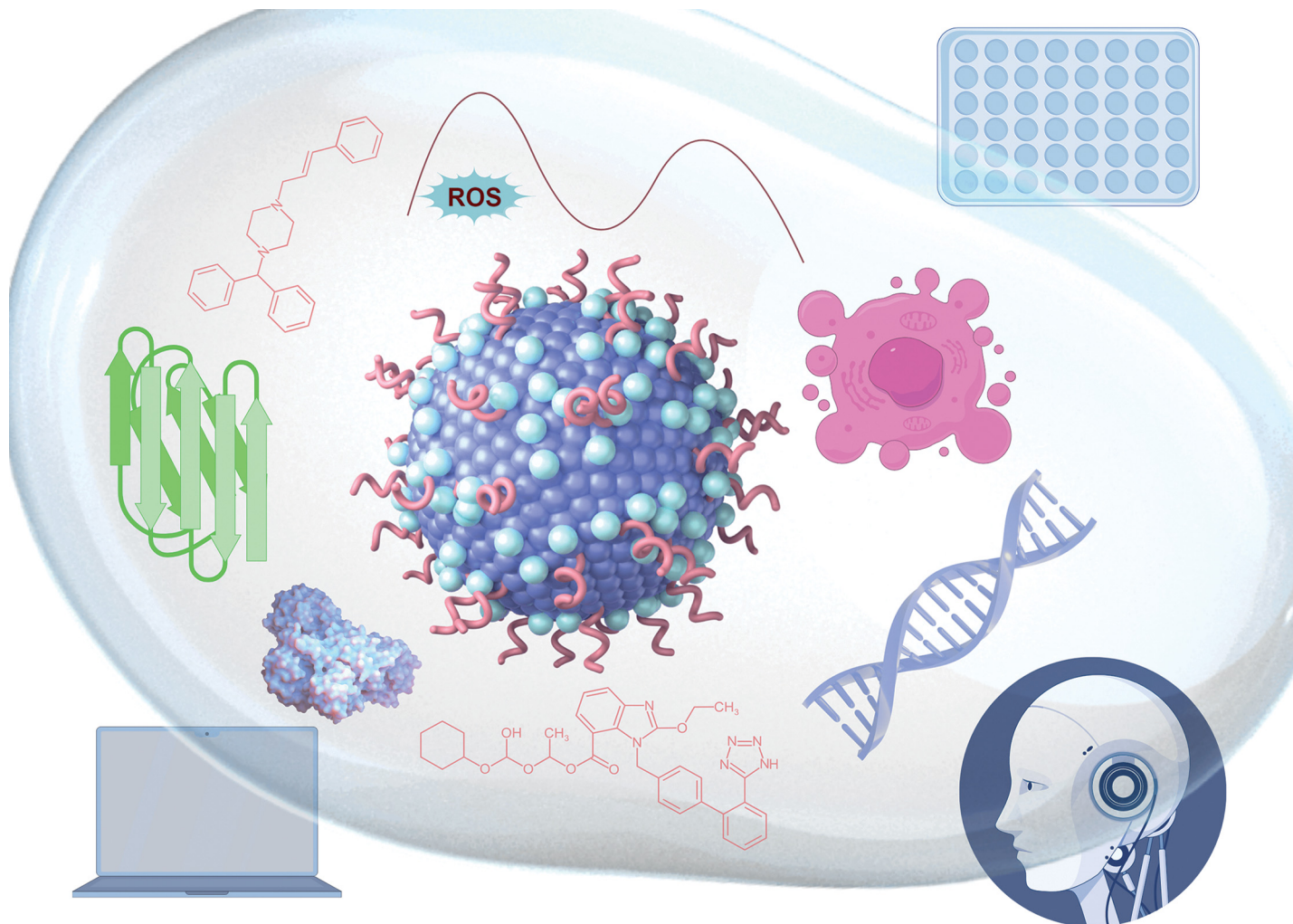


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Showcasing research from Professor Xingfa Gao's laboratory, National Center for Nanoscience and Technology, Beijing, China.

Computer-aided nanodrug discovery: recent progress and future prospects

Computational techniques, using either traditional or machine learning methods, could provide valuable insights into how the interfacial bio-nano interactions affect the therapeutic efficacy of nanodrugs and have a promising future in accelerating nanodrug discovery. Cover art credit: "Figdraw".

As featured in:



See Xingfa Gao *et al.*,
Chem. Soc. Rev., 2024, **53**, 9059.