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Diamond is our preeminent molecular and crystal structure visualization software. It combines many features that are superior to working with .NET platform. With it, you can easily view, modify and compare the results obtained using different methods. In addition, we use a number of algorithms to automatically analyze the results and build models, which allows us to quickly determine the type and nature of the data. We recommend using Diamond software to analyze information and model it for further exploration and creation of new materials.

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