



Computational Pharmaceutical Solid State Chemistry

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Computational Pharmaceutical Solid State Chemistry

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This book is the first to combine computational material science and modeling of molecular solid states for pharmaceutical industry applications.

- Provides descriptive and applied state-of-the-art computational approaches and workflows to guide pharmaceutical solid state chemistry experiments and to support/troubleshoot API solid state selection
- Includes real industrial case examples related to application of modeling methods in problem solving
- Useful as a supplementary reference/text for undergraduate, graduate and postgraduate students in computational chemistry, pharmaceutical and biotech sciences, and materials science

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