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MODELING VIRAL CAPSID ASSEMBLY

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CHARGES AT AQUEOUS INTERFACES: DEVELOPMENT OF COMPUTATIONAL APPROACHES IN DIRECT CONTACT WITH EXPERIMENT

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SOLUTE PRECIPITATE NUCLEATION: A REVIEW OF THEORY AND SIMULATION ADVANCES

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WATER IN THE LIQUID STATE: A COMPUTATIONAL VIEWPOINT

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CONSTRUCTION OF ENERGY FUNCTIONS FOR LATTICE HETEROPOLYMER MODELS: EFFICIENT ENCODINGS FOR CONSTRAINT SATISFACTION PROGRAMMING AND QUANTUM ANNEALING

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