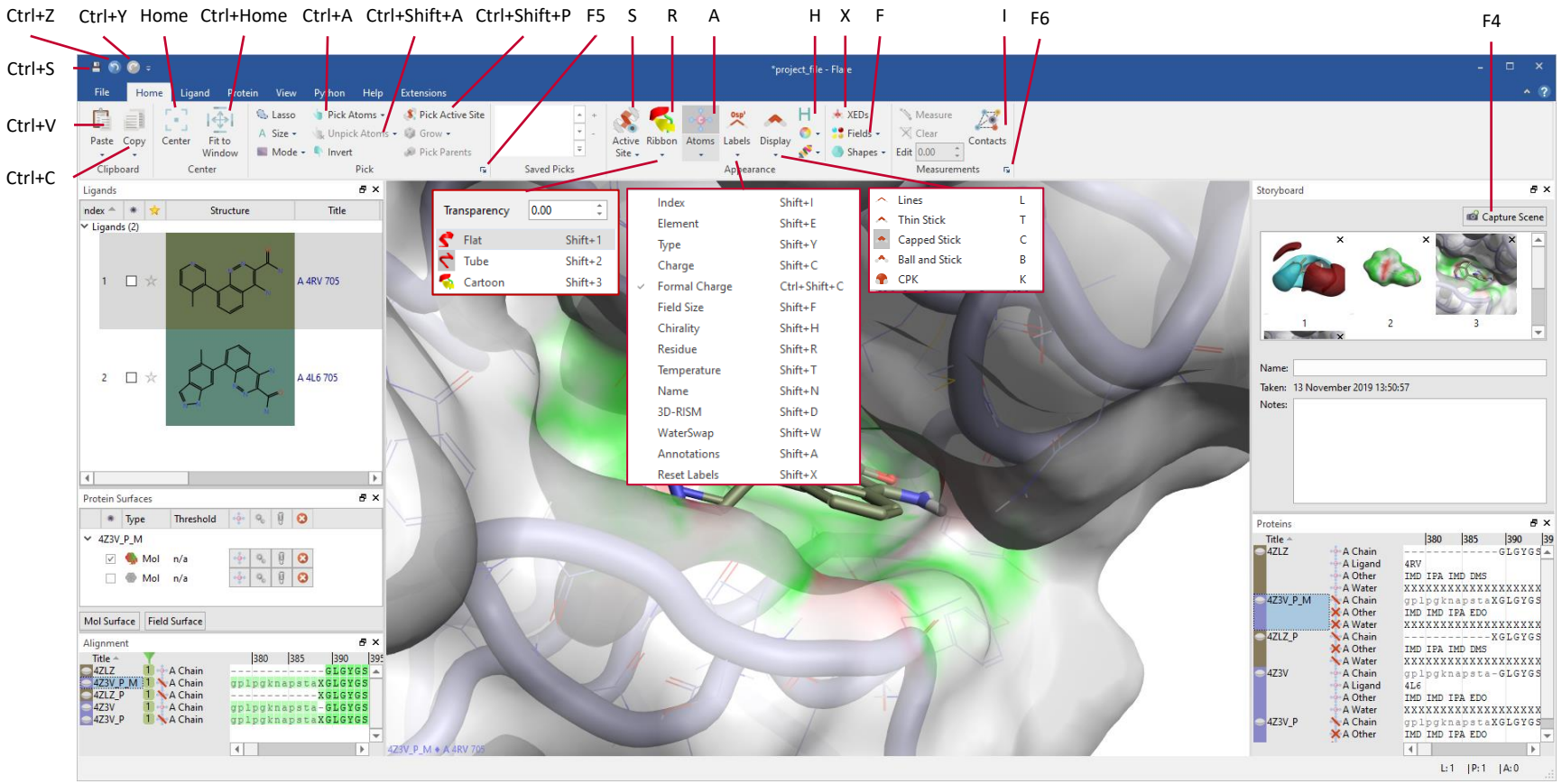
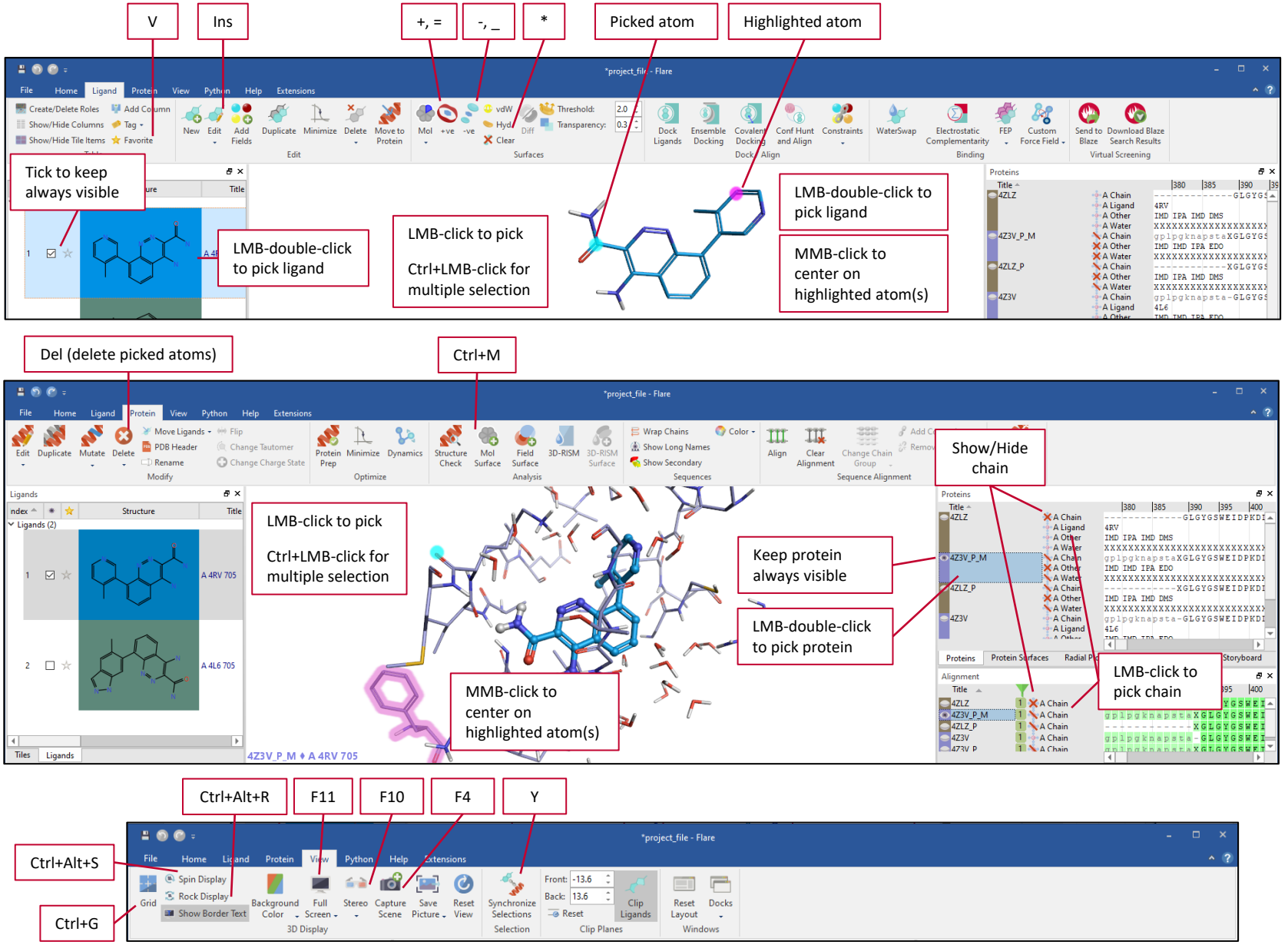


Shortcut	File operation
F1	Show manual
Ctrl+N	New Project
Ctrl+O	Open File
Ctrl+S	Save Project
Ctrl+Shift+S	Save Project As…
Alt+F4 Ctrl+Q	Exit (Windows) Quit (Linux)

Shortcut	Molecule operation
LMB-double-click on ligand atom	Pick ligand
LMB-double-click on water atom	Pick water
LMB-double-click on protein atom	Pick residue
Ctrl+C	Copy 3D
Ctrl+V	Paste

Shortcut	3D Window operation
Ctrl+G	Grid
Ctrl+Alt+R	Rock display
Ctrl+Alt+S	Spin display
F11	Full screen
Y	Synchronize ligand/ protein selection
Shift+LMB-drag	Clip front plane
Shift+RMB-drag	Clip back plane

To activate right-click menus, RMB-click on:
Atom in 3D window
Ligand in Ligands table
Protein/Chain/Residue in Proteins & Alignment Tables
Water/other molecule in Proteins table


V
Ins
+, =
-, _
*
Picked atom
Highlighted atom
Tick to keep always visible
LMB-double-click to pick ligand
LMB-click to pick
Ctrl+LMB-click for multiple selection
LMB-double-click to pick ligand
MMB-click to center on highlighted atom(s)
Del (delete picked atoms)
Ctrl+M
Show/Hide chain
LMB-click to pick
Ctrl+LMB-click for multiple selection
Keep protein always visible
LMB-double-click to pick protein
MMB-click to center on highlighted atom(s)
LMB-click to pick chain
Ctrl+Alt+R
F11
F10
F4
Y
Ctrl+Alt+S
Ctrl+G