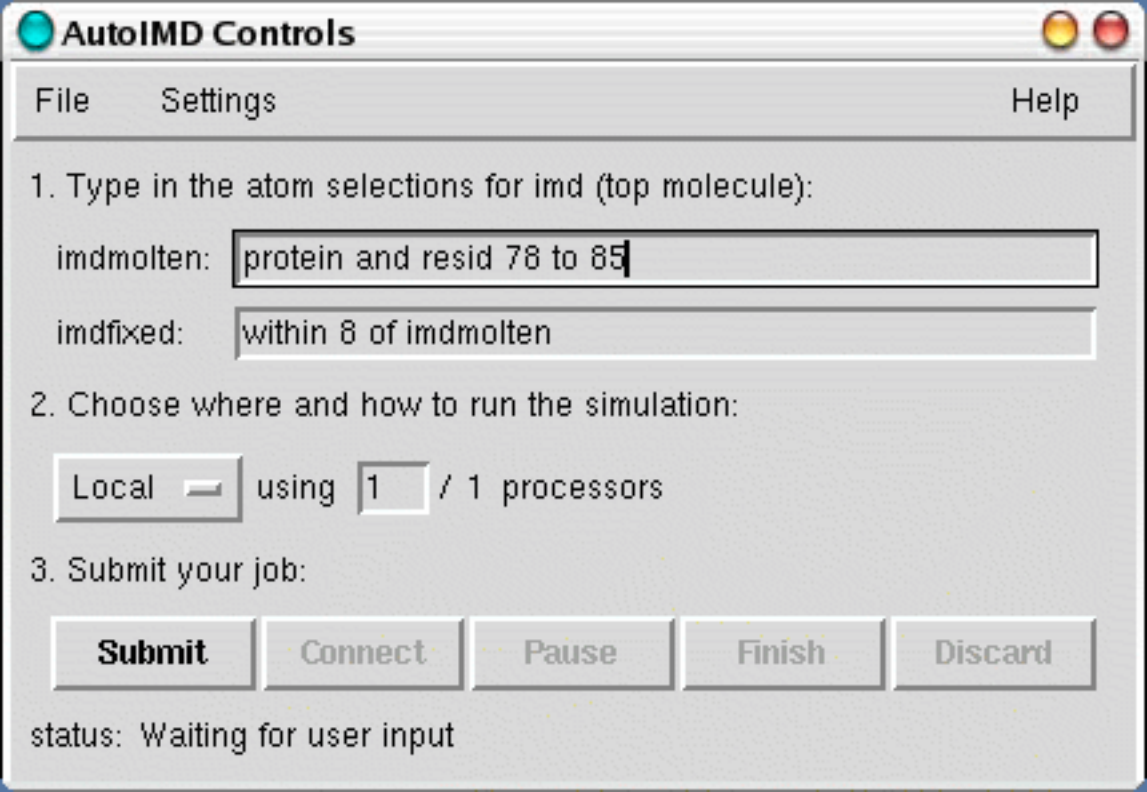
AutoIMD Controls
File
Settings
Help
1. Type in the atom selections for imd (top molecule):
imdmolten:
protein and resid 78 to 85
imdfixed:
within 8 of imdmolten
2. Choose where and how to run the simulation:
Local
using
1
/ 1 processors
3. Submit your job:
Submit
Connect
Pause
Finish
Discard
status: Waiting for user input