

Coot Crib Sheet

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1 Keyboard

1.1 Dialog Shortcuts

F5 Post Model/Fit/Refine dialog
F6 Post Go To Atom window
F7 Post Display Control Window
F8 Raster3D "Screenshot"

1.2 Previous/Next Residue

"Space" Next Residue
"Shift" "Space" Previous Residue

1.3 Closest Residue

"P" go to an atom of the closest residue (the "CA" atom if the residue has one)

1.4 Next NCS Chain

"O" - other NCS chain.

1.5 Previous/Next Rotamer

When in "Rotamer" mode, these keyboard short-cuts are available¹:

"." Next Rotamer
"," Previous Rotamer

1.6 Keyboard Chi Angles

Instead of pressing the buttons in the Chi Angles button box, you can use keyboard "1" for Chi1, "2" for Chi2 *etc.*

1.7 Keyboard Contouring

Use + or - to change the contour level

¹note: focus must be in the graphics window, not the Rotamer dialog

1.8 Keyboard Labelling

"l" to label closest atom

1.9 Keyboard Go To Residue

Ctrl-G then key in a residue number and (optionally) a chain-id and press Return

1.10 Keyboard Translation

Keypad 3 Push View (+Z translation)
Keypad . Pull View (-Z translation)

1.11 Keyboard Undo

Ctrl-Z Undo last modification
U Undo last move/navigation

1.12 Keyboard Zoom and Clip

N Zoom out
M Zoom in
D Slim clip
F Fatten clip

1.13 Crosshairs

C: cross-hairs

1.14 Skeleton

S: Generate skeleton around current point²

1.15 Continuous Rotate

I: Toggle continuous spin

²if a skeleton is being displayed

1.16 Baton Mode

B: toggle into baton rotate mode³

2 Mouse

Mouse actions are occasionally augmented with keyboard modifiers:

Left-mouse Drag	Rotate view
Ctrl Left-Mouse Drag	Translates view
Shift Left-Mouse Click	Label Atom
Right-Mouse Drag	Zoom in and out
Shift Right-Mouse Drag	Change clipping and Translate in Screen
	The movement is along orthogonal axes: up+right/down+left shifts in z, up+left/down+right changes the slab
Ctrl Shift Right-Mouse Drag	Rotate View about Screen Z
Middle-mouse Click	Centre on atom
Scroll-wheel Forward	Increase map contour level
Scroll-wheel Backward	Decrease map contour level

Intermediate (white) atoms can be dragged around by clicking on them:

Left-mouse Drag:	Move all intermediate atoms by linear shear
Left-mouse Drag	as above with
with "A" key:	non-linear shear
Left-mouse Drag	Move a single atom
with "Ctrl":	

3 Refinement Extras

Use "A" to define a residue range⁴ with a single-click. Useful in Refinement and Regularization.

- Click "Real Space Refine Zone"
- Click on an atom
- Press the "A" key

³rather than view rotate mode

⁴+/- *n* residues from the current residue