

Senior Scientist, Cheminformatics

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The department of Computational & Structural Chemistry is seeking a creative and self-motivated cheminformatics data scientist to join our team. The successful candidate will work closely with discovery scientists to drive towards a 'design first' culture, supporting project medicinal chemists with the proper predictive models and workflows. In addition to leveraging rapid advances in machine learning towards the interpretation of complex *in vitro* and *in vivo* biological data, the candidate will make informed decisions to de-risk safety liabilities and understand the mechanisms of toxicity at both lead selection and throughout the lifecycle of a discovery project.

Responsibilities include and are not limited to the following:

- Drive a 'predict first' culture in hit-to-lead and lead optimization
- Develop and apply computational tools, including machine learning
- Characterize the toxicology profile of lead molecules and mitigate potential adverse effects
- Design and implement deeper informatics methods to address specific project challenges

Education Minimum Requirement:

Ph.D. (or close to completion) in Chemistry, Mathematics, Physics, Computer Science, Biology or commensurate career experience.

Required Experience and Skills:

- Competency in programming, scripting, and databases (e.g. Python, R, MySQL)
- Background in methods for cheminformatics, statistical analysis, visualization tools, and machine learning techniques
- Strong communication skills, both written and oral
- Ability to work well in multidisciplinary teams

Preferred Experience:

- Experience with workflow applications and data analytics software (Pipeline Pilot, Knime, Spotfire, ...)
- External scientific visibility through multiple publications and/or presentations
- A basic understanding of the drug discovery processes and medicinal chemistry

To apply please use the link below

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