



Computational Chemist Centre for Protein Degradation Candidate Information

July 2026

The Institute of Cancer Research

About our organisation

We are one of the world's most influential cancer research institutes with an outstanding record of achievement dating back more than 100 years. We are world leaders in identifying cancer genes, discovering cancer drugs and developing precision radiotherapy. Together with our hospital partner The Royal Marsden, we are rated in the top four centres for cancer research and treatment worldwide.

As well as being a world-class institute, we are a college of the University of London. We came top in the league table of university research quality compiled from the Research Excellence Framework (REF 2014).

We have charitable status and rely on support from partner organisations, charities, donors and the general public.

We have more than 1000 staff and postgraduate students across three sites – in Chelsea and Sutton.

Division of Cancer Therapeutics

The Centre for Cancer Drug Discovery is a major hub within the Division of Cancer Therapeutics and a multidisciplinary 'bench to bedside' centre, comprising approximately 160 chemistry and biology scientists dedicated to the discovery and development of novel therapeutics for the treatment of cancer. We implement innovative drug discovery technologies, discover novel mechanism-based drugs, and develop these as rapidly as possible from the laboratory through to hypothesis-testing early clinical trials. These activities are carried out in highly focused multi-disciplinary project teams analogous to those in a biotechnology company, with patient benefit as the primary driver. We publish our work extensively and have a large network of collaborations with academia, biotechnology companies, and the pharmaceutical industry.

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Centre for Protein Degradation

The ICR established the Centre for Protein Degradation (CPD) in 2022 to accelerate targeted protein degradation-based drug discovery – an innovative approach utilising drug-induced degradation of harmful proteins via the cell's endogenous protein disposal system. Our aim is to discover novel molecular glue degraders and PROTACs for the treatment of the most challenging cancers. Based at the ICR's Centre for Cancer Drug Discovery, the CPD benefits from state-of-the-art drug discovery platforms – biology, medicinal and in silico chemistry, biophysics, structural biology, proteomics, bioinformatics, as well as strong cancer biology and clinical expertise within the wider ICR and the Royal Marsden Hospital. We also have established collaborations with biotechnology companies and the pharmaceutical industry.

In Silico Medicinal Chemistry, Medicinal Chemistry Team 3

The *In Silico* Medicinal Chemistry group comprises 4-6 postdoctoral and staff grade in silico scientists focussed on the application of cutting edge computational chemistry techniques and data analysis to accelerate our drug discovery. Group members work in close collaboration with chemists and biologists on projects at all stages of drug discovery, from early target validation and hit finding, through to lead optimisation and candidate selection. To achieve this, the *In Silico* Medicinal Chemistry group applies a wide range of approaches including structure- and ligand-based virtual screening, molecular dynamics, generative AI, property prediction, library design and data analysis. All necessary technologies, including high-performance computing and an integrated data-platform, are in place to make this possible. The group works closely with in-house experts in structural chemistry, structural biology, biological assays, DMPK and bioinformatics to inform their research. Many projects also involve scientific collaborations with external academic and commercial partners.

Our mission
is to make the
discoveries that
defeat cancer.

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Our values

The ICR has a highly skilled and committed workforce, with a wide variety of roles, each requiring different skills. But whether you work as a researcher, or work as part of our corporate team, your work and behaviour is underpinned by these six values. They are what bring us together as one team - as 'One ICR'.



Pursuing excellence

We aspire to excellence in everything we do, and aim to be leaders in our field.



Acting with Integrity

We promote an open and honest environment that gives credit and acknowledges mistakes, so that our actions stand up to scrutiny.



Valuing all our people

We value the contribution of all our people, help them reach their full potential, and treat everyone with kindness and respect.



Working together

We collaborate with colleagues and partners to bring together different skills, resources and perspectives.



Leading innovation

We do things differently in ways that no one else has done before, and share the expertise and learning we gain.



Making a difference

We all play our part, doing a little bit more, a little bit better, to help improve the lives of people with cancer.

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Job description

Department / division:	Centre for Protein Degradation, Division of Cancer Therapeutics
Pay grade / staff group:	Analytical Scientist 3 / Academic
Hours / duration:	Full time (35 hours per week), Monday to Friday. Fixed term contract until the 31-December-2027 in the first instance, with the potential to extend subject to funding.
Reports to:	Dr Andrea Scarpino, Lead Computational Chemist, <i>In Silico</i> Medicinal Chemistry
Main purpose of the job:	To advance the design and discovery of novel protein degraders in the ICR Centre for Protein Degradation through innovative computational chemistry methods.

Key Roles and Responsibilities

Collaborate within multidisciplinary project teams to advance projects in the Centre for Protein Degradation.
Apply structure- and ligand-based approaches to guide compound design and prioritisation.
Design, mine and interrogate compound libraries to extract actionable SAR.
Develop and deploy ternary complex modelling workflows to support the design of protein degraders.
Present scientific findings to internal project teams through effective data analysis and visualisation.
Evaluate and assess novel protein targets for ligandability and tractability across TPD modalities.
Contribute to peer-reviewed publications and conference presentations.
Maintain up-to-date knowledge of the latest developments in computational chemistry and related fields.
Contribute to expanding computational chemistry capabilities within the <i>In Silico</i> Medicinal Chemistry team.
Provide guidance to scientists from different disciplines on the appropriate use of computational tools.
Keep accurate electronic records enabling the work to be followed and reproduced by others.

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General

All staff must ensure that they familiarise themselves with and adhere to any ICR policies that are relevant to their work and that all personal and sensitive personal data is treated with the utmost confidentiality and in line with the General Data Protection Regulations

Any other duties that are consistent with the nature and grade of the post that may be required.

To work in accordance with the ICR's Values.

To promote a safe, healthy and fair environment for people to work, where bullying and harassment will not be tolerated.

This job description is a reflection of the present position and is subject to review and alteration in detail and emphasis in the light of future changes or development.

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Person specification

Education and Knowledge

PhD in computational chemistry, cheminformatics, or closely related field.	Essential*
Wide-ranging knowledge and experience in the use of computational chemistry methods.	Essential
Familiarity with concepts in medicinal chemistry, drug discovery, targeted protein degradation.	Desirable

**as a minimum requirement candidates must have submitted their thesis by the start date of their employment and awarded their PhD within the six month probationary period.*

Skills

Highly motivated with strong desire to achieve scientific excellence.	Essential
Proactive approach with excellent time management skills and ability to work on different projects simultaneously.	Essential
Excellent written and oral scientific communication.	Essential
Ability to apply structure- and ligand-based methods for drug design (e.g. docking, QSAR, pharmacophore modelling).	Essential
Ability to design, mine and interrogate compound libraries to extract actionable SAR and drive multi-parameter compound optimisation.	Essential
Familiarity with molecular modelling software packages (e.g. Schrödinger, MOE, OpenEye).	Essential
Ability to set up, run and analyse molecular dynamics simulations.	Essential
Proficiency in Python and cheminformatics packages (e.g. RDKit).	Essential
Familiarity with AI/ML methods for generative design and molecular property prediction.	Desirable

Experience

Post-PhD drug discovery experience in pharma/biotech or academia.	Essential
Experience working within multidisciplinary project teams.	Essential
Track record of creative problem solving using a wide variety of computational methods.	Essential
Experience with large-scale virtual screening for hit discovery and optimisation.	Desirable
Prior hands-on experience using computational methods for targeted protein degradation (e.g. PPI and/or ternary complex modelling).	Desirable
Experience with relative binding free energy calculations (e.g. FEP).	Desirable
Experience working with HPC clusters and job scheduling managers (e.g. SLURM).	Essential
Track record of computational chemistry publications in peer-reviewed journals.	Desirable

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Benefits

We offer a fantastic working environment, great opportunities for career development and the chance to make a real difference to defeat cancer. We aim to recruit and develop the best – the most outstanding scientists and clinicians, and the most talented professional and administrative staff.

The annual leave entitlement for full time employees is 28 days per annum on joining. This will increase by a further day after 2 years' and 5 years' service.

Staff membership to the Universities Superannuation Scheme (USS) is available. The USS is a defined benefit scheme and provides a highly competitive pension scheme with robust benefits. The rate of contributions is determined by USS and details of the costs and benefits of this scheme can be found on their website. If staff are transferring from the NHS, they can opt to remain members of the NHS Pension Scheme.

We offer a range of family friendly benefits such as flexible working, a parents' group, and a maternity mentoring scheme. Other great benefits include interest free loans for discounted season tickets for travel and bicycle purchases, access to the NHS discounts website, a free and confidential Employee Assistance Programme which offers a range of well-being, financial and legal advice services, two staff restaurants, and access to a gym and sporting facilities at our Sutton site.

Further information

You may contact Dr Andrea Scarpino for further information by emailing andrea.scarpino@icr.ac.uk. This job description is a reflection of the current position and is subject to review and alteration in detail and emphasis in the light of future changes or development.