



yes, we have a deluge of data



and sometimes, it does seem very big...



and yes, there are ways to  
crunch it fast...



Streamlining Workflows

speed up analysis using a computing cluster



# why bother?

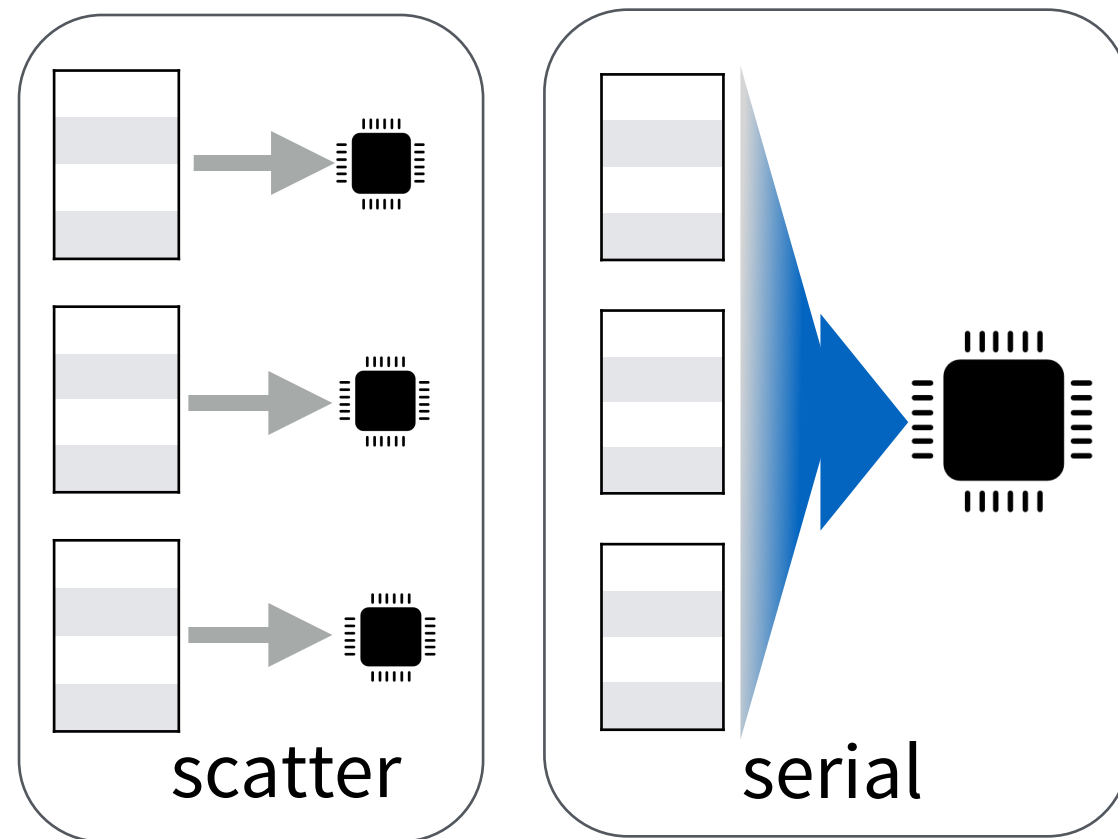
- ➔when one needs to wrangle a lot of data
- ➔and there are multiple steps involved
- ➔esp. when some of the steps can be further broken down and processed in parallel
- ➔use a computing cluster, submit a web of jobs

- ✓ Effectively process a **multi-step pipeline**, spawning it across the computing cluster
- ✓ **Reproducible and transparent**, with cleanly structured execution logs
- ✓ **Track and re-run flows**
- ✓ **Lean and Portable**, with easy installation
- ✓ Run the same pipeline in the cloud (using star cluster) OR a local machine
- ✓ Supports **multiple cluster computing platforms** (torque, lsf, sge, slurm ...)

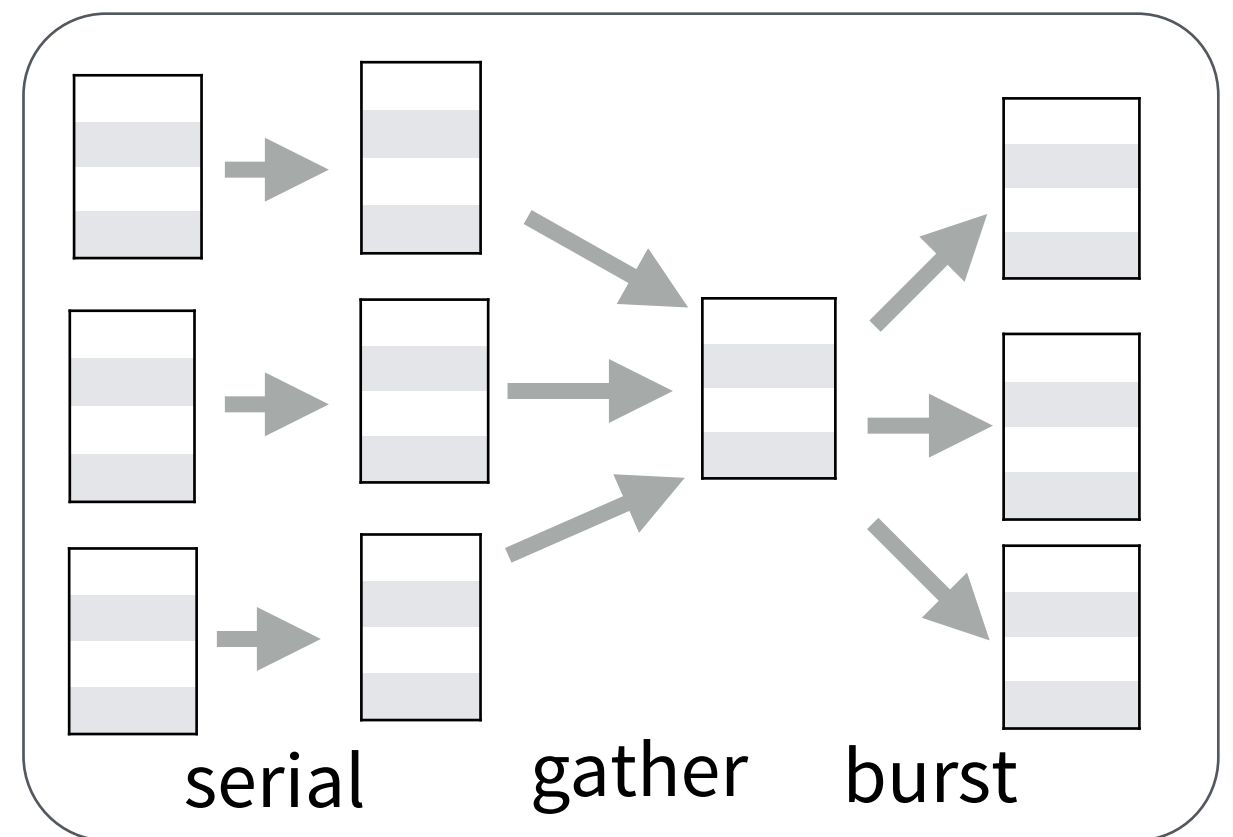


# five simple terms, defining all relationships

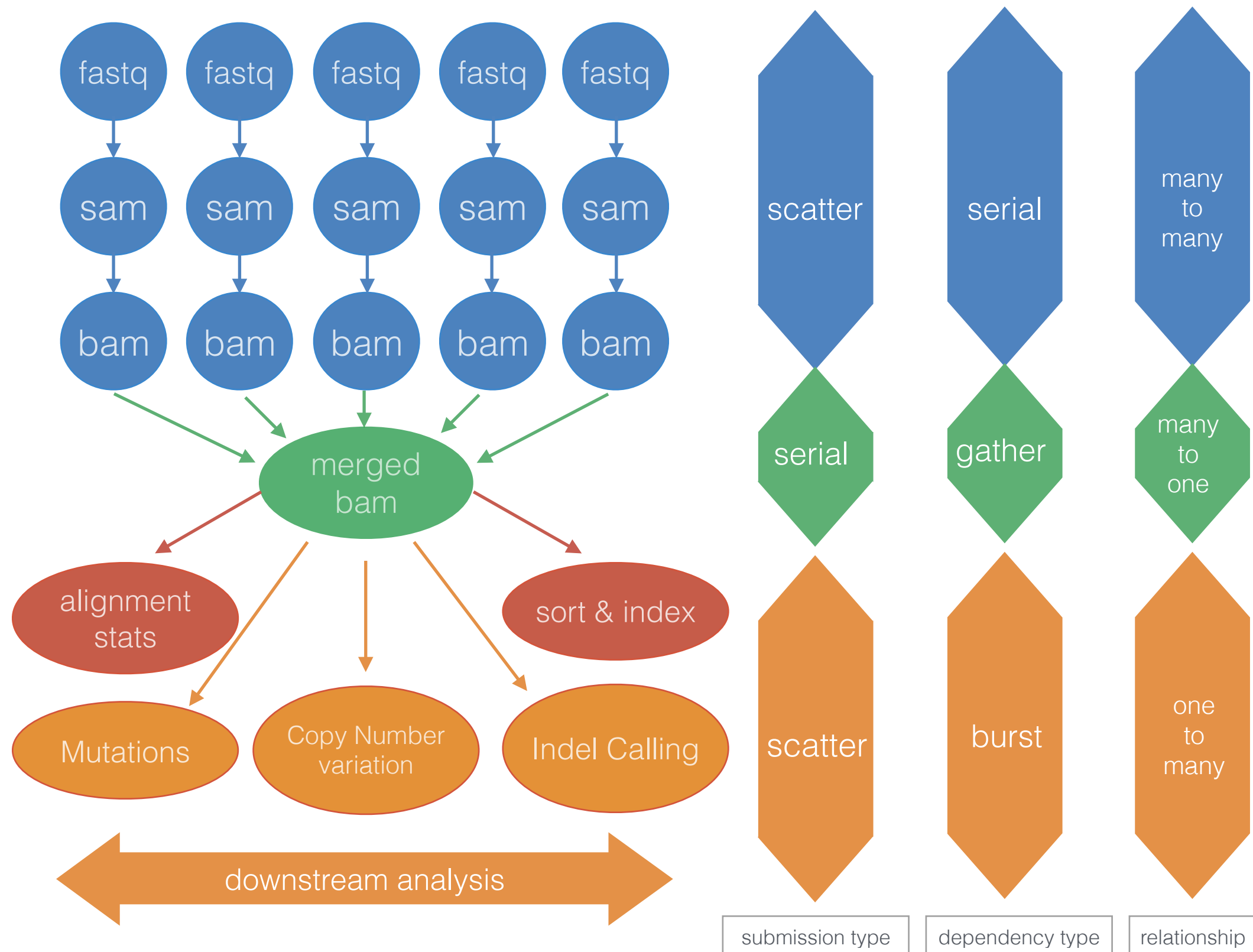
## submission types



## dependency types



# Using a genomics example flow, with flowr concepts



All attributes of this flow can  
be defined using these concepts

a simple pipeline, where

- ★ we would sleep for a few seconds
- ★ create a few small files
- ★ merge those files
- ★ get the size of the resulting merged file

# simple pipeline in bash

say Hello to  
the world

```
echo 'Hello World !'
```

wait for a few  
seconds...

```
sleep 5  
sleep 5
```

create two  
small files

```
cat $RANDOM > tmp1  
cat $RANDOM > tmp2
```

merge the  
two files

```
cat tmp1 tmp2 > tmp
```

check the size of  
the resulting file

```
du -sh tmp
```

- ★ we would sleep for a few seconds
- ★ create a few small files
- ★ merge those files
- ★ get the size of the resulting merged file



# wrap bash commands into R

say Hello to  
the world

```
hello='echo Hello World !'
```

wait for a few  
seconds...

```
sleep=c('sleep 5', 'sleep 5')
```

create two  
small files

```
tmp=c('cat $RANDOM > tmp1',  
      'cat $RANDOM > tmp2')
```

merge the  
two files

```
merge='cat tmp1 tmp2 > tmp'
```

```
size='du -sh tmp'
```

check the size of  
the resulting file

- ★ we would sleep for a few seconds
- ★ create a few small files
- ★ merge those files
- ★ get the size of the resulting merged file


# create a table of all commands

```
hello='echo Hello World !'
sleep=c('sleep 5', 'sleep 5')
tmp=c('cat $RANDOM > tmp1',
      'cat $RANDOM > tmp2')
merge='cat tmp1 tmp2 > tmp'
size='du -sh tmp'
```

```
library(flowr)
lst = list(hello=hello,
           sleep=sleep,
           tmp=tmp,
           merge=merge,
           size=size)

flowmat = to_flowmat(lst, "samp1")
```

create a  
**named** list



samplename	jobname	cmd
samp1	hello	echo Hello World !
samp1	sleep	sleep 5
samp1	sleep	sleep 5
samp1	tmp	cat \$RANDOM > tmp1
samp1	tmp	cat \$RANDOM > tmp2
samp1	merge	cat tmp1 tmp2 > tmp
samp1	size	du -sh tmp

create a  
table

a simple tab-delim table



**connect the dots...**

**flow definition decides the  
sequence of steps**

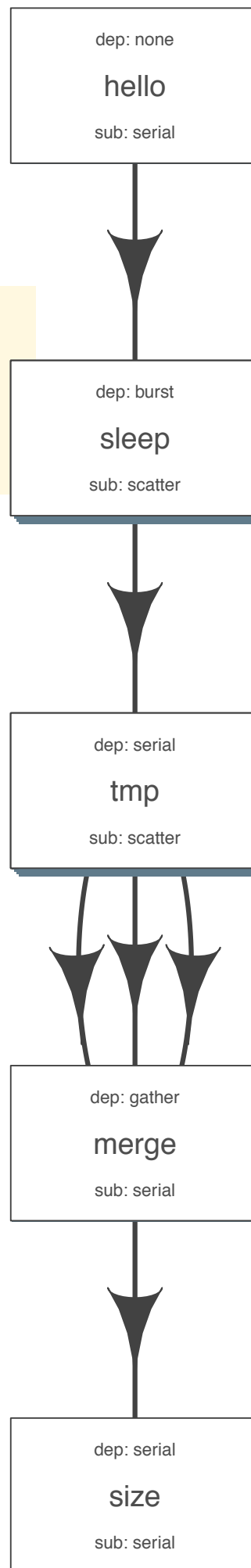
# create a flow definition

```
flowdef = to_flowdef(flowmat,  
    sub_type = c("serial", "scatter", "scatter", "serial", "serial"),  
    dep_type = c("none", "burst", "serial", "gather", "serial"),  
    platform = "local")
```

jobname	sub_type	prev_jobs	dep_type	cpu
:-----	:-----	:-----	:-----	---:
hello	serial	none	none	1
sleep	scatter	hello	burst	1
tmp	scatter	sleep	serial	1
merge	serial	tmp	gather	1
size	serial	merge	serial	1

a simple tab-delim table

```
plot_flow(flowdef)
```





**stitch a flow...**

**use a flow mat and flow def,  
to create a flow object**

# flow mat

samplename	jobname	cmd
:	:	:
samp1	hello	echo Hello World !
samp1	sleep	sleep 5
samp1	sleep	sleep 5
samp1	tmp	cat \$RANDOM > tmp1
samp1	tmp	cat \$RANDOM > tmp2
samp1	merge	cat tmp1 tmp2 > tmp
samp1	size	du -sh tmp

+

# flow def

jobname	sub_type	prev_jobs	dep_type	cpu
:	:	:	:	:
hello	serial	none	none	1
sleep	scatter	hello	burst	1
tmp	scatter	sleep	serial	1
merge	serial	tmp	gather	1
size	serial	merge	serial	1

**stitch & submit to the cluster  
(cloud or server)**

```
fobj = to_flow(flowmat, flowdef, execute = TRUE)
```

Flow is being processed. Track it from R/Terminal using:  
flowr status x=~/.flowr/runs/flowname-samp1-20151005-16-01-38-M8WniKJo  
OR from R using:  
status(x=~/.flowr/runs/flowname-samp1-20151005-16-01-38-M8WniKJo')

```
Working on: hello
|=====| 25%
Working on: sleep
|=====| 50%
Working on: merge
|=====| 100%
Working on: size
```



# submit a flow, then...

## `status()`

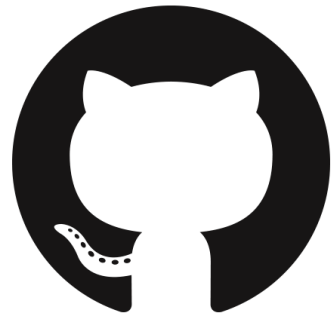
monitor the status of a single flow  
OR multiple flows

## `kill()`

kill all the associated jobs, of one or many flows

## `rerun()`

One can rerun the flow from an intermediate step



[github.com/sahilseth/flowr](https://github.com/sahilseth/flowr)

complete documentation: [docs.flowr.space](https://docs.flowr.space)

email: [sahil.seth@me.com](mailto:sahil.seth@me.com)



**Extra details**

## Flow mat

samplename	jobname	cmd
sample1	A	sleep 2 && sleep 5;echo hello
sample1	A	sleep 13 && sleep 7;echo hello
sample1	B	head -c 100000 /dev/urandom > tmp1
sample1	B	head -c 100000 /dev/urandom > tmp1
sample1	C	cat tmp1 tmp2 tmp3 > merged
sample1	D	du -sh merged
sample1	D	ls merged

- use any language to create a flow mat (a tsv file)
- cmd column defines commands to run

## Flow Definition

Define Relationships				Resource Requirements				
jobname	submission type	previous job(s)	dependency type	queue	memory	time	cpu	platform
A	scatter	none	none	medium	163185	23:00	1	lsf
B	scatter	A	serial	medium	163185	23:00	1	lsf
C	serial	B	gather	medium	163185	23:00	1	lsf
D	scatter	C	burst	medium	163185	23:00	1	lsf

- creatively define relationships using submission and dependency types
- each row describes resources for **one** step, providing full flexibility