



ms.epfl.ch: The EPFL mstoolbox

EPFL

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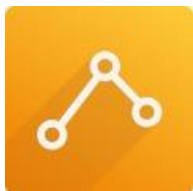
Free. Open access. No installation needed just Chrome as browser. Calculations done locally without transfer of data to our servers



Theoretical calculations

Applications for any kind of MS calculations, to generate isotopic distributions or find a molecular formula from any experimental value.

Start



Apm²s for peptides & small proteins

Application for peptide and small protein characterization. Automatic peak picking, post calibration, fragment ion assignment and resulting fragmentation map.

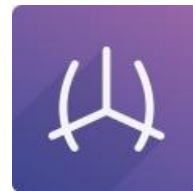
Start



Aom²s for oligonucleotides

Application for DNA/RNA characterization. Automatic peak picking, post calibration, fragment ion assignment and resulting fragmentation map.

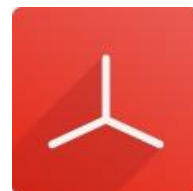
Start



MSPolyCalc for polymers

Application for interpretation of polymers mass spectra. Determination of polymer distribution, molecular formula and similarity calculation with experimental data.

Start



Complex mixtures of compounds

Application for complex data analysis. You specify groups that, when combined, generate multiple combinations that will be matched to your experimental spectra.

Start



Custom database / contaminants

Application to create your own database of compounds to be tracked in your experimental spectrum. Track your background ions with our Easycont DB.

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