Analysis V3beta2 shortcuts

Mouse events

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/Remove one to/from Selection	Ctrl/Cmd left click
Select in area and add	Ctrl/Cmd left drag
Pick one peak at position	Ctrl/Cmd shift left click
Pick peaks in area	Ctrl/Cmd shift left drag
Move single peak	Ctrl/Cmd middle 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	Sidebar
View/Edit object	Sidebar

Keyboard shortcuts

	Shortcut	Related
Graphics :		
Delete selected	Del (Fn+Backspace)	
Toggle Grid	GS	CH, TA, PC
Toggle Crosshairs	CH	GS, TA, PC
Toggle Horizontal Trace	TH	TV, TU, TD
Toggle Vertical Trace	TV	TH, TU, TD
Increase Trace Scale	TU	TH, TV, TD
Decrease Trace Scale	TD	TH, TV, TU
Create Mark at cursor position	MK	MC
Clear All Marks	MC	MK
Navigate to Peak position	FP	FN
Navigate to NmrResidue position	FN	FP
Copy PeakList	CL	CP
Copy Peaks	CP	CL
Clear Current Selected Peaks	CZ	
Refit Selected Peaks	RP	SE
Snap to Extremum Current Selected Peaks	SE	RP
Select All Peaks	Ctrl/Cmd+A	
Set Phasing Pivot to cursor position	PV	PC, PT, TR

Add Integral in 1D plot	**I1**	
Add Phasing Trace	**PT**	**PC, PV, TR**
Remove Phasing Trace	**TR**	**PC, PV, PT**
Switch Mouse Mode	**MM**	
Zoom in to Spectrum	**ZI**	**ZO, ZS, ZR**
Zoom out of Spectrum	**ZO**	**ZI, ZS, ZR**
Toggle Share Y Axis Across Display	**TA**	**GS, CH, PC**
Cycle Peak Labelling	**PL**	**PS**
Cycle Peak Symbols	**PS**	**PL**

Project :

New project	**Ctrl/Cmd+N**
Open project	**Ctrl/Cmd+O**
Load Data	**LD**
Save	**Ctrl/Cmd+S**
Save as	**SA**
Import Nef File	**IN**
Export Nef File	**EX**
Undo	**Ctrl/Cmd+Z**
Redo	**Ctrl/Cmd+Y**
Quit	**Ctrl/Cmd+Q**

Spectrum :

Load Spectrum	**LS**	**SS**
Show Spectrum Groups	**SS**	**LS**
Set Experiment Types	**ET**	
Pick 1D Peaks	**P1**	**PP**
Pick ND Peaks	**PP**	**P1**
Copy PeakList	**CL**	**CP**
Copy Peaks	**CP**	**CL**
Make Projection	**PJ**	
Print to File	**PR**	

Molecules :

Reference Chemical Shifts	**RC**

Assign :

Set up NmrResidues	**SN**
Pick and assign	**PA**
Backbone assignment	**BB**
Peak Assigner	**AA**
Assignment Inspector	**AI**
Residue information	**RI**

View :

NmrAtom Chemical Shift Table	**CT**	
NmrResidue Table	**NT**	
Peak Table	**LT**	
Integral Table	**IT**	
Restraint Table	**RT**	
Structure Table	**ST**	
Sequence Graph	**SG**	
Atom Selector	**AS**	
Chemical Shift Mapping	**CM**	
Show Sequence	**SQ**	
Notes Table	**NO**	
Python Console	**PY**	
For current strip		
Show/Hide Toolbar	**TB**	
Show/Hide Phasing Console	**PC**	**PV, PT, TR**
Move to Next Spectrum (and Hide Current)	**TabTab**	
Move to Previous Spectrum (and Hide Current)	**TabQ**	
Toggle all Spectra on Current Display	**TabA**	
Store Zoom	**ZS**	**ZI, ZO, ZR**
Restore Zoom	**ZR**	**ZI, ZO, ZS**
Flip X-Y Axis	**XY**	**XZ, YZ**
Flip X-Z Axis	**XZ**	**XY, YZ**
Flip Y-Z Axis	**YZ**	**XY, XZ**

Macro :

Define User Shortcuts	**DU**
Reserved for user macros	**Q0, Q1, ...Q9**
Reserved for user macros	**W0, W1, ...W9**
Reserved for user macros	**Y0, Y1, ...Y9**