HADOOP_INPUT_FORMAT (mrjob.job.MRJob attribute), 116	105 make_runner() (mrjob.job.MRJob method), 110 make_s3_client() (mrjob.emr.S3Filesystem method), 105
nadoop_input_format() (mrjob.job.MRJob method), 116 HADOOP_OUTPUT_FORMAT (mrjob.job.MRJob at-	make_s3_resource() (mrjob.emr.S3Filesystem method), 105
tribute), 116 nadoop_output_format() (mrjob.job.MRJob method), 116 HadoopJobRunner (class in mrjob.hadoop), 106	map_pairs() (mrjob.job.MRJob method), 115 map_version() (in module mrjob.compat), 101 mapper, <b>165</b>
	mapper() (mrjob.job.MRJob method), 106 mapper_cmd() (mrjob.job.MRJob method), 108
ncrement_counter() (mrjob.job.MRJob method), 111 inlineMRJobRunner (class in mrjob.inline), 106 NPUT (in module mrjob.step), 129 nput protocol, <b>165</b>	mapper_final() (mrjob.job.MRJob method), 107 mapper_init() (mrjob.job.MRJob method), 107 mapper_pre_filter() (mrjob.job.MRJob method), 109 mapper_raw() (mrjob.job.MRJob method), 109 md5sum() (mrjob.fs.base.Filesystem method), 126
NPUT_PROTOCOL (mrjob.job.MRJob attribute), 111 nput_protocol() (mrjob.job.MRJob method), 111 nternal protocol, 165 NTERNAL PROTOCOL (mrjob.job.MRJob attribute)	mkdir() (mrjob.fs.base.Filesystem method), 126 mr_job_script() (mrjob.job.MRJob class method), 115 MRJob (class in mrjob.job), 106
NTERNAL_PROTOCOL (mrjob.job.MRJob attribute), 111	mrjob.ami (module), 89
nternal_protocol() (mrjob.job.MRJob method), 111	mrjob.cat (module), 89 mrjob.compat (module), 100

170 Index

mrjob.conf (module), 101	$path\_to\_uri() \ (mrjob.setup.UploadDirManager \ method),$
mrjob.dataproc (module), 103	130
mrjob.emr (module), 104	paths() (mrjob.setup.WorkingDirManager method), 131
mrjob.fs.base (module), 125	pick_protocols() (mrjob.job.MRJob method), 112
mrjob.hadoop (module), 106	PickleProtocol (class in mrjob.protocol), 122
mrjob.inline (module), 106	PickleValueProtocol (class in mrjob.protocol), 122
mrjob.job (module), 106	protocol, 165
mrjob.local (module), 118	put() (mrjob.fs.base.Filesystem method), 126
mrjob.parse (module), 118	PYTHONPATH, 24, 59, 162
mrjob.protocol (module), 119	D
mrjob.retry (module), 122	R
mrjob.runner (module), 123	random_identifier() (in module mrjob.util), 133
mrjob.setup (module), 130	RapidJSONProtocol (class in mrjob.protocol), 121
mrjob.spark.runner (module), 122	RapidJSONValueProtocol (class in mrjob.protocol), 121
mrjob.step (module), 127	RawProtocol (class in mrjob.protocol), 120
mrjob.tools.diagnose (module), 95	RawValueProtocol (class in mrjob.protocol), 119
mrjob.tools.emr.audit_usage (module), 90	read_file() (in module mrjob.util), 133
mrjob.tools.emr.create_cluster (module), 91	read_input() (in module mrjob.util), 133
mrjob.tools.emr.mrboss (module), 91	reduce_pairs() (mrjob.job.MRJob method), 115
mrjob.tools.emr.report_long_jobs (module), 95	reducer, 165
mrjob.tools.emr.s3_tmpwatch (module), 96	reducer() (mrjob.job.MRJob method), 107
mrjob.tools.emr.terminate_cluster (module), 99	reducer_cmd() (mrjob.job.MRJob method), 108
mrjob.tools.emr.terminate_idle_clusters (module), 99	reducer_final() (mrjob.job.MRJob method), 108
mrjob.tools.spark_submit (module), 96	reducer_init() (mrjob.job.MRJob method), 108
mrjob.util (module), 132	reducer_pre_filter() (mrjob.job.MRJob method), 109
MRJOB_CONF, 34, 101, 162	ReprProtocol (class in mrjob.protocol), 121
MRJobRunner (class in mrjob.runner), 123	ReprValueProtocol (class in mrjob.protocol), 121
MRStep (class in mrjob.step), 127	RetryWrapper (class in mrjob.retry), 122
NI	rm() (mrjob.fs.base.Filesystem method), 126
N	run() (mrjob.job.MRJob class method), 110
name() (mrjob.setup.WorkingDirManager method), 131	run() (mrjob.runner.MRJobRunner method), 124
name_to_path() (mrjob.setup.WorkingDirManager	run_combiner() (mrjob.job.MRJob method), 115
method), 131	run_job() (mrjob.job.MRJob method), 115
name_uniquely() (in module mrjob.setup), 131	run_mapper() (mrjob.job.MRJob method), 115
	run_reducer() (mrjob.job.MRJob method), 115
O	S
OUTPUT (in module mrjob.step), 129	
output protocol, 165	S3Filesystem (class in mrjob.fs.s3), 105
OUTPUT_PROTOCOL (mrjob.job.MRJob attribute), 111	safeeval() (in module mrjob.util), 134
output_protocol() (mrjob.job.MRJob method), 111	sandbox() (mrjob.job.MRJob method), 117
D	save_current_environment() (in module mrjob.util), 134
P	save_cwd() (in module mrjob.util), 134
parse_and_save_options() (in module mrjob.util), 133	save_sys_path() (in module mrjob.util), 134
parse_legacy_hash_path() (in module mrjob.setup), 131	save_sys_std() (in module mrjob.util), 134
parse_mr_job_stderr() (in module mrjob.parse), 119	set_status() (mrjob.job.MRJob method), 111
parse_output() (mrjob.job.MRJob method), 110	shlex_split() (in module mrjob.util), 134
parse_output_line() (mrjob.job.MRJob method), 110	show_steps() (mrjob.job.MRJob method), 116
parse_s3_uri() (in module mrjob.parse), 119	SimpleJSONProtocol (class in mrjob.protocol), 121
parse_setup_cmd() (in module mrjob.setup), 132	SimpleJSONValueProtocol (class in mrjob.protocol), 121
PARTITIONER (mrjob.job.MRJob attribute), 117	SORT_VALUES (mrjob.job.MRJob attribute), 112
partitioner() (mrjob.job.MRJob method), 117	spark() (mrjob.job.MRJob method), 109
pass_arg_through() (mrjob.job.MRJob method), 113	SparkJarStep (class in mrjob.step), 128
PATH, 35, 134	SparkScriptStep (class in mrjob.step), 129
	SparkStep (class in mrjob.step), 128

Index 171

```
StandardJSONProtocol (class in mrjob.protocol), 121
StandardJSONValueProtocol (class in mrjob.protocol),
         121
step, 165
steps() (mrjob.job.MRJob method), 109
stream output() (mrjob.runner.MRJobRunner method),
strip_microseconds() (in module mrjob.util), 134
TextProtocol (class in mrjob.protocol), 120
TextValueProtocol (class in mrjob.protocol), 120
to_chunks() (in module mrjob.cat), 90
to_lines() (in module mrjob.util), 134
to_uri() (in module mrjob.parse), 119
touchz() (mrjob.fs.base.Filesystem method), 126
translate_jobconf() (in module mrjob.compat), 101
translate_jobconf_dict() (in module mrjob.compat), 101
translate_jobconf_for_all_versions() (in module mr-
         job.compat), 101
TZ, 35
U
UltraJSONProtocol (class in mrjob.protocol), 120
UltraJSONValueProtocol (class in mrjob.protocol), 121
unarchive() (in module mrjob.util), 134
unique() (in module mrjob.util), 134
UploadDirManager (class in mrjob.setup), 130
uri() (mrjob.setup.UploadDirManager method), 130
uses_yarn() (in module mrjob.compat), 101
V
version_gte() (in module mrjob.compat), 101
W
which() (in module mrjob.util), 134
WorkingDirManager (class in mrjob.setup), 130
Z
zip_dir() (in module mrjob.util), 134
```

172 Index