# Dr. NIVEDITA ACHARJEE

Contact: nivchem@gmail.com

Qualification: M.Sc., PhD. (University of Calcutta, Kolkata, India)

**Designation:** Assistant Professor (West Bengal Educational Service)

Teaching Experience (UG/PG) (in years): UG: 14 years; PG: 9 years

Research Experience (in years): 17 years

**Institutional Address**: Department of Chemistry, Durgapur Government College, Durgapur, J. N. Avenue, District-Paschim Bardhaman, West Bengal, Pin Code-713214

## **Specialization/Research Interest:**

The research area of Dr. Nivedita Acharjee is Computational Chemistry using conceptual Density Functional Theory (DFT) calculations for chemical systems. This research work involves studying the energetics of chemical reactions, radical spin traps, spectroscopy, photovoltaic technology and also structuring the plausible mechanism of chemical reactions. Presently, we are studying chemical reactions by the newly developed theoretical outlook "Molecular Electron Density Theory" (MEDT) proposed by Domingo in 2016.

## **International Research Collaborations:**

- [1] Professor Luis R Domingo, Department of Organic Chemistry, University of Valencia, Spain <a href="http://www.luisrdomingo.com/p/colaboraciones.html">http://www.luisrdomingo.com/p/colaboraciones.html</a>
- [2] Prof. H. A. Mohammad-Salim, Department of Chemistry, University of Zakho, Duhok 42001, Iraq
- [3] Prof. Khalid Abichhe, Applied Chemistry-Physics and Environment Team, Faculty of Sciences, Ibn Zohr University, B.P.8106 Cité Dakhla, Agadir, Morocco, Polydisciplinary Faculty of Taroudant, Ibn Zohr University, Hay El Mohammadi (Lastah), Taroudant, Morocco.
- [4] Prof. Mohammed Salah, Team of Chemoinformatics Research and Spectroscopy and Quantum Chemistry, Department of Chemistry, Faculty of Science, University Chouaib Doukkali, B. P. 20, 2300, El Jadida, Morocco; LS3MN2E, Department of Chemistry, Mohammed V University, Faculty of Sciences Rabat, Morocco.
- [5] Professor Thierry Prange [Laboratorie de Cristallographie et RMN Biologiques, Universite Paris Descartes, France].
- [6] Jamelah S. Al-Otaibi, Department of Chemistry, College of Science, Princess Nourah bint Abdulrahman University, Riyadh, Saudi Arabia

## **National Collaborations:**

- 1. Professor Avijit Banerji [Retired Professor, University of Calcutta and Formerly Professor Ashutosh Mukherjee Fellow, National Institute for Ayurvedic Drug Development (NRIADD), West Bengal, India
  - 2. Dr. Ani Deepthi, Department of Chemistry, University of Kerela, India
- **3.** Dr. Madhu Ganesh, Department of Chemistry, B.M.S College of Engineering, Bull Temple Road, Bengaluru 560019, India and Department of Pharmaceutical Technology, National Institute of Pharmaceutical Education & Research, Balanagar, Hyderabad-500037, Telangana, India
  - 4. Y. Sheena Mary and Y.Shyma Mary, Thushara, Kollam, Kerala, India

#### **BIBLIOMETRICS**

RESEARCH GATE LINK: https://www.researchgate.net/profile/Nivedita-Acharjee

**ORCID ID:** https://orcid.org/0000-0001-8354-8693

**SCOPUS ID:** <a href="https://www.scopus.com/authid/detail.uri?authorId=34978096200">https://www.scopus.com/authid/detail.uri?authorId=34978096200</a>

**WEB OF SCIENCE RESEARCHER ID (**AAS-2991-2020**):** (SCI indexed publications)

https://publons.com/researcher/3740315/nivedita-acharjee/

VIDWAN INFLIBNET PROFILE: <a href="https://vidwan.inflibnet.ac.in/profile/184722">https://vidwan.inflibnet.ac.in/profile/184722</a>

#### **Reviewer of International Journals:**

- Nature Communications (Nature Publishing House)
- ACS Sustainable Resource Management (American Chemical Society)
- Chemistry Select (Wiley)
- New Journal of Chemistry (Royal Society of Chemistry)
- BMC Chemistry (Springer-Nature)
- Journal of Molecular Graphics and Modelling (Elsevier)
- International Journal of Quantum Chemistry (Wiley)
- Theoretical Chemistry Accounts (Springer-Nature)
- Chemistry-Africa (Springer-Nature)
- Structural Chemistry (Springer-Nature)
- Indian Journal of Chemistry (NISCAIR)
- Indian Journal of Chemistry-Section A (NISCAIR)
- Oriental Journal of Chemistry
- Journal of Chemical and Pharmaceutical Research

# **Research Publications**

- Jamelah S Al-Otaibi, Y.Sheena Mary, Y.Shyma Mary, Asmita Mondal, Nivedita Acharjee, Deepthi S. Rajendran Nair (2024) Adsorption of toremifene on carboxylic acid functionalized N/B-doped coronene: DFT, AIM, Solvation, docking and MD analyses *Journal of Molecular Liquids (Elsevier)*, 395, 123931 IF: 6.00) https://doi.org/10.1016/j.molliq.2023.123931
- 2. Aneesh Kumar R, Jamelah S. Al-Otaibi, Y. Sheena Mary, Y. Shyma Mary, **Nivedita Acharjee**, Renjith Thomas, Leena T L (2024) Surface adsorption of adenine on pristine and B/N/O/P-doped coronene as biosensing substrate for DNA detection- DFT Study. *Journal of Molecular Liquids (Elsevier)*, 393, 123546 (**IF: 6.00**) https://doi.org/10.1016/j.molliq.2023.123546
- 3. Dr. Jamelah S Al-Otaibi, Y.Sheena Mary; Y.Shyma Mary; Asmita Mondal; **Nivedita Acharjee**; Deepthi S. Rajendran Nair (2024) Investigation of the interaction of thymine drugs with Be12O12 and Ca12O12 nanocages: a quantum chemical study, *Spectrochimica Acta Part A: Molecular and Biomolecular Spectroscopy (Elsevier)*, 308 (2024) 123728 (**IF: 4.4**) https://doi.org/10.1016/j.saa.2023.123728
- 4. Bhaskar Mondal, Luis R. Domingo, Haydar A. Mohammad-Salim, Nivedita Acharjee (2024) Unveiling [3+2] cycloaddition reactions of diphenyl diazomethane to thiobenzophenone and cycloaliphatic thioketones in the light of molecular electron density theory, *Computational and Theoretical Chemistry (Elsevier)*, 114660 (**IF: 2.80**) <a href="https://doi.org/10.1016/j.comptc.2024.114660">https://doi.org/10.1016/j.comptc.2024.114660</a>
- Amine Rafik, Hamza El Hadki, Abdeljabbar Jaddi, Najia Komiha, Oum Keltoum Kabbaj, Nivedita Acharjee, Mohammed Salah, Mohammad Azam, khadija Marakchi (2024) A new investigation into the regioselectivity and stereoselectivity of the reaction between di-methyl nitrone and fluorinated alkenes: DFT, independent gradient model, and molecular docking, *Computational and Theoretical Chemistry (Elsevier)*, 1234, 114511 (IF: 2.80) <a href="https://doi.org/10.1016/j.comptc.2024.114511">https://doi.org/10.1016/j.comptc.2024.114511</a>
- 6. Moulay Driss MELLAOUI, Khalid Abbiche, **Nivedita Acharjee**, Haydar Mohammad-Salim, Abdallah Imjjad, Rachid Boutiddar, Khadija Marakchi, Souad El Issami, Hanane Zejli (2024)Unveiling the mechanism and selectivity of the [3 + 2] cycloaddition reactions of nitrone with acetylene derivatives leading to anticancer 4-isoxazoline derivatives from the MEDT perspective, *Computational and Theoretical Chemistry (Elsevier)*, 1237, 114619 (**IF: 2.80**), <a href="https://doi.org/10.1016/j.comptc.2024.114619">https://doi.org/10.1016/j.comptc.2024.114619</a>
- 7. Haitham K.R. Al-Sharifi; Ani Deepthi; Nivedita Acharjee; M S Swapna; S. Sankararaman (2024) A Window to Energy Efficiency: Tryptanthrin-Malononitrile Knoevenagel Adduct as an IR Window Material, *Journal of Materials Science: Materials in Electronics (Springer)* (**IF** = **2.440**) https://doi.org/10.1007/s10854-024-13018-0
- 8. Jamelah S Al-Otaibi, Fowzia S Alamro, Aljawhara H Almugrin, Y Sheena Mary, **Nivedita Acharjee** (2024) Adsorption, sensor properties, AIM analysis, docking of an anticancer drug benzamide on nanocones: SERS, solvent effects and DFT investigation, *Journal of Computational Biophysics and Chemistry* (World Scientific), (**IF** = **2.440**) <a href="https://doi.org/10.1142/S2737416524500091">https://doi.org/10.1142/S2737416524500091</a>
- 9. Abdellatif Rafik, Fatima Lakhdar, afid Zouihri, Taoufiq Guedira, **Nivedita Acharjee**, Mohammad Shahidul Islam, Mohammed Salah, Abdellah Zeroual (2024) Exploring nonlinear optical properties in a hybrid dihydrogen phosphate system: an experimental and theoretical approach, *Journal of Molecular Modeling* (Springer) (**IF** = **2.20**) 30:151, <a href="https://doi.org/10.1007/s00894-024-05936-x">https://doi.org/10.1007/s00894-024-05936-x</a>
- 10. Asmita Mondal, **Nivedita Acharjee\***, Haydar Mohammad-Salim, Mrinmoy Chakraborty (2024) A Molecular Electron Density Theory Study to understand intramolecular [3+2] cycloaddition reactions of azides and

- diazoalkanes, Structural Chemistry (Springer), <a href="https://doi.org/10.1007/s11224-023-02270-5">https://doi.org/10.1007/s11224-023-02270-5</a> (IF = 1.795)
- 11. Mohammed Salah, Oumaima El Alaoui El Abdallaoui, Abdellah Zeroual, **Nivedita Acharjee** and Mohammed El idrissi (2024) Insight into a new discovery of SARS-CoV-2 inhibitor activated through Chloroquine derivatives, *Current Chemistry Letters*, 13(1), 49-60. <a href="https://doi.org/10.5267/j.ccl.2023.8.010">https://doi.org/10.5267/j.ccl.2023.8.010</a>

- 12. Jamelah S Al-Otaibi, Y.Sheena Mary, Y.Shyma Mary, **Nivedita Acharjee**, Deepthi S. Rajendran Nair, Yathirajan HS (2023) An investigation on spectroscopic, wavefunction dependent reactivity, docking and anti-Covid-19 ability of Flupentixol dihydrochloride: DFT and MD simulations at different temperatures, *Journal of Molecular Liquids (Elsevier)*, 392, 123177 (**IF: 6.00**) <a href="https://doi.org/10.1016/j.molliq.2023.123177">https://doi.org/10.1016/j.molliq.2023.123177</a>
- 13. Luis R Domingo\*, **Nivedita Acharjee**, (2023) "Unveiling the Mg(II) Promoted [3+2] Cycloaddition Reaction of Mesitonitrile Oxide to Baylis-Hilman Adduct from the Molecular Electron Density Theory Perspective", *New Journal of Chemistry (Royal Society of Chemistry)*, 47, 2495-2506 (**IF** = **3.925**) https://doi.org/10.1039/D2NJ05424H
- 14. Jamelah S.Al-Otaibi\*, Y. Sheena Mary, Y. Shyma Mary, **Nivedita Acharjee**, Sreejit Soman (2023) Spectroscopic, reactivity analysis and docking studies of 3-(adamantan-1-yl)-4-(4-fluorophenyl)-1-[(4-phenylpiperazin-1-yl)methyl]-4,5-dihdyro-1H-1,2,4-triazole—5-thione: DFT and MD simulations, *Journal of Molecular Structure (Elsevier)*, 134418. (**IF = 3.841**) https://doi.org/10.1016/j.molstruc.2022.134418
- 15. Jamelah S. Al-Otaibi, Y.Sheena Mary, Y.Shyma Mary, Asmita Mondal, Nivedita Acharjee, Jineet kumar Gawad, Insights on adsorption properties of a DNA base, guanine on nano metal cages (Ag24/Au24/Cu24): DFT, SERS, NCI and solvent effects, *Journal of Molecular Structure* (Elsevier), 135541. (**IF** = **3.841**) <a href="https://doi.org/10.1016/j.molstruc.2023.135541">https://doi.org/10.1016/j.molstruc.2023.135541</a>
- 16. P. Vennila, Jamelah S. Al-Otaibi, G. Venkatesh, Y. Sheena Mary, V. Raj, **Nivedita Acharjee** & P. Tamilselvi (2023): Structural, Spectral, Molecular Docking, and Molecular Dynamics Simulations of Phenylthiophene-2-Carboxylate Compounds as Potential Anticancer Agents, *Polycyclic Aromatic Compounds (Taylor and Francis)* (**IF** = **3.744**) https://doi.org/10.1080/10406638.2023.2172052
- 17. A. Khannous a, S. Jorio, **N. Acharjee**, M. Azam, K. Min, M. Salah (2023) Unveiling the atropoisomerism induced facial selectivity and regioselectivity in the [3 + 2] cycloaddition reaction of benzonitrile oxide with 5-methylenehydantoin from the molecular electron density theory perspective, *Computational and Theoretical Chemistry* (*Elsevier*), Volume 1227, September 2023, 114243 (**IF** = **2.800**) <a href="https://doi.org/10.1016/j.comptc.2023.114243">https://doi.org/10.1016/j.comptc.2023.114243</a>
- 18. Bhaskar Mondal; Luis R. Domingo; Haydar A. Mohammad-Salim; **Nivedita Acharjee\*** (2023), Unveiling the mechanism, regiochemistry and substituent effects of the [3+2] cycloaddition reactions of C, N-diaryl nitrile imine to ethylene derivatives from the molecular electron density theory perspective, *Computational and Theoretical Chemistry (Elsevier)* https://doi.org/10.1016/j.comptc.2023.114393 (**IF** = **2.800**)
- 19. Jamelah S Al-Otaibi\*, Y.Sheena Mary, Y.Shyma Mary, **Nivedita Acharjee**, Abdulaziz A. Al-Saadi, Maria Cristina Gamberini (2023) Influence of pyramidal M20 (M=Cu, Ag, Au) clusters on SERS and non-covalent interactions towards tuberculosis drug Pretomanid (PTD): DFT study, *Journal of Computational Biophysics and Chemistry* (accepted) (**IF** = **2.440**)
- 20. Moulay Driss Mellaoui, **Nivedita Acharjee\***, Abdallah Imjjad, Jamal Koubachi, Abdellatif El Hammadi, Hassan Bourzi, Souad El Issami, Hanane Zejli, Majdi Hochlaf\*, and Khalid Abbiche\* (2023) Unveiling the mechanism, selectivity, solvent and temperature effects on the [3+2] cycloaddition reaction of N-methyl-C-(2-furyl) nitrone with maleimide derivatives from the molecular electron density theory perspective, *Theoretical Chemistry Accounts (Springer)*, 142, 33 (**IF** = **1.700**) <a href="https://doi.org/10.1007/s00214-023-02972-3">https://doi.org/10.1007/s00214-023-02972-3</a>
- 21. Asmita Mondal, Luis R Domingo, **Nivedita Acharjee**\* (2023) Revealing the influence of tether length on the intramolecular [3+2] cycloaddition reactions of nitrones from the molecular electron density theory perspective, *Journal of Physical Organic Chemistry* (*Wiley*) (**IF: 1.800**). <a href="https://doi.org/10.1002/poc.4574">https://doi.org/10.1002/poc.4574</a>
- 22. Barsali Banerjee, **Nivedita Acharjee**\*, Debnath Palit (2023) Revealing the cyclization selectivity in intramolecular [3+2] cycloaddition reactions of allenic nitrones from the molecular electron density theory perspective, *Structural Chemistry (Springer)*, (**IF** = **1.795**) https://doi.org/10.1007/s11224-023-02175-3
- 23. Asmita Mondal, **Nivedita Acharjee\***, Debnath Palit (2023) Unveiling [3+2] cycloaddition reactions of pyridinium bis(methoxycarbonyl)methylides and pyridinium dicyanomethylides with cyclooctyne for indolizine synthesis from the molecular electron density theory perspective, *Structural Chemistry (Springer)*, (**IF** = **1.795**) <a href="https://doi.org/10.1007/s11224-023-02193-1">https://doi.org/10.1007/s11224-023-02193-1</a>
- 24. Haydar Mohammad-Salim, Asmita Mondal, Jesus Vicente de Julián-Ortiz, **Nivedita Acharjee** (2023), Diastereoselective green synthesis of pyrrolo[1,2-a]quinolines via [3+2] cycloaddition reaction: insights from molecular electron density theory, *Theoretical Chemistry* Account (Springe)s, 142, 129, (**IF:1.7**), <a href="https://doi.org/10.1007/s00214-023-03068-8">https://doi.org/10.1007/s00214-023-03068-8</a>

- 25. Asmita Mondal, **Nivedita Acharjee\*** (2023) Unveiling the exclusive stereo- and site selectivity in [3+2] cycloaddition reactions of a tricyclic strained alkene with nitrile oxides from the molecular electron density theory perspective, *Chemistry of Heterocyclic Compounds Special Issue on "Cycloaddition reactions: Recent Progress (Springer)"* (**IF** = **1.490**) https://doi.org/10.1007/s10593-023-03176-0
- 26. Asmita Mondal, Hayder A Mohammad-Salim, **Nivedita Acharjee\*** (2023), Unveiling substituent effects in [3+2] cycloaddition reactions of benzonitrile N-oxide and benzylideneanilines from the molecular electron density theory perspective, *Scientific Radices*, Volume 2, page-88-105, <a href="https://doi.org/10.58332/scirad2023v2i1a05">https://doi.org/10.58332/scirad2023v2i1a05</a>

- 27. Jamelah S. Al-Otaibi\*, Y.Sheena Mary, Y.Shyma Mary, Asmita Mondal, **Nivedita Acharjee**, S. Balachandar (2022) Investigation of the adsorption of a DNA based purine derivative on N/B-doped coronene and coronene by means of DFT and NCI interaction analysis, *Journal of Molecular Liquids (Elsevier)*, 367, 120373 (**IF** = **6.633**) <a href="https://doi.org/10.1016/j.molliq.2022.120373">https://doi.org/10.1016/j.molliq.2022.120373</a>
- 28. Jamelah S. Al-Otaibi\*, Y. Sheena Mary, Y. Shyma Mary, Nivedita Acharjee, S.Balachandar, H.S.Yathirajan (2022), Insights into solvation effects, spectroscopic, Hirshfeld Surface Analysis, reactivity analysis and anti-Covid-19 ability of doxylamine succinate: Experimental, DFT, MD and docking simulations, *Journal of Molecular Liquids (Elsevier)*, 361, 119609 (IF = 6.633) <a href="https://doi.org/10.1016/j.molliq.2022.119609">https://doi.org/10.1016/j.molliq.2022.119609</a>
- 29. Jamelah S. Al-Otaibi\*, Y.Sheena Mary, Y.Shyma Mary, Asmita Mondal, **Nivedita Acharjee** (2022) DFT analysis on the adsorption of melamine in Ga12-N12/P12 nanocages: solvent effects, SERS analysis, reactivity properties, Journal of Biomolecular Structure and Dynamics (*Taylor and Francis*) (**IF** = **5.235**) https://doi.org/10.1080/07391102.2022.2137700
- 30. **Nivedita Acharjee\***, Asmita Mondal, Mrinmoy Chakraborty (2022) "Unveiling intramolecular [3+2] cycloaddition reactions of C, N-disubstituted nitrones with a molecular electron density theory perspective", *New Journal of Chemistry (Royal Society of Chemistry)*, 2022, 46, 7721-7735 (**IF** = **3.925**) <a href="https://doi.org/10.1039/d2nj00888b">https://doi.org/10.1039/d2nj00888b</a>
- 31. Luis R Domingo\*, Mar-Rioz Gutiérrez, **Nivedita Acharjee** (2022) "A Molecular Electron Density Theory Study of the Lewis Acid Catalyzed [3+2] Cycloaddition Reactions of Nitrones with Nucleophilic Ethylenes", *European Journal of Organic Chemistry* (*Wiley*) 2022, e202101417 (**IF** = **3.261**) <a href="https://doi.org/10.1002/ejoc.202101417">https://doi.org/10.1002/ejoc.202101417</a>
- 32. Luis R. Domingo\*, Mar Ríos-Gutiérrez, **Nivedita Acharjee** (2022) "Unveiling the cb-type Intramolecular [3+2] Cycloaddition Reactions of Fluorinated Azomethine Ylides to Ester Carbonyls with a Molecular Electron Density Theory Perspective", *ChemistrySelect* (*Wiley*), 7, e202201845(**IF** = **2.307**) https://doi.org/10.1002/slct.202201845
- 33. Jamelah S. Al-Otaibi\*, Y. Sheena Mary, Y. Shyma Mary, Sreejit Soman, **Nivedita Acharjee**, B. Narayana (2022), Theoretical and experimental investigation of a pyrazole derivative- solvation effects, reactivity analysis and MD simulations, *Chemical Physics Letters* (*Elsevier*), 139469. (**IF** = **2.719**) https://doi.org/10.1016/j.cplett.2022.139469
- 34. Ani Deepthi, **Nivedita Acharjee**, Sruthi S. L., Meenakshy C. B (2022) An overview of nitrile imine based [3+2] cycloadditions over half a decade, *Tetrahedron* (*Elsevier*), 116, 132812 (**IF** = **2.388**) <a href="https://doi.org/10.1016/j.tet.2022.132812">https://doi.org/10.1016/j.tet.2022.132812</a>
- 35. Jamelah S. Al-Otaibi\*, Y. Sheena Mary, Y. Shyma Mary, **Nivedita Acharjee**, David G.Churchill (2022), Spectroscopic studies of 5-fluoro-1H-pyrimidine-2,4-dione adsorption on nanorings, solvent effects and SERS analysis, *Computational and Theoretical Chemistry(Elsevier)*, (**IF** = **2.292**) <a href="https://doi.org/10.1016/j.comptc.2022.113873">https://doi.org/10.1016/j.comptc.2022.113873</a>
- 36. Jamelah S. Al-Otaibi\*, Y. Sheena Mary, Y. Shyma Mary, Asmita Mondal, **Nivedita Acharjee**, David G.Churchill (2022), Sumanene as a delivery carrier for methimazole drug: DFT, AIM, SERS and solvent effects, *Computational and Theoretical Chemistry*(*Elsevier*),1215,113811 (**IF** = **2.292**) <a href="https://doi.org/10.1016/j.comptc.2022.113811">https://doi.org/10.1016/j.comptc.2022.113811</a>
- 37. Jamelah S. Al-Otaibi, Y. Sheena Mary\*, Y. Shyma Mary, **Nivedita Acharjee** (2022) Quantum mechanical investigation into the adsorption pattern of clomipramine and methotrimeprazine HCl with graphene and fullerene, *Polycyclic Aromatic Hydrocarbons (Taylor and Francis)* (**IF** = **2.195**) <a href="https://doi.org/10.1080/10406638.2022.2041678">https://doi.org/10.1080/10406638.2022.2041678</a>
- 38. Said Kerraj, Mohammed Salah\*, Samir Chtita, Mohammed El Idrissi, Said Belaaouad, Moutaabid Mohammed, **Nivedita Acharjee**, Najia Komiha (2022), Theoretical study of photovoltaic performances of Ru, Rh and Ir half sandwich complexes containing N,N chelating ligands in Dye- Sensitized Solar Cells (DSSCs). DFT and TD-DFT investigation, *Computational and Theoretical Chemistry* (*Elsevier*), 1209,

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- 39. **Nivedita Acharjee\***, H. A. M. Salim, Mrinmoy Chakraborty (2022)" Unveiling the synthesis of spirocyclic, tricyclic and bicyclic triazolooxazines from intramolecular [3+2] azide-alkyne cycloadditions with a molecular electron density theory perspective", *Structural Chemistry* (*Springer*), 33, 555-570 (**IF =1.795**) https://doi.org/10.1007/s11224-021-01870-3
- 40. **Nivedita Acharjee\***, H. A. M. Salim, Mrinmoy Chakraborty (2022) "Unveiling the regioselective synthesis of antiviral 5-isoxazol-5-yl-2' deoxyuridines with a molecular electron density theory perspective", *Journal of the Serbian Chemical Society*, 87, 707-721 (**IF** = **1.240**) https://doi.org/10.2298/JSC211014106A
- 41. Ali Barhoumi, Mohammed Salah, Abdellah Zeroual, Mrinmoy Chakraborty, **Nivedita Acharjee\*** (2022), Unveiling the reactions of triethylphosphite and its diethylamino substituted derivatives tocarbon tetrachloride with a molecular electron density theory perspective, *Indian Journal of Chemistry*, 61, 127-135. <a href="http://op.niscpr.res.in/index.php/IJC/article/view/60707">http://op.niscpr.res.in/index.php/IJC/article/view/60707</a>
- 42. Jamelah S. Al-Otaibi\*, Y.Sheena Mary, Y.Shyma Mary, **Nivedita Acharjee**, D. G. Churchill (2023) Theoretical study of Glycoluril by highly symmetrical magnesium oxide Mg 12 O 12 nanostructure: Adsorption, Detection, SERS enhancement and electrical conductivity study, *Journal of Molecular Modeling (Springer)* 28, 332 (**IF** = **2.172**)
- 43. Jamelah S. Al-Otaibi\*, Y.Sheena Mary, Y.Shyma Mary, **Nivedita Acharjee**, Zakir Ullah (2022) DFT, solvation effects, reactivity and SERS analysis on structural, optical, and vibrational properties of a biomolecule of pyrimidine derivative adsorbed on metal clusters of Ag/Au/Cu, *Journal of the Indian Chemical Society* (*Elsevier*), 99, 100753, <a href="https://doi.org/10.1016/j.jics.2022.100753">https://doi.org/10.1016/j.jics.2022.100753</a>

- 44. Luis R Domingo\*, **Nivedita Acharjee** (2021), Unveiling the Substituent Effects in the Stereochemistry of [3+2] Cycloaddition Reactions of Aryl- and Alkyldiazomethylphosphonates with Norbornadiene within a MEDT Perspective, *ChemistrySelect* (*Wiley*), 6, 10722-10733 (**IF** = 2.109) <a href="http://dx.doi.org/10.1002/slct.202102942">http://dx.doi.org/10.1002/slct.202102942</a>
- 45. **Nivedita Acharjee\***, H. A. M. Salim, Mrinmoy Chakraborty (2021) Unveiling [3+2] cycloaddition reactions of benzonitrile oxide and diphenyl diazomethane to cyclopentene and norbornene: A molecular electron density theory perspective, *Theoretical Chemistry Accounts (Springer)*, 140, 113 (**IF** = **1.702**). https://doi.org/10.1007/s00214-021-02811-3
- 46. Luis R. Domingo\*, **Nivedita Acharjee**\* (2021) Unveiling the Chemo- and Regioselectivity of the [3+2] Cycloaddition Reaction between 4-Chlorobenzonitrile Oxide and β-Aminocinnamonitrile with a MEDT Perspective, *ChemistrySelect (Wiley)*, 6, 4521-4532 (**IF = 2.109**) (**Top downloaded paper of 2021-2022 recognized by Wiley**) <a href="https://doi.org/10.1002/slct.202100978">https://doi.org/10.1002/slct.202100978</a>
- 47. **Nivedita Acharjee\***, H. A. M. Salim, Mrinmoy Chakraborty, Madhuri P Rao, Madhu Ganesh (2021) Unveiling the high regioselectivity and stereoselectivity within the synthesis of spirooxindolenitropyrrolidine: A molecular electron density theory perspective, *Journal of PhysicalOrganic Chemistry (Wiley)*, e4189 (**IF = 2.391**) (**Top cited paper of 2021-2022 recognized by Wiley**) https://doi.org/10.1002/poc.4189
- 48. H. A. M. Salim, **Nivedita Acharjee\***, Luis R. Domingo, H. H. Abdallah (2021) A Molecular Electron Density Theory Study for [3+2] Cycloaddition Reactions of 1-Pyrroline-1-oxide with Disubstituted Acetylenes Leading to Bicyclic 4-isoxazolines, *International Journal of Quantum Chemistry (Wiley)*, 121, e2653 (**IF** = **2.444**) https://doi.org/10.1002/qua.26503
- 49. A. Bazzi, K. Abbiche\*, S. Izzaouihda, **Nivedita Acharjee\***, H. Zejli, K. Marakchi, N. Komiha, S. El Issami, L. Bazzi, M. Hilali (2021) Inhibition efficiency and adsorption mechanism of 4-aminobenzoic acid for copper corrosion in nitric acid medium: a combined experimental andtheoretical investigation, *Structural Chemistry* (Springer), (**IF** = 1.887) https://doi.org/10.1007/s11224-021-01784-0
- 50. H. A. M. Salim\*, **Nivedita Acharjee**, H. H. Abdallah (**2021**) Insights into the mechanism andregioselectivity of the [3 + 2] cycloaddition reactions of cyclic nitrone to nitrile functions with a molecular electron density theory perspective, *Theoretical Chemistry Accounts*, 140, 1 (**IF** = **1.702**) https://doi.org/10.1007/s00214-020-02703-y
- 51. Luis R. Domingo \*, Mar Ríos-Gutiérrez, **Nivedita Acharjee** (**2021**) Unveiling the UnexpectedReactivity of Electrophilic Diazoalkanes in [3+2] Cycloaddition Reactions within MolecularElectron Density Theory, *Chemistry* (*MDPI*), 3(1), 74-93. https://doi.org/10.3390/chemistry3010006
- 52. **Nivedita Acharjee\***, Sourav Mondal (**2021**), A molecular electron density theory study tounderstand the strain promoted [3+2] cycloaddition (SPAAC) reaction of benzyl azide and cyclooctyne, *Indian Journal of*

- 53. Luis R Domingo\*, **Nivedita Acharjee (2020)** "Unravelling the Strain-Promoted [3+2] Cycloaddition Reactions of Phenyl Azide with Cycloalkynes from the Molecular Electron Density Theory Perspective", *New Journal of Chemistry (Royal Society of Chemistry)*, 44, 13633-13643(**IF** = **3.591**) https://doi.org/10.1039/D0NJ02711A
- 54. Luis R Domingo\*, **Nivedita Acharjee**\*, H. A. M. Salim (**2020**) Understanding the Reactivity of Trimethylsilyldiazoalkanes participating in [3+2] Cycloaddition Reactions towards Diethylfumarate with a Molecular Electron Density Theory Perspective, *Organics (MDPI)*, "*SpecialIssue on Cycloaddition Reactions in Organic Synthesis*", 1(1), 3-18. <a href="https://doi.org/10.3390/org1010002">https://doi.org/10.3390/org1010002</a>
- 55. **Nivedita Acharjee\* (2020)** Unravelling the regio- and stereoselective synthesis of bicyclic N,O-nucleoside analogues within the molecular electron density theory perspective, *Structural Chemistry (Springer)*, 31, 2147 (**IF = 1.887**) <a href="https://doi.org/10.1007/s11224-020-01569-x">https://doi.org/10.1007/s11224-020-01569-x</a>
- 56. Luis R Domingo\*, **Nivedita Acharjee** (**2020**) Unveiling the High Reactivity of Strained Dibenzocyclooctyne in [3+2] Cycloaddition Reactions with Diazoalkanes through the Molecular Electron Density Theory, *Journal of Physical Organic Chemistry* (*Wiley*), 33, e4100. (**IF** = **2.391**) https://doi.org/10.1002/poc.4100
- 57. Luis R Domingo, **Nivedita Acharjee\*** (2020), A Molecular Electron Density Theory Study of the Grignard Reagent Mediated Regioselective Direct Synthesis of 1,5-disubstituted-1,2,3-Triazoles, *Journal of Physical Organic Chemistry* (*Wiley*) 33, e4062, (**IF** = 2.391) <a href="https://doi.org/10.1002/poc.4062">https://doi.org/10.1002/poc.4062</a>
- 58. **Nivedita Acharjee\***, Avijit Banerji (**2020**), A Molecular electron density theory study to understand the interplay of theory and experiment in nitrone-enone cycloaddition, *Journal of Chemical Sciences (Springer)*, 2020, 132, 65. (**IF** = **1.573**) <a href="https://doi.org/10.1007/s12039-020-01766-5">https://doi.org/10.1007/s12039-020-01766-5</a>
- 59. **Nivedita Acharjee\*** (2020) Theoretical analysis of the regio- and stereoselective synthesis of spiroisoxazolines, *Journal of Molecular Modeling (Springer)*, 26, 117. (**IF** = 1.810) <a href="https://doi.org/10.1007/s00894-020-04372-x">https://doi.org/10.1007/s00894-020-04372-x</a>
- 60. K.Abbiche, **Nivedita Acharjee\***, M. Salah, M. Hilali, A.Laknifli, N. Komiha, and K. Marakchi (2020) Unveiling the mechanism and selectivity of [3+2] cycloaddition reactions of benzonitrile oxide to ethyl transcinnamate, ethyl crotonate and trans-2-penten-1-ol through DFT analysis, *Journal of Molecular Modeling* (*Springer*), 26, 279 (**IF = 1.810**) https://doi.org/10.1007/s00894-020-04547-6
- 61. **Nivedita Acharjee\*** (2020), Understanding the regio- and diastereoselective synthesis of a potent antinociceptive isoxazolidine from C-(pyridin-3-yl)-N-phenyl nitrone in the light of molecular electron density theory, *Journal of Serbian Chemical Society*, 85(6), 765-779. (**IF** = 1.240) https://doi.org/10.2298/JSC190914136A
- 62. **Nivedita Acharjee\***, Avijit Banerji, Mira Jhawar, Ambar Barman (**2020**), Comparative DFT analysis of ELF topology and global properties of allyl and allenyl type three atom components (TACs), *Journal of The Indian Chemical Society*, 97, 753-759. [Selected Cover page of May 2020 issue]
- 63. **Nivedita Acharjee,** Tuhin Ghosh\* (**2020**), Structure analysis of sulfated polysaccharides extracted from Scinaia interrupta: Experimental and density functional theory studies, *Asian Journal of Chemistry*, Volume 32(7), 1589-1596. <a href="https://doi.org/10.14233/ajchem.2020.22609">https://doi.org/10.14233/ajchem.2020.22609</a>
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- 65. Luis R Domingo\*, Mar-RiozGutiérrez, **Nivedita Acharjee\*** (**2019**), A Molecular Electron Density Theory Study of the Chemoselectivity, Regioselectivity, and Diastereofacial Selectivity in the Synthesis of an Anticancer Spiroisoxazoline derived from α-Santonin, *Molecules*, 24, 832. (**IF** = **3.267**) https://doi.org/10.3390/molecules24050832
- 66. **Nivedita Acharjee\* (2019)**, Understanding the geometry and [3+2] cycloadditions of nitrile imine in terms of molecular electron density theory, *Indian Journal of Chemistry*, 58A, 645-652. (**IF=0.489**) <a href="https://op.niscpr.res.in/index.php/IJCA/article/download/23411/465478047">https://op.niscpr.res.in/index.php/IJCA/article/download/23411/465478047</a>
- 67. Avijit Banerji, Rina Saha, **Nivedita Acharjee\*** (2019), Reaction intermediates for microwave irradiation assisted synthesis of N-Aryl Enaminoketones and their subsequent conversion to quinolines: An experimental and DFT study, *Asian Journal of Research in Chemistry*; 12(2):103-111. https://doi.org/10.5958/0974-4150.2019.00023.3

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68. Luis R Domingo, Nivedita Acharjee\* (2018), [3+2] Cycloaddition Reaction of C-Phenyl-N-methyl

- Nitrone to Acyclic-Olefin-Bearing-Electron-Donating Substituent: A Molecular Electron Density Theory Study, *Chemistry Select*, 3, 8373. (**IF** = **1.811**) <a href="https://doi.org/10.1002/slct.201801528">https://doi.org/10.1002/slct.201801528</a>
- 69. **Nivedita Acharjee\*** (2018), Solvent effects on cycloaddition reactions of potent spin trappingprobe Ntert-butylmethanimine-N-oxide: A DFTstudy, *Journal of Theoretical andComputational Chemistry*, Vol. 17. No. 4, 1850027 (23 pages) (**IF** = **0.848**) <a href="https://doi.org/10.1142/S021963361850027X">https://doi.org/10.1142/S021963361850027X</a>
- 70. Sutapa Mandal, Kaustav Kumar Maiti, Avijit Banerji, Thierry Prange, Alain Neuman, **Nivedita Acharjee\*** (2018) Experimental and DFT studies for substituent effects oncycloadditions of C,N-disubstituted nitrones to cinnamoyl piperidine, *Indian Journal of Chemistry*, *Section* B, 57B, 108. (**IF** = 0.388) http://nopr.niscair.res.in/bitstream/123456789/43414/1/IJCB%2057B%281%29%20108-119.pdf

- 71. Sutapa Mandal, **Nivedita Acharjee\*** (**2017**), DFT study for radical capture by mitochondria oxidotoxin protective ionic and non-ionic amphiphilic α-phenyl-N-t-butyl nitrone derivatives; *Indian Journal of Chemistry Section A*, 56A, 9. http://nopr.niscair.res.in/bitstream/123456789/39745/1/IJCA%2056A%281%29%209- 20.pdf
- 72. Sutapa Mandal, **Nivedita Acharjee\***, Bhaskar Chakraborty, Avijit Banerji (**2017**) DFT analysis for the regio- and stereospecific synthesis of antibacterial isoxazolidine, *Journal of the Indian Chemical Society*, *94*, pp. 535-542.

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- 73. **Nivedita Acharjee\*** (2014), Acyclic and cyclic nitrone cycloadditions to 1-cinnamoyl-1- piperidine: A DFT study. *Journal of Theoretical and Computational Chemistry*, 13, 1450071. (**IF** = **0.848**) https://doi.org/10.1142/S0219633614500710
- 74. **Nivedita Acharjee\*** (**2014**), DFT interpretations for cycloadditions of an electron deficient C- aryl-N-phenyl nitrone. *Journal of Theoretical and Computational Chemistry*, 13, 1450007. (**IF** = **0.848**) <a href="https://doi.org/10.1142/S0219633614500072">https://doi.org/10.1142/S0219633614500072</a>

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- 75. Nivedita Acharjee, Avijit Banerji\*, Thierry Prangé (2012), Remarkable influence of mildLewis acid catalysts on cycloadditions leading to tetrasubstituted isoxazolidines: DFT analysis augmented by X-Ray and NMR studies. *Monatshefte für Chemie, chemical monthly*, 143, 1687. https://doi.org/10.1007/s00706-012-0828-z
- 76. **Nivedita Acharjee**, Avijit Banerji\*, (**2011**), DFT interpretation of 1,3-dipolar cycloaddition reaction of *C*,*N*-diphenyl nitrone to methyl crotonate in terms of reactivity indices, interaction energy and activation parameters. *Computational and Theoretical Chemistry*, 967, 50. https://doi.org/10.1016/j.comptc.2011.03.040
- 77. **Nivedita Acharjee**, Avijit Banerji\*, Thierry Prangé (**2011**), Catalytic effects of Lewis acids on1,3-dipolar cycloaddition of *C*-(4-chlorophenyl)-*N*-phenyl nitrone to benzylidene acetophenone, *Indian Journal of Chemistry Section B*, 50B, 835.
- 78. **Nivedita Acharjee**, Avijit Banerji\*, Biswajit Gayen (**2011**), Cycloaddition approach to five-membered heterocycles Some recent experimental and theoretical aspects. Invited Reviewon 150<sup>th</sup> Birth anniversary Commemoration of Acharya Prafulla Chandra Ray, *Journal of the Indian Chemical Society*, 88, 1857.
- 79. **Nivedita Acharjee**, Avijit Banerji\*, Thierry Prangé (**2010**), DFT study of 1,3-dipolarcycloadditions of *C,N* disubstituted aldonitrones to chalcone evidenced by NMR and X-Rayanalysis. *Monatshefte für Chemie, chemical monthly*, 141, 1213. https://doi.org/10.1007/s00706-010-0393-2
- 80. **Nivedita Acharjee,** Tapas Kumar Das, Avijit Banerji, Manas Banerjee\*, Thierry Prangé (**2010**)(**2010**), A DFT based exploration augmented by X-ray and NMR of the stereoselectivity in the 1,3-dipolar cycloaddition of 1-Pyrroline-1-oxide to methyl cinnamate and benzylideneacetophenone. *Journal of Physical Organic Chemistry*, 83, 1187. (**IF =1.536**) <a href="https://doi.org/10.1002/poc.1690">https://doi.org/10.1002/poc.1690</a>
- 81. **Nivedita Acharjee,** Avijit Banerji\*, (**2010**), DFT study of the configuration and conformation of *C*-aryl-*N*-phenyl nitrones and their reactivities as 1,3-dipoles to methyl and ethylcrotonates, *Indian Journal of Chemistry Section A*, 49A, 1444.
- 82. **Nivedita Acharjee,** Avijit Banerji\*, Manas Banerjee\*, Tapas Kr. Das (**2009**), Experimental andtheoretical spectroscopic investigations of *C*-aryl-*N*-methyl nitrones., *Indian Journal of Chemistry Section A*, 48A, 1627.
- 83. Avijit Banerji\*, **Nivedita Acharjee**, (**2009**) Frontier Molecular orbitals of *C*-aryl- *N*-methylnitrones– A DFT study. *Journal of Indian Chemical Society*, 86, 1068.
- 84. Avijit Banerji\*, **Nivedita Acharjee**, Pizush Kanti Biswas (**2009**) 1,3-Dipolar cycloadditions. Part XV: Systematic spectroscopic investigations of *C* aryl- *N*-methyl nitrones, *Journal of the Indian Chemical Society*, 86, 63.

## **Book Chapter**

85. Luis R Domingo\*, **Nivedita Acharjee** (2020), Molecular Electron Density Theory-A New Theoretical Outlook on Organic Chemistry *Frontiers in Computational Chemistry-Volume 5*, Published by Bentham and Science Pvt. Ltd. https://doi.org/10.2174/9789811457791120050007

## **Research Articles in Edited Volumes**

- 86. **Mandal S, Acharjee N\***, Chakraborty B "DFT studies for generation of stereospecific antibacterial cycloadductfrom symmetrically substituted reactant.", "Chemistry on its Way-Impacts on the Environment", (Ed: Samiran Mandal), Saldiha College (ISBN: 978-93-84191-17-7); [Published in2017]
- 87. **Mandal S**, **Acharjee N\***, Banerji A, "Conceptual DFT interpretations for diastereoselective synthesis of tetrasubstituted isoxazoidines", (Ed: Kamala Mitra) Dept. of Chemistry, Prasanta Chandra MahalanobishMahavidyalaya (ISBN: 978-93-8310-31-8) [**Published in 2017**]

# **General Articles/Periodicals**

- 88. Asmita Mondal, Avijit Banerjee\*, **Nivedita Acharjee** (2022) Understanding [3+2] cycloaddition reactions from the molecular electron density perspective: A new theoretical outlook on organic reactions, Edu. Chem. Sci. Tech., Vol 10, August 2022 (Periodical published by the Indian Chemical Society)
- 89. Nivedita Acharjee (2019): Impact of solar radiation on human health, CURE (Congregation for Urban and Rural Environment), Environmental Bulletin 2019, Volume XXV, ISSN: 2319-1333
- 90. **Nivedita Acharjee (2019):** Beauty of "Golden Ratio" in paintings, architecture, music, nature and technology, *Communique*, Academic Journal of Durgapur Government College, *ISSN No. 0973-8584*.
- 91. Nivedita Acharjee\* (2018), Impact of smog on public health, CURE (Congregation for Urban and Rural Environment), Environmental Bulletin 2018, Volume XXIV, ISSN: 2319-1333
- 92. **Nivedita Acharjee\*** (2018), "Scientific Story of our laughter, crying and stress": *Everyman's Science*, Journal published by Indian Science Congress Association, *ISSN*: 0531-495X
- 93. **Nivedita Acharjee\*** (2017), The Scientific Face of our emotions, *Communique*, Academic Journal of Durgapur Government College, *ISSN No. 0973-8584*.
- 94. **Nivedita Acharjee\*** (2017), Effect of particulate matter on Human Health and Environment, CURE (Congregation for Urban and Rural Environment), Environmental Bulletin 2017, Volume XXIII, ISSN: 2319-1333
- 95. **Nivedita Acharjee\*** (2015), "Molecular Processes of Aging": Vol. XXI, Environmental Bulletin 2015, *CURE* (*Congregation for urban and rural environment*), *ISSN*: 2319-1333
- 96. **Nivedita Acharjee\*** (2014), "The Aging Process: Accelerators and Decelerators": Aug13- Sep13 issue, Everyman's Science, Journal published by Indian Science Congress association, *ISSN*: 0531-495X

## Sponsored Projects

- 1. **Dr. Nivedita Acharjee**, Coordinator, DBT STAR COLLEGE SCHEME for Durgapur Government College sponsored by Department of Biotechnology, Government of India, Project No. HRD-11011/24/2022-HRD-DBT dated 29-09-2022 (Period: 29-09-2022 to 28-09-2025)] (Rs. 120.38469 Lakhs)
- 2. **Dr. Nivedita Acharjee** (PI), UGC Minor research Project No. F. PSW-111/11-12 (ERO) "Experimental and theoretical DFT investigations of the effects of mild Lewis acid catalysts on the cycloaddition reactions of acyclic and cyclic aldonitrones to α,β-unsaturated ketones and esters" (Period: 8.11.2011 to 7.11.2013)] (Rs. 1.35 Lakhs)

## Guidance of Ph.D. Scholar and M.Sc. Project Dissertations

- 1. Asmita Mondal (2022, Durgapur Government College) [enrolled for Ph.D.]
- 2. Barsali Banerji (2023, Durgapur Government College) [enrolled for Ph.D.]
- 3. Bhaskar Mondal (2023, Durgapur Government College) [enrolled for Ph.D.]
- 4. Soubhik Chakraborty (2023, *Durgapur Government College*) N-t-Butyl- aryl Nitrones as Potent Spin Traps: DFT Analaysis of Electron Localization Function Topology, Local Selectivity, Reactivity and Solvent Effects [M.Sc.]
- 5. Sangram Singh (2023) DFT Calculations for the Analysis of Selectivity, Reactivity and Solvent Effects of 5,5-Dimethyl-1-Pyrroline-N-Oxide (DMPO) Nitrones Towards Radical Capture )As Spin Traps) [M.Sc]
- 6. Shreyasee Nath (2022, Durgapur Government College): DFT calculations for the analysis of frontier molecular orbitals and reactivity indices of tricyclic pyrrolidine-fused benzo[b] thiophene 1,1-dioxide(M.Sc.)

- 7. Subarna Khan (2021, *Durgapur Government College*): Energy profile study associated with the[3+2] cycloaddition reaction of 1-pyrroline-1-oxide with nvinyl nucleobases[M.Sc.]
- 8. Asmita Mondal (2020, Durgapur Government College): Chemical profiling and theoretical analysis of polysaccharides in rice bran [M.Sc.]
- 9. Sourav Mondal (2019, Durgapur Government College): Solvent effects on [3+2] cycloadditionreactions of N-tert-butylmethanimine-N-oxide [M.Sc.]
- 10. Monirul Islam (*2019, Durgapur Government College*); 1,3-dipolar cycloaddition reactions: classification, selectivity and catalysis [M.Sc.]
- 11. Mira Jhawar (2018, Durgapur Government College): Allyl type 1,3-dipoles: DFT study augmented with ELF topological analysis [M.Sc.]
- 12. Ambar Barman (*2018, Durgapur Government College*): Allenyl type 1,3-dipoles: DFT study augmented with ELF topological analysis [M.Sc.]
- 13. Suranjana Sen (2017, Durgapur Government College): Experimental and Theoretical Studies On Synthesis and Spectroscopy of C-(4-chlorophenyl)-N-phenyl nitrone [M.Sc.]
- 14. Jayanta Mondal (2017, Durgapur Government College): Synthesis, global properties and vibrational analysis of C-(4-bromophenyl)-N-phenyl nitrone [M.Sc.]
- 15. Nasrin Jahan (2017, Durgapur Government College): Synthesis, vibrational analysis and cycloaddition reactions of chalcone to C-aryl-N-phenyl nitrones :Experimental andtheoretical studies [M.Sc.]
- 16. Akash Mitra (2016, Durgapur Government College): Effect of substituents on the frontier molecular orbital energies and global properties of Isoxazolidines: AM1 study [M.Sc.]

# **Awards & Recognitions**

- 1. **Prof. Sadhan Basu Memorial award (by University of Calcutta)** for excellence in the performance in post graduate examination, 2005.
- **2. University Gold Medal** from University of Calcutta for 1<sup>st</sup> class 1<sup>st</sup> position in M.Sc. examination in the year 2005.
- 3. Umesh Chandra Memorial Cash Prize, Manoranjan Banerjee Memorial award, Pradip Kr. Ghosh Memorial Prize, Kananbala Sriti Purashkar for first class first position in Chemistry Hons from Lady Brabourne College, Kolkata.
- **4. Best poster presentation** in UGC sponsored National Level Seminar on "*Modern Trends in Chemical Sciences*" during November 25-26, 2016 at Department of Chemistry, Prasanta Chandra Mahalanobish Mahavidyalaya.
- **5. Best oral presentation** in organic Chemistry in 2<sup>nd</sup> UGC sponsored National seminar on advanced spectroscopy, theoretical chemistry, synthesis, reactivity and structure evaluation, organized by University of Burdwan, Burdwan, 20-22<sup>nd</sup> February 2009.
- **6.** Dr. Nivedita Acharjee has contributed the first research paper of the newly launched journal "*Organics*" by MDPI in 2020 in "*Special Issue on Cycloaddition Reactions in Organic Synthesis*",1(1), 3-18. <a href="https://doi.org/10.3390/org1010002">https://doi.org/10.3390/org1010002</a>.
- 7. Dr. Nivedita Acharjee has contributed an MEDT research paper to analyse steric effects in 32CAreactions to the Special Issue "2020 Profile Papers by Chemistry' Editorial Board Members" of theWeb of Science indexed journal "Chemistry-Switzerland" by MDPI in 2021 (https://doi.org/10.3390/chemistry3010006)
- **8. National Scholarship** under National Scholarship Scheme for B.Sc. (III year Hons.) Examination 2003 in Chemistry
- **9. Certificate of Merit and National scholarship** from CBSE for being among top 0.1% studentsin All India Secondary School examination, 1998.

### **Invited Lectures**

- **1.** "How to develop and upload and upload E-content in 4 quadrants?"; Workshop organized by Internal Quality Assurance Cell, Durgapur Government College on March 14, **2019**
- 2. "Career Counselling and Placement after Secondary and Senior Secondary level"; Special Programme organized by Assembly of God Church School, Okhra, West Bengal on 1st November 2017.
- 3. "Gender Equality for Sustainable Development"; NSS Special Camp, Organized by NSS Unit, Durgapur Government College in 2016
- **4.** "Gender Equality: Indispensable yet overlooked"; Awareness Programme on International Women's Day; organized by Cell for Gender Sensitization and Protection of Sexual Harassment against Women, Durgapur

## Abstracts in seminar/conference

- 1. Equinox 2022, The Annual International Conference on the occasion of celebrating Century Year of University of Delhi, Paper entitled "Unveiling [3+2] Cycloaddition Reactions from the Molecular Electron Density Theory Perspective", August 11-13, 2022, Presented by PhD student Asmita Mondal.
- 2. International Seminar on Recent Advances in Chemistry and Material Science (RACMS-2022) on the occasion of 161<sup>st</sup> Birth Anniversary Celebration of Acharya P. C. Ray, Paper entitled "Unveiling the mechanism and substituent effects in the synthesis of fused tricyclic 1,2,3-triazolines and 1,2,3-triazoles from intramolecular [3+2] cycloadditions within the molecular electron density theory framework", July 30-31, 2022, Presented by PhD student Asmita Mondal.
- 3. International Research Scholars' meet on the occasion of 99th Foundation Day celebration of Indian Chemical Society, Paper entitled "Unveiling intramolecular [3+2] cycloaddition reactions of nitrones and azides from the molecular electron density theory perspective", May 9, 2022, Presented by PhD student Asmita Mondal and received award of excellence.
- 4. International Seminar on "Advancements in Chemical Sciences: Issues and Challenges", organized by Asansol B.B. College, Asansol, February 4-5, 2020. Paper entitled: Molecular electron density theory study to analyse [3+2] cycloaddition reactions leading to biologically active five-membered heterocycles", Presented by Dr. Nivedita Acharjee
- 5. International Seminar on "Innovation, Expansion, Impacts and Challenges in Chemical and Biological Sciences" organized by Surendranath College, Kolkata, January 8-9, 2020, Paper entitled: N-t-butyl-α-aryl nitrones as potent spin traps of biologically relevant free radicals: DFT analysis of electron localization function topology, local selectivity, reactivity, and solvent effects", Presentedby Sourav Mondal (M.Sc. Student)
- 6. International Seminar on "Innovation, Expansion, Impacts and Challenges in Chemical and Biological Sciences" organized by Surendranath College, Kolkata, **January 8-9, 2020**, Paper entitled: Understanding the chemo-, regio and steroselective synthesis of an anticancer spiroisoxazoline derived from α-santonin in the light of molecular electron density theory", **Presented by Nivedita Acharjee**
- 7. National Seminar on "Modern Era of Biosciences: Lab to land" organized by Departments of Conservation Biology and Zoology, Durgapur Government College, **December 05, 2019**, Paper entitled: DFT study of the Electronic Structure and Reactivity of the Nitrone moiety as potent radical spin trap for hydroxyl radical to fight the aging process", Presented by Nivedita Acharjee
- 8. National Seminar on Chemistry for Sustainable Development organized by Sidho-Kanho-Birsha University in collaboration with Indian Chemical Society; during November 26-27, 2019, Paper entitled: "DFT study of the selectivity and reactivity of potent radical spin trap N-t-butyl-α- phenyl nitrone (PBN) in terms of global & local indices and electron localization function topological analysis", Presented by Sourav Mondal (M.Sc. Student).
- 9. DAE and CSIR sponsored International Level Conference on "Non-linear Dynamics and its applications" during March 12-14, 2019 at Dept. of Physics, Durgapur Government College; Paper entitled" Understanding the pathway of [3 + 2] cycloaddition reactions of C-phenyl-N- methyl nitrone in terms of BET (Bonding Evolution Theory) and QTAIM (Quantum Theory of atoms in Molecules) analyses", Presented by Nivedita Acharjee
- 10. DAE and CSIR sponsored International Level Conference on "Non-linear Dynamics and its applications" during March 12-14, 2019 at Dept. of Physics, Durgapur Government College; Paper entitled" Molecular Electron Density Theory: A computational chemistry approach tounderstand the reactivity of organic reactions", Presented by Nivedita Acharjee
- 11. DST, DRDO and DAE sponsored National Level Seminar on "Non-linear Dynamics and its applications" during February 7-9, 2017 at Dept. of Physics, Durgapur Government College; Paper entitled "Density Functional Theory applications for interpretations of chemical reactions", Presented by Nivedita Acharjee
- 12. UGC sponsored National Level Seminar on "Modern Trends in Chemical Sciences" during November 25-26, 2016 at Dept. of Chemistry, Prasanta Chandra Mahalanobish Mahavidyalaya; Paper entitled "DFT Study for nitrone cycloadditions to benzylidene acetone in terms of reactivity indices, interaction energies and activation parameters", Presented by Nivedita Acharjee (Awardedbest poster presentation)
- 13. UGC Sponsored National Seminar on "A Journey Through Recent Developments In Chemistry" organized jointly by Department of Chemistry, Acharya B. N. Seal College Cooch Behar, West Bengal and Department of Chemistry, University of North Bengal, Darjeeling, West Bengal, March 1st and 2nd 2012; Paper entitled "Accelerative and stereoselective influence of mild Lewis acid coordination in dipolar cycloadditions: Experimental and DFT study", Presented by Nivedita Acharjee
- 14. Chemical Research Society of India Eastern Zonal Meeting 2011 and Celebration of the International Year of Chemistry 2011, July 22-24, 2011; Poster no. 4, Page: 22, organized by Department of Chemistry, University of North Bengal, Darjeeling, West Bengal; Paper Entitled "DFT study evidenced by X-Ray and NMR analysis for the cycloaddition reactions of C,N disubstituted aldonitrones in terms of reactivity indices, interaction energies and activationparameters", Presented by Nivedita Acharjee
- 15. 2<sup>nd</sup> UGC sponsored National seminar on advanced spectroscopy, theoretical chemistry, synthesis, reactivity and

structure evaluation, organized by University of Burdwan, Burdwan, 20-22<sup>nd</sup> February **2009**; Paper entitled: "Studies on Nitrone Cycloadditions to  $\alpha$ ,  $\beta$ - unsaturated ketones" (Awarded best oral presentation in organic chemistry); Presented by Nivedita Acharjee

16. 45th Annual convention of chemists 2008; Paper entitled "Systematic spectroscopic investigations of C-aryl- N-methyl nitrones" Presented by Nivedita Acharjee
Symposium on "Acharya Prafulla Chandra Ray and Chemistry Today" organized jointly by Indian Chemical Society and Department of Chemistry, University of Calcutta, 2007; Paper entitled "A cycloaddition approach to masked triketomethanes" Presented by Nivedita Acharjee