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Collana	Lecture Notes in Bioinformatics ; ; 4453
Disciplina	572.8
Soggetti	Molecular biology - Computer simulation
Nota di bibliografia	Includes bibliographical references and author index.
Nota di contenuto	QNet: A Tool for Querying Protein Interaction Networks Pairwise Global Alignment of Protein Interaction Networks by Matching Neighborhood Topology Reconstructing the Topology of Protein Complexes Network Legos: Building Blocks of Cellular Wiring Diagrams An Efficient Method for Dynamic Analysis of Gene Regulatory Networks and in silico Gene Perturbation Experiments A Feature-Based Approach to Modeling Protein-DNA Interactions Network Motif Discovery Using Subgraph Enumeration and Symmetry- Breaking Nucleosome Occupancy Information Improves de novo Motif Discovery Framework for Identifying Common Aberrations in DNA Copy Number Data Estimating Genome-Wide Copy Number Using Allele Specific Mixture Models GIMscan: A New Statistical Method for Analyzing Whole-Genome Array CGH Data Production- Passage-Time Approximation: A New Approximation Method to Accelerate the Simulation Process of Enzymatic Reactions Shift- Invariant Adaptive Double Threading: Learning MHC II - Peptide Binding Reconstructing the Phylogeny of Mobile Elements Beyond Galled Trees - Decomposition and Computation of Galled Networks Variational Upper Bounds for Probabilistic Phylogenetic Models Heuristics for the Gene-Duplication Problem: A ?(n) Speed-Up for the

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	Local Search Support Vector Training of Protein Alignment Models
	Tools for Simulating and Analyzing RNA Folding Kinetics Multiple
	Sequence Alignment Based on Profile Alignment of Intermediate
	Sequences Connectedness Profiles in Protein Networks for the
	Analysis of Gene Expression Data Multivariate Segmentation in the
	Analysis of Transcription Tiling Array Data A Bayesian Model That
	Links Microarray mRNA Measurements to Mass Spectrometry Protein
	Measurements Rearrangements in Genomes with Centromeres Part I:
	Translocations Identification of Deletion Polymorphisms from
	Haplotypes Free Energy Estimates of All-Atom Protein Structures
	Using Generalized Belief Propagation Minimizing and Learning
	Energy Functions for Side-Chain Prediction Protein Conformational
	Flexibility Analysis with Noisy Data Deterministic Pharmacophore
	Detection Via Multiple Flexible Alignment of Drug-Like Molecules
	Design of Compact, Universal DNA Microarrays for Protein Binding
	Microarray Experiments Improved Ranking Functions for Protein and
	Modification-Site Identifications Peptide Retention Time Prediction
	Yields Improved Tandem Mass Spectrum Identification for Diverse
	Chromatography Conditions A Fast and Accurate Algorithm for the
	Quantification of Peptides from Mass Spectrometry Data Association
	Mapping of Complex Diseases with Ancestral Recombination Graphs:
	Models and Efficient Algorithms An Efficient and Accurate Graph-
	Based Approach to Detect Population Substructure RB-Finder: An
	Improved Distance-Based Sliding Window Method to Detect
	Recombination Breakpoints Comparative Analysis of Spatial Patterns
	of Gene Expression in Drosophila melanogaster Imaginal Discs.
Sommario/riassunto	This book constitutes the refereed proceedings of the 11th Annual
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	Biology, RECOMB 2007, held in Oakland, CA, USA in April 2007. The 37
	revised full papers presented were carefully reviewed and selected from
	just under 170 submissions. As the top conference in computational
	molecular biology, RECOMB addresses all current issues in algorithmic,
	theoretical, and experimental bioinformatics.