

# Renana Gershoni Poranne

Assistant Professor

Schulich Faculty of Chemistry – Technion, Israel

✉ rporanne@technion.ac.il

🌐 <https://poranne-group.github.io/>

## List of Publications and Software

October 2024

Notes: 1) \* denotes corresponding author. 2) Titles in red are clickable links.

### Peer-Reviewed Publications.....

42. S. Garhwal, S. Raje, K. Mlodzikowska-Pienko, T. S. Mohammed, R. Rafaeli, N. Fridman, L. J. W. Shimon, [R. Gershoni-Poranne\\*](#), and G. de Ruiter\*.  
*N<sub>2</sub> Dissociation vs. Reversible 1,2-Methyl Migration in PCNHCP Cobalt(I) Complexes in the Stereoselective Isomerization (E/Z) of Allyl Ethers*  
JACS Au **2024**, In Press.  
⇒ **Selected for a Supplementary Front Cover**
41. Y. Zhu, Z. Zhou, Z. Wei, A. Tsybizova, [R. Gershoni-Poranne\\*](#), and M. A. Petrukhina\*  
*What a Difference an Electron Makes: Structural Response of Saddle-Shaped Tetraphenylene to One and Two Electron Uptake*  
ChemistryEurope **2024**, 2, e202400055.  
⇒ **Selected for a Front Cover**  
⇒ **Highlighted in ChemistryViews**
40. S. Chakraborty, E. Mayo Yanes, and [R. Gershoni-Poranne\\*](#)  
*Hetero-Polycyclic Aromatic Systems: A Data-Driven Investigation of Structure-Property Relationships*  
Beilstein Journal of Organic Chemistry **2024**, 20, 1817.
39. A. Wahab and [R. Gershoni-Poranne\\*](#)  
*COMPAS-3: a dataset of peri-condensed polybenzenoid hydrocarbons*  
Physical Chemistry Chemical Physics **2024**, 26, 15344.
38. S. Jalife, A. Tsybizova, [R. Gershoni-Poranne\\*](#), and J. I. Wu\*  
*Modulating Paratropicity in Heteroarene-Fused Expanded Pentalenenes*  
Organic Letters **2024**, 26, 1293.
37. E. Mayo Yanes, S. Chakraborty, and [R. Gershoni-Poranne\\*](#)  
*COMPAS-2: a Dataset of Cata-Condensed Hetero-Polycyclic Aromatic Systems*  
Scientific Data **2024**, 11, 97.
36. Z. Yang, R. Nandi, A. Orieshyna, [R. Gershoni-Poranne](#), S. Zhang\*, and N. Amdursky\*  
*Light-Triggered Enhancement of Fluorescence Efficiency in Organic Cages*  
Journal of Physical Chemistry Letters **2023**, 15, 136.
35. T. Weiss, E. Mayo Yanes, S. Chakraborty, A. M. Bronstein\* and [R. Gershoni-Poranne\\*](#)  
*Guided Diffusion for Inverse Molecular Design*  
Nature Computational Science **2023**, 3, 873.  
⇒ **Selected for a Front Cover**  
⇒ **Highlighted in News & Views**

34. M. Pennachio, Z. Wei, R. G. Clevenger, K. V. Kilway, A. Tsybizova, R. Gershoni-Poranne,\* and M. A. Petrukhina\*  
*Repercussions of Multi-Electron Uptake by a Twistacene: A Reduction-Induced Double Dehydrogenative Annulation*  
 Organic Chemistry Frontiers **2023**, 10, 5823.  
 ⇒ **Selected for an Inside Front Cover**
33. R. Gershoni-Poranne\* and A. Tsybizova  
*A Crowning Achievement: The First Solution-Phase Synthesis of Circumcoronenes*  
 Angewandte Chemie Int. Ed. **2023**, 62, e202305289.
32. T. Weiss, A. Wahab, A. M. Bronstein and R. Gershoni-Poranne\*  
*Interpretable Deep-Learning Unveils Structure-Property Relationships in Polybenzenoid Hydrocarbons*  
 Journal of Organic Chemistry **2023**, 88, 9645.  
 ⇒ **Selected for a Front Cover**
31. M. Pennachio, Z. Zhou, Z. Wei, A. Tsybizova, R. Gershoni-Poranne\*, and M. A. Petrukhina\*  
*Interplay of Charge and Aromaticity Upon Chemical Reduction of p-Quinquephenyl with Alkali Metals*  
 Organometallics **2023**, 42, 2492.
30. S. Fite, A. Wahab, E. Paenurk, Z. Gross and R. Gershoni-Poranne\*  
*Text-Based Representations with Interpretable Machine Learning Reveal Structure-Property Relationships of Polybenzenoid Hydrocarbons*  
 Journal of Physical Organic Chemistry **2022**, 36, e4458.  
 ⇒ **Invited contribution for the special issue on Excited State Aromaticity and Antiaromaticity**
29. A. Wahab, L. Pfuderer, E. Paenurk, and R. Gershoni-Poranne\*  
*The COMPAS Project: A Computational Database of Polycyclic Aromatic Systems. Phase 1: cata-Condensed Polybenzenoid Hydrocarbons*  
 Journal of Chemical Information and Modeling **2022**, 62, 3704.  
 ⇒ **Selected for a Front Cover**
28. Z. Zhou, D. T. Egger, C. Hu, M. Pennachio, Z. Wei, R. K. Kawade, Ö. Üngör, R. Gershoni-Poranne\*, M. A. Petrukhina\*, and I. V. Alabugin\*  
*Localized Antiaromaticity Hot-spot Drives Reductive Dehydrogenative Cyclizations in Bis- and Mono-Helicenes*  
 Journal of the American Chemical Society **2022**, 144, 12321.  
 ⇒ **Selected for a Front Cover**
27. E. Paenurk\* and R. Gershoni-Poranne\*  
*Simple and Efficient Visualization of Aromaticity: Bond Currents Calculated from NICS Values*  
 Physical Chemistry Chemical Physics **2022**, 24, 8631.  
 ⇒ **Selected for a Front Cover**
26. R. Thenarukandiyil, E. Paenurk, A. Wong, N. Fridman, A. Karton, R. Carmieli, G. Ménard, R. Gershoni-Poranne\*, and G. de Ruiter\*  
*Extensive Redox Non-Innocence in Iron Bipyridine-Diimine Complexes: a Combined Spectroscopic and Computational Study*  
 Inorganic Chemistry **2021**, 60, 18296.
25. Z. Zhou, Ö. Üngör, Z. Wei, M. Shatruk\*, A. Tsybizova, R. Gershoni-Poranne\*, and M. A. Petrukhina\*  
*Tuning Magnetic Interactions Between Triphenylene Radicals by Variation of Crystal Packing in Structures with Alkali Metal Counterions*  
 Inorganic Chemistry **2021**, 60, 14844.

24. G. Markert, E. Paenurk, and R. Gershoni-Poranne\*  
*Prediction of Spin Density, Baird-Antiaromaticity, and Singlet-Triplet Energy Gap in Triplet-State Polybenzenoid Systems from Simple Structural Motifs*  
 Chemistry - A European Journal **2021**, 27, 6923.  
 ⇒ **Selected for a Cover Feature**  
 ⇒ **Denoted as a Hot Paper**
23. E. Paenurk, S. Feusi, and R. Gershoni-Poranne\*  
*Predicting Bond-currents in Polybenzenoid Hydrocarbons with an Additivity Scheme*  
 Journal of Chemical Physics **2021**, 154, 024110.  
 ⇒ **Invited contribution for the Issue Honoring Women in Chemical Physics and Physical Chemistry**
22. M. A. Hope, T. Nakamura, P. Ahlawat, A. Mishra, M. Cordova, F. Jahanbakhshi, M. Mladenović, R. Runjhun, L. Merten, A. Hinderhofer, B. I. Carlsen, D. J. Kubicki, R. Gershoni-Poranne, T. Schneeberger, L. C. Carbone, Y. Liu, S. M. Zakeeruddin, J. Lewinski, A. Hagfeldt, F. Schreiber, U. Rothlisberger, M. Grätzel\*, J. V. Milić\*, and L. Emsley\*  
*Nanoscale Phase Segregation in Supramolecular  $\pi$ -Templating for Hybrid Perovskite Photovoltaics from NMR Crystallography*  
 Journal of the American Chemical Society **2021**, 143, 1529.
21. T. Schnitzer, E. Paenurk, N. Trapp, R. Gershoni-Poranne, and H. Wennemers\*  
*Peptide-Metal Frameworks with Metal Strings Guided by Dispersion Interactions*  
 Journal of the American Chemical Society **2021**, 143, 644.
20. A. Wahab, F. Fleckenstein, S. Feusi, and R. Gershoni-Poranne\*  
*Predi-XY: A Python program for automated generation of NICS-XY-Scans based on an Additivity Scheme*  
 Electronic Structure **2020**, 2, 047002.  
 ⇒ **Invited contribution for the Emerging Leaders issue**  
 ⇒ **Selected as Editor's Choice paper**
19. E. Solel, D. Pappo, O. Reany, T. Mejuch, R. Gershoni-Poranne, M. Botoshansky, A. Stanger, and E. Keinan\*  
*Flat corannulene: when a transition state becomes a stable molecule*  
 Chemical Science **2020**, 11, 13015.  
 ⇒ **Selected for a Cover Feature**
18. S. Eichenberger, M. Hönig, M. J. R. Richter, R. Gershoni-Poranne\* and E. M. Carreira\*  
*Ring-fused cyclobutanes via cycloisomerization of alkylidenecyclopropane acylsilanes*  
 Chemical Science **2020**, 11, 5294.
17. M. A. Ruiz-Preciado, D. J. Kubicki, A. Hofstetter, L. McGovern, M. H. Futscher, A. Ummadisingu, R. Gershoni-Poranne, S. M. Zakeeruddin, B. Ehrler, L. Emsley\*, J. V. Milić\*, and M. Grätzel\*  
*Supramolecular Modulation of Hybrid Perovskite Solar Cells via Bifunctional Halogen Bonding Revealed by Two-Dimensional  $^{19}\text{F}$  Solid-State NMR Spectroscopy*  
 Journal of the American Chemical Society **2020**, 142, 1645.
16. Z. Zhou, R. K. Kawade, Z. Wei, F. Kuriakose, Ö. Üngör, M. Jo, M. Shatruk, R. Gershoni-Poranne\*, M. A. Petrukhina\*, and I. V. Alabugin\*  
*Negative charge as a lens for concentrating antiaromaticity: using pentagonal "defect" and helicene strain for cyclizations*  
 Angewandte Chemie Int. Ed. **2020**, 59, 1256.

15. P. Finkelstein and R. Gershoni-Poranne\*  
*An Additivity Scheme for Aromaticity: The Heteroatom Case*  
ChemPhysChem **2019**, 20, 1508.
14. J. V. Milić, C., N. Hellou, F. Isenrich, R. Gershoni-Poranne, D. Neshchadin, S. Egloff, N. Trapp, L. Ruhlmann, C. Boudon, G. Gescheidt, J. Crassous, and F. Diederich\*  
*Light-Responsive Pyrazine-Based Systems: Probing Aromatic Diarylethene Photocyclization*  
Journal of Physical Chemistry C **2018**, 122, 19100.
13. R. Gershoni-Poranne\*, A. P. Rahalkar, and A. Stanger\*  
*The Predictive Power of Aromaticity: Quantitative Correlation between Aromaticity and Ionization Potentials and HOMO-LUMO Gaps in Oligomers of Benzene, Pyrrole, Furan, and Thiophene*  
Physical Chemistry Chemical Physics **2018**, 20, 14808.
12. R. Gershoni-Poranne\*  
*Piecing it Together: An Additivity Scheme for Aromaticity using NICS-XY-Scans*  
Chemistry – A European Journal **2018**, 24, 4165.
11. S. Künzi, R. Gershoni-Poranne, and P. Chen\*  
*Mechanistic Studies on the Nickel-Catalyzed Cyclopropanation with Lithiomethyltrimethylammonium Triflate*  
Organometallics **2019**, 38, 1928.
10. P. Chen\* and R. Gershoni-Poranne  
*Response to “Covalent Bonding and Charge Shift Bonds: Comment on ‘The Carbon–Nitrogen Bonds in Ammonium Compounds Are Charge Shift Bonds’”*  
Chemistry – A European Journal **2017**, 23, 18325.
9. E. Paenurk, R. Gershoni-Poranne, and P. Chen\*  
*Trends in Metallophilic Bonding in Pd-Zn and Pd-Cu Complexes*  
Organometallics **2017**, 36, 4854.
8. R. Gershoni-Poranne and P. Chen\*  
*The C-N Bonds in Ammoniums are Charge Shift Bonds*  
Chemistry – A European Journal **2017**, 23, 4659.
7. R. Gershoni-Poranne and A. Stanger\*  
*Magnetic Criteria of Aromaticity*  
**Invited Review** Chemical Society Reviews **2015**, 44, 6597.
6. M. Schaffroth, R. Gershoni-Poranne, A. Stanger\*, and U. H. F. Bunz\*  
*Tetraazacenes Containing Four-membered Rings in Different Oxidation States. Are They Aromatic? A Computational Study*  
Journal of Organic Chemistry **2014**, 79, 11644.
5. R. Gershoni-Poranne and A. Stanger\*  
*The NICS-XY-Scan: Identification of Local and Global Ring Currents in Multi-Ring Systems*  
Chemistry – A European Journal **2014**, 20, 5673.
4. R. Gershoni-Poranne, C. M. Gibson, P. W. Fowler, and A. Stanger\*  
*Concurrence between Current Density, Nucleus-Independent Chemical Shifts, and Aromatic Stabilization Energy: The Case of Isomeric [4]- and [5]Phenylenes*  
Journal of Organic Chemistry **2013**, 78, 7544.

3. R. Gershoni-Poranne and A. Stanger\*  
*An MO-Based Identification of Charge-Shift Bonds*  
ChemPhysChem **2012**, *13*, 2377.
2. M. Standera, R. Haefliger, R. Gershoni-Poranne, A. Stanger, G. Jeschke, J. D. van Beek, and A. D. Schlüter\*  
*Evidence for Fully Conjugated Double-Stranded Cycles*  
Chemistry – A European Journal **2011**, *17*, 12163.
1. R. Gershoni-Poranne, D. Pappo, E. Solel, and E. Keinan\*  
*Corannulene Ethers Via Ullmann Condensation*  
Organic Letters **2009**, *11*, 5146.

### Preprint / Under Review.....

3. G. I. Warren, K. Mlodzikowska-Pienko, S. Jalife, I. S. Demachkie, J. I. Wu, M. M. Haley, and R. Gershoni-Poranne\*  
*Effects of Benzoheterocyclic Annulation on the s-Indacene Core: a Computational Analysis*  
Submitted **2024**, currently under review.
2. A. Wahab and R. Gershoni-Poranne\*  
*The  $xTB$  fractional occupation number enables identification of polyradical character in polybenzenoid hydrocarbons*  
Submitted **2024**, currently under review.
1. Y. Davidson, A. Philipp, S. Chakraborty, A. M. Bronstein, and R. Gershoni-Poranne\*  
*How Local is 'Local'? Deep Learning Reveals Locality of the Induced Magnetic Field of Polycyclic Aromatic Hydrocarbons*  
Submitted **2024**, currently under review.

### Book Chapters.....

1. R. Gershoni-Poranne\* and A. Stanger\*  
*Chapter 4: NICS – Nucleus Independent Chemical Shifts*  
in Aromaticity: Modern Computational Methods and Applications, **2021**  
Edited by I. Fernandez.

### Software.....

Notes: All of our software is freely available to download from the *Poranne Group Repository*.

2. BC-Wizard  
Python package implementing the NICS2BC method for calculating bond-currents from NICS values.
1. Predi-XY  
Python package implementing an additivity scheme for rapid generation of NICS-XY-Scans for polycyclic aromatic systems.