

# Joseph Rebehmed

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2830465/publications.pdf>

Version: 2024-02-01

23  
papers

470  
citations

840776

11  
h-index

713466

21  
g-index

23  
all docs

23  
docs citations

23  
times ranked

890  
citing authors

| #  | ARTICLE  | IF  | CITATIONS |
|----|--|-----|-----------|
| 1  | Shaking the $\hat{I}^2$ -Bulges. IEEE/ACM Transactions on Computational Biology and Bioinformatics, 2022, 19, 14-18.   | 3.0 | 3         |
| 2  | Current status of PTMs structural databases: applications, limitations and prospects. Amino Acids, 2022, 54, 575-590.  | 2.7 | 9         |
| 3  | Editorial: Advances in Molecular Docking and Structure-Based Modelling. Frontiers in Molecular Biosciences, 2022, 9, 839826.   | 3.5 | 2         |
| 4  | Discrete analyses of protein dynamics. Journal of Biomolecular Structure and Dynamics, 2020, 38, 2988-3002.  | 3.5 | 20        |
| 5  | Combining In Silico Phylogenetic and Threading Approaches to Assist the in vitro Protein Engineering of BVMO Enzymes. Biophysical Journal, 2020, 118, 45a.   | 0.5 | 1         |
| 6  | In silico prediction of protein flexibility with local structure approach. Biochimie, 2019, 165, 150-155.  | 2.6 | 15        |
| 7  | A topology-based investigation of protein interaction sites using Hydrophobic Cluster Analysis. Biochimie, 2019, 167, 68-80.   | 2.6 | 12        |
| 8  | Investigation of the impact of PTMs on the protein backbone conformation. Amino Acids, 2019, 51, 1065-1079.  | 2.7 | 17        |
| 9  | iPBAvizu: a PyMOL plugin for an efficient 3D protein structure superimposition approach. Source Code for Biology and Medicine, 2019, 14, 5.  | 1.7 | 14        |
| 10 | Dynamics and deformability of $\hat{I}^{\pm}$ , $\beta$ 10- and $\beta$ €-helices. Archives of Biological Sciences, 2018, 70, 21-31.   | 0.5 | 10        |
| 11 | Enhancing the Activity of a <i>Dietzia</i> sp. D5 Baeyer-Villiger Monooxygenase towards Cyclohexanone by Saturation Mutagenesis. ChemistrySelect, 2017, 2, 7169-7177.  | 1.5 | 7         |
| 12 | The respective roles of polar/nonpolar binary patterns and amino acid composition in protein regular secondary structures explored exhaustively using hydrophobic cluster analysis. Proteins: Structure, Function and Bioinformatics, 2016, 84, 624-638. | 2.6 | 12        |
| 13 | Enzymatic Synthesis of Galactosylated Serine/Threonine Derivatives by $\hat{I}^2$ -Galactosidase from <i>Escherichia coli</i> . International Journal of Molecular Sciences, 2015, 16, 13714-13728.  | 4.1 | 4         |
| 14 | Protein flexibility in the light of structural alphabets. Frontiers in Molecular Biosciences, 2015, 2, 20.   | 3.5 | 71        |
| 15 | PTM-SD: a database of structurally resolved and annotated posttranslational modifications in proteins. Database: the Journal of Biological Databases and Curation, 2014, 2014, bau041-bau041.  | 3.0 | 44        |
| 16 | Expanding the SRI domain family: A common scaffold for binding the phosphorylated C-terminal domain of RNA polymerase II. FEBS Letters, 2014, 588, 4431-4437.  | 2.8 | 19        |
| 17 | Efficient Protein Structure Alignment Methods Based on a Structural Alphabet. Biophysical Journal, 2014, 106, 54a.   | 0.5 | 0         |
| 18 | Cis-trans isomerization of omega dihedrals in proteins. Amino Acids, 2013, 45, 279-289.  | 2.7 | 60        |

| #  | ARTICLE  | IF  | CITATIONS |
|----|--|-----|-----------|
| 19 | Î²â€Bulges: Extensive structural analyses of Î²â€sheets irregularities. <i>Protein Science</i> , 2013, 22, 1366-1378.  | 7.6 | 27        |
| 20 | Evolution study of the Baeyerâ€Villiger monoxygenases enzyme family: Functional importance of the highly conserved residues. <i>Biochimie</i> , 2013, 95, 1394-1402.          | 2.6 | 19        |
| 21 | 2D and 3D QSAR studies of diarylpyrimidine HIV-1 reverse transcriptase inhibitors. <i>Journal of Computer-Aided Molecular Design</i> , 2008, 22, 831-841.                      | 2.9 | 10        |
| 22 | Flexible computational docking studies of new aminoglycosides targeting RNA 16S bacterial ribosome site. <i>European Journal of Medicinal Chemistry</i> , 2008, 43, 1648-1656. | 5.5 | 26        |
| 23 | Molecular modeling studies of N-substituted pyrrole derivativesâ€Potential HIV-1 gp41 inhibitors. <i>Biorganic and Medicinal Chemistry</i> , 2008, 16, 3039-3048.             | 3.0 | 68        |