Resource Summary Report

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Molecular Dynamics (MD) Simulations, step by step protocol

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Protocol Information

URL: https://dx.doi.org/10.17504/protocols.io.mztc76n

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Summary: To MD analysis for one of our previous results of docking, for example to analyze the drug/target, almost 40 steps are required to finalize the MD results.The following protocol is according to our own experience in this project and was improved so that the errors were solved during the projects.

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