## MarvinSketch cheat sheet

## General

| Wheel up and wheel down | Scrolls canvas vertically |
| :---: | :---: |
| Shift + Wheel up and wheel down | Scrolls canvas horizontally |
| Ctrl + Wheel up and wheel down | Zooms canvas in and out |
| Arrow keys or Ctrl <br> + Arrow keys | Scrolls canvas in the proper direction if no object is selected on the canvas |
| Arrow keys | Moves the selected object if an item is selected on the canvas. You can scroll the canvas with Ctrl + Arrow Keys in this case |
| Shift + Arrow keys | Move the selected object on the canvas in greater units |
| Del | Removes the selected element |
| Ctrl + A | Select All |
| Ctrl + Shift +1 | Invert selection |
| $\mathrm{Ctrl}+\mathrm{C} \text { or } \mathrm{Ctrl}+$ <br> Insert | Copy |
| Ctrl + D | Copy As OLE |
| $\mathrm{Ctrl}+\mathrm{K}$ | Copy As... |
| Ctrl + L | Copy As Smiles |
| $\mathrm{Ctrl}+\mathrm{M}$ | Copy As MRV |
| $\begin{aligned} & \text { Ctrl + X or Ctrl + } \\ & \text { Shift + Del } \end{aligned}$ | Cut |
| $\begin{aligned} & \text { Ctrl + V or Ctrl + } \\ & \text { Shift + Insert } \end{aligned}$ | Paste |
| Ctrl + Y | Redo |
| Ctrl +Z or Alt + <br> Backspace | Undo |
| Ctrl + L | Copy as SMILES |
| $\mathrm{Ctrl}+\mathrm{O}$ | File open (if available) |
| Ctrl +S | Save to file (if available) |
| Ctrl + Shift + S | Save as... (if available) |
| Ctrl + P | Print (if available) |
| Ctrl +E | Display Periodic Table dialog |
| $\mathrm{Ctrl}+\mathrm{N}$ | Create a new window |
| Ctrl + Del | Clear Desk |
| $\mathrm{Ctrl}+\mathrm{W}$ | Close current window |


| Ctrl + Q | Exit from the application |
| :---: | :---: |
| Ctrl + G | Create a Group |
| Ctrl +2 | Clean in 2D |
| Ctrl +3 | Clean in 3D |
| $\mathrm{Ctrl}+\mathrm{F}$ | Select conformer |
| Ctrl + T | Opens the Template Library. |
| $\mathrm{Ctrl}+\mathrm{R}$ | Checks and corrects chemical structures. |
| Ctrl + 1 | Places the Analysis box on the canvas. |
| Ctrl + Shift + N | You can view the name of the current structure, and enter a new name to be imported. |
| Ctrl + Shift + M | Open MarvinSpace |
| Ctrl + Shift + F9 | Guidelines |
| Shift + F9 | Grid |
| F5 | Exit transformation mode and return to Sketching mode. |
| F6 | Switch on the Zoom mode. |
| F7 | 3D rotation |
| F11 | Sets the visibility of the main menubar. |
| Space | Changes transformation mode from Drag to Rotate in 2D, Rotate in 2D to Rotate in 3D, while Rotate in 3D to Drag. |
| - | Negative charge |
| + | Positive charge |
| 1 | Single bond |
| 2 | Double bond |
| 3 | Triple bond |
| 4 | Aromatic bond |
| 5 | Single up bond (active only in 2D) |
| 6 | Single down bond (active only in 2D) |
| 7 | Single up or down bond |
| 12 | Single or double bond |
| 14 | Single or aromatic bond |
| 24 | Double or aromatic bond |
| 0 | Any bond |
| * | Any atom |


| Q | Heteroatom |
| :--- | :--- |
| C | Carbon |
| N | Nitrogen |
| H | Hydrogen |

For the other elements, type the mark of the element, for example, Cl for chlorine (case insensitive).

Au,Ag,Pt...: Atom List can be defined by typing chemical symbols separated by commas (case insensitive).
!Au,Ag,Pt...: NOT List can be defined by starting the atom list with an exclamation mark (case insensitive).

Source: MarvinSketch docs

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More information: defkey.com/marvinsketch-
shortcuts

