

A

View **Model building**

## Model building

## Imports

- import volumes
- import atomic structure
- import sequence

## Preprocess map

- xmipp3 - create 3d mask
- xmipp3 - local MonoRes
- xmipp3 - localdeblur sharpening
- xmipp3 - extract unit cell

## Initial model

- chimera - model from template

## Rigid fitting

- powerfit\_scipion - powerfit
- chimera - chimera rigid fit

## Flexible fitting

- phenix - real space refine
- ccp4 - coot refinement
- ccp4 - refrac

## Validation

- phenix - emringer
- phenix - molprobtity
- phenix - validation\_cryoem

## Tools-Calculators

- phenix - superpose pdbs
- atomstructutils - operator
- atomstructutils - convert\_sym
- chimera - chimera operate
- chimera - chimera restore session
- chimera - contacts

## Others

- xmipp3 - 3d bionotes

## Exports

- export to EMDB

B

Protocol Run: PowerfitProtRigidFit

Protocol: powerfit\_scipion - powerfit [Cite](#) [Help](#)

**Run**

Run name: powerfit\_scipion - powerfit [Edit](#) Comment: [Edit](#)

Run mode: ☒ Continue ☐ Restart [?](#) Host: localhost [?](#)

Parallel Threads: 4 [?](#) Use queue? ☐ Yes ☒ No [?](#)

Wait for: [?](#)

**Expert Level** ☐ Normal ☒ Advanced

**Input**

Input atomic structure to fit [Search](#) [Delete](#) [View](#) [?](#)

Input volume [Search](#) [Delete](#) [View](#) [Edit](#) [?](#)

Resolution (Å): 6.0 [?](#)

Angular step: 10.0 [?](#)

Number of models: 10 [?](#)

Apply Laplacian: ☐ Yes ☒ No [?](#)

Apply core weight: ☐ Yes ☒ No [?](#)

Other parameters for Powerfit: [?](#)

[Close](#) [Save](#) [Execute](#)