

# Martin Tschöppe

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2341080/publications.pdf>

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6  
papers

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1478505

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1872680

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docs citations

6  
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citing authors

#	ARTICLE	IF	CITATIONS
1	Efficient and automated quantum chemical calculation of rovibrational nonresonant Raman spectra. Journal of Chemical Physics, 2022, 156, 124102.	3.0	9
2	<i>Ab initio</i> calculation of rovibrational states for non-degenerate double-well potentials: <i>cis</i> $\leftrightarrow$ <i>trans</i> isomerization of HOPO. Journal of Chemical Physics, 2020, 152, 174306.	3.0	11
3	Neural network approach for the dynamics on the normally hyperbolic invariant manifold of periodically driven systems. Physical Review E, 2020, 101, 022219.	2.1	8
4	Toward a fully automated calculation of rovibrational infrared intensities for semi-rigid polyatomic molecules. Journal of Chemical Physics, 2020, 152, 244104.	3.0	29
5	High-Level Rovibrational Calculations on Ketenimine. Frontiers in Chemistry, 2020, 8, 623641.	3.6	13
6	Invariant Manifolds and Rate Constants in Driven Chemical Reactions. Journal of Physical Chemistry B, 2019, 123, 2070-2086.	2.6	23