

Ctrl+Z Ctrl+Y Ctrl+H X Ctrl+F = Shift -

Hover mouse to change Element:
N/O/S/F..

Atom RMB-click

LM rotate, RM translate, MM scales. Shift-LM/RM moves coordinates. Drag on an atom to draw.

Shortcut	Action
Shift + </>	Zoom in/out
RMB-click on white space	Display menu
RMB-click on any atom	Display long menu
L/RMB	Left/Right mouse button

Shortcut	Action
LMB-drag on white space	Rotate all objects
Shift+ LMB-drag on white space	Rotate molecule relative to Ref/Prot
RMB-drag on white space	Translate all objects
Shift+RMB-drag on white space	Translate molecule relative to Ref/Prot

Shortcut	Action
LMB-click on atom + atom selected in drawing widget	Change atom type
LMB-click on atom + ring selected in drawing widget	Growing a ring
RMB-drag on atom	Move selected atom
Ctrl+RMB-drag	Z-clip the display

The screenshot displays the Molecule Editor interface. The main window shows a 3D ball-and-stick model of a complex organic molecule. Annotations with red lines point to specific features:

- "LMB-click to change bond order: 1/2/3" points to a bond in the lower part of the molecule.
- "Hover mouse to change bond order: 1/2/3" points to a bond in the upper part of the molecule.
- "LMB-drag to sprout a new atom or bond" points to a bond in the upper part of the molecule.
- "LMB-drag to rotate the smaller group" and "Shift-LMB drag to rotate the bigger group" point to a bond in the upper part of the molecule.

A context menu is open over the molecule, titled "Bond RMB-click". The menu items are:

- Undo (Ctrl+Z)
- Redo Change Element (Ctrl+Y)
- Select All (Ctrl+A)
- Invert Selection (Ctrl+I)
- Paste (Ctrl+V)
- Recenter View (Home)
- Delete (Del)
- Change Bond Order (highlighted)
 - Single 1
 - Double 2
 - Triple 3
- Save Image As...

On the left side, the "Elements" panel shows a grid of chemical elements (H, C, N, O, S, P, F, Cl, Br, I, B, Si) and rings (benzene, furan, pyrrole, thiophene). Below it, the "Select" mode is set to "Select Mode" (indicated by a radio button), with "Rotate Mode" and "Select Mode" options. A note says "Hold Ctrl" and "Select Mode".

On the right side, the "Actions" panel contains buttons for "Add H", "Del H", "Charge for pH7", "Add Fields", "Minimize", and "Optimize Alignment". Below these are icons for "Ref", "Prot", "ASite", "Ribb", "HBnds", "Meas", and "Clear".

At the bottom of the window, a status bar reads: "LM rotate, RM translate, MM scales. Shift-LM/RM moves coordinates. Drag on an atom to draw."

L/RMB	Left/Right mouse button	Shortcut	Action
		Ctrl+A	Select all molecule
		Ctrl+I	Invert selection
		Shift+RMB-drag on any atom of selection	Translate a selection of the molecule