

# Cheat-Sheet for tools-for-g16.bash

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## Introduction

This accompanies the repository `polyluxus/tools-for-g16.bash`.

Various bash scripts to aid the use of the quantum chemistry software package Gaussian 16.

## Preliminary notes

The notation in brackets [ ] indicate optional arguments/inputs; arguments in angles < > require human input; a bar | indicates alternatives.

The following abbreviations will be used:

`opt` Short for option(s)

`ARG` String type argument

`INT` Positive integer

`NUM` Whole number

`FLT` Floating point number

`DUR` Duration in format [[HH:]MM:]SS

## Installation & Configuration

General settings for the scripts can be found in the file `g16.tools.rc`. Alternatively, settings can be stored in `.g16.toolsrc`, which always has precedence. Every script will check three different directories in the order 1. installation directory 2. user's home 3. parent working directory. It will load the last configuration file it finds.

Setting files can be generated with the `configure/configure.sh` script.

## g16.prepare.sh

This tool reads in a file containing a set of cartesian coordinates and writes a Gaussian inputfile with predefined keywords. The script interfaces to Xmol format, Turbomole/ GFN-xTB coord format, too.

Usage: `g16.prepare.sh [opt] <file>`

- T <FLT> Temperature (kelvin)
- P <FLT> Pressure (atmosphere)
- r <ARG> Add ARG to route section
- R <ARG> Specific route section ARG
- l <INT> Load predefined route section
- l list Show all predefined route sections
- t <ARG> Adds ARG to end of file

- C <ARG> Specify caption/title of job;  
Replacements: %F input filename; %f input filename without .xyz; %s like %f, also filtering start; %j jobname.
- j <ARG> Jobname
- j %s Jobname is filename filtering start.xyz
- f <ARG> Filename of generated input
- c <NUM> Charge
- M <INT> Multiplicity (> 1)
- U <INT> Unpaired electrons (> 0)
- m <INT> Memory (megabyte)
- p <INT> Processors
- d <INT> disksize via MaxDisk (megabyte)
- s Silence script (incremental)
- h Help file

## g16.testroute.sh

This tool parses a Gaussian 16 inputfile and tests the route section for syntax errors with the Gaussian 16 utility `testrt`.

- s Silence script (incremental)
- h Help file

## g16.dissolve.sh

This tool reads in a Gaussian 16 inputfile and adds relevant keywords for solvent corrections.

Usage: `g16.dissolve.sh [opt] <file>`

- o <ARG> Adds option ARG to the `scrfl` keyword.
- S <ARG> Adds option `solvent=ARG` to the `scrfl` option list.
- O Runs an optimisation (preserves or adds OPT)
- r <ARG> Add ARG to route section
- t <ARG> Adds ARG to end of file
- m <INT> Memory (megabyte)
- p <INT> Processors
- d <INT> disksize via MaxDisk (megabyte)
- s Silence script (incremental)
- h Help file

## g16.freqinput.sh

This tool reads in a Gaussian 16 inputfile and adds relevant keywords for a frequency calculation.

Usage: `g16.freqinput.sh [opt] <file>`

- o <ARG> Adds option ARG to the `freq` keyword.
- R Adds option `ReadFC` to the `freq` option list.

- T <FLT> Temperature (kelvin)
- P <FLT> Pressure (atmosphere)
- r <ARG> Add ARG to route section
- t <ARG> Adds ARG to end of file
- m <INT> Memory (megabyte)
- p <INT> Processors
- d <INT> disksize via MaxDisk (megabyte)
- s Silence script (incremental)
- h Help file

## g16.submit.sh

This tool parses and then submits a Gaussian 16 inputfile to a queueing system.

Usage: `g16.submit.sh [opt] <file>`

- m <INT> Memory (megabyte)
- p <INT> Processors
- d <INT> disksize via MaxDisk (megabyte)
- w <DUR> Walltime limit
- e <ARG> Specify an environment variable ARG in format <VAR=value>
- j <INT> Wait for job with ID INT
- H Submit with status hold (PBS) or PSUSP (BSUB)
- k Only create (keep) the jobscript, do not submit it.
- Q <ARG> Queue for which job script should be created (`pbs-gen/bsub-rwth`)
- P <ARG> Account to project (BSUB); if ARG is `default/0/'` presets are overwritten.
- u <ARG> set user email address (BSUB); if ARG is `default/0/'` presets are overwritten.
- s Silence script (incremental)
- h Help file

## g16.getenergy.sh

This tool finds energy statements from Gaussian 16 calculations.

Usage: `g16.getenergy.sh [opt] [<file(s)>]`

If no files given, it finds energy statements from all log files in the current directory.

- i <ARG> Specify input suffix if processing directory
- o <ARG> Specify output suffix if processing directory
- s Silence script (incremental)
- h Help file

## g16.getfreq.sh

This tool summarises a frequency calculation and extracts the thermochemistry data.

Usage: `g16.getfreq.sh` [`opt`] `<file(s)>`

`-v` Incrementally increase verbosity  
`-V <INT>` Set level of verbosity directly, (0-4)  
`-c` Separate values by comma (`-V0` or `-V1`)  
`-O <ARG>` Write summary to file instead of screen  
`-s` Silence script (incremental)  
`-h` Help file

## g16.chk2xyz.sh

A tool to convert a checkpoint file to an xyz file. This formats the `chk` first to a `fchk`.

Usage: `g16.chk2xyz.sh` [`-s`] `-h` | `-f` | `<chk-file(s)>`

`-f` Formats all checkpointfiles that are found in the current directory  
`-s` Silence script (incremental)  
`-h` Help file

## Author, Bugs, and the Rest

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