



Version 2

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# LC-MS/MS Label-Free Proteomic Data Analysis Parameters V.2

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Works for me

[dx.doi.org/10.17504/protocols.io.bfshjnb6](https://dx.doi.org/10.17504/protocols.io.bfshjnb6)

VU Biomolecular Multimodal Imaging Center

Human BioMolecular Atlas Program (HuBMAP) Method Development Community



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## ABSTRACT

List of parameters and settings for searching label free proteomic data in MaxQuant version 1.6.7.

- 1 Label-free proteomic samples were searched using MaxQuant version 1.6.7.
- 2 *Group Specific Parameter settings included:*
  - Standard
  - Multiplicity = 1
  - Variable modifications: Oxidation (M); Acetylation (Protein N-term), Carbamidomethyl (C)
  - Fixed modifications: Carbamidomethyl (C)
  - Max number of modifications per peptide: 5
  - Instrument settings: Orbitrap (default settings)
  - Digestion: Specific, Trypsin /P
  - Max missed cleavages: 2
  - LFQ: None
- 3 *Global Parameters included:*
  - Database: UniProt Human Proteome (UP000005640, 9606) Reviewed, downloaded July 30, 2019
  - Min peptide length: 7
  - Max peptide mass: 4600 Da
  - MS/MS analyzer: Default settings
  - Identification: Default settings