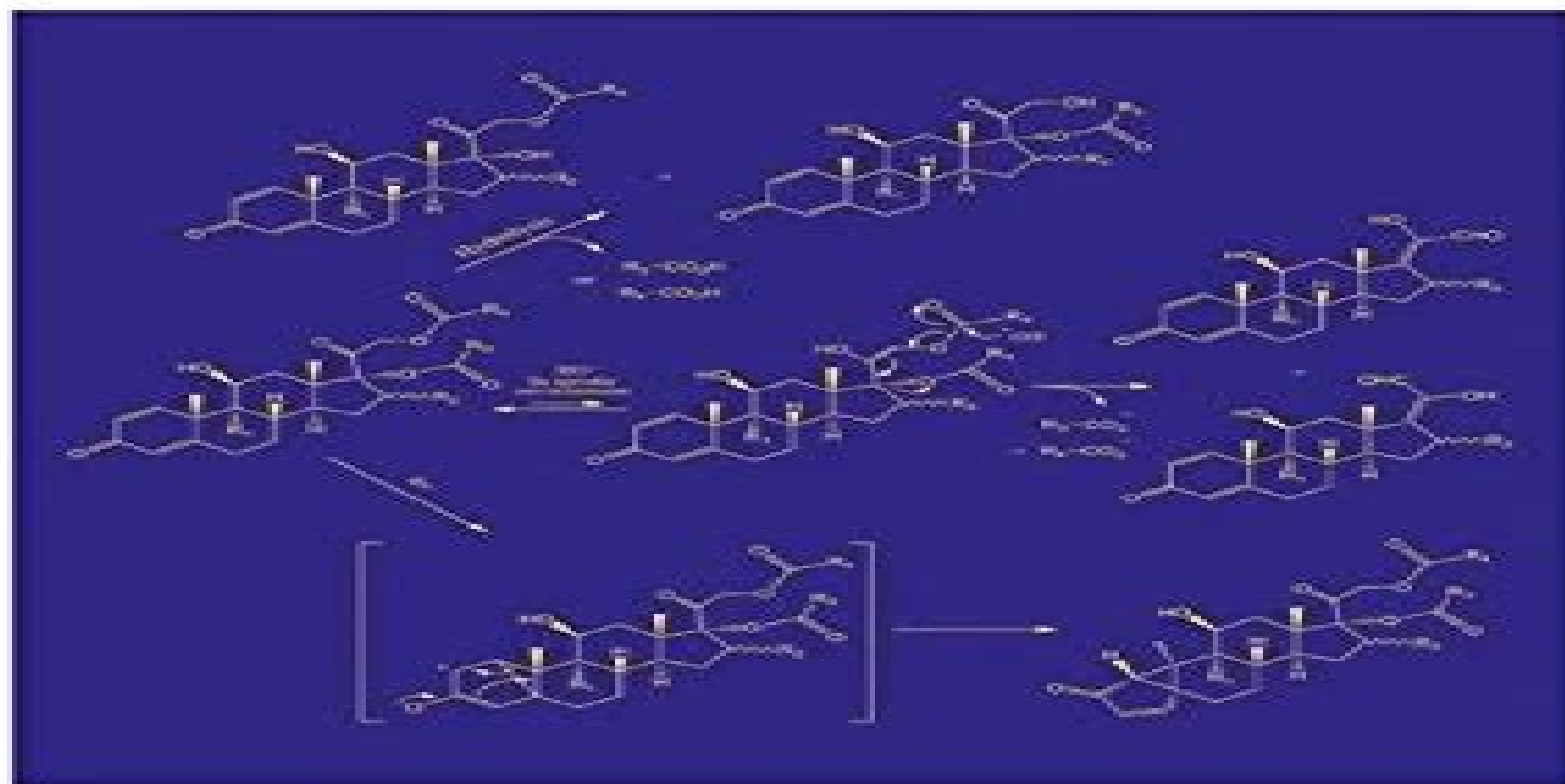


Min Li

Organic Chemistry of Drug Degradation



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Min Li



Organic Chemistry Of Drug Degradation Rsc Rsc Drug Discovery:

Organic Chemistry of Drug Degradation Min Li, 2015-10-20 The vast majority of drugs are organic molecular entities. A clear understanding of the organic chemistry of drug degradation is essential to maintaining the stability, efficacy, and safety of a drug product throughout its shelf life. During analytical method development, stability testing, and pharmaceutical manufacturing troubleshooting activities, one of the frequently occurring and usually challenging events would be the identification of drug degradants and understanding of drug degradation mechanisms and pathways. This book is written by a veteran of the pharmaceutical industry who has first-hand experience in drug design and development, drug degradation mechanism studies, analytical development, and manufacturing process troubleshooting and improvement. The author discusses various degradation pathways with an emphasis on the mechanisms of the underlying organic chemistry, which should aid greatly in the efforts of degradant identification, formulation development, analytical development, and manufacturing process improvement. Organic reactions that are significant in drug degradation will first be reviewed and then illustrated by examples of drug degradation reported in the literature. The author brings the book to a close with a final chapter dedicated to the strategy for rapid elucidation of drug degradants with regard to the current regulatory requirements and guidelines. One chapter that should be given special attention is Chapter 3, Oxidative Degradation. Oxidative degradation is one of the most common degradation pathways but perhaps the most complex one. This chapter employs more than sixty drug degradation case studies with in-depth discussion in regard to their unique degradation pathways. With the increasing regulatory requirements on the quality and safety of pharmaceutical products, in particular with regard to drug impurities and degradants, the book will be an invaluable resource for pharmaceutical and analytical scientists who engage in formulation development, analytical development, stability studies, degradant identification, and support of manufacturing process improvement. In addition, it will also be helpful to scientists engaged in drug discovery and development as well as in drug metabolism studies.

Biotherapeutics Lyn H. Jones, Andrew J. McKnight, 2013 Biotherapeutics are often considered to be beyond the reach of the medicinal chemist, but this book demonstrates that chemistry has an essential role in the future success of this area.

Comprehensive Biomarker Discovery and Validation for Clinical Application Péter Horvatovich, Rainer Bischoff, 2013 This book focuses on proteomics biomarker discovery and validation procedures from the clinical perspective. It provides an overview of current technology and the challenges encountered throughout the process. This covers all key stages from biomarker discovery and validation through to registration with the European and US regulatory authorities (EMA and FDA). All the important elements such as patient selection, sample handling, data processing, and statistical analysis are described in detail, and the reader is introduced to each topic with well-described examples or guidelines for best practice. Case studies are also included to demonstrate clinical applications. Individual chapters explain the best performing techniques for profiling complex body fluids and biomarker discovery. This includes the application of different LC-MS

profiling platforms and affinity array for screening complex body fluids Future developments needed to improve the success rate of translating biomarker discovery into useful clinical tests are also discussed Common pitfalls and success stories are described as are the limitations of the various technologies involved Broad and interdisciplinary in approach this book provides an excellent source of information for industrial and academic researchers Emerging Drugs and Targets for Parkinson's Disease Ana Martinez, Carmen Gil, 2013 This exemplary new book reviews some of the most outstanding examples of new drugs currently in pharmaceutical development or new targets under the validation process to try to reach the Parkinson's drug market in the next few years as potential disease modifying drugs *Green Chemistry Strategies for Drug Discovery* Emily A. Peterson, Julie B. Manley, 2015-06-30 The incorporation of Green Chemistry is a relatively new phenomenon in the drug discovery discipline since the scale that chemists operate on in drug discovery is smaller than those of process and manufacturing chemistry The necessary metrics are more difficult to obtain in drug discovery due to the diversity of reactions conducted However pharmaceutical companies are realizing that incorporation of green chemistry techniques at earlier stages of drug development can speed the development of a drug candidate Edited by experts who have pioneered green chemistry efforts within their own institutions this book provides a practical guide for both academic and industrial labs wanting to know where to start with introducing greener approaches for greatest return on investment The Editors have taken a comprehensive approach to the topic covering the entire drug discovery process from molecule conception through synthesis formulation and toxicology with specific examples and case studies where green chemistry strategies have been implemented Currently employed as well as emerging techniques for performing greener drug discovery chemistry are addressed as well as cutting edge topics like biologics discovery Moreover important surrounding issues such as intellectual property are included This book will serve as a practical guide for both academic and industrial chemists who work across the breadth of the drug discovery discipline Ultimately readers will learn how to incorporate green chemistry strategies into their everyday workflow without slowing down their science **Pain Therapeutics** Charlotte Allerton, 2013 This book presents a contemporary review of the field of pain therapeutics including the historical medicines which still dominate standard of care treatments as well as the new mechanisms and combinations reformulations that have dominated the regulatory approvals over the last decade In addition this book provides a deep review of the key biological mechanisms currently under investigation for their utility into the treatment of pain such as ion channels opiates and others Additional discussion highlights the current challenges of pain research covering a range of topics from difficulties in identifying new targets and pre clinical models to the current regulatory and commercial challenges This background sets the scene for recent scientific developments in pain research such as the drive for genetic validation of targets and the derivation of human cell platforms from stem cells Finally the book covers the discovery and development stories of two pain products approved in the last decade These case studies for Lyrica and the Butrans patch will give insight into the discovery

and development challenges and successes for both an oral and non oral product **Allosterism in Drug Discovery** Dario Doller, 2016-11-24 Although the concept of allosterism has been known for over half a century its application in drug discovery has exploded in recent years The emergence of novel technologies that enable molecular level ligand receptor interactions to be studied in unprecedented detail has driven this trend This book written by the leaders in this young research area describes the latest developments in allosterism for drug discovery Bringing together research in a diverse range of scientific disciplines Allosterism in Drug Discovery is a key reference for academics and industrialists interested in understanding allosteric interactions The book provides an in depth review of research using small molecules as chemical probes and drug candidates that interact allosterically with proteins of relevance to life sciences and human disease Knowledge of these interactions can then be applied in the discovery of the novel therapeutics of the future This book will be useful for people working in all disciplines associated with drug discovery in academia or industry as well as postgraduate students who may be working in the design of allosteric modulators **Computational Approaches to Nuclear**

Receptors Pietro Cozzini, Glen E. Kellogg, 2012 Nuclear receptors NR are ligand induced activated transcription factors that are involved in numerous biological processes Since the 1990 s when the first structures were determined by means of X ray diffraction the number of NR structures has increased considerably Moreover several omics projects genomics pharmacogenomics and proteomics have opened up great opportunities for the discovery of new targets the characterization of abnormal protein patterns the selection of tailored drugs and the evaluation of drug efficacy even with a lack of structural data Furthermore structure based drug design computational methods for in silico screening and nanobiotechnology based tools are simplifying this time consuming and money intensive research of lead compounds and possibly new drugs Biological interactions such as those that occur between a protein and ligand are concerted events where flexible molecules interact Thus understanding flexibility of large molecules or biological complexes is of primary importance to help define the right model to approximate the reality for drug discovery virtual screening food safety analysis etc NRs are known as flexible targets with many structural similarities in particular for their Ligand Binding Domain these similarities could be assumed to share behavioural qualities that belong to this class of compounds Thus to supply a possible complete and exhaustive answer to questions about the behaviour of NRs their interactions with new potential drugs endocrine disruptors such as animal and human food toxins food additives or industry residuals it is mandatory to approach the problem from a different point of view a molecular modelling approach steered synthesis and in vitro and in vivo tests etc The aim of this book is to provide a state of the art review on investigations into Nuclear Receptors **Macrocycles in Drug Discovery** Jeremy I Levin, 2015 This series provides a comprehensive resource for postgraduate students and for scientists in academia or industry wanting to learn topics outside their own areas of expertise

Traditional Chinese Medicine James David Adams (Jr.), Eric J. Lien (ed.), 2013 A volume that will provide a mechanism based approach to understanding Chinese medicine for both Western and

Eastern scientists *Ion Channel Drug Discovery* Brian Cox, Martin Gosling, 2014-09-24 A rapidly growing field this book covers the recent advances in screening technology ion channel structure and modelling with up to date case histories

Accounts in Drug Discovery Joel Barrish, Percy Carter, Peter Cheng, Robert Zahler, 2010-09-30 Accounts in Drug Discovery describes recent case studies in medicinal chemistry with a particular emphasis on how the inevitable problems that arise during any project can be surmounted or overcome The Editors cover a wide range of therapeutic areas and medicinal chemistry strategies including lead optimization starting from high throughput screening hits as well as rational structure based design The chapters include follow ons and next generation compounds that aim to improve upon first generation agents This volume surveys the range of challenges commonly faced by medicinal chemistry researchers including the optimization of metabolism and pharmacokinetics toxicology pharmaceutics and pharmacology including proof of concept in the clinic for novel biological targets The case studies include medicinal chemistry stories on recently approved and marketed drugs but also chronicle near misses i e exemplary compounds that may have proceeded well into the clinic but for various reasons did not result in a successful registration As the vast majority of projects fail prior to registration much can be learned from such narratives By sharing a wide range of drug discovery experiences and information across the community of medicinal chemists in both industry and academia we believe that these accounts will provide insights into the art of medicinal chemistry as it is currently practiced and will help to serve the needs of active medicinal chemists **Third**

SCI-RSC Medicinal Chemistry Symposium Robert Wilson Lambert, 1986 *Synthetic Methods in Drug Discovery* David C Blakemore, Paul M Doyle, Yvette M Fobian, 2016-07-15 Building on key reactions presented in Volume 1 Synthetic Methods in Drug Discovery Volume 2 covers a range of important reaction types including organometallic chemistry fluorination approaches and asymmetric methods as well as new and exciting areas such as Csp² Csp³ couplings catalytic amide bond forming reactions hydrogen borrowing chemistry and methods to access novel motifs and monomers This book provides both academic and industrial perspectives on key reactions giving the reader an excellent overview of the techniques used in modern synthesis Reaction types are conveniently framed in the context of their value to industry and the challenges and limitations of methodologies are discussed with relevant illustrative examples Moreover key opportunities in expanding chemical space are presented including the increasingly important syntheses that introduce three dimensional molecular shape Edited and authored by leading scientists from both academia and industry this book will be a valuable reference for all chemists involved in drug discovery as well as postgraduate students in medicinal chemistry Chemistry and Industry, 2008 **Synthetic Methods in Drug Discovery** David C. Blakemore, Paul M. Doyle, Yvette M. Fobian, 2016

Fragment-Based Drug Discovery Steven Howard, Chris Abell, 2015-06-17 Fragment based drug discovery is a rapidly evolving area of research which has recently seen new applications in areas such as epigenetics GPCRs and the identification of novel allosteric binding pockets The first fragment derived drug was recently approved for the treatment of melanoma It is

hoped that this approval is just the beginning of the many drugs yet to be discovered using this fascinating technique This book is written from a Chemist s perspective and comprehensively assesses the impact of fragment based drug discovery on a wide variety of areas of medicinal chemistry It will prove to be an invaluable resource for medicinal chemists working in academia and industry as well as anyone interested in novel drug discovery techniques MInd, the Meetings Index ,1995

Exploiting Chemical Diversity for Drug Discovery Paul A Bartlett,Michael Entzeroth,2007-10-31 Conceptual and technological advances in chemistry and biology have transformed the drug discovery process Evolutionary pressure among the diverse scientific and engineering disciplines that contribute to the identification of biologically active compounds has resulted in synergistic improvements at every step in the process Exploiting Chemical Diversity for Drug Discovery encompasses the many components of this transformation and presents the current state of the art of this critical endeavour From the theoretical and operational considerations in generating a collection of compounds to screen to the design and implementation of high capacity and high quality assays that provide the most useful biological information this book provides a comprehensive overview of modern approaches to lead identification Beginning with an introductory overview subsequent chapters address topics that include the design of chemical libraries and methods for optimizing their diversity automated and accelerated chemistry high throughput assay design and detection techniques and strategies for data analysis and property optimization Written by experts in the field both academic and industrial and illustrated in full colour this book provides an excellent overview for current practitioners and will also serve as a stimulating resource for future generations Researchers in organic and medicinal chemistry the biological and pharmacological sciences as well as those interested in allied computational and engineering disciplines will value the comprehensive and up to date coverage Drug Discovery from Natural Products Olga Genilloud,Francisca Vicente,2012 An integrated review of the most recent trends in natural products drug discovery and key lead candidates that are outstanding for their chemistry and biology in novel drug development

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