!Vámonos a experimentar!

```
load $2,3-dibromopentane
load $(2R,3S)-2,3-dibromopentane
load $(2S,3R)-2,3-dibromopentane

set zoomlarge false
load files "$2,3-dibromopentane"
           "$(2S,3R)-2,3-dibromopentane"
select 1.1; translateSelected {-4 0 0}
select 2.1; translateSelected {4 0 0}
frame *
compare {1.1} {2.1} rotate translate
compare {1.1} {2.1} atoms {!_H} rotate
translate
prompt compare({1.1}, {2.1}, “isomer”)
```

http://chemapps.stolaf.edu/jmol/jsmol

simple2.htm	general
jsmetest2.htm	2D->3D
jmol-flot-energy.htm	la rotación enlace
jsmol.htm	más capacidades

http://chemagic.org/molecules/mini.html

http://www.chemtube3d.com