

# Analysis V3<sub>beta2</sub> shortcuts

## Mouse events

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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

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	Shortcut	Related
<i>Graphics :</i>		
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 or +	ZO, ZS, ZR
Zoom out of Spectrum	ZO or -	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
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*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

<b>Restraint Table</b>	<b>RT</b>	
<b>Structure Table</b>	<b>ST</b>	
<b>Sequence Graph</b>	<b>SG</b>	
<b>Atom Assigner</b>	<b>AS</b>	
<b>Chemical Shift Mapping</b>	<b>CM</b>	
<b>Notes Table</b>	<b>NO</b>	
<b>Python Console</b>	<b>Space-Space</b>	
<i>For current strip</i>		
<b>Show/Hide Toolbar</b>	<b>TB</b>	
<b>Show/Hide Phasing Console</b>	<b>PC</b>	<b>PV, PT, TR</b>
<b>Move to Next Spectrum (and Hide Current)</b>	<b>TabTab</b>	
<b>Move to Previous Spectrum (and Hide Current)</b>	<b>TabQ</b>	
<b>Toggle all Spectra on Current Display</b>	<b>TabA</b>	
<b>Store Zoom</b>	<b>ZS</b>	<b>ZI, ZO, ZR</b>
<b>Restore Zoom</b>	<b>ZR</b>	<b>ZI, ZO, ZS</b>
<b>Flip X-Y Axis</b>	<b>XY</b>	<b>XZ, YZ</b>
<b>Flip X-Z Axis</b>	<b>XZ</b>	<b>XY, YZ</b>
<b>Flip Y-Z Axis</b>	<b>YZ</b>	<b>XY, XZ</b>
 <i>Macro :</i>		
<b>Define User Shortcuts</b>	<b>DU</b>	
<b>Reserved for user macros</b>	<b>Q0, Q1, ...Q9</b>	
<b>Reserved for user macros</b>	<b>W0, W1, ...W9</b>	
<b>Reserved for user macros</b>	<b>Y0, Y1, ...Y9</b>	