

# Rollin A King

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2956718/publications.pdf>

Version: 2024-02-01

9  
papers

5,264  
citations

1039880

9  
h-index

1474057

9  
g-index

11  
all docs

11  
docs citations

11  
times ranked

6204  
citing authors

#	ARTICLE	IF	CITATIONS
1	Advances in molecular quantum chemistry contained in the Q-Chem 4 program package. <i>Molecular Physics</i> , 2015, 113, 184-215.	0.8	2,561
2	<sc>Psi4</sc> 1.1: An Open-Source Electronic Structure Program Emphasizing Automation, Advanced Libraries, and Interoperability. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 3185-3197.	2.3	961
3	Psi4: an open-source <i>ab initio</i> electronic structure program. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2012, 2, 556-565.	6.2	838
4	P<sc>SI4</sc> 1.4: Open-source software for high-throughput quantum chemistry. <i>Journal of Chemical Physics</i> , 2020, 152, 184108.	1.2	440
5	PSI3: An open-source Ab Initio electronic structure package. <i>Journal of Computational Chemistry</i> , 2007, 28, 1610-1616.	1.5	258
6	P<sc>si</sc>4N<sc>um</sc>P<sc>y</sc>: An Interactive Quantum Chemistry Programming Environment for Reference Implementations and Rapid Development. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 3504-3511.	2.3	106
7	Coupled cluster methods including triple excitations for excited states of radicals. <i>Journal of Chemical Physics</i> , 2005, 122, 054110.	1.2	64
8	Quantum Chemistry Common Driver and Databases (QCDB) and Quantum Chemistry Engine (QCE<sc>ngine</sc>): Automation and interoperability among computational chemistry programs. <i>Journal of Chemical Physics</i> , 2021, 155, 204801.	1.2	15
9	Competition Between $\pi$ - $\pi$ and $C\ddot{=}H/\ddot{=}H$ Interactions: A Comparison of the Structural and Electronic Properties of Alkoxy-Substituted 1,8-Bis((propyloxyphenyl)ethynyl)naphthalenes. <i>Chemistry - A European Journal</i> , 2015, 21, 19168-19175.	1.7	9