

Antreas C Kalli

List of Publications by Year in descending order

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Version: 2024-02-01

40
papers

1,603
citations

293460

24
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371746

37
g-index

50
all docs

50
docs citations

50
times ranked

2123
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----|-----------|
| 1 | Pocket delipidation induced by membrane tension or modification leads to a structurally analogous mechanosensitive channel state. <i>Structure</i> , 2022, 30, 608-622.e5. | 1.6 | 16 |
| 2 | Organization and Dynamics of the Red Blood Cell Band 3 Anion Exchanger SLC4A1: Insights From Molecular Dynamics Simulations. <i>Frontiers in Physiology</i> , 2022, 13, 817945. | 1.3 | 6 |
| 3 | Characterization of the membrane interactions of phospholipase C β 3 reveals key features of the active enzyme. <i>Science Advances</i> , 2022, 8, . | 4.7 | 7 |
| 4 | The guidance and adhesion protein FLRT2 dimerizes in cis via dual small-X3-small transmembrane motifs. <i>Structure</i> , 2022, 30, 1354-1365.e5. | 1.6 | 4 |
| 5 | Modeling of full-length Piezo1 suggests importance of the proximal N-terminus for dome structure. <i>Biophysical Journal</i> , 2021, 120, 1343-1356. | 0.2 | 23 |
| 6 | Molecular dynamics simulations of Piezo1 channel opening by increases in membrane tension. <i>Biophysical Journal</i> , 2021, 120, 1510-1521. | 0.2 | 33 |
| 7 | Allosteric activation of T β cell antigen receptor signaling by quaternary structure relaxation. <i>Cell Reports</i> , 2021, 36, 109375. | 2.9 | 23 |
| 8 | Multi-scale simulations of the T cell receptor reveal its lipid interactions, dynamics and the arrangement of its cytoplasmic region. <i>PLoS Computational Biology</i> , 2021, 17, e1009232. | 1.5 | 13 |
| 9 | Sphingomyelinase Disables Inactivation in Endogenous PIEZO1 Channels. <i>Cell Reports</i> , 2020, 33, 108225. | 2.9 | 47 |
| 10 | Distortion of the bilayer and dynamics of the BAM complex in lipid nanodiscs. <i>Communications Biology</i> , 2020, 3, 766. | 2.0 | 32 |
| 11 | RBCs prevent rapid PIEZO1 inactivation and expose slow deactivation as a mechanism of dehydrated hereditary stomatocytosis. <i>Blood</i> , 2020, 136, 140-144. | 0.6 | 23 |
| 12 | Multiple lipid binding sites determine the affinity of PH domains for phosphoinositide-containing membranes. <i>Science Advances</i> , 2020, 6, eaay5736. | 4.7 | 44 |
| 13 | Inter-domain dynamics in the chaperone SurA and multi-site binding to its outer membrane protein clients. <i>Nature Communications</i> , 2020, 11, 2155. | 5.8 | 48 |
| 14 | Membrane Recognition and Binding by the Phosphatidylinositol Phosphate Kinase PIP5K1A: A Multiscale Simulation Study. <i>Structure</i> , 2019, 27, 1336-1346.e2. | 1.6 | 22 |
| 15 | Force Sensing by Piezo Channels in Cardiovascular Health and Disease. <i>Arteriosclerosis, Thrombosis, and Vascular Biology</i> , 2019, 39, 2228-2239. | 1.1 | 147 |
| 16 | Molecular Simulations of Intact Anion Exchanger 1 Reveal Specific Domain and Lipid Interactions. <i>Biophysical Journal</i> , 2019, 117, 1364-1379. | 0.2 | 16 |
| 17 | β 1 integrin is a sensor of blood flow direction. <i>Journal of Cell Science</i> , 2019, 132, . | 1.2 | 41 |
| 18 | Modes of Interaction of Pleckstrin Homology Domains with Membranes: Toward a Computational Biochemistry of Membrane Recognition. <i>Journal of Molecular Biology</i> , 2018, 430, 372-388. | 2.0 | 42 |

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|----|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----|-----------|
| 19 | Distinctive phosphoinositide- and Ca ²⁺ -binding properties of normal and cognitive performance-linked variant forms of KIBRA C2 domain. <i>Journal of Biological Chemistry</i> , 2018, 293, 9335-9344. | 1.6 | 8 |
| 20 | Structural Lipids Enable the Formation of Functional Oligomers of the Eukaryotic Purine Symporter UapA. <i>Cell Chemical Biology</i> , 2018, 25, 840-848.e4. | 2.5 | 64 |
| 21 | Interaction of the human erythrocyte Band 3 anion exchanger 1 (AE1, SLC4A1) with lipids and glycophorin A: Molecular organization of the Wright (Wr) blood group antigen. <i>PLoS Computational Biology</i> , 2018, 14, e1006284. | 1.5 | 37 |
| 22 | Interactions of the EphA2 Kinase Domain with PIPs in Membranes: Implications for Receptor Function. <i>Structure</i> , 2018, 26, 1025-1034.e2. | 1.6 | 33 |
| 23 | Dynamic interactions between a membrane binding protein and lipids induce fluctuating diffusivity. <i>Science Advances</i> , 2017, 3, e1601871. | 4.7 | 59 |
| 24 | Structure and lipid-binding properties of the kindlin-3 pleckstrin homology domain. <i>Biochemical Journal</i> , 2017, 474, 539-556. | 1.7 | 40 |
| 25 | Effects of Periplasmic Chaperones and Membrane Thickness on BamA-Catalyzed Outer-Membrane Protein Folding. <i>Journal of Molecular Biology</i> , 2017, 429, 3776-3792. | 2.0 | 63 |
| 26 | The Integrin Receptor in Biologically Relevant Bilayers: Insights from Molecular Dynamics Simulations. <i>Journal of Membrane Biology</i> , 2017, 250, 337-351. | 1.0 | 29 |
| 27 | Multiscale Simulations Suggest a Mechanism for the Association of the Dok7 PH Domain with PIP-Containing Membranes. <i>PLoS Computational Biology</i> , 2016, 12, e1005028. | 1.5 | 24 |
| 28 | On the interpretation of reflectivity data from lipid bilayers in terms of molecular-dynamics models. <i>Acta Crystallographica Section D: Structural Biology</i> , 2016, 72, 1227-1240. | 1.1 | 10 |
| 29 | Band 3, the human red cell chloride/bicarbonate anion exchanger (AE1, SLC4A1), in a structural context. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2016, 1858, 1507-1532. | 1.4 | 154 |
| 30 | Interactions of Pleckstrin Homology Domains with Membranes: Adding Back the Bilayer via High-Throughput Molecular Dynamics. <i>Structure</i> , 2016, 24, 1421-1431. | 1.6 | 68 |
| 31 | Anomalous Dynamics of a Lipid Recognition Protein on a Membrane Surface. <i>Scientific Reports</i> , 2016, 5, 18245. | 1.6 | 38 |
| 32 | Association of Peripheral Membrane Proteins with Membranes: Free Energy of Binding of GRP1 PH Domain with Phosphatidylinositol Phosphate-Containing Model Bilayers. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 1219-1224. | 2.1 | 47 |
| 33 | Molecular Dynamics Simulations of the Bacterial UraA H ⁺ -Uracil Symporter in Lipid Bilayers Reveal a Closed State and a Selective Interaction with Cardiolipin. <i>PLoS Computational Biology</i> , 2015, 11, e1004123. | 1.5 | 40 |
| 34 | Interactions of peripheral proteins with model membranes as viewed by molecular dynamics simulations. <i>Biochemical Society Transactions</i> , 2014, 42, 1418-1424. | 1.6 | 31 |
| 35 | Interactions of Phosphatase and Tensin Homologue (PTEN) Proteins with Phosphatidylinositol Phosphates: Insights from Molecular Dynamics Simulations of PTEN and Voltage Sensitive Phosphatase. <i>Biochemistry</i> , 2014, 53, 1724-1732. | 1.2 | 42 |
| 36 | Interactions of the Auxilin-1 PTEN-like Domain with Model Membranes Result in Nanoclustering of Phosphatidyl Inositol Phosphates. <i>Biophysical Journal</i> , 2013, 105, 137-145. | 0.2 | 24 |

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|----|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----|-----------|
| 37 | Conformational Changes in Talin on Binding to Anionic Phospholipid Membranes Facilitate Signaling by Integrin Transmembrane Helices. PLoS Computational Biology, 2013, 9, e1003316. | 1.5 | 30 |
| 38 | A Helix Heterodimer in a Lipid Bilayer: Prediction of the Structure of an Integrin Transmembrane Domain via Multiscale Simulations. Structure, 2011, 19, 1477-1484. | 1.6 | 39 |
| 39 | Multiscale simulations suggest a mechanism for integrin inside-out activation. Proceedings of the National Academy of Sciences of the United States of America, 2011, 108, 11890-11895. | 3.3 | 62 |
| 40 | The Structure of the Talin/Integrin Complex at a Lipid Bilayer: An NMR and MD Simulation Study. Structure, 2010, 18, 1280-1288. | 1.6 | 57 |