

# Florian Mittendorfer

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4824770/publications.pdf>

Version: 2024-02-01

32

papers

899

citations

516710

16

h-index

454955

30

g-index

33

all docs

33

docs citations

33

times ranked

1699

citing authors

#	ARTICLE	IF	CITATIONS
1	Adsorption of Unsaturated Hydrocarbons on Pd(111) and Pt(111): A DFT Study. <i>Journal of Physical Chemistry B</i> , 2003, 107, 12287-12295.	2.6	148
2	Nickel Carbide as a Source of Grain Rotation in Epitaxial Graphene. <i>ACS Nano</i> , 2012, 6, 3564-3572.	14.6	77
3	Disorder and Defect Healing in Graphene on Ni(111). <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 136-139.	4.6	65
4	Carbon in palladium catalysts: A metastable carbide. <i>Journal of Chemical Physics</i> , 2010, 132, 024711.	3.0	63
5	Adsorption of water at the SrO surface of Ruthenates. <i>Nature Materials</i> , 2016, 15, 450-455.	27.5	63
6	A first-principles study of bulk oxide formation on Pd(100). <i>Journal of Chemical Physics</i> , 2009, 131, 054701.	3.0	53
7	Artificially lattice-mismatched graphene/metal interface: Graphene/Ni/Ir(111). <i>Physical Review B</i> , 2013, 87, .	3.2	53
8	Growth of an Ultrathin Zirconia Film on Pt <sub>3</sub> Zr Examined by High-Resolution X-ray Photoelectron Spectroscopy, Temperature-Programmed Desorption, Scanning Tunneling Microscopy, and Density Functional Theory. <i>Journal of Physical Chemistry C</i> , 2015, 119, 2462-2470.	3.1	46
9	Pt $\times$ mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>3</mml:mn></mml:msub></mml:math>Zr(0001): A substrate for growing well-ordered ultrathin zirconia films by oxidation. <i>Physical Review B</i> , 2012, 86, .	3.2	41
10	Self-Organized Growth, Structure, and Magnetism of Monatomic Transition-Metal Oxide Chains. <i>Physical Review Letters</i> , 2016, 117, 046101.	7.8	32
11	Tuning the Growth Orientation of Epitaxial Films by Interface Chemistry. <i>Physical Review Letters</i> , 2012, 108, 066101.	7.8	27
12	Water adsorption at zirconia: from the ZrO <sub>2</sub> (111)/Pt <sub>3</sub> Zr(0001) model system to powder samples. <i>Journal of Materials Chemistry A</i> , 2018, 6, 17587-17601.	10.3	24
13	One-dimensional Oxide-Metal Hybrid Structures: Site-Specific Enhanced Reactivity for CO Oxidation. <i>ChemPhysChem</i> , 2010, 11, 2506-2509.	2.1	20
14	Quasiliquid Layer Promotes Hexagonal Boron Nitride (h-BN) Single-Domain Growth: h-BN on Pt(110). <i>ACS Nano</i> , 2019, 13, 7083-7090.	14.6	19
15	High Chemical Activity of a Perovskite Surface: Reaction of CO with mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow /><mml:msub><mml:mrow><mml:mi>Sr</mml:mi></mml:mrow><mml:mi>3</mml:mi></mml:msub></mml:mrow><mml:mathvariant="normal">O</mml:mi></mml:mrow><mml:mrow><mml:mi>7</mml:mi></mml:mrow></mml:mrow></mml:msub></mml:mrow></mml:mrow>. <i>Physical Review Letters</i> , 2014, 113, 116101.	14.6	18
16	Metal Adatoms and Clusters on Ultrathin Zirconia Films. <i>Journal of Physical Chemistry C</i> , 2016, 120, 9920-9932.	3.1	18
17	A full monolayer of superoxide: oxygen activation on the unmodified Ca <sub>3</sub> Ru <sub>2</sub> O <sub>7</sub> (001) surface. <i>Journal of Materials Chemistry A</i> , 2018, 6, 5703-5713.	10.3	17
18	Low-dimensional surface oxides in the oxidation of Rh particles. <i>Journal of Physics Condensed Matter</i> , 2010, 22, 393001.	1.8	13

#	ARTICLE	IF	CITATIONS
19	Role of Precursor Carbides for Graphene Growth on Ni(111). <i>Scientific Reports</i> , 2018, 8, 2662.	3.3	13
20	Point defects at cleaved surfaces. <i>Physical Review B</i> , 2014, 90, .	3.2	12
21	Ordered hydroxyls on Ca <sub>3</sub> Ru <sub>2</sub> O <sub>7</sub> (001). <i>Nature Communications</i> , 2017, 8, 23.	12.8	12
22	Adsorption of hydrogen on stable and metastable Ir(100) surfaces. <i>Surface Science</i> , 2017, 656, 66-76.	1.9	9
23	Monitoring the Interaction of CO with Graphene Supported Ir Clusters by Vibrational Spectroscopy and Density Functional Theory Calculations. <i>Journal of Physical Chemistry C</i> , 2018, 122, 4281-4289.	3.1	9
24	Many-electron calculations of the phase stability of ZrO <sub>2</sub> polymorphs. <i>Physical Review Research</i> , 2020, 2, .	3.6	8
25	A density-functional theory study of the adsorption of CO molecules on Au/Ni(111). <i>Journal of Physics Condensed Matter</i> , 2006, 18, 10825-10835.	1.8	8
26	Thermodynamic modelling of the partially ordered solid solution Hf <sub>5-x</sub> Nb <sub>x</sub> Ge <sub>4</sub> supported by ab initio calculations. <i>Solid State Sciences</i> , 2007, 9, 159-165.	3.2	8
27	Oxygen-Stabilized Rh Adatoms: 0D Oxides on a Vicinal Surface. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2747-2751.	4.6	5
28	Single-domain h-BN on Pt(110): Electronic structure, correlation, and bonding. <i>Physical Review Research</i> , 2020, 2, .	3.6	5
29	First principles studies of the electronic and structural properties of the rutile VO <sub>2</sub> (110) surface and its oxygen-rich terminations. <i>Journal of Physics Condensed Matter</i> , 2021, 33, 475002.	1.8	4
30	Oxygen-rich tetrahedral surface phase on high-temperature rutile single crystals. <i>Physical Review Materials</i> , 2021, 5, .	2.4	3
31	Adsorption of CO on the Ca <sub>3</sub> Ru <sub>2</sub> O <sub>7</sub> (001) Åsurface. <i>Surface Science</i> , 2019, 680, 18-23.	1.9	2
32	Adsorption of a superoxo O <sub>2</sub> - Åspecies on the pure and Ca-doped Sr <sub>3</sub> Ru <sub>2</sub> O <sub>7</sub> (001) Åsurface. <i>Surface Science</i> , 2019, 680, 24-31.	1.9	2