NAME
parallel_alternatives - Alternatives to GNU parallel

DIFFERENCES BETWEEN GNU Parallel AND ALTERNATIVES

There are a lot programs that share functionality with GNU parallel. Some of these are specialized tools, and while GNU parallel can emulate many of them, a specialized tool can be better at a given task. GNU parallel strives to include the best of the general functionality without sacrificing ease of use.

parallel has existed since 2002-01-06 and as GNU parallel since 2010. A lot of the alternatives have not had the vitality to survive that long, but have come and gone during that time.

GNU parallel is actively maintained with a new release every month since 2010. Most other alternatives are fleeting interests of the developers with irregular releases and only maintained for a few years.

SUMMARY LEGEND

The following features are in some of the comparable tools:

Inputs
I1. Arguments can be read from stdin
I2. Arguments can be read from a file
I3. Arguments can be read from multiple files
I4. Arguments can be read from command line
I5. Arguments can be read from a table
I6. Arguments can be read from the same file using #! (shebang)
I7. Line oriented input as default (Quoting of special chars not needed)

Manipulation of input
M1. Composed command
M2. Multiple arguments can fill up an execution line
M3. Arguments can be put anywhere in the execution line
M4. Multiple arguments can be put anywhere in the execution line
M5. Arguments can be replaced with context
M6. Input can be treated as the complete command line

Outputs
O1. Grouping output so output from different jobs do not mix
O2. Send stderr (standard error) to stderr (standard error)
O3. Send stdout (standard output) to stdout (standard output)
O4. Order of output can be same as order of input
O5. Stdout only contains stdout (standard output) from the command
O6. Stderr only contains stderr (standard error) from the command
O7. Buffering on disk
O8. No temporary files left if killed
O9. Test if disk runs full during run
O10. Output of a line bigger than 4 GB

Execution
E1. Running jobs in parallel
E2. List running jobs
E3. Finish running jobs, but do not start new jobs
E4. Number of running jobs can depend on number of cpus
E5. Finish running jobs, but do not start new jobs after first failure
E6. Number of running jobs can be adjusted while running
E7. Only spawn new jobs if load is less than a limit

Remote execution
R1. Jobs can be run on remote computers
R2. Basefiles can be transferred
R3. Argument files can be transferred
R4. Result files can be transferred
R5. Cleanup of transferred files
R6. No config files needed
R7. Do not run more than SSHD's MaxStartups can handle
R8. Configurable SSH command
R9. Retry if connection breaks occasionally

Semaphore
S1. Possibility to work as a mutex
S2. Possibility to work as a counting semaphore

Legend
- = no
x = not applicable
ID = yes

As every new version of the programs are not tested the table may be outdated. Please file a bug report if you find errors (See REPORTING BUGS).

parallel:
I1 I2 I3 I4 I5 I6 I7
M1 M2 M3 M4 M5 M6
O1 O2 O3 O4 O5 O6 O7 O8 O9 O10
E1 E2 E3 E4 E5 E6 E7
R1 R2 R3 R4 R5 R6 R7 R8 R9
S1 S2

DIFFERENCES BETWEEN xargs AND GNU Parallel
Summary (see legend above):
I1 I2 - - - - -
- M2 M3 - - -
- O2 O3 - O5 O6
E1 - - - - - -
- - - - x - - -
- -

xargs offers some of the same possibilities as GNU parallel.
xargs deals badly with special characters (such as space, \, ' and "). To see the problem try this:

touch important_file
touch 'not important_file'
ls not* | xargs rm
mkdir -p "My brother's 12" records
ls | xargs rmdir
touch 'c:\windows\system32\clfs.sys'
echo 'c:\windows\system32\clfs.sys' | xargs ls -1

You can specify -0, but many input generators are not optimized for using GNU NUL as separator but are optimized for newline as separator. E.g. awk, ls, echo, tar -v, head (requires using -z), tail (requires using -z), sed (requires using -z), perl (-0 and `0 instead of \n), locate (requires using -0), find (requires using -print0), grep (requires using -z or -Z), sort (requires using -z).

GNU parallel's newline separation can be emulated with:
cat | xargs -d "\n" -n1 command

xargs can run a given number of jobs in parallel, but has no support for running number-of-cpu-cores jobs in parallel.

xargs has no support for grouping the output, therefore output may run together, e.g. the first half of a line is from one process and the last half of the line is from another process. The example Parallel grep cannot be done reliably with xargs because of this. To see this in action try:

parallel perl -e '$a="1"."\{\}"x10000000;print \$a,\\"\n\"'; \n    > { } :::: a b c d e f g h
# Serial = no mixing = the wanted result
# 'tr -s a-z' squeezes repeating letters into a single letter
parallel perl -f -e 'perl -ne \051=S|=1; for(split//)\{ print; select($a,$a,$a,0.100);\}\' \n    > { } :::: a b c d e f g h
parallel perl -e 'perl -n l grep 1 | tr -s a-z
# Compare to 8 jobs in parallel
parallel -kP8 -nI grep 1 :::: a b c d e f g h | tr -s a-z
parallel perl -f -e 'perl -ne \051=S|=1; for(split//)\{ print; select($a,$a,$a,0.100);\}\' \n    > { } :::: a b c d e f g h
parallel perl -e 'perl -n l grep 1 | tr -s a-z
parallel perl -f -e 'perl -ne \051=S|=1; for(split//)\{ print; select($a,$a,$a,0.100);\}\' \n    > { } :::: a b c d e f g h

Or try this:

slow_seq() {
    echo Count to "$@
    seq "$@" |
    perl -ne \051=1; for(split//)\{ print; select($a,$a,$a,0.100);\}"}
export -f slow_seq
# Serial = no mixing = the wanted result
seq 8 | xargs -n 1 -P1 -I {} bash -c slow_seq {}'
# Compare to 8 jobs in parallel
seq 8 | parallel -P8 slow_seq {}
seq 8 | xargs -n 1 -P8 -I {} bash -c slow_seq {}'

xargs has no support for keeping the order of the output, therefore if running jobs in parallel using xargs the output of the second job cannot be postponed till the first job is done.

xargs has no support for running jobs on remote computers.

xargs has no support for context replace, so you will have to create the arguments.

If you use a replace string in xargs (-I) you can not force xargs to use more than one argument.
Quoting in `xargs` works like `-q` in GNU parallel. This means composed commands and redirection require using `bash -c`.

```bash
s l s  |  parallel "wc {} >\{}.wc"
ls | parallel "echo {}; ls {}|wc"
```

becomes (assuming you have 8 cores and that none of the filenames contain space, " or ").

```bash
ls | xargs -d "\n" -P8 -I {} bash -c "wc {}
ls | xargs -d "\n" -P8 -I {} bash -c "echo {}; ls {}|wc"
```

A more extreme example can be found on: https://unix.stackexchange.com/q/405552/

https://www.gnu.org/software/findutils/

**DIFFERENCES BETWEEN find -exec AND GNU Parallel**

Summary (see legend above):

```
- - - x - x -
- M2 M3 - - - -
- O2 O3 O4 O5 O6
- - - - - - -
- - - - - - -
x x
find -exec offers some of the same possibilities as GNU parallel.
find -exec only works on files. Processing other input (such as hosts or URLs) will require creating these inputs as files. find -exec has no support for running commands in parallel.
https://www.gnu.org/software/findutils/ (Last checked: 2019-01)
```

**DIFFERENCES BETWEEN make -j AND GNU Parallel**

Summary (see legend above):

```
- - - - - - -
- O1 O2 O3 - x O6
E1 - - - E5 -
- - - - - - -
- -
make -j can run jobs in parallel, but requires a crafted Makefile to do this. That results in extra quoting to get filenames containing newlines to work correctly.
make -j computes a dependency graph before running jobs. Jobs run by GNU parallel does not depend on each other.
(Very early versions of GNU parallel were coincidentally implemented using make -j).
https://www.gnu.org/software/make/ (Last checked: 2019-01)
```

**DIFFERENCES BETWEEN ppss AND GNU Parallel**

Summary (see legend above):

```
I1 I2 - - - I7
M1 - M3 - - M6
```
GNU Parallel alternatives

O1 - - x - -
E1 E2 ?E3 E4 - -
R1 R2 R3 R4 - - ?R7 ??
--

ppss is also a tool for running jobs in parallel.

The output of ppss is status information and thus not useful for using as input for another command. The output from the jobs are put into files.

The argument replace string ($ITEM) cannot be changed. Arguments must be quoted - thus arguments containing special characters (space "","&!") may cause problems. More than one argument is not supported. Filenames containing newlines are not processed correctly. When reading input from a file null cannot be used as a terminator. ppss needs to read the whole input file before starting any jobs.

Output and status information is stored in ppss_dir and thus requires cleanup when completed. If the dir is not removed before running ppss again it may cause nothing to happen as ppss thinks the task is already done. GNU parallel will normally not need cleaning up if running locally and will only need cleaning up if stopped abnormally and running remote (--cleanup may not complete if stopped abnormally). The example Parallel grep would require extra postprocessing if written using ppss.

For remote systems PPSS requires 3 steps: config, deploy, and start. GNU parallel only requires one step.

EXAMPLES FROM ppss MANUAL

Here are the examples from ppss’s manual page with the equivalent using GNU parallel:

1$ ./ppss.sh standalone -d /path/to/files -c 'gzip '

1$ find /path/to/files -type f | parallel gzip

2$ ./ppss.sh standalone -d /path/to/files -c 'cp "$ITEM" /destination/dir '

2$ find /path/to/files -type f | parallel cp {} /destination/dir

3$ ./ppss.sh standalone -f list-of-urls.txt -c 'wget -q ' 

3$ parallel -a list-of-urls.txt wget -q

4$ ./ppss.sh standalone -f list-of-urls.txt -c 'wget -q "$ITEM"'

4$ parallel -a list-of-urls.txt wget -q {} 

5$ ./ppss config -C config.cfg -c 'encode.sh ' -d /source/dir 

-m 192.168.1.100 -u ppss -k ppss-key.key -S ./encode.sh 

-n nodes.txt -o /some/output/dir --upload --download; 

./ppss deploy -C config.cfg 

./ppss start -C config

5$ # parallel does not use configs. If you want a different username put it in nodes.txt: user@hostname

find source/dir -type f | 

parallel --sshloginfile nodes.txt --trc {.}.mp3 lame -a {} -o

--
GNU Parallel alternatives

DIFFERENCES BETWEEN pexec AND GNU Parallel

Summary (see legend above):
I1 I2 - I4 I5 --
M1 - M3 - - M6
O1 O2 O3 - O5 O6
E1 -- E4 - E6 -
R1 ---- R6 ---
S1 -
pexec is also a tool for running jobs in parallel.

EXAMPLES FROM pexec MANUAL

Here are the examples from pexec's info page with the equivalent using GNU parallel:

1$ pexec -o sqrt-%s.dat -p "$(seq 10)" -e NUM -n 4 -c --
   'echo "scale=10000;sqrt($NUM)" | bc'

1$ seq 10 | parallel -j4 'echo "scale=10000;sqrt({})" | \
   bc > sqrt-{}.dat'

2$ pexec -p "$(ls myfiles*.ext)" -i %s -o %s.sort -- sort

2$ ls myfiles*.ext | parallel sort {} ">{}.sort"

3$ pexec -f image.list -n auto -e B -u star.log -c --
   'fistar $B.fits -f 100 -F id,x,y,flux -o $B.star'

3$ parallel -a image.list \
   'fistar {}.fits -f 100 -F id,x,y,flux -o {}.star' 2>star.log

https://github.com/louwrentius/PPSS
GNU Parallel alternatives

4$ pexec -r *.png -e IMG -c -o - -- \
   'convert $IMG $(IMG%.png).jpeg ; "echo $IMG: done"

4$ ls *.png | parallel 'convert {} {} .jpeg; echo {}: done'

5$ pexec -r *.png -i %s -o %s.jpeg -c 'pngtopnm | pnmtojpeg'

5$ ls *.png | parallel 'pngtopnm < {} | pnmtojpeg > {}.jpg'

6$ for p in *.png ; do echo ${p%.png} ; done | 
   pexec -f -i %s.png -o %s.jpg -c 'pngtopnm | pnmtojpeg'

6$ ls *.png | parallel 'pngtopnm < {} | pnmtojpeg > {}.jpg'

7$ LIST=$(for p in *.png; do echo ${p%.png}; done)
   pexec -r $LIST -i %s.png -o %s.jpg -c 'pngtopnm | pnmtojpeg'

7$ ls *.png | parallel 'pngtopnm < {} | pnmtojpeg > {}.jpg'

8$ pexec -n 8 -r *.jpg -y unix -e IMG -c \
   'pexec -j -m blockread -d $IMG | \ 
    jpegtopnm | pnmscale 0.5 | pnmtojpeg | \ 
    pexec -j -m blockwrite -s th_$IMG'

8$ # Combining GNU B<parallel> and GNU B<sem>.
   ls *jpg | parallel -j8 'sem --id blockread cat {} | jpegtopnm |' \
   'pnmscale 0.5 | pnmtojpeg | sem --id blockwrite cat > th_{}';

   # If reading and writing is done to the same disk, this may be 
   # faster as only one process will be either reading or writing:
   ls *jpg | parallel -j8 'sem --id diskio cat {} | jpegtopnm |' \
   'pnmscale 0.5 | pnmtojpeg | sem --id diskio cat > th_{}';

https://www.gnu.org/software/pexec/

DIFFERENCES BETWEEN xjobs AND GNU Parallel

xjobs is also a tool for running jobs in parallel. It only supports running jobs on your local computer.

xjobs deals badly with special characters just like xargs. See the section DIFFERENCES BETWEEN xargs AND GNU Parallel.

EXAMPLES FROM xjobs MANUAL

Here are the examples from xjobs's man page with the equivalent using GNU parallel:

1$ ls -l *.zip | xjobs unzip

1$ ls *.zip | parallel unzip

2$ ls -l *.zip | xjobs -n unzip

2$ ls *.zip | parallel unzip >/dev/null

3$ find . -name '*.bak' | xjobs gzip
GNU Parallel alternatives

3$ find . -name '*.bak' | parallel gzip

4$ ls -l *.jar | sed 's/\(.*\)/\1 \1.idx/' | xjobs jar tf

4$ ls *.jar | parallel jar tf () '()' {}.idx

5$ xjobs -s script

5$ cat script | parallel

6$ mkfifo /var/run/my_named_pipe; xjobs -s /var/run/my_named_pipe &
    echo unzip 1.zip >> /var/run/my_named_pipe;
    echo tar cf /backup/myhome.tar /home/me >> /var/run/my_named_pipe

6$ mkfifo /var/run/my_named_pipe;
    cat /var/run/my_named_pipe | parallel &
    echo unzip 1.zip >> /var/run/my_named_pipe;
    echo tar cf /backup/myhome.tar /home/me >> /var/run/my_named_pipe

https://www.maier-komor.de/xjobs.html (Last checked: 2019-01)

DIFFERENCES BETWEEN prll AND GNU Parallel

prll is also a tool for running jobs in parallel. It does not support running jobs on remote computers.

prll encourages using BASH aliases and BASH functions instead of scripts. GNU parallel supports scripts directly, functions if they are exported using export -f, and aliases if using env_parallel.

prll generates a lot of status information on stderr (standard error) which makes it harder to use the stderr (standard error) output of the job directly as input for another program.

EXAMPLES FROM prll's MANUAL

Here is the example from prll's man page with the equivalent using GNU parallel:

1$ prll -s 'mogrify -flip $1' *.jpg

1$ parallel mogrify -flip ::: *.jpg

https://github.com/exzombie/prll (Last checked: 2019-01)

DIFFERENCES BETWEEN dxargs AND GNU Parallel

dxargs is also a tool for running jobs in parallel.

dxargs does not deal well with more simultaneous jobs than SSHD's MaxStartups. dxargs is only built for remote run jobs, but does not support transferring of files.


DIFFERENCES BETWEEN mdm/middleman AND GNU Parallel

middleman(mdm) is also a tool for running jobs in parallel.

EXAMPLES FROM middleman's WEBSITE

Here are the shellscripts of https://web.archive.org/web/20110728064735/http://mdm.berlios.de/usage.html ported to GNU parallel:
GNU Parallel alternatives

1$ seq 19 | parallel buffon -o - | sort -n > result
cat files | parallel cmd
find dir -execdir sem cmd () \\

https://github.com/cklin/mdm (Last checked: 2019-01)

DIFFERENCES BETWEEN xapply AND GNU Parallel

xapply can run jobs in parallel on the local computer.

EXAMPLES FROM xapply's MANUAL

Here are the examples from xapply's man page with the equivalent using GNU parallel:

1$ xapply '(cd %1 && make all)' */
1$ parallel 'cd () && make all' ::: */

2$ xapply -f 'diff %1 ../version5/%1' manifest | more
2$ parallel diff () ../version5/{} < manifest | more

3$ xapply -p/dev/null -f 'diff %1 %2' manifest1 checklist1
3$ parallel --link diff {1} {2} ::: manifest1 checklist1

4$ xapply 'indent' * .c
4$ parallel indent ::: * .c

5$ find ~ksb/bin -type f ! -perm -111 -print | \
xapply -f -v 'chmod a+x' -
5$ find ~ksb/bin -type f ! -perm -111 -print | \
parallel -v chmod a+x

6$ find */ -... | fmt 960 1024 | xapply -f -i /dev/tty 'vi' -
6$ sh <(find */ -... | parallel -s 1024 echo vi)

6$ find */ -... | parallel -s 1024 -Xuj1 vi

7$ find ... | xapply -f -5 -i /dev/tty 'vi' - - - - -
7$ sh <(find ... | parallel -n5 echo vi)

7$ find ... | parallel -n5 -uj1 vi

8$ xapply -fn "" /etc/passwd
8$ parallel -k echo < /etc/passwd

9$ tr ': ' '\012' < /etc/passwd | \\
DIFFERENCES BETWEEN AIX apply AND GNU Parallel

**apply** can build command lines based on a template and arguments - very much like GNU **parallel**. **apply** does not run jobs in parallel. **apply** does not use an argument separator (like :::); instead the template must be the first argument.

**EXAMPLES FROM IBM's KNOWLEDGE CENTER**

Here are the examples from IBM's Knowledge Center and the corresponding command using GNU **parallel**:

To obtain results similar to those of the ls command, enter:

```
1$ apply echo *
1$ parallel echo ::: *
```

To compare the file named a1 to the file named b1, and the file named a2 to the file named b2, enter:

```
2$ apply -2 cmp a1 b1 a2 b2
2$ parallel -N2 cmp ::: a1 b1 a2 b2
```

To run the who command five times, enter:

```
3$ apply -0 who 1 2 3 4 5
3$ parallel -N0 who ::: 1 2 3 4 5
```

To link all files in the current directory to the directory /usr/joe, enter:

```
4$ apply 'ln %1 /usr/joe' *
4$ parallel ln {} /usr/joe ::: *
```


**DIFFERENCES BETWEEN paexec AND GNU Parallel**

**paexec** can run jobs in parallel on both the local and remote computers.

**paexec** requires commands to print a blank line as the last output. This means you will have to write a wrapper for most programs.

**paexec** has a job dependency facility so a job can depend on another job to be executed successfully. Sort of a poor-man's **make**.
EXAMPLES FROM paexec’s EXAMPLE CATALOG

Here are the examples from paexec’s example catalog with the equivalent using GNU parallel:

1_div_X_run

```
1$ ../../paexec -s -l -c "`pwd`/1_div_X_cmd" -n +1 <<EOF [...] 
1$ parallel echo {} '|' `pwd`/1_div_X_cmd <<EOF [...] 
```

all_substr_run

```
2$ ../../paexec -lp -c "`pwd`/all_substr_cmd" -n +3 <<EOF [...] 
2$ parallel echo {} '|' `pwd`/all_substr_cmd <<EOF [...] 
```

cc_wrapper_run

```
3$ ../../paexec -c "env CC=gcc CFLAGS=-O2 `pwd`/cc_wrapper_cmd" 
   -n 'host1 host2' 
   -t '/usr/bin/ssh -x' <<EOF [...] 
3$ parallel echo {} '|' "env CC=gcc CFLAGS=-O2 `pwd`/cc_wrapper_cmd" 
   -S host1,host2 <<EOF [...] 
   # This is not exactly the same, but avoids the wrapper 
   parallel gcc -O2 -c -o {}.o {} 
   -S host1,host2 <<EOF [...] 
```

toupper_run

```
4$ ../../paexec -lp -c "`pwd`/toupper_cmd" -n +10 <<EOF [...] 
4$ parallel echo {} '|' `.toupper_cmd <<EOF [...] 
   # Without the wrapper: 
   parallel echo {} '|' awk (print toupper("$0"))' <<EOF [...] 
```

https://github.com/cheusov/paexec

DIFFERENCES BETWEEN map(sitaramc) AND GNU Parallel

Summary (see legend above):

I1 - I4 - (I7)  
M1 (M2) M3 (M4) M5 M6 
- O2 O3 - O5 - N/A N/A O10 
E1 ------ 
-------- 
--

(I7): Only under special circumstances. See below.  
(M2+M4): Only if there is a single replacement string.  
map rejects input with special characters: 

```
echo "The Cure" > My\ brother\'s\ 12\"\ records 
```
ls | map 'echo %; wc %'

It works with GNU parallel:

ls | parallel 'echo {}; wc {}'

Under some circumstances it also works with map:

ls | map 'echo % works %'

But tiny changes make it reject the input with special characters:

ls | map 'echo % does not work "%"'

This means that many UTF-8 characters will be rejected. This is by design. From the web page: "As such, programs that quietly handle them, with no warnings at all, are doing their users a disservice."

map delays each job by 0.01 s. This can be emulated by using parallel --delay 0.01.

map prints '+' on stderr when a job starts, and '-' when a job finishes. This cannot be disabled. parallel has --bar if you need to see progress.

map's replacement strings (% %D %B %E) can be simulated in GNU parallel by putting this in ~/.parallel/config:

--rpl '%'
--rpl '%D $=_Q::{:dirname($_)};'
--rpl '%B s:.*/::*;s:./[^/./]+$::*;'
--rpl '%E s::*\::'

map does not have an argument separator on the command line, but uses the first argument as command. This makes quoting harder which again may affect readability. Compare:

map -p 2 'perl -ne '/^\S+\s+\S+$/' and print $ARGV,"\n"' *

parallel -q perl -ne '/\S+\s+\S+$/' and print $ARGV,"\n" :: *

map can do multiple arguments with context replace, but not without context replace:

parallel --xargs echo 'BEGIN{"'}END' :: 1 2 3

map "echo 'BEGIN{"'}END'" 1 2 3

map has no support for grouping. So this gives the wrong results:

parallel perl -e '\$a="1\"x10000000;\print \$a,\"\n\"'} '*() \ :: a b c d e f
ls -l a b c d e f
parallel -kP4 -n1 grep 1 :: a b c d e f > out.par
map -n1 -p 4 'grep 1' a b c d e f > out.map-unbuf
map -n1 -p 4 'grep --line-buffered 1' a b c d e f > out.map-linebuf
map -n1 -p 1 'grep --line-buffered 1' a b c d e f > out.map-serial
ls -l out* md5sum out*
EXAMPLES FROM map's WEBSITE

Here are the examples from map's web page with the equivalent using GNU parallel:

```bash
1$ ls *.gif | map convert % %B.png        # default max-args: 1
1$ ls *.gif | parallel convert {} {} .png

2$ map "mkdir %B; tar -C %B -xf %" *.tgz  # default max-args: 1
2$ parallel 'mkdir {}; tar -C {} -xf {}' ::: *.tgz

3$ ls *.gif | map cp % /tmp             # default max-args: 100
3$ ls *.gif | parallel -X cp {} /tmp

4$ ls *.tar | map -n 1 tar -xf %
4$ ls *.tar | parallel tar -xf

5$ map "cp % /tmp" *.tgz
5$ parallel cp {} /tmp ::: *.tgz

6$ map "du -sm /home/%/mail" alice bob carol
6$ parallel "du -sm /home/{}/mail" ::: alice bob carol
or if you prefer running a single job with multiple args:
6$ parallel -Xj1 "du -sm /home/{}/mail" ::: alice bob carol

7$ cat /etc/passwd | map -d: 'echo user %1 has shell %?'
7$ cat /etc/passwd | parallel --colsep: 'echo user {1} has shell {7}''

8$ export MAP_MAX_PROCS=$(( `nproc` / 2 ))
8$ export PARALLEL=-j50%
```

https://github.com/sitaramc/map (Last checked: 2020-05)

DIFFERENCES BETWEEN ladon AND GNU Parallel

ladon can run multiple jobs on files in parallel.

ladon only works on files and the only way to specify files is using a quoted glob string (such as \\*.jpg). It is not possible to list the files manually.

As replacement strings it uses FULLPATH DIRNAME BASENAME EXT RELDIR RELPATH

These can be simulated using GNU parallel by putting this in ~/.parallel/config:

```bash
--rpl 'FULLPATH $_=Q($_);chomp($_=qx{readlink -f $_}})''
--rpl 'DIRNAME $_=Q{:dirname($_);}chomp($_=qx{readlink -f $_}})''
--rpl 'BASENAME s::*;s:\.[^/]+$::'  
--rpl 'EXT s::*\.'
```
GNU Parallel alternatives

ladon deals badly with filenames containing " and newline, and it fails for output larger than 200k:

ladon '/*' -- seq 36000 | wc

EXAMPLES FROM ladon MANUAL

It is assumed that the '--rpl's above are put in ~/.parallel/config and that it is run under a shell that supports "*" globbing (such as zsh):

1$ ladon "**/*.txt" -- echo RELPATH
   1$ parallel echo RELPATH ::: **/*.txt

2$ ladon "~/Documents/**/*.pdf" -- shasum FULLPATH >hashes.txt
   2$ parallel shasum FULLPATH ::: ~/Documents/**/*.pdf >hashes.txt

3$ ladon -m thumbs/RELDIR "**/*.jpg" -- convert FULLPATH \        -thumbnail 100x100^ -gravity center -extent 100x100 \ thumbs/RELPATH
   3$ parallel mkdir -p thumbs/RELDIR\; convert FULLPATH \        -thumbnail 100x100^ -gravity center -extent 100x100 \ thumbs/RELPATH ::: **/*.jpg

4$ ladon "~/Music/*.wav" -- lame -V 2 FULLPATH DIRNAME/BASENAME.mp3
   4$ parallel lame -V 2 FULLPATH DIRNAME/BASENAME.mp3 ::: ~/Music/*.wav

https://github.com/danielgtaylor/ladon (Last checked: 2019-01)

DIFFERENCES BETWEEN jobflow AND GNU Parallel

Summary (see legend above):

I1 - - - - I7
   -- M3 - -(M6)
O1 O2 O3 - O5 O6 (O7) - - O10
E1 - - - E6 -
-----------

jobflow can run multiple jobs in parallel.

Just like xargs output from jobflow jobs running in parallel mix together by default. jobflow can buffer into files with -buffered (placed in /run/shm), but these are not cleaned up if jobflow dies unexpectedly (e.g. by Ctrl-C). If the total output is big (in the order of RAM+swap) it can cause the system to slow to a crawl and eventually run out of memory.

Just like xargs redirection and composed commands require wrapping with bash -c.
Input lines can at most be 4096 bytes.

jobflow is faster than GNU parallel but around 6 times slower than parallel-bash.

jobflow has no equivalent for --pipe, or --sshlogin.

jobflow makes it possible to set resource limits on the running jobs. This can be emulated by GNU parallel using bash's ulimit:

```
jobflow -limits=mem=100M,cpu=3,fsize=20M,nofiles=300 myjob
```

```
parallel 'ulimit -v 102400 -t 3 -f 204800 -n 300 myjob'
```

EXAMPLES FROM jobflow README

```
1$ cat things.list | jobflow -threads=8 -exec ./mytask {}
```

```
1$ cat things.list | parallel -j8 ./mytask {}
```

```
2$ seq 100 | jobflow -threads=100 -exec echo {}
```

```
2$ seq 100 | parallel -j100 echo {}
```

```
3$ cat urls.txt | jobflow -threads=32 -exec wget {}
```

```
3$ cat urls.txt | parallel -j32 wget {}
```

```
4$ find . -name '*.bmp' | \ 
   jobflow -threads=8 -exec bmp2jpeg {.}.bmp {.}.jpg
```

```
4$ find . -name '*.bmp' | \ 
   parallel -j8 bmp2jpeg {.}.bmp {.}.jpg
```

```
5$ seq 100 | jobflow -skip 10 -count 10
```

```
5$ seq 100 | parallel --filter '{1} > 10 and {1} <= 20' echo
```

```
5$ seq 100 | parallel echo '{= $_[>10 and $_[<=20 or skip() =]}'
```

https://github.com/rofl0r/jobflow (Last checked: 2022-05)

DIFFERENCES BETWEEN gargs AND GNU Parallel

gargs can run multiple jobs in parallel.

Older versions cache output in memory. This causes it to be extremely slow when the output is larger than the physical RAM, and can cause the system to run out of memory.

See more details on this in man parallel_design.

Newer versions cache output in files, but leave files in $TMPDIR if it is killed.

Output to stderr (standard error) is changed if the command fails.

EXAMPLES FROM gargs WEBSITE

```
1$ seq 12 -1 1 | gargs -p 4 -n 3 "sleep (0); echo (1) (2)"
```
1$ seq 12 -1 1 | parallel -P 4 -n 3 "sleep (1); echo (2) (3)"

2$ cat t.txt | gargs --sep "\s+" -p 2 "echo '{0}:{1}-{2}' full-line: '{3}'"

2$ cat t.txt | parallel --colsep "\s+" -P 2 "echo '{1}:{2}-{3}' full-line: '{0}'"

https://github.com/brentp/gargs

DIFFERENCES BETWEEN orgalorg AND GNU Parallel

orgalorg can run the same job on multiple machines. This is related to --onall and --nonall.

orgalorg supports entering the SSH password - provided it is the same for all servers. GNU parallel advocates using ssh-agent instead, but it is possible to emulate orgalorg's behavior by setting SSHPASS and by using --ssh "sshpass ssh".

To make the emulation easier, make a simple alias:

```
alias par_emul="parallel -j0 --ssh 'sshpass ssh' --nonall --tag --lb"
```

If you want to supply a password run:

```
SSHPASS=`ssh-askpass`
```

or set the password directly:

```
SSHPASS=P4$$w0rd!
```

If the above is set up you can then do:

```
orgalorg -o frontend1 -o frontend2 -p -C uptime
par_emul -S frontend1 -S frontend2 uptime

orgalorg -o frontend1 -o frontend2 -p -C top --bid 1
par_emul -S frontend1 -S frontend2 top --bid 1

orgalorg -o frontend1 -o frontend2 -p -er /tmp -n 
/md5sum /tmp/bigfile" -S bigfile
par_emul -S frontend1 -S frontend2 --basefile bigfile
--workdir /tmp md5sum /tmp/bigfile
```

orgalorg has a progress indicator for the transferring of a file. GNU parallel does not.

https://github.com/reconquest/orgalorg

DIFFERENCES BETWEEN Rust parallel AND GNU Parallel

Rust parallel focuses on speed. It is almost as fast as xargs, but not as fast as parallel-bash. It implements a few features from GNU parallel, but lacks many functions. All these fail:

```
# Read arguments from file
parallel -a file echo
# Changing the delimiter
parallel -d _ echo :: a_b_c_
```

These do something different from GNU parallel
DIFFERENCES BETWEEN Rush AND GNU Parallel

rush (https://github.com/shenwei356/rush) is written in Go and based on gargs.

Just like GNU parallel rush buffers in temporary files. But opposite GNU parallel rush does not clean up, if the process dies abnormally.

 rush has some string manipulations that can be emulated by putting this into ~/.parallel/config (/ is used instead of %, and $ is used instead of ^ as that is closer to bash's $(var%postfix):

```
--rpl '(): s:\(.\[^/]+\)'*$$:'
--rpl '(:%([^]+)?) s:\$1\(.\[^/]+\)'*$$:'
--rpl './%(\[^/]+\)\$)' s:\.*\$1\(.\[^/]+\)'*$$:'
--rpl '(/:\$1\(\[^/]+\)\(\[^/]+\)\$)' s:\.*\$1\(\[^/]+\)'*$$:'
--rpl '(@(.*)) /\$1/ and $=_$1;'
EXAMPLES FROM rush’s WEBSITE

Here are the examples from rush’s website with the equivalent command in GNU parallel.

1. Simple run, quoting is not necessary
   1$ seq 1 3 | rush echo {}
   1$ seq 1 3 | parallel echo {}

2. Read data from file (`-i`)
   2$ rush echo {} -i data1.txt -i data2.txt
   2$ cat data1.txt data2.txt | parallel echo {}

3. Keep output order (`-k`)
   3$ seq 1 3 | rush 'echo {}' -k
   3$ seq 1 3 | parallel -k echo {}

4. Timeout (`-t`)
   4$ time seq 1 | rush 'sleep 2; echo {}' -t 1
   4$ time seq 1 | parallel --timeout 1 'sleep 2; echo {}'

5. Retry (`-r`)
   5$ seq 1 | rush 'python unexisted_script.py' -r 1
   5$ seq 1 | parallel --retries 2 'python unexisted_script.py'

   Use -u to see it is really run twice:
   5$ seq 1 | parallel -u --retries 2 'python unexisted_script.py'

6. Dirname (`{/}`) and basename (`{%}`) and remove custom suffix (`{^suffix}`)
   6$ echo dir/file_1.txt.gz | rush 'echo // {%} {^_1.txt.gz}'
   6$ echo dir/file_1.txt.gz | parallel --plus echo // // {%} {^_1.txt.gz}

7. Get basename, and remove last (`.{}`) or any (`{[:]}`) extension
   7$ echo dir.d/file.txt.gz | rush 'echo . {>({:.})
   7$ echo dir.d/file.txt.gz | parallel 'echo . {>({:.})

8. Job ID, combine fields index and other replacement strings
   8$ echo 12 file.txt dir/s_1.fq.gz | rush 'echo job (#): {2} {2.} {3:^_1}'}
8$ echo 12 file.txt dir/s_1.fq.gz | parallel --colsep ' ' 'echo job #: {2} (2:) (3:/{_1})'

9. Capture submatch using regular expression (`[@regexp]`)  
9$ echo read_1.fq.gz | rush 'echo {@(\.+)_\d}'
9$ echo read_1.fq.gz | parallel 'echo {@(\.+)_\d}'

10. Custom field delimiter (`-d`)  
10$ echo a=b=c | rush 'echo {1} {2} {3}' -d =
10$ echo a=b=c | parallel -d = echo {1} {2} {3}

11. Send multi-lines to every command (`-n`)  
11$ seq 5 | rush -n 2 -k 'echo {}'; echo'
11$ seq 5 | parallel -n 2 -k 
    'echo {=-1 $_={join}\n',@arg{1..$#arg} =}; echo'
11$ seq 5 | rush -n 2 -k 'echo {}'; echo' -J '
11$ seq 5 | parallel -n 2 -k 'echo {}'; echo'

12. Custom record delimiter (`-D`), note that empty records are not used.  
12$ echo a b c d | rush -D " " -k 'echo {}'
12$ echo a b c d | parallel -d " " -k 'echo {}'
12$ echo abcd | rush -D " " -k 'echo {}'

Cannot be done by GNU Parallel

12$ cat fasta.fa
>seq1
tag
>seq2
cat
gat
>seq3
attac
a
cat

12$ cat fasta.fa | rush -D "">"
    'echo FASTA record #: name: {1} sequence: {2}' -k -d "\n"
# rush fails to join the multiline sequences

12$ cat fasta.fa | (read -n1 ignore_first_char;
parallel -d '-> --colsep \n' echo FASTA record {#: \
    name: {1} sequence: '{=2 $_=join'',@arg[2..$#arg]=}'}
)

13. Assign value to variable, like `awk -v` (`-v`)

13$ seq 1 | rush 'echo Hello, {fname} {lname}!' -v 
    'fname=Wei -v lname=Shen'

13$ seq 1 | 
    parallel -N0 \
    'fname=Wei; lname=Shen; echo Hello, ${fname} ${lname}!'

13$ for var in a b; do \
13  seq 1 3 | rush -k -v \
    'echo var: $var, data: {}'; \
13$ done

In GNU parallel you would typically do:

13$ seq 1 3 | parallel -k echo var: (1), data: (2) :: a b :::

If you really want the var:

13$ seq 1 3 | parallel -k var=${1} 'echo var: $var, data: {}' :: a b :::

If you really want the for-loop:

13$ for var in a b; do \n    export var; \n    seq 1 3 | parallel -k 'echo var: $var, data: {}'; \n    done

Contrary to rush this also works if the value is complex like:

My brother's 12" records

14. Preset variable (-v), avoid repeatedly writing verbose replacement strings

14$ # naive way
    echo read_1.fq.gz | rush 'echo :^_1 {^_1}_2.fq.gz'

14$ echo read_1.fq.gz | parallel 'echo :%_1 {_%1}_2.fq.gz'

14$ # macro + removing suffix
    echo read_1.fq.gz | rush -v p='{:^_1}' 'echo {p} {p}_2.fq.gz'

14$ echo read_1.fq.gz | parallel 'p=:{_%1}; echo $p ${p}_2.fq.gz'

14$ # macro + regular expression
    echo read_1.fq.gz | rush -v p='[@(\.+?)]d' 'echo {p} {p}_2.fq.gz'

14$ echo read_1.fq.gz | parallel 'p=@(\.+?)_d'; echo $p ${p}_2.fq.gz'
Contrary to rush GNU parallel works with complex values:

```
14$ echo "My brother's 12"read_1.fq.gz" |
    parallel 'p=\(@(.+?)_d\); echo $p ${p}_2.fq.gz'
```

15. Interrupt jobs by `Ctrl-C`, rush will stop unfinished commands and exit.

```
15$ seq 1 20 | rush 'sleep 1; echo {}'
    ^C
15$ seq 1 20 | parallel 'sleep 1; echo {}'
    ^C
```

16. Continue/resume jobs (`-c`). When some jobs failed (by execution failure, timeout, or canceling by user with `Ctrl + C`), please switch flag `-c/--continue` on and run again, so that `rush` can save successful commands and ignore them in NEXT run.

```
16$ seq 1 3 | rush 'sleep {}; echo {}' -t 3 -c
    cat successful_cmds.rush
16$ seq 1 3 | rush 'sleep {}; echo {}' -t 3 -c
```

Multi-line jobs:

```
16$ seq 1 3 | rush 'sleep {}; echo {}';
    echo finish {}' -t 3 -c finished.rush
    cat finished.rush
16$ seq 1 3 | rush 'sleep {}; echo {}';
    echo finish {}' -t 3 -c finished.rush
```

```
16$ seq 1 3 |
    parallel --joblog mylog --timeout 2 'sleep {}; echo {}';
    echo finish {}' -t 3 -c finished.rush
16$ seq 1 3 |
    parallel --joblog mylog --retry-failed 'sleep {}; echo {}';
    echo finish {}'
```

17. A comprehensive example: downloading 1K+ pages given by three URL list files using `phantomjs save_page.js` (some page contents are dynamically generated by Javascript, so `wget` does not work). Here I set max jobs number (`-j`) as `20`, each job has a max running time (`-t`) of `60` seconds and `3` retry changes (`-r`). Continue flag `-c` is also switched on, so we can continue unfinished jobs. Luckily, it’s accomplished in one run :) 

```
17$ for f in $(seq 2014 2016); do 
    /bin/rm -rf $f; mkdir -p $f; 
    cat $f.html.txt | rush -v d=$f -d=
        'phantomjs save_page.js "{}" > {d}/{3}.html' 
        -j 20 -t 60 -r 3 -c; 
done
```
GNU parallel can append to an existing joblog with `+':

```
17$ rm mylog
    for f in $(seq 2014 2016); do
        /bin/rm -rf $f; mkdir -p $f;
        cat $f.html.txt | parallel -j20 --timeout 60 --retries 4 --joblog +mylog \
               --colsep = \phantomjs save_page.js {1}={2}={3} '>' $f/{3}.html
    done
```

18. A bioinformatics example: mapping with `bwa`, and processing result with `samtools`:

```
18$ ref=ref/xxx.fa
threads=25
    ls -d raw.cluster.clean.mapping/* \ 
        | rush -v ref=$ref -v j=$threads -v p='/{}/{}' \ 
        'bwa mem -t {j} -M -a {ref} {p}_1.fq.gz {p}_2.fq.gz >{p}.sam;\ 
        samtools view -bS {p}.sam > {p}.bam; \ 
        samtools sort -T {p}.tmp -@ {j} {p}.bam -o {p}.sorted.bam; \ 
        samtools index {p}.sorted.bam; \ 
        samtools flagstat {p}.sorted.bam > {p}.sorted.bam.flagstat; \ 
        /bin/rm {p}.bam {p}.sam;' \ 
        -j 2 --verbose -c -C mapping.rush
```

GNU parallel would use a function:

```
18$ ref=ref/xxx.fa
    export ref
    thr=25
    export thr
    bwa_sam() { 
        p="$1"
        bam="$p".bam
        sam="$p".sam
        sortbam="$p".sorted.bam
        bwa mem -t $thr -M -a $ref ${p}_1.fq.gz ${p}_2.fq.gz > "$sam"
        samtools view -bS "$sam" > "$bam"
        samtools sort -T ${p}.tmp -@ $thr "$bam" -o "$sortbam"
        samtools index "$sortbam"
        samtools flagstat "$sortbam" > "$sortbam.flagstat"
        /bin/rm ${p}bam ${p}sam;'
    }
    export -f bwa_sam
    ls -d raw.cluster.clean.mapping/* | parallel -j 2 --verbose --joblog mylog bwa_sam
```

Other rush features

** rush has:**

* awk -v like custom defined variables (-v)

With GNU parallel you would simply set a shell variable:

```
parallel 'v={}; echo "$v"' :::: foo 
echo foo | rush -v v={} 'echo (v)'
```

Also rush does not like special chars. So these do not work:
GNU Parallel alternatives

echo does not work | rush -v v="'echo (v)'"
echo "My brother's 12" records | rush -v v=() 'echo (v)'

Whereas the corresponding GNU parallel version works:
parallel 'v=\"'; echo "$\v\"' ::: works
parallel 'v={}; echo "$\v\"' ::: "My brother's 12" records"

* Exit on first error(s) (-e)
  This is called --halt now,fail=1 (or shorter: --halt 2) when used with GNU parallel.

* Settable records sending to every command (-n, default 1)
  This is also called -n in GNU parallel.

* Practical replacement strings
  {\} remove any extension
  With GNU parallel this can be emulated by:
  parallel --plus echo '{/\.*}' ::: foo.ext.bar.gz

  {^suffix}, remove suffix
  With GNU parallel this can be emulated by:
  parallel --plus echo '{.%} bar.gz' ::: foo.ext.bar.gz

  {@regexp}, capture submatch using regular expression
  With GNU parallel this can be emulated by:
  parallel --rpl '{@(.*).gz}' /$$1/ and $$2=' \n  echo '{@\d_.).*}lfoo.gz

  {%}, {%:}, basename without extension
  With GNU parallel this can be emulated by:
  parallel echo '{= s:.*/;s/\..*/=}' ::: dir/foo.bar.gz

  And if you need it often, you define a --rpl in $HOME/.parallel/config:
  --rpl '%{%} s:.*/;' s/\..*/'
  --rpl '%{^} s:.*/;' s/\..*/'

  Then you can use them as:
  parallel echo {%} {%:} ::: dir/foo.bar.gz

* Preset variable (macro)
  E.g.
  echo foosuffix | rush -v p={^suffix} 'echo {p}_new_suffix'

  With GNU parallel this can be emulated by:
  echo foosuffix |
  parallel --plus 'p={%suffix}; echo ${p}_new_suffix'

  Opposite rush GNU parallel works fine if the input contains double space, ',' and ':
  echo "1'6'\" foosuffix" |
  parallel --plus 'p={%suffix}; echo "${p}"_new_suffix'
* Commands of multi-lines

While you can use multi-lined commands in GNU parallel, to improve readability GNU parallel discourages the use of multi-line commands. In most cases it can be written as a function:

```
seq 1 3 |
  parallel --timeout 2 --joblog my.log 'sleep {}; echo {}; 
   echo finish {}'
```

Could be written as:

```
doit() {
  sleep "$1"
  echo "$1"
  echo finish "$1"
}
export -f doit
seq 1 3 | parallel --timeout 2 --joblog my.log doit
```

The failed commands can be resumed with:

```
seq 1 3 |
  parallel --resume-failed --joblog my.log 'sleep {}; echo {}; 
   echo finish {}'
```

https://github.com/shenwei356/rush

DIFFERENCES BETWEEN ClusterSSH AND GNU Parallel

ClusterSSH solves a different problem than GNU parallel.

ClusterSSH opens a terminal window for each computer and using a master window you can run the same command on all the computers. This is typically used for administrating several computers that are almost identical.

GNU parallel runs the same (or different) commands with different arguments in parallel possibly using remote computers to help computing. If more than one computer is listed in -S GNU parallel may only use one of these (e.g. if there are 8 jobs to be run and one computer has 8 cores).

GNU parallel can be used as a poor-man’s version of ClusterSSH:

```
parallel --nonall -S server-a,server-b do_stuff foo bar
```

https://github.com/duncs/clusterssh

DIFFERENCES BETWEEN coshell AND GNU Parallel

coshell only accepts full commands on standard input. Any quoting needs to be done by the user.

Commands are run in sh so any bash/tcsh/zsh specific syntax will not work.

Output can be buffered by using -d. Output is buffered in memory, so big output can cause swapping and therefore be terrible slow or even cause out of memory.

https://github.com/gdm85/coshell (Last checked: 2019-01)

DIFFERENCES BETWEEN spread AND GNU Parallel

spread runs commands on all directories.

It can be emulated with GNU parallel using this Bash function:

```
spread() {
  _cmds() {
    perl -e '$"=" "$" "$;print @ARGV" \"cd \" "$'
  }
```

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This works except for the --exclude option.

(Last checked: 2017-11)

**DIFFERENCES BETWEEN pyargs AND GNU Parallel**

*pyargs* deals badly with input containing spaces. It buffers stdout, but not stderr. It buffers in RAM. {} does not work as replacement string. It does not support running functions.

*pyargs* does not support composed commands if run with --lines, and fails on pyargs traceroute
gnu.org fsf.org.

**Examples**

```bash
seq 5 | pyargs -P50 -L seq
seq 5 | parallel -P50 --lb seq

seq 5 | pyargs -P50 --mark -L seq
seq 5 | parallel -P50 --lb - --tagstring OUTPUT'[(= $_=$job->replaced())=]" seq
# Similar, but not precisely the same
seq 5 | parallel -P50 --lb --tag seq

seq 5 | pyargs -P50 --mark command
# Somewhat longer with GNU Parallel due to the special
# --mark formatting
cmd="$(echo "command" | parallel --shellquote)"

wrap_cmd()
{
    echo "MARK $cmd $@="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="="=

https://github.com/robertblackwell/pyargs (Last checked: 2019-01)

**DIFFERENCES BETWEEN concurrently AND GNU Parallel**

*concurrently* runs jobs in parallel.

The output is prepended with the job number, and may be incomplete:

```
$ concurrently 'seq 100000' | (sleep 3; wc -l)
7165
```

When pretty printing it caches output in memory. Output mixes by using test MIX below whether or
not output is cached.

There seems to be no way of making a template command and have concurrently fill that with different args. The full commands must be given on the command line.

There is also no way of controlling how many jobs should be run in parallel at a time - i.e. "number of jobslots". Instead all jobs are simply started in parallel.

https://github.com/kimmobrunfeldt/concurrently (Last checked: 2019-01)

DIFFERENCES BETWEEN map(soveran) AND GNU Parallel

map does not run jobs in parallel by default. The README suggests using:

```
... | map t 'sleep $t && say done &'
```

But this fails if more jobs are run in parallel than the number of available processes. Since there is no support for parallelization in map itself, the output also mixes:

```
seq 10 | map i 'echo start-$i && sleep 0.$i && echo end-$i &'
```

The major difference is that GNU parallel is built for parallelization and map is not. So GNU parallel has lots of ways of dealing with the issues that parallelization raises:

- Keep the number of processes manageable
- Make sure output does not mix
- Make Ctrl-C kill all running processes

EXAMPLES FROM maps WEBSITE

Here are the 5 examples converted to GNU Parallel:

```
1$ ls *.c | map f 'foo $f'
1$ ls *.c | parallel foo

2$ ls *.c | map f 'foo $f; bar $f'
2$ ls *.c | parallel 'foo {}; bar {}'

3$ cat urls | map u 'curl -O $u'
3$ cat urls | parallel curl -O

4$ printf "1\n1\n1\n" | map t 'sleep $t && say done'
4$ printf "1\n1\n1\n" | parallel 'sleep () && say done'
4$ parallel 'sleep {} && say done' ::: 1 1 1

5$ printf "1\n1\n1\n" | map t 'sleep $t && say done &'
5$ printf "1\n1\n1\n" | parallel -j0 'sleep {} && say done'
5$ parallel -j0 'sleep {} && say done' ::: 1 1 1
```

https://github.com/soveran/map (Last checked: 2019-01)

DIFFERENCES BETWEEN loop AND GNU Parallel

loop mixes stdout and stderr:

```
loop 'ls /no-such-file' >/dev/null
```

loop's replacement string $ITEM does not quote strings:
echo 'two spaces' | loop 'echo $ITEM'

```
loop cannot run functions:

myfunc() { echo joe; }
export -f myfunc
loop 'myfunc this fails'
```

**EXAMPLES FROM loop's WEBSITE**

Some of the examples from https://github.com/Miserlou/Loop/ can be emulated with GNU parallel:

```
# A couple of functions will make the code easier to read
$ loopy() {
    yes | parallel -uN0 -j1 "$@"
}
$ export -f loopy
$ time_out() {
    parallel -uN0 -q --timeout "$@" :: 1
}
$ match() {
    perl -0777 -ne 'grep /"$1"/,$_ and print or exit 1'
}
$ export -f match

$ loop 'ls' --every 10s
$ loopy --delay 10s ls

$ loop 'touch $COUNT.txt' --count-by 5
$ loopy touch '($= $=seq()*5 =)'.txt

$ loop --until-contains 200 --\n    ./get_response_code.sh --site mysite.biz
$ loopy --halt now,success=1 \n    './get_response_code.sh --site mysite.biz | match 200'

$ loop './poke_server' --for-duration 8h
$ time_out 8h loopy ./poke_server

$ loop './poke_server' --until-success
$ loopy --halt now,success=1 ./poke_server

$ cat files_to_create.txt | loop 'touch $ITEM'
$ cat files_to_create.txt | parallel touch {}

$ loop 'ls' --for-duration 10min --summary
# --joblog is somewhat more verbose than --summary
$ time_out 10m loopy --joblog my.log ./poke_server; cat my.log

$ loop 'echo hello'
$ loopy echo hello

$ loop 'echo $COUNT'
# GNU Parallel counts from 1
$ loopy echo (#)
```
# Counting from 0 can be forced
$ loopy echo '($_.=seq()-1 =)' 

$ loop 'echo $COUNT' --count-by 2
$ loopy echo '($._=2*(seq()-1) =)'

$ loop 'echo $COUNT' --count-by 2 --offset 10
$ loopy echo '($._=10+2*(seq()-1) =)'

$ loop 'echo $COUNT' --count-by 1.1
# GNU Parallel rounds 3.3000000000000003 to 3.3
$ loopy echo '($._=1.1*(seq()-1) =)'

$ loop 'echo $COUNT $ACTUALCOUNT' --count-by 2
$ loopy echo '($._=2*(seq()-1) =) (#)'

$ loop 'echo $COUNT' --num 3 --summary
# --joblog is somewhat more verbose than --summary
$ seq 3 | parallel --joblog my.log echo; cat my.log

$ loop 'ls -foobartatz' --num 3 --summary
# --joblog is somewhat more verbose than --summary
$ seq 3 | parallel --joblog my.log -N0 ls -foobartatz; cat my.log

$ loop 'echo $COUNT' --count-by 2 --num 50 --only-last
# Can be emulated by running 2 jobs
$ seq 49 | parallel echo '($._=2*(seq()-1) =)' >/dev/null
$ echo 50 | parallel echo '($._=2*(seq()-1) =)'

$ loop 'date' --every 5s
$ loopy --delay 5s date

$ loop 'date' --for-duration 8s --every 2s
$ time_out 8s loopy --delay 2s date

$ loop 'date -u' --until-time '2018-05-25 20:50:00' --every 5s
$ seconds=$(`date -d 2019-05-25T20:50:00 +%s` - `date +%s`)
$ time_out $seconds loopy --delay 5s date -u

$ loop 'echo $RANDOM' --until-contains "666"
$ loopy --halt now,success=1 'echo $RANDOM | match 666'

$ loop 'if (( RANDOM % 2 )); then
  (echo "TRUE"; true);
else
  (echo "FALSE"; false);
fi' --until-success
$ loopy --halt now,success=1 'if (( $RANDOM % 2 )); then
  (echo "TRUE"; true);
else
  (echo "FALSE"; false);
fi'
GNU Parallel alternatives

$ loop 'if (( RANDOM % 2 )); then
  (echo "TRUE"); true);
else
  (echo "FALSE"); false);
fi' --until-error
$ loopy --halt now,fail=1 'if (( $RANDOM % 2 )); then
  (echo "TRUE"); true);
else
  (echo "FALSE"); false);
fi'

$ loop 'date' --until-match "(\d{4})"
$ loopy --halt now,success=1 'date | match [0-9][0-9][0-9][0-9]'

$ loop 'echo $ITEM' --for red,green,blue
$ parallel echo :: red green blue

$ cat /tmp/my-list-of-files-to-create.txt | loop 'touch $ITEM'
$ cat /tmp/my-list-of-files-to-create.txt | parallel touch

$ ls | loop 'cp $ITEM $ITEM.bak'; ls
$ ls | parallel cp {} {} .bak; ls

$ loop 'echo $ITEM | tr a-z A-Z' -i
$ parallel 'echo {} | tr a-z A-Z'
# Or more efficiently:
$ parallel --pipe tr a-z A-Z

$ loop 'echo $ITEM' --for "`ls`"
$ parallel echo {} ::: "`ls`"

$ ls | loop './my_program $ITEM' --until-success;
$ ls | parallel --halt now,success=1 ./my_program {}

$ ls | loop './my_program $ITEM' --until-fail;
$ ls | parallel --halt now,fail=1 ./my_program {}

$ ./deploy.sh;
  loop 'curl -sw "%(http_code)" http://coolwebsite.biz' \
     --every 5s --until-contains 200;
./announce_to_slack.sh
$ ./deploy.sh;
  loopy --delay 5s --halt now,success=1 \n  'curl -sw "%(http_code)" http://coolwebsite.biz | match 200'; \
./announce_to_slack.sh

$ loop "ping -c 1 mysite.com" --until-success; ./do_next_thing
$ loopy --halt now,success=1 ping -c 1 mysite.com; ./do_next_thing

$ ./create_big_file -o my_big_file.bin;
  loop 'ls' --until-contains 'my_big_file.bin';
./upload_big_file my_big_file.bin
# inotifywait is a better tool to detect file system changes.
# It can even make sure the file is complete
# so you are not uploading an incomplete file

$ inotifywait -q -r -mre MOVED_TO -e CLOSE_WRITE --format %w%f . |
  grep my_big_file.bin

$ ls | loop 'cp $ITEM $ITEM.bak'
$ ls | parallel cp {} {} .bak

$ loop './dothing.sh' --every 15s --until-success --num 5
$ parallel --retries 5 --delay 15s ::: ./dothing.sh

https://github.com/Miserlou/Loop/ (Last checked: 2018-10)

DIFFERENCES BETWEEN loriikeet AND GNU Parallel

lorikeet can run jobs in parallel. It does this based on a dependency graph described in a file, so this is similar to make.


DIFFERENCES BETWEEN spp AND GNU Parallel

spp can run jobs in parallel. spp does not use a command template to generate the jobs, but requires jobs to be in a file. Output from the jobs mix.

https://github.com/john01dav/spp (Last checked: 2019-01)

DIFFERENCES BETWEEN paral AND GNU Parallel

paral prints a lot of status information and stores the output from the commands run into files. This means it cannot be used the middle of a pipe like this

```
paral "echo this" "echo does not" "echo work" | wc
```

Instead it puts the output into files named like `out_{#}_command.out.log`. To get a very similar behaviour with GNU parallel use

```
--results 'out_{#}_\{s/[\w\-a-z_0-9]/g;s/\s+/_/g=_}.log' --eta
```

paral only takes arguments on the command line and each argument should be a full command. Thus it does not use command templates.

This limits how many jobs it can run in total, because they all need to fit on a single command line.

paral has no support for running jobs remotely.

EXAMPLES FROM README.markdown

The examples from README.markdown and the corresponding command run with GNU parallel (```
--results 'out_{#}_\{s/[\w\-a-z_0-9]/g;s/\s+/_/g=_}.log' --eta
```) is omitted from the GNU parallel command):

```
1$ paral "command 1" "command 2 --flag" "command arg1 arg2"
1$ parallel ::: "command 1" "command 2 --flag" "command arg1 arg2"

2$ paral "sleep 1 && echo c1" "sleep 2 && echo c2" 
  "sleep 3 && echo c3" "sleep 4 && echo c4" "sleep 5 && echo c5"
2$ parallel ::: "sleep 1 && echo c1" "sleep 2 && echo c2" 
  "sleep 3 && echo c3" "sleep 4 && echo c4" "sleep 5 && echo c5"
# Or shorter:
parallel "sleep {} && echo c{}" ::: {1..5}

3$ paral -n=0 "sleep 5 && echo c5" "sleep 4 && echo c4" 
```
"sleep 3 && echo c3" "sleep 2 && echo c2" "sleep 1 && echo c1"
3$ parallel ::: "sleep 5 && echo c5" "sleep 4 && echo c4"
   "sleep 3 && echo c3" "sleep 2 && echo c2" "sleep 1 && echo c1"
# Or shorter:
parallel -j0 "sleep () && echo c()" ::: 5 4 3 2 1

4$ paral -n=1 "sleep 5 && echo c5" "sleep 4 && echo c4"
   "sleep 3 && echo c3" "sleep 2 && echo c2" "sleep 1 && echo c1"
4$ parallel -j1 "sleep () && echo c()" ::: 5 4 3 2 1

5$ paral -n=2 "sleep 5 && echo c5" "sleep 4 && echo c4"
   "sleep 3 && echo c3" "sleep 2 && echo c2" "sleep 1 && echo c1"
5$ parallel -j2 "sleep () && echo c()" ::: 5 4 3 2 1

6$ paral -n=5 "sleep 5 && echo c5" "sleep 4 && echo c4"
   "sleep 3 && echo c3" "sleep 2 && echo c2" "sleep 1 && echo c1"
6$ parallel -j5 "sleep () && echo c()" ::: 5 4 3 2 1

7$ paral -n=1 "echo a && sleep 0.5 && echo b && sleep 0.5 && 
   echo c && sleep 0.5 && echo d && sleep 0.5 && 
   echo e && sleep 0.5 && echo f && sleep 0.5 && 
   echo g && sleep 0.5 && echo h"
7$ parallel ::: "echo a && sleep 0.5 && echo b && sleep 0.5 && 
   echo c && sleep 0.5 && echo d && sleep 0.5 && 
   echo e && sleep 0.5 && echo f && sleep 0.5 && 
   echo g && sleep 0.5 && echo h"

https://github.com/amattr/paral (Last checked: 2019-01)

DIFFERENCES BETWEEN concur AND GNU Parallel
concur is built to run jobs in parallel using a client/server model.

EXAMPLES FROM README.md
The examples from README.md:

1$ concur 'echo job {}' on slot {}: {}' : arg1 arg2 arg3 arg4
1$ parallel 'echo job {}' on slot {}: {}' ::: arg1 arg2 arg3 arg4

2$ concur 'echo job {}' on slot {}: {}' :: file1 file2 file3
2$ parallel 'echo job {}' on slot {}: {}' :: file1 file2 file3

3$ concur 'echo {}' < input_file
3$ parallel 'echo {}' < input_file

concur deals badly empty input files and with output larger than 64 KB.

https://github.com/mmstick/concurr (Last checked: 2019-01)

DIFFERENCES BETWEEN lesser-parallel AND GNU Parallel
lesser-parallel is the inspiration for parallel --embed. Both lesser-parallel and parallel --embed
define bash functions that can be included as part of a bash script to run jobs in parallel.
GNU Parallel alternatives

lesser-parallel implements a few of the replacement strings, but hardly any options, whereas parallel--embed gives you the full GNU parallel experience.

https://github.com/kou1okada/lesser-parallel (Last checked: 2019-01)

DIFFERENCES BETWEEN npm-parallel AND GNU Parallel

npm-parallel can run npm tasks in parallel.

There are no examples and very little documentation, so it is hard to compare to GNU parallel.

https://github.com/spion/npm-parallel (Last checked: 2019-01)

DIFFERENCES BETWEEN machma AND GNU Parallel

machma runs tasks in parallel. It gives time stamped output. It buffers in RAM.

EXAMPLES FROM README.md

The examples from README.md:

1$ # Put shorthand for timestamp in config for the examples
echo '---rpl "\"{time}\$_=::strftime("%Y-%m-%d %H:%M:%S", localtime())"" \n> ~/.parallel/machma
echo '---line-buffer --tagstring "(#) (time) ()"' \n>> ~/.parallel/machma

2$ find . -iname '*.jpg' |
   machma -- mogrify -resize 1200x1200 -filter Lanczos {}
find . -iname '*.jpg' |
   parallel --bar --Jmachma mogrify -resize 1200x1200 \n   -filter Lanczos {}

3$ cat /tmp/ips | machma -p 2 -- ping -c 2 -q {}
3$ cat /tmp/ips | parallel -j2 --Jmachma ping -c 2 -q {}

4$ cat /tmp/ips |
   machma -- sh -c 'ping -c 2 -q $0 > /dev/null && echo alive' {}
4$ cat /tmp/ips |
   parallel --Jmachma 'ping -c 2 -q {} > /dev/null && echo alive'

5$ find . -iname '*.jpg' |
   machma --timeout 5s -- mogrify -resize 1200x1200 \n   -filter Lanczos {}
5$ find . -iname '*.jpg' |
   parallel --timeout 5s --bar mogrify -resize 1200x1200 \n   -filter Lanczos {}

6$ find . -iname '*.jpg' -print0 |
   machma --null -- mogrify -resize 1200x1200 -filter Lanczos {}
6$ find . -iname '*.jpg' -print0 |
   parallel --null --bar mogrify -resize 1200x1200 \n   -filter Lanczos {}

https://github.com/fd0/machma (Last checked: 2019-06)
DIFFERENCES BETWEEN interlace AND GNU Parallel

Summary (see legend above):
- I2 I3 I4 - - -
- M1 - M3 - - M6
- O2 O3 - ---- x x
E1 E2 - ----
---------
--

**interlace** is built for network analysis to run network tools in parallel.

**interface** does not buffer output, so output from different jobs mixes.

The overhead for each target is \(O(n^2)\), so with 1000 targets it becomes very slow with an overhead in the order of 500ms/target.

EXAMPLES FROM interlace’s WEBSITE

Using **prips** most of the examples from https://github.com/codingo/Interlace can be run with GNU parallel:

Blocker commands.txt:
```bash
mkdir -p _output_//_target_/scans/_blocker_
  nmap _target_ -oA _output_//_target_/scans/_target_-nmap
interlace -tL ./targets.txt -cL commands.txt -o $output
```
```
parallel -a targets.txt \
  mkdir -p $output/()/scans/; nmap () -oA $output/()/scans/()-nmap
```

Blocks commands.txt:
```bash
_block:nmap_
  mkdir -p _target_/output/scans/
  nmap _target_ -oN _target_/output/scans/_target_-nmap
_block:nmap_
  nikto --host _target_
interlace -tL ./targets.txt -cL commands.txt

_nmap() { 
  mkdir -p $1/output/scans/
  nmap $1 -oN $1/output/scans/$1-nmap
}
export -f _nmap
parallel ::: _nmap "nikto --host" ::: targets.txt
```

Run Nikto Over Multiple Sites
```
interlace -tL ./targets.txt -threads 5 \
  -c "nikto --host _target_ > ./_target_-nikto.txt" -v
```
```
parallel -a targets.txt -P5 nikto --host {} \
  > ./{}-nikto.txt
```
GNU Parallel alternatives

Run Nikto Over Multiple Sites and Ports

interlace -tL ./targets.txt -threads 5 -c \
  "nikto --host _target_:__port_ > ./_target__port__nikto.txt" \
  -p 80,443 -v

parallel -P5 nikto --host [1]:[2] \> ./[1]-[2]-nikto.txt \
::: targets.txt ::: 80 443

Run a List of Commands against Target Hosts

commands.txt:
  nikto --host _target_:__port_ > _output_/target-nikto.txt 
  sslscan _target_:__port_ > _output_/target-sslscan.txt 
  testssl.sh _target_:__port_ > _output_/target-testssl.txt 
interlace -t example.com -o ~/Engagements/example/ \ 
  -cL ./commands.txt -p 80,443

parallel --results ~/Engagements/example/[2]:[3]{1} {1} (2):[3] \
::: "nikto --host" sslscan testssl.sh ::: example.com ::: 80 443

CIDR notation with an application that doesn't support it

interlace -t 192.168.12.0/24 -c "vhostscan _target_ \ 
  -oN _output_/target-vhosts.txt" -o ~/scans/ -threads 50 

prips 192.168.12.0/24 | 
  parallel -P50 vhostscan {} -oN ~/scans/{}-vhosts.txt

Glob notation with an application that doesn't support it

interlace -t 192.168.12.* -c "vhostscan _target_ \ 
  -oN _output_/target-vhosts.txt" -o ~/scans/ -threads 50

# Glob is not supported in prips 
prips 192.168.12.0/24 | 
  parallel -P50 vhostscan {} -oN ~/scans/{}-vhosts.txt

Dash (-) notation with an application that doesn't support it

interlace -t 192.168.12.1-15 -c \ 
  "vhostscan _target_ -oN _output_/target-vhosts.txt" \ 
  -o ~/scans/ -threads 50

# Dash notation is not supported in prips 
prips 192.168.12.1 192.168.12.15 | 
  parallel -P50 vhostscan {} -oN ~/scans/{}-vhosts.txt

Threading Support for an application that doesn't support it

interlace -tL ./target-list.txt -c \ 
  "vhostscan -t _target_ -oN _output_/target-vhosts.txt" \ 
  -o ~/scans/ -threads 50

cat ./target-list.txt |
GNU Parallel alternatives

```
parallel -P50 vhostscan -t {} -oN ~/scans/{}-vhosts.txt

alternatively

./vhosts-commands.txt:
    vhostscan -t $target -oN _output_/_target_-vhosts.txt
interlace -cl ./vhosts-commands.txt -tL ./target-list.txt \ 
    -threads 50 -o ~/scans

./vhosts-commands.txt:
    vhostscan -t "$1" -oN "$2"
parallel -P50 ./vhosts-commands.txt {} -oN ~/scans/{}-vhosts.txt \ 
    :::: ./target-list.txt

Exclusions

interlace -t 192.168.12.0/24 -e 192.168.12.0/26 -c \ 
    "vhostscan _target_ -oN _output_/_target_-vhosts.txt" \ 
    -o ~/scans/ -threads 50

prips 192.168.12.0/24 | grep -xv -Ff <(prips 192.168.12.0/26) | 
parallel -P50 vhostscan {} -oN ~/scans/{}-vhosts.txt

Run Nikto Using Multiple Proxies

interlace -tL ./targets.txt -pL ./proxies.txt -threads 5 -c \ 
    "nikto --host _target_:port_ -useproxy _proxy_ > \ 
    ./_target_:port_:nikto.txt" -p 80,443 -v

parallel -j5 \ 
    "nikto --host {1}:{2} -useproxy {3} > ./{1}-{2}-nikto.txt" \ 
    :::: ./targets.txt :::: 80 443 :::: ./proxies.txt

https://github.com/codingo/Interlace (Last checked: 2019-09)

DIFFERENCES BETWEEN otonvm Parallel AND GNU Parallel

I have been unable to get the code to run at all. It seems unfinished.

https://github.com/otonvm/Parallel (Last checked: 2019-02)

DIFFERENCES BETWEEN k-bx par AND GNU Parallel

par requires Haskell to work. This limits the number of platforms this can work on.

par does line buffering in memory. The memory usage is 3x the longest line (compared to 1x for parallel --lb). Commands must be given as arguments. There is no template.

These are the examples from https://github.com/k-bx/par with the corresponding GNU parallel command.

```
par "echo foo; sleep 1; echo foo; sleep 1; echo foo" \ 
    "echo bar; sleep 1; echo bar; sleep 1; echo bar" & & echo "success"
parallel --lb ::: "echo foo; sleep 1; echo foo; sleep 1; echo foo" \ 
    "echo bar; sleep 1; echo bar; sleep 1; echo bar" & & echo "success"

par "echo foo; sleep 1; foofoo" \ 
    "echo bar; sleep 1; echo bar; sleep 1; echo bar" & & echo "success"
parallel --lb --halt 1 ::: "echo foo; sleep 1; foofoo" 
```
"echo bar; sleep 1; echo bar; sleep 1; echo bar" && echo "success"

par "PARPREFIX=[fooechoer] echo foo" "PARPREFIX=[bar] echo bar"
parallel --lb --colsep , --tagstring {1} {2} \\
::: "[fooechoer],echo foo" "[bar],echo bar"

par --succeed "foo" "bar" && echo 'wow'
parallel "foo" "bar"; true && echo 'wow'

https://github.com/k-bx/par (Last checked: 2019-02)

DIFFERENCES BETWEEN parallelshell AND GNU Parallel

parallelshell does not allow for composed commands:

# This does not work
parallelshell 'echo foo;echo bar' 'echo baz;echo quuz'

Instead you have to wrap that in a shell:

parallelshell 'sh -c "echo foo;echo bar"' 'sh -c "echo baz;echo quuz"'

It buffers output in RAM. All commands must be given on the command line and all commands are started in parallel at the same time. This will cause the system to freeze if there are so many jobs that there is not enough memory to run them all at the same time.

https://github.com/keithamus/parallelshell (Last checked: 2019-02)

DIFFERENCES BETWEEN shell-executor AND GNU Parallel

shell-executor does not allow for composed commands:

# This does not work
sx 'echo foo;echo bar' 'echo baz;echo quuz'

Instead you have to wrap that in a shell:

sx 'sh -c "echo foo;echo bar"' 'sh -c "echo baz;echo quuz"'

It buffers output in RAM. All commands must be given on the command line and all commands are started in parallel at the same time. This will cause the system to freeze if there are so many jobs that there is not enough memory to run them all at the same time.


DIFFERENCES BETWEEN non-GNU par AND GNU Parallel

par buffers in memory to avoid mixing of jobs. It takes 1s per 1 million output lines.

par needs to have all commands before starting the first job. The jobs are read from stdin (standard input) so any quoting will have to be done by the user.

Stdout (standard output) is prepended with o:. Stderr (standard error) is sendt to stdout (standard output) and prepended with e:.

For short jobs with little output par is 20% faster than GNU parallel and 60% slower than xargs.

https://github.com/UnixJunkie/PAR
DIFFERENCES BETWEEN fd AND GNU Parallel

*fd* does not support composed commands, so commands must be wrapped in `sh -c`.

It buffers output in RAM.

It only takes file names from the filesystem as input (similar to `find`).

https://github.com/sharkdp/fd (Last checked: 2019-02)

DIFFERENCES BETWEEN lateral AND GNU Parallel

*lateral* is very similar to `sem`: It takes a single command and runs it in the background. The design means that output from parallel running jobs may mix. If it dies unexpectly it leaves a socket in `~/.lateral/socket.PID`.

*lateral* deals badly with too long command lines. This makes the *lateral* server crash:

```
lateral run echo `seq 100000| head -c 1000k`
```

Any options will be read by *lateral* so this does not work (*lateral* interprets the `-l`):

```
lateral run ls -l
```

Composed commands do not work:

```
lateral run pwd ';' ls
```

Functions do not work:

```
myfunc() { echo a; }
export -f myfunc
lateral run myfunc
```

Running *emacs* in the terminal causes the parent shell to die:

```
echo '#!/bin/bash' > mycmd
echo emacs -nw >> mycmd
chmod +x mycmd
lateral start
lateral run ./mycmd
```

Here are the examples from https://github.com/akramer/lateral with the corresponding GNU *sem* and GNU *parallel* commands:

```
1$ lateral start
    for i in $(cat /tmp/names); do
        lateral run -- some_command $i
done
lateral wait

1$ for i in $(cat /tmp/names); do
    sem some_command $i
done
    sem --wait

1$ parallel some_command :::: /tmp/names

2$ lateral start
```
for i in $(seq 1 100); do
    lateral run -- my_slow_command < workfile$i > /tmp/logfile$i
done
lateral wait

2$ for i in $(seq 1 100); do
    sem my_slow_command < workfile$i > /tmp/logfile$i
done
sem --wait

2$ parallel 'my_slow_command < workfile{} > /tmp/logfile{}' \ 
   ::: (1..100)

3$ lateral start -p 0 # yup, it will just queue tasks
for i in $(seq 1 100); do
    lateral run -- command_still_outputs_but_wont_spam inputfile$i
done
# command output spam can commence
lateral config -p 10; lateral wait

3$ for i in $(seq 1 100); do
    echo "command inputfile$i" >> joblist
done
parallel -j 10 :::: joblist

3$ echo 1 > /tmp/njobs
parallel -j /tmp/njobs command inputfile() \ 
   ::: (1..100)
echo 10 >/tmp/njobs
wait

https://github.com/akramer/lateral (Last checked: 2019-03)

DIFFERENCES BETWEEN with-this AND GNU Parallel

The examples from https://github.com/amritb/with-this.git and the corresponding GNU parallel command:

with -v "$\{cat myurls.txt\}" "curl -L this"
parallel curl -L ::: myurls.txt

with -v "$\{cat myregions.txt\}"
    "aws --region=\{\} ec2 describe-instance-status"
parallel aws --region={} ec2 describe-instance-status \ 
   ::: myregions.txt

with -v "$\{ls\}" "kubectl --kubeconfig=\{this get pods\"
ls | parallel kubectl --kubeconfig={} get pods

with -v "$\{ls | grep config\}" "kubectl --kubeconfig=\{this get pods\"
ls | grep config | parallel kubectl --kubeconfig={} get pods

with -v "$\{echo (1..10)\}" "echo 123"
parallel -N0 echo 123 ::: (1..10)
Stderr is merged with stdout. **with-this** buffers in RAM. It uses 3x the output size, so you cannot have output larger than 1/3rd the amount of RAM. The input values cannot contain spaces. Composed commands do not work.

**with-this** gives some additional information, so the output has to be cleaned before piping it to the next command.

https://github.com/amritb/with-this.git (Last checked: 2019-03)

**DIFFERENCES BETWEEN Tollef's parallel (moreutils) AND GNU Parallel**

Summary (see legend above):

- - - - I4 - - I7
- - M3 - - M6
- O2 O3 - O5 O6 - x x
E1 - - - - E7
- x x x x x x x x
- -

**EXAMPLES FROM Tollef's parallel MANUAL**

**Tollef** parallel sh -c "echo hi; sleep 2; echo bye" -- 1 2 3

**GNU** parallel "echo hi; sleep 2; echo bye" ::: 1 2 3

**Tollef** parallel -j 3 ufraw -o processed -- *.NEF

**GNU** parallel -j 3 ufraw -o processed ::: *.NEF

**Tollef** parallel -j 3 -- ls df "echo hi"

**GNU** parallel -j 3 ::: ls df "echo hi"

(Last checked: 2019-08)

**DIFFERENCES BETWEEN rargs AND GNU Parallel**

Summary (see legend above):

I1 - - - - - I7
- - M3 M4 - -
- O2 O3 - O5 O6 - O8 -
E1 - - E4 - -
- - - - - - - - -
- -

**rargs** has elegant ways of doing named regexp capture and field ranges.

With **GNU parallel** you can use **--rpl** to get a similar functionality as regexp capture gives, and use **join** and **@arg** to get the field ranges. But the syntax is longer. This:

```
--rpl 'r\{r\(\d+\)\..\(\d+\)\}"$_=join"$opt::colsep",@arg\[\$$1..\$$2\]'
```

would make it possible to use:

```
{1r3..6}
```

for field 3..6.

For full support of {n..m:s} including negative numbers use a dynamic replacement string like this:
PARALLEL=--rpl"\'(r((?-\d+)?)\.\.(?-\d+)?((:[^\}]*))?)\')
   $a = defined $2 ? $2 : 0 ? 1+$#arg+$2 : $2 : 1;
   $b = defined $4 ? $4 : 0 ? 1+$#arg+$4 : $4 : $#arg+1;
   $s = defined $6 ? $7 : " ";
   $_ = join $s,@arg[$a..$b]'

export PARALLEL

You can then do:

head /etc/passwd | parallel --colsep : echo ..=(1r..) ..3=(1r..3) \\
   4..=(1r4..) 2..4=(1r2..4) 3..3=(1r3..3) ..3:=(1r..3:)

EXAMPLES FROM rargs MANUAL

ls *.bak | rargs -p '(.*)\.bak' mv {0} {1}
ls *.bak | parallel mv {} .

cat download-list.csv | rargs -p '(?P<url>.*),(?P<filename>.*)' wget {url} -O {filename}
   cat download-list.csv | parallel --csv wget {1} -O {2}
   # or use regexps:
   cat download-list.csv |
   parallel --rpl '{url} s/,.*//' --rpl '{filename} s/.*?,//' wget {url} -O {filename}

   cat /etc/passwd | rargs -d: echo -e 'id: "\{1\}\t name: "\{5\}\t rest: "\{6..::\}"'
   cat /etc/passwd |
   parallel -q --colsep : echo -e 'id: "\{1\}\t name: "\{5\}\t rest: "\{=6 \\
   $._=join":","@arg[6..#arg]=")''

https://github.com/lotabout/rargs (Last checked: 2020-01)

DIFFERENCES BETWEEN threader AND GNU Parallel

Summary (see legend above):

<table>
<thead>
<tr>
<th></th>
<th>I1</th>
<th>M1</th>
<th>O1</th>
<th>E1</th>
</tr>
</thead>
<tbody>
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<td>-</td>
<td>-</td>
<td>-</td>
</tr>
<tr>
<td></td>
<td></td>
<td>M3</td>
<td>0</td>
<td>-</td>
</tr>
<tr>
<td></td>
<td></td>
<td>M6</td>
<td>0</td>
<td>-</td>
</tr>
<tr>
<td></td>
<td></td>
<td>M3</td>
<td>0</td>
<td>-</td>
</tr>
<tr>
<td></td>
<td></td>
<td>M6</td>
<td>N/A</td>
<td>N/A</td>
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</tbody>
</table>

Newline separates arguments, but newline at the end of file is treated as an empty argument. So this runs 2 jobs:

    echo two_jobs | threader -run 'echo "$THREADID"'

threader ignores stderr, so any output to stderr is lost. threader buffers in RAM, so output bigger than the machine's virtual memory will cause the machine to crash.

https://github.com/voodooEntity/threader (Last checked: 2020-04)
DIFFERENCES BETWEEN runp AND GNU Parallel

Summary (see legend above):

<table>
<thead>
<tr>
<th>I1</th>
<th>I2</th>
<th>M1</th>
<th>M6</th>
<th>O1</th>
<th>O2</th>
<th>O3</th>
<th>O5</th>
<th>O6</th>
<th>N/A</th>
<th>N/A</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
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<td></td>
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<td></td>
</tr>
</tbody>
</table>

(M3): You can add a prefix and a postfix to the input, so it means you can only insert the argument on the command line once.

runp runs 10 jobs in parallel by default. runp blocks if output of a command is > 64 Kbytes. Quoting of input is needed. It adds output to stderr (this can be prevented with -q)

Examples as GNU Parallel

```sh
base='https://images-api.nasa.gov/search'
query='jupiter'
desc='planet'
type='image'
url="$base?q=$query&description=$desc&media_type=$type"

# Download the images in parallel using runp
curl -s $url | jq -r .collection.items[].href | \n   runp -p 'curl -s' | jq -r .[] | grep large | \n   runp -p 'curl -s -L -O'

time curl -s $url | jq -r .collection.items[].href | \n   runp -g 1 -q -p 'curl -s' | jq -r .[] | grep large | \n   runp -g 1 -q -p 'curl -s -L -O'

# Download the images in parallel
curl -s $url | jq -r .collection.items[].href | \n   parallel curl -s | jq -r .[] | grep large | \n   parallel curl -s -L -O

time curl -s $url | jq -r .collection.items[].href | \n   parallel -j 1 curl -s | jq -r .[] | grep large | \n   parallel -j 1 curl -s -L -O
```

Run some test commands (read from file)

```sh
# Create a file containing commands to run in parallel.
cat << EOF > /tmp/test-commands.txt
sleep 5
sleep 3
blah    # this will fail
ls $PWD  # PWD shell variable is used here
EOF

# Run commands from the file.
runp /tmp/test-commands.txt > /dev/null
```
parallel -a /tmp/test-commands.txt > /dev/null

Ping several hosts and see packet loss (read from stdin)

```bash
# First copy this line and press Enter
runp -p 'ping -c 5 -W 2' -s '|' grep loss
localhost
1.1.1.1
8.8.8.8
# Press Enter and Ctrl-D when done entering the hosts

# First copy this line and press Enter
parallel ping -c 5 -W 2 '{}' | grep loss
localhost
1.1.1.1
8.8.8.8
# Press Enter and Ctrl-D when done entering the hosts
```

Get directories' sizes (read from stdin)

```bash
echo -e "$HOME
/etc
/tmp" | runp -q -p 'sudo du -sh'
echo -e "$HOME
/etc
/tmp" | parallel sudo du -sh
# or:
parallel sudo du -sh :::
"$HOME
/etc
/tmp"
```

Compress files

```bash
find . -iname '*.txt' | runp -p 'gzip --best'
find . -iname '*.txt' | parallel gzip --best
```

Measure HTTP request + response time

```bash
export CURL="curl -w 'time_total: %{time_total}\n"
CURL="$CURL -o /dev/null -s https://golang.org/"
perl -wE 'for (1..10) { say $ENV{CURL} }' | runp -q # Make 10 requests
perl -wE 'for (1..10) { say $ENV{CURL} }' | parallel # or:
parallel -N0 "$CURL" ::: (1..10)
```

Find open TCP ports

```bash
cat << EOF > /tmp/host-port.txt
localhost 22
localhost 80
localhost 81
127.0.0.1 443
127.0.0.1 444
scanme.nmap.org 22
scanme.nmap.org 23
scanme.nmap.org 443
EOF
```

```bash
cat /tmp/host-port.txt |
```
```bash
1$ cat /tmp/host-port.txt |
```
GNU Parallel alternatives

```
runp -q -p 'netcat -v -w2 -z' 2>&1 | egrep '(succeeded!|open)\$
```

```
# --colsep is needed to split the line
1$ cat /tmp/host-port.txt |
    parallel --colsep ' netcat -v -w2 -z 2>&1 |
        egrep '(succeeded!|open)\$'
# or use uq for unquoted:
1$ cat /tmp/host-port.txt |
    parallel netcat -v -w2 -z {=uq=} 2>&1 |
        egrep '(succeeded!|open)\$'
```

https://github.com/jreisinger/runp (Last checked: 2020-04)

DIFFERENCES BETWEEN papply AND GNU Parallel

Summary (see legend above):

- - - I4 - - -
M1 - M3 - - M6
- - O3 - O5 - - N/A N/A O10
E1 - - E4 - - -
----------
--

papply does not print the output if the command fails:

```
$ papply 'echo %F; false' foo
"echo foo; false" did not succeed
```

papply's replacement strings (%F %d %f %n %e %z) can be simulated in GNU parallel by putting this in ~/.parallel/config:

```
--rpl '%F'
--rpl '%d $_=Q(::dirname($));'
--rpl '%f $s:.*/\:.;'
--rpl '%n $s:.*/\:.;s:\.[^/\.]+$:;'
--rpl '%e $s:.*\.:;'
--rpl '%z $=_""'
```

papply buffers in RAM, and uses twice the amount of output. So output of 5 GB takes 10 GB RAM.

The buffering is very CPU intensive: Buffering a line of 5 GB takes 40 seconds (compared to 10 seconds with GNU parallel).

Examples as GNU Parallel

```
1$ papply gzip *.txt
1$ parallel gzip ::: *.txt
```

```
2$ papply "convert %F %n.jpg" *.png
2$ parallel convert () (.).jpg ::: *.png
```

https://pypi.org/project/papply/ (Last checked: 2020-04)
DIFFERENCES BETWEEN async AND GNU Parallel

Summary (see legend above):
--- I4 - I7
------ M6
- O2 O3 - O5 O6 - N/A N/A O10
E1 - E4 - E6 -
----------
S1 S2

async is very similar to GNU parallel's --semaphore mode (aka sem). async requires the user to start a server process.

The input is quoted like -q so you need bash -c "...;..." to run composed commands.

Examples as GNU Parallel

1$ S="/tmp/example_socket"

1$ ID=myid

2$ async -s="$S" server --start

2$ # GNU Parallel does not need a server to run

3$ for i in {1..20}; do
    # prints command output to stdout
    async -s="$S" cmd -- bash -c "sleep 1 && echo test $i"
    done

3$ for i in {1..20}; do
    # prints command output to stdout
    sem --id "$ID" -j100% "sleep 1 && echo test $i"
    # GNU Parallel will only print job when it is done
    # If you need output from different jobs to mix
    # use -u or --line-buffer
    sem --id "$ID" -j100% --line-buffer "sleep 1 && echo test $i"
    done

4$ # wait until all commands are finished
async -s="$S" wait

4$ sem --id "$ID" --wait

5$ # configure the server to run four commands in parallel
async -s="$S" server -j4

5$ export PARALLEL=-j4

6$ mkdir "/tmp/ex_dir"
    for i in {21..40}; do
        # redirects command output to /tmp/ex_dir/file*
        async -s="$S" cmd -o "/tmp/ex_dir/file$i" -- \
        bash -c "sleep 1 && echo test $i"
6$ mkdir "/tmp/ex_dir"
   for i in {21..40}; do
     # redirects command output to /tmp/ex_dir/file*
     sem --id "$ID" --result '/tmp/my-ex/file-{=$_=""=}'"$i" \ 
     "sleep 1 && echo test $i"
   done

7$ sem --id "$ID" --wait

7$ async -s="$S" wait

8$ # stops server
   async -s="$S" server --stop

8$ # GNU Parallel does not need to stop a server

https://github.com/ctbur/async/ (Last checked: 2020-11)

DIFFERENCES BETWEEN pardi AND GNU Parallel

Summary (see legend above):
I1 I2 - - - I7
M1 - - - M6
O1 O2 O3 O4 O5 - O7 - - O10
E1 - E4 - -
--------
--

pardi is very similar to parallel --pipe --cat: It reads blocks of data and not arguments. So it cannot insert an argument in the command line. It puts the block into a temporary file, and this file name (%IN) can be put in the command line. You can only use %IN once.

It can also run full command lines in parallel (like: cat file | parallel).

EXAMPLES FROM pardi test.sh

1$ time pardi -v -c 100 -i data/decoys.smi -ie .smi -oe .smi \ 
   -o data/decoys_std_pardi.smi \ 
   -w '{(standardiser -i %IN -o %OUT 2>&1) > /dev/null}'

1$ cat data/decoys.smi |
   time parallel -N 100 --pipe --cat \ 
   '{(standardiser -i {} -o {}) 2>&1) > /dev/null; cat {}; rm {}}' \ 
   > data/decoys_std_pardi.smi

2$ pardi -n 1 -i data/test_in.types -o data/test_out.types \ 
   -d 'r:^#atoms:' -w 'cat %IN > %OUT'

2$ cat data/test_in.types | parallel -n 1 -k --pipe --cat \ 
   --regexp --recstart '^#atoms' 'cat {}' > data/test_out.types

3$ pardi -c 6 -i data/test_in.types -o data/test_out.types \
GNU Parallel alternatives

DIFFERENCES BETWEEN bthread AND GNU Parallel

Summary (see legend above):

- - - I4 - - -
- - - - - M6
O1 - O3 - - O7 O8 - -
E1 - - - - - - - - - -
--

bthread takes around 1 sec per MB of output. The maximal output line length is 1073741759.

You cannot quote space in the command, so you cannot run composed commands like sh -c "echo a; echo b".

https://gitlab.com/netikras/bthread (Last checked: 2021-01)

DIFFERENCES BETWEEN simple_gpu_scheduler AND GNU Parallel

Summary (see legend above):

I1 - - - - I7
M1 - - - - M6
- O2 O3 - - O6 - x x O10
E1 - - - - - - - - - -
--

EXAMPLES FROM simple_gpu_scheduler MANUAL

1$ simple_gpu_scheduler --gpus 0 1 2 < gpu_commands.txt

1$ parallel -j3 --shuf 
    CUDA_VISIBLE_DEVICES='{:1 $_=slot()-1 =} {=ug;=}' < gpu_commands.txt

https://github.com/UnixJunkie/pardi (Last checked: 2021-01)
2$ simple_hypersearch "python3 train_dnn.py --lr {lr} --batch_size {bs}" \
   -p lr 0.001 0.0005 0.0001 -p bs 32 64 128 | \
   simple_gpu_scheduler --gpus 0,1,2

2$ parallel --header : --shuf -j3 -v \ 
   CUDA_VISIBLE_DEVICES='(={1 $=slot()-1 =}')' \ 
   python3 train_dnn.py --lr {lr} --batch_size {bs} \ 
   ::: lr 0.001 0.0005 0.0001 ::: bs 32 64 128

3$ simple_hypersearch \
   "python3 train_dnn.py --lr {lr} --batch_size {bs}" \ 
   --n-samples 5 -p lr 0.001 0.0005 0.0001 -p bs 32 64 128 | \
   simple_gpu_scheduler --gpus 0,1,2

3$ parallel --header : --shuf \ 
   CUDA_VISIBLE_DEVICES='(={1 $=slot()-1; seq() > 5 and skip() =}')' \ 
   python3 train_dnn.py --lr {lr} --batch_size {bs} \ 
   ::: lr 0.001 0.0005 0.0001 ::: bs 32 64 128

4$ touch gpu.queue \
   tail -f -n 0 gpu.queue | simple_gpu_scheduler --gpus 0,1,2 & \
   echo "my_command_with | and stuff > logfile" >> gpu.queue

4$ touch gpu.queue \
   tail -f -n 0 gpu.queue | \
   parallel -j3 CUDA_VISIBLE_DEVICES='(={1 $=slot()-1; seq() > 5 and skip() =}')' & \
   # Needed to fill job slots once \
   # Add jobs \
   echo "my_command_with | and stuff > logfile" >> gpu.queue \
   # Needed to flush output from completed jobs \
   seq 3 | parallel echo true >> gpu.queue

https://github.com/ExpectationMax/simple_gpu_scheduler (Last checked: 2021-01)

**DIFFERENCES BETWEEN parasweep AND GNU Parallel**

*parasweep* is a Python module for facilitating parallel parameter sweeps.

A *parasweep* job will normally take a text file as input. The text file contains arguments for the job. Some of these arguments will be fixed and some of them will be changed by *parasweep*.

It does this by having a template file such as template.txt:

```
Xval: {x}
Yval: {y}
FixedValue: 9
# x with 2 decimals
DecimalX: {x:.2f}
TenX: ${x*10}
RandomVal: {r}
```

and from this template it generates the file to be used by the job by replacing the replacement strings.

Being a Python module *parasweep* integrates tighter with Python than GNU *parallel*. You get the parameters directly in a Python data structure. With GNU *parallel* you can use the JSON or CSV
GNU Parallel alternatives

output format to get something similar, but you would have to read the output.

**parasweep** has a filtering method to ignore parameter combinations you do not need.

Instead of calling the jobs directly, **parasweep** can use Python's Distributed Resource Management Application API to make jobs run with different cluster software.

**GNU parallel** **--tmpl** supports templates with replacement strings. Such as:

```
Xval: (x)
Yval: (y)
FixedValue: 9
# x with 2 decimals
DecimalX: {$=x$_=sprintf("%.2f",$_) =}
TenX: {$=x$_=$_*10 =}
RandomVal: {$=1$_=rand() =}
```

that can be used like:

```
parallel --header : --tmpl my.tmpl={#}.t myprog {#}.t \\
::: x 1 2 3 ::: y 1 2 3
```

Filtering is supported as:

```
parallel --filter '{1} > {2}' echo ::: 1 2 3 ::: 1 2 3
```

https://github.com/eviatarbach/parasweep (Last checked: 2021-01)

**DIFFERENCES BETWEEN** parallel-bash **AND GNU Parallel**

Summary (see legend above):

```
I1 I2 ----- 
-- M3 -- M6
- O2 O3 - O5 O6 - O8 x O10
E1 ------- 
---------- 
--
```

**parallel-bash** is written in pure bash. It is really fast (overhead of ~0.05 ms/job compared to GNU parallel's 3-10 ms/job). So if your jobs are extremely short lived, and you can live with the quite limited command, this may be useful.

It works by making a queue for each process. Then the jobs are distributed to the queues in a round robin fashion. Finally the queues are started in parallel. This works fine, if you are lucky, but if not, all the long jobs may end up in the same queue, so you may see:

```
$ printf "%b\n" 1 1 1 4 1 1 1 4 1 1 1 4 | 
  time parallel -P4 sleep {}
(7 seconds)
$ printf "%b\n" 1 1 1 4 1 1 1 4 1 1 1 4 | 
  time ./parallel-bash.bash -p 4 -c sleep {}
(12 seconds)
```

Because it uses bash lists, the total number of jobs is limited to 167000..265000 depending on your environment. You get a segmentation fault, when you reach the limit.

Ctrl-C does not stop spawning new jobs. Ctrl-Z does not suspend running jobs.
EXAMPLES FROM parallel-bash

```
1$ some_input | parallel-bash -p 5 -c echo
1$ some_input | parallel -j 5 echo
2$ parallel-bash -p 5 -c echo < some_file
2$ parallel -j 5 echo < some_file
3$ parallel-bash -p 5 -c echo <<< 'some string'
3$ parallel -j 5 -c echo <<< 'some string'
4$ something | parallel-bash -p 5 -c echo {} {}
4$ something | parallel -j 5 echo {} {}
```


DIFFERENCES BETWEEN bash-concurrent AND GNU Parallel

bash-concurrent is more an alternative to make than to GNU parallel. Its input is very similar to a Makefile, where jobs depend on other jobs.

It has a nice progress indicator where you can see which jobs completed successfully, which jobs are currently running, which jobs failed, and which jobs were skipped due to a depending job failed. The indicator does not deal well with resizing the window.

Output is cached in tempfiles on disk, but is only shown if there is an error, so it is not meant to be part of a UNIX pipeline. If bash-concurrent crashes these tempfiles are not removed.

It uses an O(n*n) algorithm, so if you have 1000 independent jobs it takes 22 seconds to start it.

https://github.com/themattrix/bash-concurrent (Last checked: 2021-02)

DIFFERENCES BETWEEN spawntool AND GNU Parallel

Summary (see legend above):

```
I1 - - - - - -
M1 - - - - M6
- O2 O3 - O5 O6 - x x O10
E1 - - - - - - - -
--------
-
```

spawn reads a full command line from stdin which it executes in parallel.

http://code.google.com/p/spawntool/ (Last checked: 2021-07)

DIFFERENCES BETWEEN go-pssh AND GNU Parallel

Summary (see legend above):

```
- - - - - - - -
M1 - - - - - -
```
go-pssh does ssh in parallel to multiple machines. It runs the same command on multiple machines similar to --nonall.

The hostnames must be given as IP-addresses (not as hostnames).

Output is sent to stdout (standard output) if command is successful, and to stderr (standard error) if the command fails.

EXAMPLES FROM go-pssh

1$ go-pssh -l <ip>,<ip> -u <user> -p <port> -P <passwd> -c "<command>"

1$ parallel --nonall "<command>"

2$ go-pssh scp -f host.txt -u <user> -p <port> -P <password> \
   --basefile /local/file_or_directory \
   --wd /remote/directory

2$ parallel --nonall --slf host.txt \
   --basefile /local/file_or_directory/. \
   --wd /remote/directory \
   --ssh 'sshpass -p <password> ssh -p <port> -l <user>' true

3$ go-pssh scp -l <ip>,<ip> -u <user> -p <port> -P <password> \
   --basefile /local/file_or_directory \
   --wd /remote/directory

3$ parallel --nonall "<command>"

https://github.com/xuchenCN/go-pssh (Last checked: 2021-07)

DIFFERENCES BETWEEN go-parallel AND GNU Parallel

Summary (see legend above):

I1 I2 - - - I7
-- M3 -- M6
- O2 O3 - O5 - - x x - O10
E1 -- E4 - -
----------
--

go-parallel uses Go templates for replacement strings. Quite similar to the \(= perl expr =\)
replacement string.

EXAMPLES FROM go-parallel

1$ go-parallel -a ./files.txt -t 'cp {{.Input}} {{.Input | dirname |
   dirname}}'

1$ parallel -a ./files.txt cp () '{= $_=::dirname::dirname($_) =}'
2$ go-parallel -a ./files.txt -t 'mkdir -p {{.Input}} {{noExt .Input}}'

2$ parallel -a ./files.txt echo mkdir -p {}

3$ go-parallel -a ./files.txt -t 'mkdir -p {{.Input}} {{noExt .Input}}'

3$ parallel -a ./files.txt echo mkdir -p {}

https://github.com/mylanconnolly/parallel (Last checked: 2021-07)

DIFFERENCES BETWEEN p AND GNU Parallel

Summary (see legend above):
--- I4 -- N/A
----- M6
- O2 O3 O5 O6 - x x - O10
E1 ------
----------
--

p is a tiny shell script. It can color output with some predefined colors, but is otherwise quite limited.
It maxes out at around 116000 jobs (probably due to limitations in Bash).

EXAMPLES FROM p

Some of the examples from p cannot be implemented 100% by GNU parallel: The coloring is a bit different, and GNU parallel cannot have --tag for some inputs and not for others.

The coloring done by GNU parallel is not exactly the same as p.

1$ p -bc blue "ping 127.0.0.1" -uc red "ping 192.168.0.1" \ -rc yellow "ping 192.168.1.1" -t example "ping example.com"

1$ parallel --lb -j0 --color --tag ping \ :::: 127.0.0.1 192.168.0.1 192.168.1.1 example.com

2$ p "tail -f /var/log/httpd/access_log" \ -bc red "tail -f /var/log/httpd/error_log"

2$ cd /var/log/httpd;
parallel --lb --color --tag tail -f :::: access_log error_log

3$ p tail -f "some file" \
p tail -f "other file with space.txt"

3$ parallel --lb tail -f :::: 'some file' "other file with space.txt"

4$ p -t project1 "hg pull project1" -t project2 \ "hg pull project2" -t project3 "hg pull project3"

4$ parallel --lb hg pull :::: project{1..3}

Todo

http://code.google.com/p/push/ (cannot compile)
https://github.com/krashanoff/parallel
https://github.com/Nukesor/pueue
https://github.com/JeKeLiM/simple_distribute_job
https://github.com/reggi/pkgrun - not obvious how to use
https://github.com/benoror/better-npm-run - not obvious how to use
https://github.com/bahmutov/with-package
https://github.com/flesler/parallel
https://github.com/Julian/Verge
https://manpages.ubuntu.com/manpages/xenial/man1/tsp.1.html
https://vicerveza.homeunix.net/~viric/soft/ts/
https://github.com/chapmanjacobd/que

TESTING OTHER TOOLS

There are certain issues that are very common on parallelizing tools. Here are a few stress tests. Be warned: If the tool is badly coded it may overload your machine.

MIX: Output mixes

Output from 2 jobs should not mix. If the output is not used, this does not matter; but if the output is used then it is important that you do not get half a line from one job followed by half a line from another job.

If the tool does not buffer, output will most likely mix now and then.

This test stresses whether output mixes.

#!/bin/bash

paralleltool="parallel -j0"

cat <&-EOF > mycommand
#!/bin/bash

# If a, b, c, d, e, and f mix: Very bad
perl -e 'print STDOUT "a"x3000_000,"' "
perl -e 'print STDERR "b"x3000_000,"' "
perl -e 'print STDOUT "c"x3000_000,"' "
perl -e 'print STDERR "d"x3000_000,"' "
perl -e 'print STDOUT "e"x3000_000,"' "
perl -e 'print STDERR "f"x3000_000,"' "
echo
echo >&2
EOF
chmod +x mycommand
# Run 30 jobs in parallel

```
seq 30 |
$paralleltool ./mycommand > (tr -s abcdef) 2> (tr -s abcdef >&2)
```

# 'a c e' and 'b d f' should always stay together
# and there should only be a single line per job

**STDERRMERGE: Stderr is merged with stdout**

Output from stdout and stderr should not be merged, but kept separated.

This test shows whether stdout is mixed with stderr.

```
#!/bin/bash
paralleltool="parallel -j0"

cat <<<EOF > mycommand
#!/bin/bash

echo stdout
echo stderr >&2
echo stdout
echo stderr >&2
EOF
chmod +x mycommand

# Run one job
echo |
   $paralleltool ./mycommand > stdout 2> stderr
cat stdout
cat stderr
```

**RAM: Output limited by RAM**

Some tools cache output in RAM. This makes them extremely slow if the output is bigger than physical memory and crash if the output is bigger than the virtual memory.

```
#!/bin/bash

paralleltool="parallel -j0"

cat <EOF > mycommand
#!/bin/bash

# Generate 1 GB output
yes "$perl -e 'print "c"x30_000'" | head -c 1G
EOF
chmod +x mycommand

# Run 20 jobs in parallel
# Adjust 20 to be > physical RAM and < free space on /tmp
seq 20 | time $paralleltool ./mycommand | wc -c
DISKFULL: Incomplete data if /tmp runs full

If caching is done on disk, the disk can run full during the run. Not all programs discover this. GNU Parallel discovers it, if it stays full for at least 2 seconds.

#!/bin/bash

paralleltool="parallel -j0"

# This should be a dir with less than 100 GB free space
smalldisk=/tmp/shm/parallel

TMPDIR="$smalldisk"
export TMPDIR

max_output()
{
    # Force worst case scenario:
    # Make GNU Parallel only check once per second
    sleep 10
    # Generate 100 GB to fill $TMPDIR
    # Adjust if /tmp is bigger than 100 GB
    yes | head -c 100G >$TMPDIR/$$
    # Generate 10 MB output that will not be buffered due to full disk
    perl -e 'print "X"x10_000_000' | head -c 10M
    echo This part is missing from incomplete output
    sleep 2
    rm $TMPDIR/$$
    echo Final output
}

export -f max_output
seq 10 | $paralleltool max_output | tr -s X

CLEANUP: Leaving tmp files at unexpected death

Some tools do not clean up tmp files if they are killed. If the tool buffers on disk, they may not clean up, if they are killed.

#!/bin/bash

paralleltool=parallel

ls /tmp >/tmp/before
seq 10 | $paralleltool sleep &
pid=$!
# Give the tool time to start up
sleep 1
# Kill it without giving it a chance to cleanup
kill -9 $!
# Should be empty: No files should be left behind
diff <(ls /tmp) /tmp/before

SPCCHAR: Dealing badly with special file names.

It is not uncommon for users to create files like:

    My brother's 12" *** record (costs $$$).jpg
Some tools break on this.

```bash
#!/bin/bash
paralleltool=parallel

touch "My brother's 12\" *** record (costs \$\$\$).jpg"
ls My*jpg | $paralleltool ls -l
```

**COMPOSED: Composed commands do not work**
Some tools require you to wrap composed commands into `bash -c`.

```bash
echo bar | $paralleltool echo foo';' echo {}
```

**ONEREP: Only one replacement string allowed**
Some tools can only insert the argument once.

```bash
echo bar | $paralleltool echo () foo {}
```

**INPUTSIZE: Length of input should not be limited**
Some tools limit the length of the input lines artificially with no good reason. GNU parallel does not:

```bash
perl -e 'print "foo."."x"x100_000_000' | parallel echo {}
```

GNU parallel limits the command to run to 128 KB due to `execve(1):

```bash
perl -e 'print "x"x131_000' | parallel echo {} | wc
```

**NUMWORDS: Speed depends on number of words**
Some tools become very slow if output lines have many words.

```bash
#!/bin/bash
paralleltool=parallel

cat <<-EOF > mycommand
#!/bin/bash

# 10 MB of lines with 1000 words
yes "`seq 1000``" | head -c 10M
EOF
chmod +x mycommand

# Run 30 jobs in parallel
seq 30 | time $paralleltool -j0 ./mycommand > /dev/null
```

**4GB: Output with a line > 4GB should be OK**

```bash
#!/bin/bash

paralleltool="parallel -j0"

cat <<-EOF > mycommand
#!/bin/bash
```
perl -e '\$a="a"x1000_000; for(1..5000) { print \$a }'
EOF
chmod +x mycommand

# Run 1 job
seq 1 | $paralleltool ./mycommand | LC_ALL=C wc

**AUTHOR**

When using GNU parallel for a publication please cite:


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Parts of the manual concerning xargs compatibility is inspired by the manual of xargs from GNU findutils 4.4.2.

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**DEPENDENCIES**

GNU parallel uses Perl, and the Perl modules Getopt::Long, IPC::Open3, Symbol, IO::File, POSIX, and File::Temp. For remote usage it also uses rsync with ssh.

**SEE ALSO**

find(1), xargs(1), make(1), pexec(1), ppss(1), xjobs(1), prll(1), dxargs(1), mdm(1)