

1. Record Nr.	UNINA9910452586103321
Autore	Callaway Gloria
Titolo	The early years curriculum : a view from outdoors / / Gloria Callaway
Pubbl/distr/stampa	London : , : David Fulton Publishers, , 2005
ISBN	0-203-82095-9 1-136-75391-5
Descrizione fisica	1 online resource (89 p.)
Disciplina	951.9504/3
Soggetti	Outdoor education Electronic books.
Lingua di pubblicazione	Inglese
Formato	Materiale a stampa
Livello bibliografico	Monografia
Note generali	Includes index.
Nota di contenuto	section 1. The weekly routine -- section 2. Principles into practice -- section 3. The grown-ups -- section 4. The curriculum in action -- section 5. Conclusions.
Sommario/riassunto	Based on an account of Cornwall's Education Action Zone project, this book explores how to plan, implement and assess a rigorous outdoor early years curriculum that complements classroom-based learning and meets the Early Learning Goals.

2. Record Nr.	UNINA9910782320303321
Autore	Lee Wen Ho
Titolo	Computer simulation of shaped charge problems [[electronic resource] /] / Wen Ho Lee
Pubbl/distr/stampa	Singapore ; ; Hackensack, NJ, : World Scientific, c2006
ISBN	1-281-90896-7 9786611908966 981-270-713-1
Descrizione fisica	1 online resource (220 p.)
Disciplina	623.4/545
Soggetti	Shaped charges - Computer simulation Shaped charges - Mathematical models Flow visualization - Computer simulation Penetration mechanics - Computer simulation Lagrange equations
Lingua di pubblicazione	Inglese
Formato	Materiale a stampa
Livello bibliografico	Monografia
Note generali	Description based upon print version of record.
Nota di bibliografia	Includes bibliographical references and indexes.
Nota di contenuto	Contents; A Personal Introduction; Chapter 1 Small Molecules for Chemogenomics-based Drug Discovery Edgar Jacoby, Ansgar Schuffenhauer, Kamal Azzaoui, Maxim Popov, Sigmar Dressler, Meir Glick, Jeremy Jenkins, John Davies and Silvio Roggo; 1. Introduction; 2. Compound Categories; 2.1. Natural products and derivatives; 2.2. Primary metabolites, co-substrates, co-factors, and marketed drugs; 2.3. Peptides and peptido-mimetics; 2.4. Diversity oriented synthesis molecules; 3. Designing Comprehensive Chemogenomics Screening Compound Collections 4. Essential Properties and Selection Processes along the Discovery Pipeline5. Molecular Information Systems and Annotated Compound Libraries; 6. Conclusion; Acknowledgements; References; Chapter 2 Mapping the Chemogenomic Space Jordi Mestres; 1. The Chemogenomic Space; 2. Annotation and Classification Schemes for Proteins; 2.1. Enzymes; 2.2. Receptors; 2.2.1. Channel receptors; 2.2.2. G Protein-coupled receptors; 2.2.3. Nuclear receptors; 3. Structural Representativity of Protein Families; 4. Annotation and Classification

Schemes for Molecules; 5. Mapping the Molecule-Protein Space
 6. Exploiting the Chemogenomic Space 7. Conclusions; References;
 Chapter 3 Natural Product Scaffolds and Protein Structure Similarity
 Clustering (PSSC) as Inspiration Sources for Compound Library Design
 in Chemogenomics and Drug Development Frank J. Dekker, Stefan
 Wetzel and Herbert Waldmann; 1. Introduction; 2. Biological Relevance
 in Compound Library Design; 2.1. Compound libraries as sources for
 small molecule modulators of protein function; 2.2. Annotated
 libraries; 2.3. Natural products as inspiration sources for library design;
 2.4. Library design based on privileged structures
 3. Natural Product Inspired Compound Library Synthesis 4. Target
 Clustering as Strategy in Drug Discovery; 4.1. Target clustering; 4.2.
 Target clustering based on structural and functional similarity; 5. PSSC
 as Guiding Principle for Compound Library Design; 5.1. Protein
 structure similarity clustering (PSSC); 5.2. PSSC based reanalysis of the
 development of leukotriene A4 hydrolase inhibitors; 5.3. PSSC based
 reanalysis of the development of nuclear hormone receptor ligands
 5.4. Application of PSSC for de novo ligand development for the protein
 cluster Cdc25A phosphataseacetylcholinesterase-11-hydroxysteroid
 dehydrogenase5.5. Position of the PSSC concept in drug development
 and chemogenomics; 6. Conclusions; Acknowledgments; References;
 Chapter 4 A Reductionist Approach to Chemogenomics in the Design of
 Drug Molecules and Focused Libraries Roger Crossley and Martin Slater;
 1. Introduction; 2. Molecular Recognition and Vicinity AnalysisTM; 3.
 Thematic AnalysisTM; Examples of Themes; 4. Family B and C GPCRs; 5.
 Classification of GPCRs; 6. Pharmacophore Maps
 7. Library Design Using Thematic AnalysisTM

Sommario/riassunto

Devoted to the subject of shape charge design using numerical
 methods, this book offers the defense and commercial industries
 unique material not contained in any other single volume. The coverage
 of the Lagrangian and Eulerian methods as well as the equation of state
 provides first hand help to engineers working on shape charge
 problems. The book includes detailed descriptions of oil-well
 perforation not available from any other sources and, coupled with the
 material flow physics discussed in Chapters 2 and 3 and Appendix B,
 readers can design the fuel rod configurations for a nuclear reactor