

















Signal Tool Set

Lets you graphically work with a chromatogram.




-  Align the x-axis of multiple signals
-  Align the y-axis of multiple signals
-  Reset the alignment of your signals
-  Create a 3D overlay of signals
-  Mirror signals
-  Subtract signals
-  Integrate all chromatograms
-  Smooth all signals

The following cursor tools are for working with alignment markers and the signal displays.










-  Set a time reference point
-  Move a time reference point
-  Delete a time reference point
-  Delete object from chromatogram display
-  Display signals overlaid
-  Display signals separated
-  Display the signals in full scale
-  Display all signals in the same scale

Spectral Tool Set

Lets you perform spectral evaluation tasks.

-  Set spectral options
-  Open a spectral library
-  Save a spectral library

The following cursor tools are for spectral selection tasks.

-  Select spectrum
-  Select peak apex spectrum
-  Select average spectrum
-  Average spectra across a peak
-  Select peak for purity evaluation
-  Select 1st reference spectrum
-  Select 2nd reference spectrum
-  Zoom in on MS window
-  Extract ion chromatogram