1. Record Nr. UNISA996393282603316

Titolo	Miserere cleri: or, The factions of the church [[electronic resource]] : Being a short view of the pernicious consequences of the clergy's intermedling with affairs of state
Pubbl/distr/stampa	London : , : Printed for W. Boreham, at the Angel in Pater-noster-row., 1718
Descrizione fisica	35, [5] p
Altri autori (Persone)	DefoeDaniel <1661?-1731.>
Soggetti	Church and state
Lingua di pubblicazione	Inglese
Formato	Materiale a stampa
Livello bibliografico	Monografia
Note generali	A re-issue of The pernicious consequences of the clergy's intermedling with affairs of state, with very minor amendments. Cf. Furbank and Owens, Defoe De-attributions. Sometimes attributed to Daniel Defoe (Moore, Novak ("perhaps"). Attribution disputed by Furbank and Owens, Defoe de-attributions. Reproduction of original in the British Library.
Sommario/riassunto	eebo-0018

2.

Record Nr.	UNISALENTO991004131819707536
Autore	Truini, Dario
Titolo	Gli strumenti del pensiero : guida alla conoscenza dei meccanismi logici e allo studio razionale dei problemi / Dario Truini
Pubbl/distr/stampa	Milano : F. Angeli, 1991
ISBN	8820430614
Edizione	[3. ed.]
Descrizione fisica	138 p. ; 22 cm
Collana	Trend ; 15
Disciplina	301.1
Soggetti	Decisione Ragionamento
Lingua di pubblicazione	Italiano
Formato	Materiale a stampa
Livello bibliografico	Monografia

3. Record Nr. UNINA9911019659603321

Titolo	Ab initio methods in quantum chemistry . Part I / / edited by K.P Lawley
Pubbl/distr/stampa	Chichester [West Sussex] ; ; New York, : Wiley, c1987
ISBN	9786612347221 9781282347229 1282347225 9780470142936 0470142936 9780470143377 0470143371
Descrizione fisica	1 online resource (568 p.)
Collana	Advances in chemical physics ; ; 67
Altri autori (Persone)	LawleyK. P
Disciplina	541.2/8 541.28 541.305 541/.08
Soggetti	Quantum chemistry Quantum theory
Lingua di pubblicazione	Inglese
Formato	Materiale a stampa
Livello bibliografico	Monografia
Note generali	Description based upon print version of record.
Nota di bibliografia	Includes bibliographies and indexes.
Nota di contenuto	AB INITIO METHODS IN QUANTUM CHEMISTRY-Part I; CONTENTS; EXCITED-STATE POTENTIALS; MOLECULAR PROPERTY DERIVATIVES; TRANSITION STRUCTURE COMPUTATIONS AND THEIR ANALYSIS; OPTIMIZATION OF EQUILIBRIUM GEOMETRIES AND TRANSITION STRUCTURES; RELATIVISTIC QUANTUM CHEMISTRY; EFFECTIVE HAMILTONIANS AND PSEUDO-OPERATORS AS TOOLS FOR RIGOROUS MODELLING; MOLECULAR CALCULATIONS WITH THE DENSITY FUNCTIONAL FORMALISM; BASIS SETS; THE COUPLED PAIR . APPROXIMATION; AUTHOR INDEX; SUBJECT INDEX
Sommario/riassunto	The Advances in Chemical Physics series provides the chemical physics and physical chemistry fields with a forum for critical, authoritative evaluations of advances in every area of the discipline. Filled with cutting-edge research reported in a cohesive manner not found

elsewhere in the literature, each volume of the Advances in Chemical Physics series serves as the perfect supplement to any advanced graduate class devoted to the study of chemical physics.