



Delivering a flexible, cost-effective and high quality computer-aided drug design consultancy service

At CADD Matters, we specialise in providing computational chemistry/chemoinformatics solutions tailored to meet your business needs.

We offer a wide range of services including:

Scientific

- ❖ Project Support for any stage of the drug discovery pipeline. We are highly experienced in the application of both protein, ligand-based and ADMET modelling approaches or a combination thereof; to kinases, proteases, ion channels and GPCRs.
- ❖ Advice and guidance on setting up a computational chemistry/ informatics platform in-house (software, hardware & costs).
- ❖ Evaluation of existing and novel computational software ensuring your tools meet your needs.
- ❖ We can enhance existing corporate collections, build and maintain databases and mine large datasets.

Supporting Services

- ❖ Training of chemists and computational chemists. This includes delivering a hand-on-training program for desktop tools for chemists. Such as 3D visualisation tools, property calculations through to seminars on specified computational chemistry methods e.g. Virtual screening, SBDD and property calculations.
- ❖ Professional guidance and interview support with recruitment for computational chemistry vacancies.
- ❖ Generation of powerpoint presentations, images and videos to promote your scientific competitive edge for marketing purposes.
- ❖ Cover for maternity leave, illness and vacant positions. We can also provide additional resources to complement your existing team at times of peak demand.

Our Business Model

- ❖ We can work on-site and/or remotely.
- ❖ We have access to vendor's software including MOE from CCG as well as freeware tools to support your business either on short term or longer term projects.
- ❖ We charge on a simple daily basis.

Don't hesitate to contact us for an initial consultation at info@cadd-matters.com

