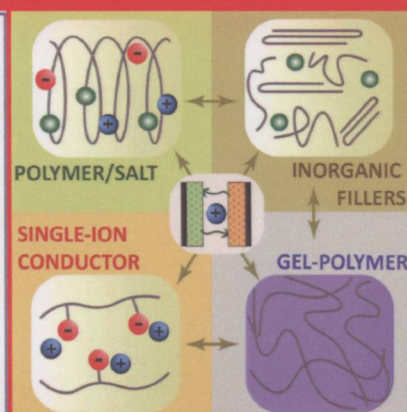
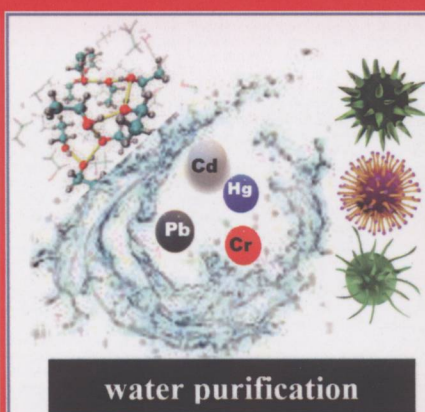
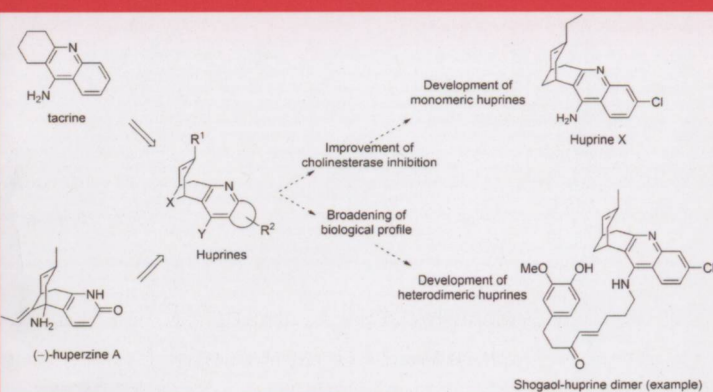
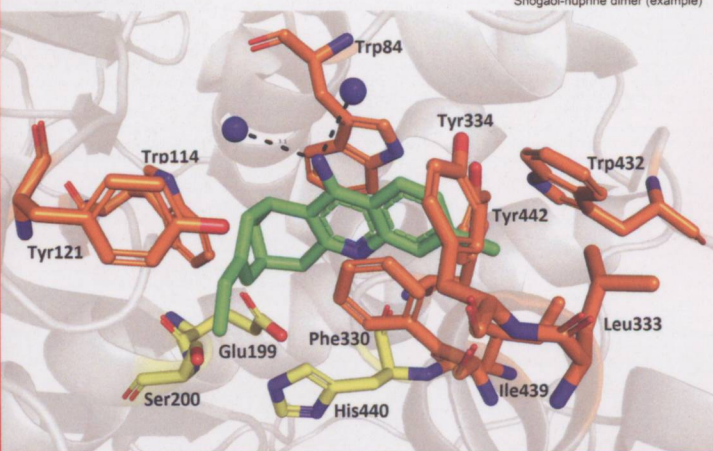


Успехи химии



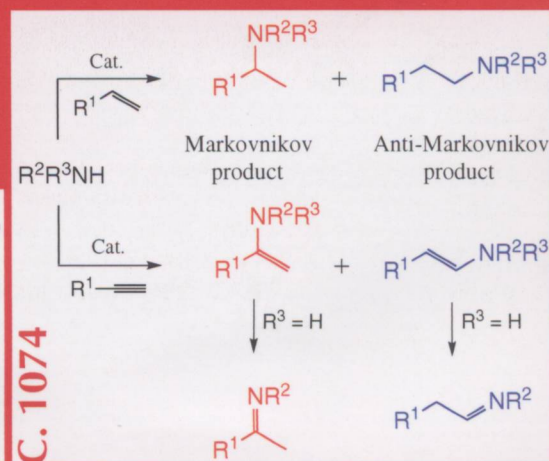
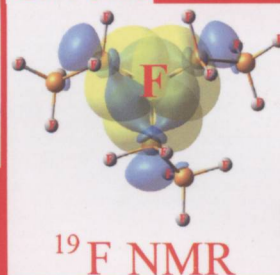
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C. 999

C. 1040



Обзорный журнал по химии

Том 89

Номер 10

2020

стр. 999–1155

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Huprines — an insight into the synthesis and biological properties

999

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The overlapping of tacrine and (–)-huperzine A templates yielded a family of highly potent cholinesterase inhibitors, so-called huprines. A relatively easy access to these compounds led to the development of dozens of huprine derivatives allowing to draw structure–activity relationship mainly for acetylcholinesterase and butyrylcholinesterase enzymes, but also with application to other biological targets of interest. An extension of their pharmacological profile is commonly associated with huprine scaffold binding to some other pharmacophores that yield high-molecular-weight heterodimers. The main purpose in developing the huprine family is related to Alzheimer's disease therapy. However, these compounds are also interesting lead structures in the treatment of other disorders, such as Myasthenia gravis, African trypanosomiasis, malaria, and prion diseases. The present review provides a rationale behind the development of huprines, detailed synthetic routes leading to different classes of huprines, and a thorough discussion of their potential pharmacological applications.

Bibliography — 174 references.

Computational aspects of ¹⁹F NMR

1040

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This review covers most recent advances in computation of ¹⁹F NMR chemical shifts and spin–spin coupling constants involving ¹⁹F nucleus calculated at different levels of theory. It deals mainly with electronic and stereochemical effects influencing these parameters, with a special emphasis on practical applications of such calculations. For ¹⁹F NMR chemical shifts, a good deal of attention is focused on less common solid state studies, in addition to much more popular liquid state data. For spin–spin coupling constants, the main interest is concentrated on the most popular ¹⁹F–¹H, ¹⁹F–¹³C and ¹⁹F–¹⁹F couplings providing marked stereochemical behaviour.

Bibliography — 149 references.

Catalysis and regioselectivity in hydrofunctionalization reactions of unsaturated carbon bonds. Part II. Hydroamination

1074

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This review continues consideration of the regioselectivity problem in the catalyzed hydrofunctionalization of unsaturated organic compounds and addresses hydroamination of unsaturated hydrocarbons. Particular Sections deal with reactions of alkenes, alkynes, allenes and dienes. It is shown that the selectivity of hydroamination depends on the natures of the reactants and the catalyst. Reaction conditions are described; in some cases, the reaction mechanisms are discussed. Reactions in which divergent regioselectivity is possible are noted.

Bibliography — 249 references.

