Analysis V3 shortcuts

Mouse events

Action	Shortcut
Context menu	Right click
Pan	Left drag
Zoom into area	Shift right drag
Zoom into area	Middle drag
Zoom in/out around cursor	Scroll wheel in display
Zoom in/out for selected axis	Scroll wheel in axis active area
Select new	Left click
Add one to Selection	Ctrl left click
Select in area and add	Shift left drag
Pick one peak at position	Ctrl shift left click
Pick peaks in area	Ctrl shift left drag
Move module	Drag&drop purple bar
Move whole window	Alt Left drag
Pop out module (manage as windows)	Double click on Purple bar
Make new object	Double click <New> in Sidebar
View/Edit object	Double click object in Sidebar

Keyboard shortcuts

Action	Shortcut
Graphics :	
Delete selected	Del (Fn+Backspace)
Toggle Grid	GS
Toggle Crosshairs	CH
Toggle Horizontal Trace	TH
Toggle Vertical Trace	TV
Increase Trace Scale	TU
Decrease Trace Scale	TD
Create Mark at cursor position	MK
Clear All Marks	MC
Navigate to Peak position	FP
Navigate to NmrResidue position	FN
Copy Assignments	CA
Project :	
New project	PN
Open project	PO
Load Spectrum	LS
Load Data	LD

Save	**PS**
Save as	**SA**
Undo	**Ctrl+Z**
Redo	**Ctrl+Y**
Quit	**QT**
Spectrum :	
Spectrum Groups	***SS***
Set Experiment Types	**ET**
Pick Peaks	**PP**
Peak Integration (1D)	**IT**
Make Projection	**PJ**
Toggle Phasing Console	**PC**
Add Phasing Traces	**PT**
Remove Phasing Traces	**PR**
Molecules :	
Create new molecule	**CM**
Reference Chemical Shifts	**RC**
Assign :	
Set up NmrResidues	**SN**
Pick and assign	**PA**
Backbone assignment	**BB**
Peak Assigner	**AA**
Residue information	**RI**
View :	
New Blank Display	**ND**
NmrAtom Chemical Shift Table	**CT**
NmrResidue Table	**NT**
Peak Table	**LT**
Sequence Graph	**SG**
Atom Selector	**AS**
Python Console	**PY**
Macro :	
Run macro	**RM**
Reserved for user macros	**Q1, Q2, ...Q0**
Reserved for user macros	**W1, W2, …W0**
Reserved for user macros	**U1, U2, …U0**