Universiteit
Leiden
The Netherlands

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

Nattino, F.

Citation
Nattino, F. (2015, October 28). *Ab initio molecular dynamics calculations on reactions of molecules with metal surfaces*. Retrieved from https://hdl.handle.net/1887/35980

Version:	Corrected Publisher's Version
License:	Licence agreement concerning inclusion of doctoral thesis in the Institutional Repository of the University of Leiden
Downloaded from:	https://hdl.handle.net/1887/35980

Note: To cite this publication please use the final published version (if applicable).

Cover Page

Universiteit Leiden

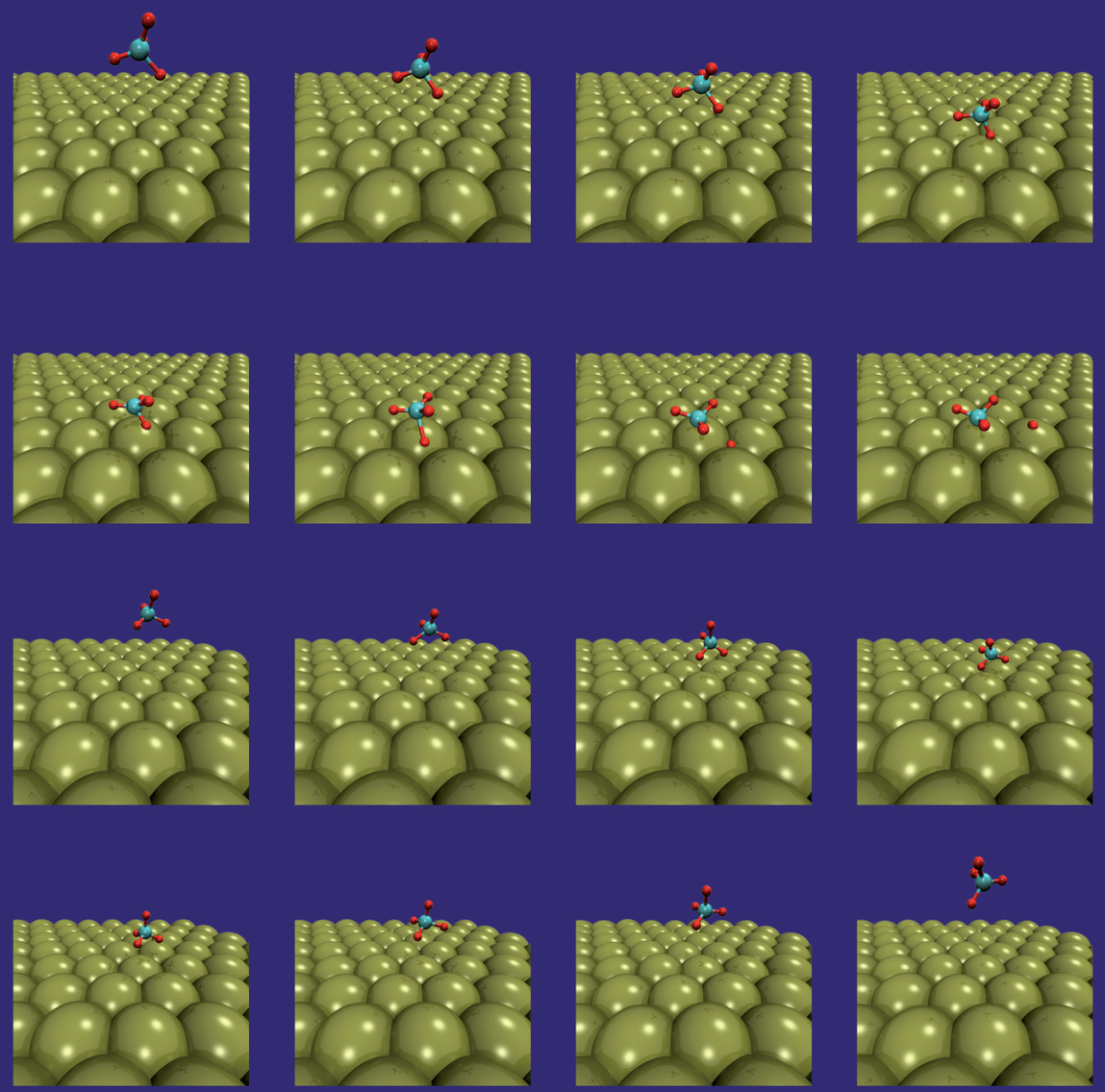
The handle http://hdl.handle.net/1887/35980 holds various files of this Leiden University dissertation

Author: Nattino, Francesco
Title: Ab initio molecular dynamics calculations on reactions of molecules with metal surfaces
Issue Date: 2015-10-28

AIMD Calculations on Reactions of Molecules with Metal Surfaces

Francesco Nattino

Ab Initio Molecular Dynamics Calculations on Reactions of Molecules with Metal Surfaces



Francesco Nattino