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

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	