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Ab initio molecular dynamics calculations on reactions of molecules with metal surfaces

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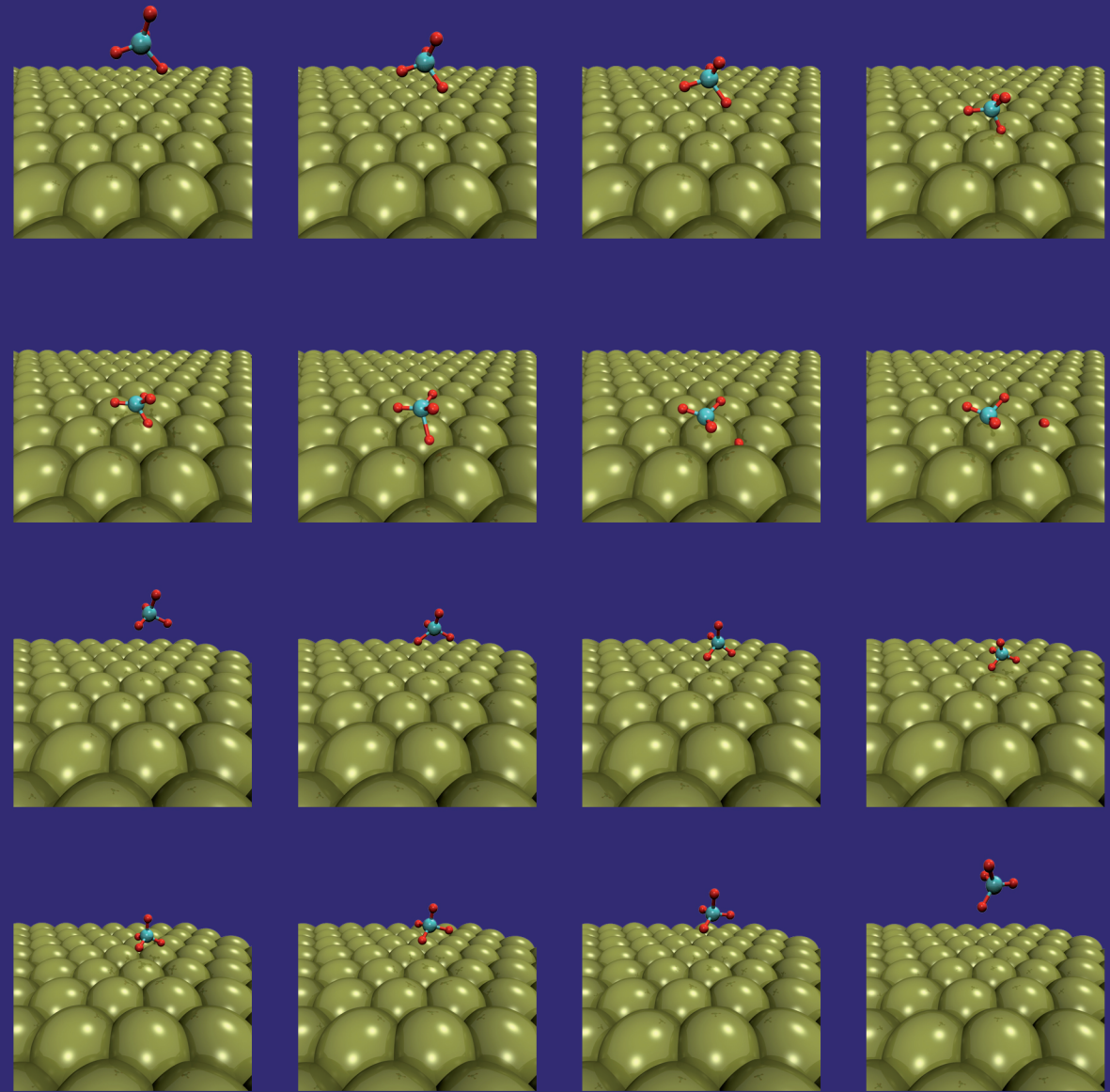
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AIMD Calculations on Reactions of Molecules with Metal Surfaces

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