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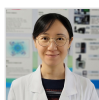
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Roberto Gobetto, Alceo Macchioni, Maurizio Peruzzini, Riccardo Pettinari, Valerio Zanotti

Article 100495

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Organic Chemistry

Research article *Open access*

Optical property investigations of coumarin and indene diketone structure dyes: Experiment and calculation

Chao-jun Hua, Wen-jie Niu, Yu-jin Li

Article 100257

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optical properties of these coumarin derivatives (**4a-4d**), such as the influence of different solvent to the fluorescence intensity and absorption intensity, and the influence of different substituent groups and the concentration of the solution are investigated. The results show that the fluorescence intensity of the compounds are the largest in the DMF (N,N-dimethylformamide). Specifically speaking, the absorption spectrum is at approximately 308 nm, the fluorescence spectrum is at nearly 435 nm in the DMF, the Stokes shift approaches 127 nm. In our case the fluorescence intensity decreases with the increase of the concentration of the solution, which is called concentration quenching. Additionally, the affection of solvent was also confirmed by TD-DFT (Time-dependent density functional theory) calculations.

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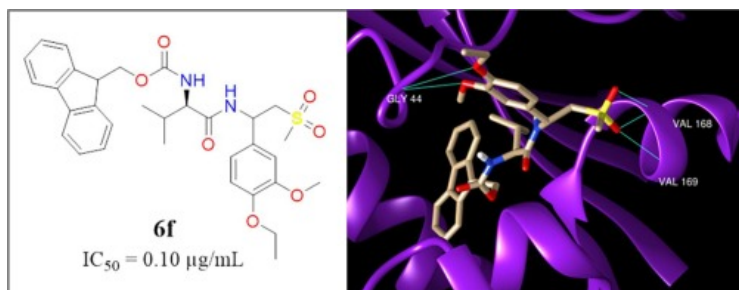
Vanillin containing 9H-fluoren sulfone scaffolds: Synthesis, biological evaluation and molecular docking study

Hanuman Narode, Manoj Gayke, Rajesh S. Bhosale, Gyanchander Eppa, ... Jhillu Singh Yadav

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Base and metal free true recyclable medium for Knoevenagel condensation reaction in SDS-ionic liquid-aqueous micellar composite system

Sourav Chakraborty, Abhijit Rudra Paul, Swapan Majumdar

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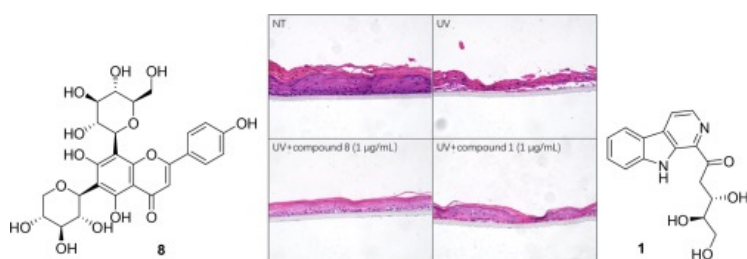
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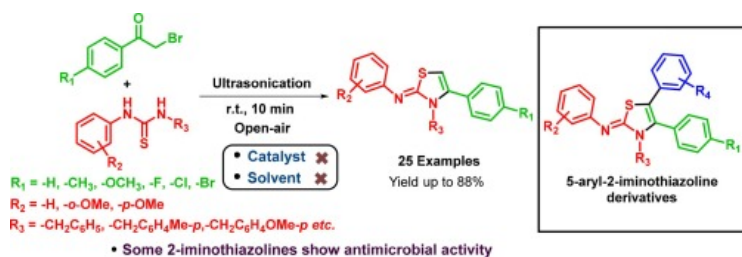
Yi-Fang Chen, Dong-Dong Zhang, Dong-Bao Hu, Xiao-Nian Li, ... Yue-Hu Wang

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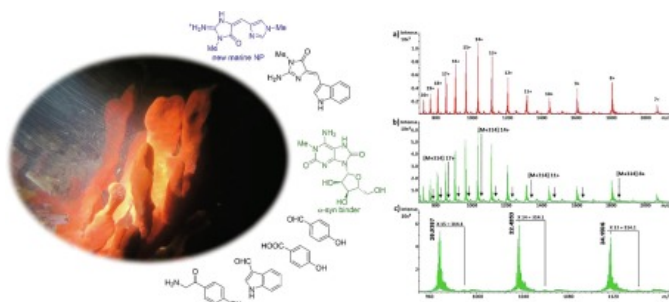
Chandan Bodhak, Subhro Mandal, Pia Dey, Samir Kumar Mukherjee, Animesh Pramanik

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PEG-cored phosphorus dendrimers: Synthesis and functionalization

Hanna Dib, Jérémy Rebière, Cyrille Rebout, Omar Alami, ... Anne-Marie Caminade

Article 100304

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Abstract

Abstract

Dendrimers are macromolecular hyperbranched nano-structures, which properties are mainly driven by the type of their terminal functions. This paper reports the synthesis of dendritic structures built from difunctional polyethylene glycols (PEGs) used as core. Three different PEG lengths have been used: PEG₃₀₀ (ca 6 CH₂CH₂-O units), PEG₁₀₀₀ (ca 22 CH₂CH₂-O units), and PEG₂₀₀₀ (ca 45 CH₂CH₂-O units). The first level of branching units is based on a dialdehyde of a dipiperazine triazine derivative. The second and third level of branching units are constituted of phosphorhydrazone linkages. The growing of the dendritic branches was found extremely slow with both PEG₁₀₀₀ and PEG₂₀₀₀ used as core, and was stopped early, whereas this problem was not observed with PEG₃₀₀ used as core. These PEG-cored dendritic structures were finally functionalized on the surface with two types of nitrogen ligands.

Short communication [Open access](#)Thermal decomposition behavior of MnO₂/Mg-Al layered double hydroxide after removal and recovery of acid gas

Tomohito Kameda, Tanya Kurutach, Yuriko Takahashi, Shogo Kumagai, ... Toshiaki Yoshioka

Article 100310

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Abstract

Abstract

Acidic gases such as HCl, SO₂, and NO_x are compounds that are generated through the incineration of waste products. Hence, an efficient method of removing such gases is highly desired. In this study, we present a novel method for treatment of acidic

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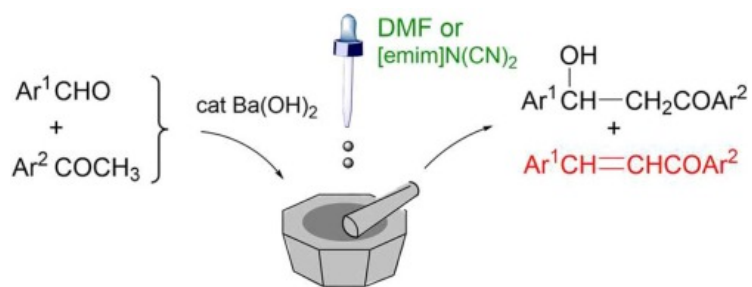
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Medicinal significance of novel coumarin analogs: Recent studies

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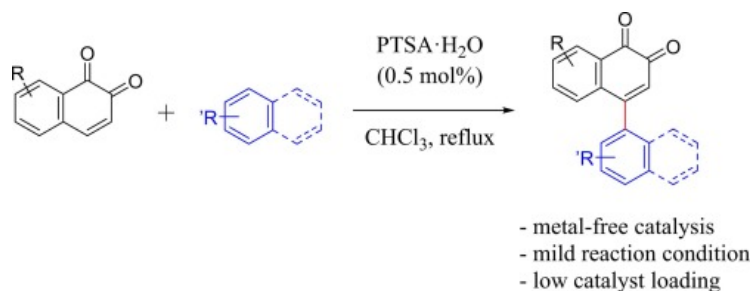
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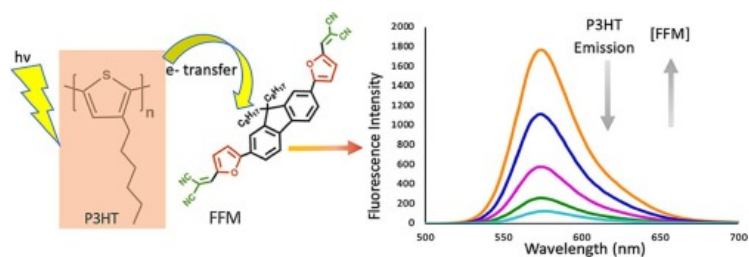
Synthesis and properties of fluorene based small molecule acceptors containing aromatic malononitrile functionalities

Sarah E. Brock, Dariia Yehorova, Brycelyn M. Boardman

Article 100320

[View PDF](#) [Article preview](#) Abstract **Graphical abstract**

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Research article *Open access*Synthesis of benzoyl esters of β -amyrin and lupeol and evaluation of their antibiofilm and antidiabetic activities

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antimicrobial activities. Biofilm inhibitory potential of the compounds on *S. aureus*, *E. coli* and *C. albicans* were performed at

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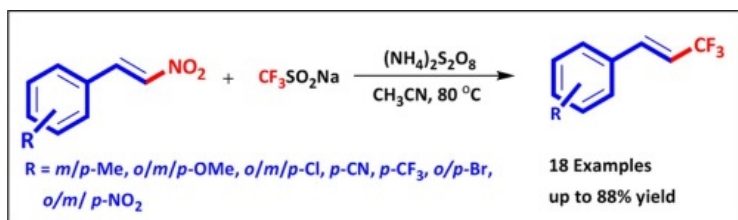
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Vitamin C-Catalyzed Hantzsch reaction under microwave condition: A greener access to 1,4-Dihydropyridines

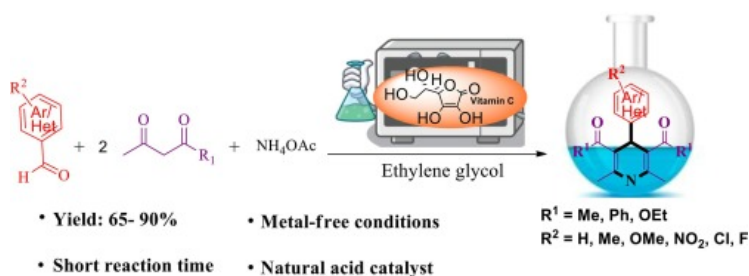
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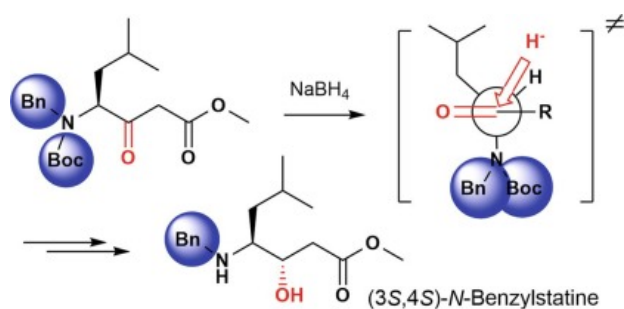
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This work presented the microwave assisted synthesis of six new 2'-hydroxychalcones and their characterization based on FTIR, UV-Vis, ¹H NMR, and mass spectral analysis. Quantum chemical studies confirmed the structures of prepared chalcones. Antioxidant, *in vitro* antimicrobial and *in silico* antiviral studies have been performed to evaluate their biological performance. Results of molecular docking of prepared 2'-hydroxychalcones against SARS-CoV-2 (7BQY) main protease disclosed their inhibition which is comparable to standard, remdesivir and better than hydroxychloroquine (HCQ). ADMET prediction revealed them to be non-carcinogenic and relatively safe.

Research article [Open access](#)**Highly stereoselective multigram scale synthesis of (3*S*,4*S*)-Statine and (3*S*,4*S*)-*N*-Benzylstatine**

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Chiranjeevi Donthulachitti, Yelukala Ramakrishna, Ravikumar Shekunti, Chandra Kiran Neella

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$R_1=OH$ & $R_2=H$ ($U=lyxo, y$)
(P = Protecting Group)

Research article [Open access](#)An isatin aldol adduct as a precursor to α, α' -difunctionalized methyl vinyl ketones

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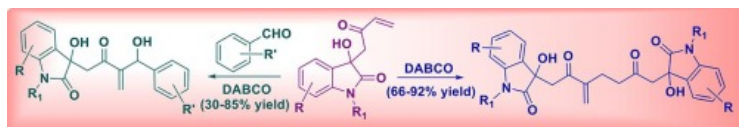
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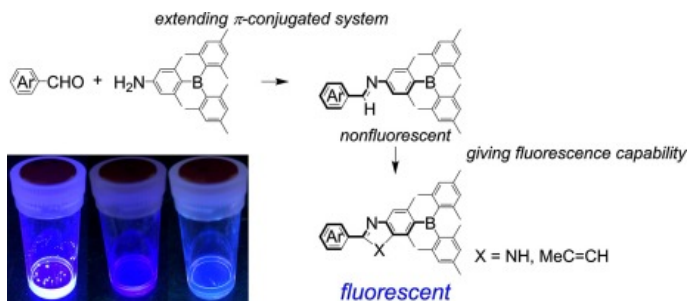
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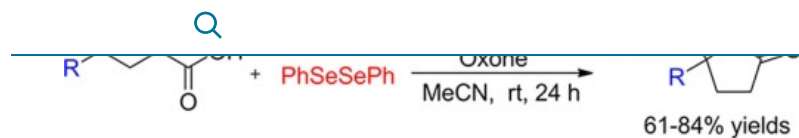
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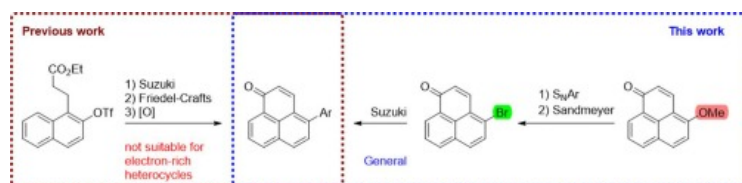
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Citric acid mediated simple and stereoselective synthesis of *o*-linked glycosides by Ferrier rearrangement

P.I. Ramesh, Mohamad Saif Ali, Subhash Ghosh, Madhu Babu Tatina

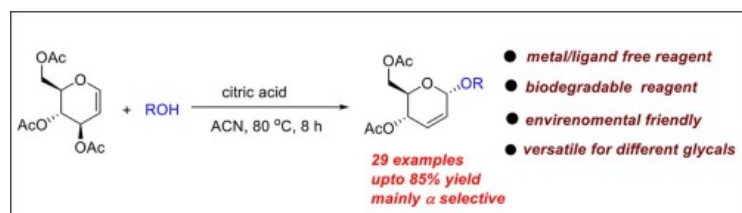
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Citric acid is versatile reagent for many glycols with good yield and selectivity well suited for the synthesis of pseudodisaccharides.



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Coagulation factor Xa (FXa), a serine endopeptidase is a common coagulation factor activated as a result of the initiation of both intrinsic and extrinsic blood coagulation pathways. Hence, FXa has been regarded as an important pharmaceutical target for the treatment of thrombotic disorders. In this study, we reported the design and synthesis of pyrazolyl piperidine analogs **4(a–h)** as a new class of anticoagulant drug candidates. Among the synthesized analogs **4(a–h)**, compound **4a** consisting of the 4-chlorophenyl substitution displayed the highest *in vitro* FXa inhibition activity with an IC₅₀ value of 13.4 nM. The PT and aPTT assay indicated that compound **4a** showed good anticoagulation activity compared to Heparin. Furthermore, docking studies suggested that the synthesized analogs displayed binding modes similar to the cocrystallized Rivaroxaban ligand. In addition, *in-silico* ADMET and DFT studies were carried out for all the designed compounds. Together, our study suggests that the compound (**4a**) displayed anti-coagulant activity through the inhibition of FXa

[Review article](#) [Open access](#)**Preparation of glycerol derivatives by entered of glycerol in different chemical organic reactions: A review**

Fatima Alashek, Mohammad Keshe, Ghader Alhassan

Article 100359

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Glycerol is now used in a large variety of applications because the unique structure, and renewability feature, and chemical and physical properties, and because it is physiologically innocuous.

The use of glycerol for advanced organic syntheses and the conversion of glycerol into value-added chemicals is now emerging as a fascinating challenge. such as Oxidation, Esterification, Chlorination, Dehydration, Hydrogenolysis, Polymerization, Etherification, glycerol is nowadays considered one of the most relevant platform chemicals.

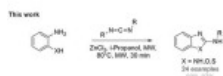
[Short communication](#) [Open access](#)**Microwave assisted one pot synthesis of 2-alkyl amino benzimidazoles, 2-alkyl amino benzoxazoles and 2-alkyl amino benzthiazoles by using various carbodiimides**

Thirupathi Rapolu, Polkam Naveen, K.V.P. Pavan Kumar, Krishnakanth Reddy Leleti, Raghubabu korapolu

Article 100351

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An efficient triple cascade process for synthesis of novel disperse dyes from lawsone: A modification of natural colorant

Paresh N. Patel, Nilam C. Patel, Dipen H. Desai

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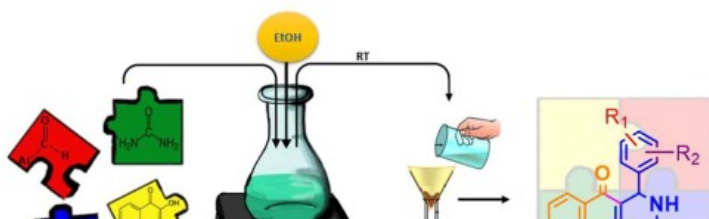
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The present studies highlight the novel approach of triple cascade reaction for the synthesis of dihydropyrimidone in a very effective way. Various novel dihydropyrimidone derivatives are synthesis by this process. Structures of all these derivatives are established by various spectroscopic techniques.

Research article [Open access](#)

In situ detection of an unstable *C,N*-Au(III) chelate by ¹⁵N NMR methods

Jostein Lund, Helgi Freyr Jónsson, Anne Fiksdahl

Article 100360

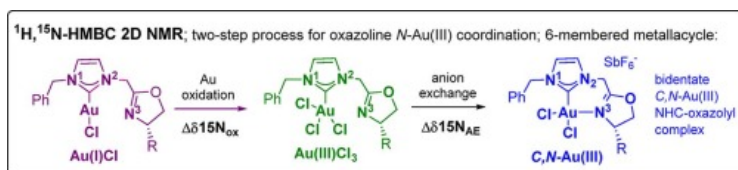
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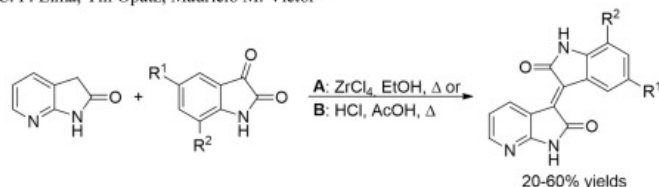


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Synthesis of new substituted 7-azaisoindigos

Cintia M. C. F. Lima, Till Opatz, Mauricio M. Victor*



Research article Open access

Asymmetric silane reduction of ketones and β -Keto esters catalyzed by a chiral azolium/iridium system in the presence of a base in methanol at room temperature

Toshiya Matsuki, Hiro Teramoto, Ryo Ichihara, Kazuo Inui, Satoshi Sakaguchi

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Abstract

Abstract

Asymmetric silane reduction of ketones catalyzed by a system comprising a chiral azolium salt **1** and $[\text{Ir}(\text{cod})_2]\text{BF}_4$ (**2**) in the presence of a base in methanol was achieved at room temperature within 2 h. Implementing a “pre-mixing” experimental procedure improved the efficiency of the catalytic reaction. Specifically, after treating a mixture of **1** and **2** with $(\text{EtO})_2\text{MeSiH}$ (**3a**) in THF, the ketone substrate was allowed to react in the presence of K_2CO_3 in methanol at room temperature to afford the corresponding optically active alcohol in high yield and enantioselectivity. The reaction parameters were fully evaluated. The presence of **3a** in the pre-mixing step and that of a base in the reduction step were found to be essential for success. The developed method was applied to the stereoselective reduction of β -keto esters. Additionally, an iridium species derived from the reaction of **1** with $[\text{IrCl}(\text{cod})]_2$ and AgBF_4 was investigated.

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Thermophysical properties of hexapropyl and hexabutyl phosphoramides in *n*-dodecane

G. Jegan, Satyabrata Mishra, B. Sreenivasulu, Puspaltha Rajesh, ... N. Sivaraman

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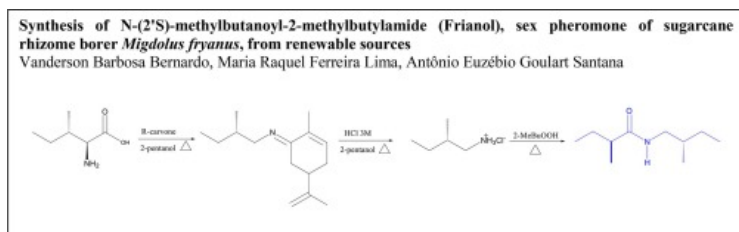
Synthesis of N-(2'S)-methylbutanoyl-2-methylbutylamide (Frianol), sex pheromone of sugarcane rhizome borer *Migdolus fryanus*, from renewable sources

Vanderson Barbosa Bernardo, Maria Raquel Ferreira Lima, Antônio Euzébio Goulart Santana

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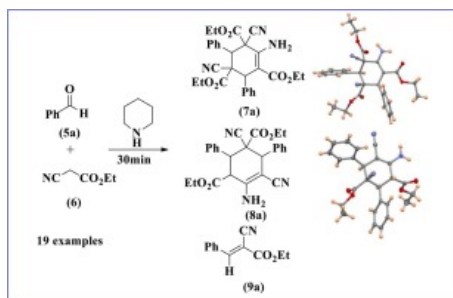
A new finding in the old Knoevenagel condensation reaction

Attimogee Shivamurthy Harisha, Kuppuswamy Nagarajan, S.Saravanan, Venkat Manohar, ... Tayur Narasingarow Guru Row

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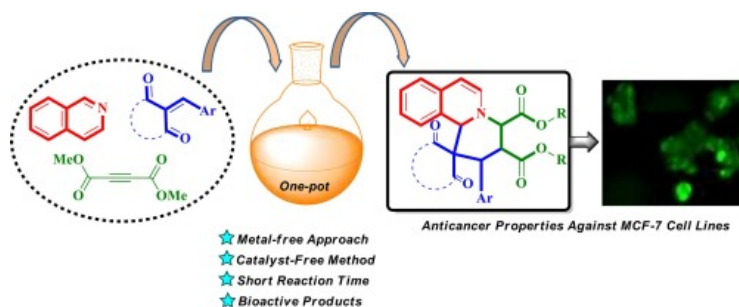
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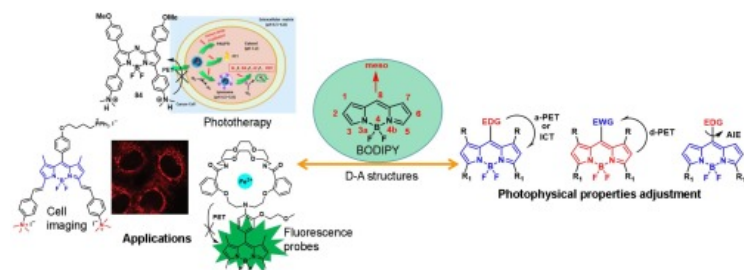
Photophysical properties regulation and applications of BODIPY-based derivatives with electron donor-acceptor system

Fang-Zhou Li, Zhiwei Wu, Changwei Lin, Qiang Wang, Gui-Chao Kuang

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An approach for preparing 3,3-disubstituted oxindole from acyclic tetrasubstituted aldehyde: Total synthesis of (-)-coerulescine & (-)-coixspirolactam A

M. Shahnawaz Ali, Mizzanoor Rahaman, Jawad Bin Belayet, Sharif A. Asad, M. Mahmum Hossain

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[\(-\)-Coeruleosine
\(6 steps, 29.4% yield\)](#)

H

[\(-\)-Coixisprolactam A
\(5 steps, 24.7% yield\)](#)Research article [Open access](#)

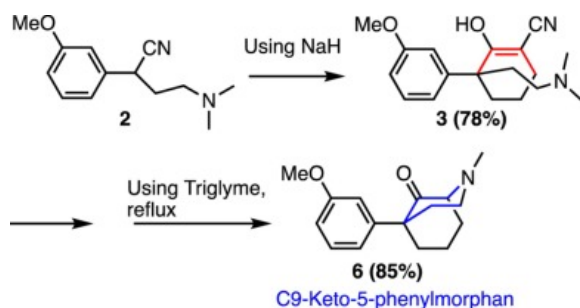
Optimized synthesis of enantiomeric C9-keto-5-phenylmorphans, essential intermediates for novel MOR agonists and antagonists

Agnieszka Sulima, Jason A. Deck, Muneaki Kurimura, Arthur E. Jacobson, Kenner C. Rice

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Short communication [Open access](#)

Radical character quenches luminescence in the all-hydrocarbon radical benzoBDPA

Adrian R. Birge, Matthew J. Piper, Karl J. Painter, Graham T. Sazama

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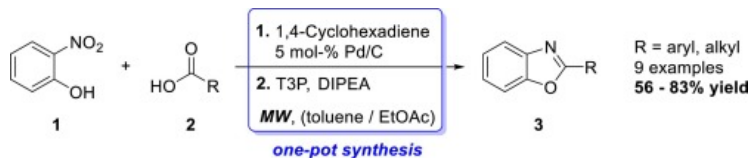
Abstract

We have synthesized and characterized a benzannulated analog, bBDPA, of the well-known stable radical bis(diphenylene)phenylallyl (BDPA). Benzannulation gives rise to fluorescence in the closed-shell precursor molecule bBDPAH, but luminescence is quenched in the radical species. We detail the synthesis, characterization by absorption, emission, and EPR spectroscopy, and examine the molecular orbital landscape dictating the properties of the radical. We find that benzannulation alters the energy of the radical's frontier molecular orbitals differently than in luminescent radicals brought about by incorporation of heteroatoms into their frameworks.

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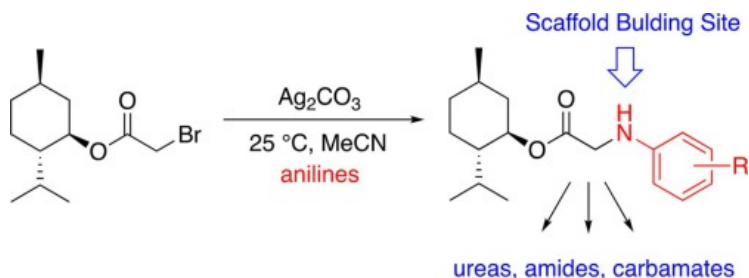
Synthesis of aniline-based menthol glycinates and derivatives

Joseph E. Mathews, Jacob C. Hood, Douglas A. Klumpp

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Highly efficient method for the synthesis of substituted benzimidazoles using sodium metabisulfite adsorbed on silica gel

Manoj Kumar, Sushil K. Pandey, Neha Chaudhary, Anand Mishra, Deepshikha Gupta

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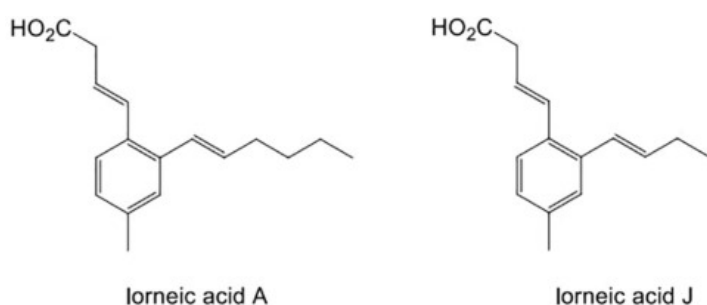
The first total synthesis of lorneic acid J and an alternative synthesis to lorneic acid A

Brandon Sarrels, Wisam Morrar, Claudia G. Lucero

Article 100408

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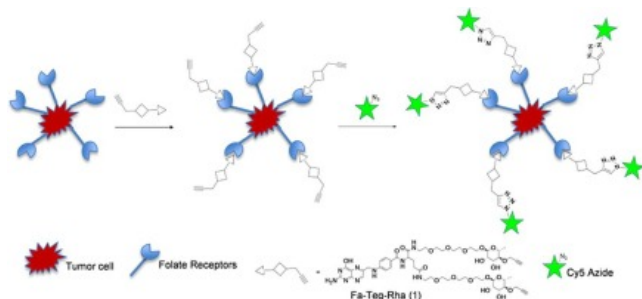
A clickable folic acid-rhamnose conjugate for selective binding to cancer cells

Shubham Parashar, Vatika Gupta, Rakesh Bhatnagar, Amina Kausar

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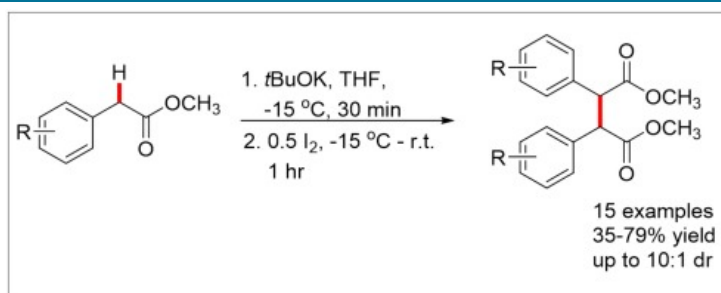
Iodine-promoted oxidative homocoupling of methyl 2-arylacrylates under equilibrium conditions using potassium *tert*-butoxide as a base

Ahmed Y. Nuriye, Joseph Barr

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One pot three components solvent free synthesis of 4-substituted phenyl-2-(sulfanyl/oxo) pyrimidine-5-carboxylate derivatives

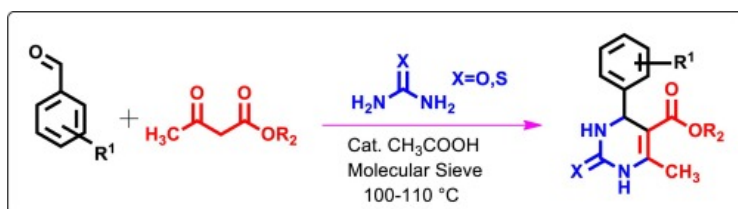
Sourav Handique, Priyanka Sharma

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Synthesis of phenol derivatives from cyclopentenones with arynes as key intermediates

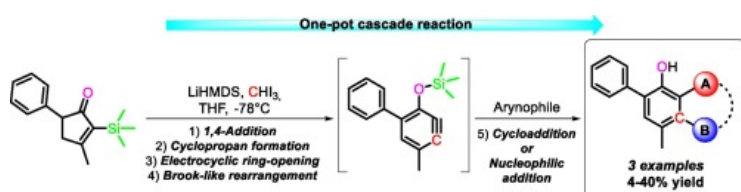
Dario Telese, Markus Staudt, Lennart Bunch

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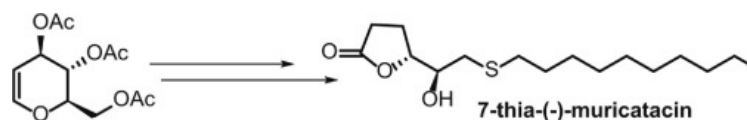
[View PDF](#)[Article preview](#) **Abstract****Abstract**

A facile method for the reduction of carboxylic acid group of Bezafibrate, an approved drug, is described. The selective reduction of carboxylic acid group to corresponding alcohol was carried out by activation of the carboxylic acid moiety via mixed anhydride followed by the addition of stoichiometric amount of NaBH_4 and methanol to obtain the first alcohol variant of Bezafibrate. The reaction was completed in 5–10 min in excellent yield and purity. The new alcohol derivative was characterized by spectroscopic methods. This is the first report on this new molecule.

Research article [Open access](#)**First total synthesis of 7-thia-(-)-muricatacin**

Uxía Gómez Bouzó, Irene Sánchez-Sanz, Maria González, Generosa Gómez, Yagamare Fall

Article 100431

[View PDF](#)[Article preview](#) **Abstract****Graphical abstract****Graphical abstract**Research article [Open access](#)**Oxa-Michael/Ugi-Smiles cascade reaction with α,β -unsaturated aldehydes**

Aubrey Thessing, Robert Ayres, Joshua Jones, Haleigh Bauer, ... Sarah B. Luesse

Article 100416

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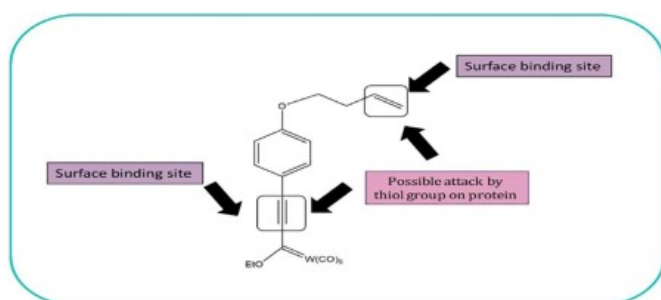
Design of a novel Fischer carbene complex which can facilitate thiol mediated site-specific protein immobilization

Namrata Ray

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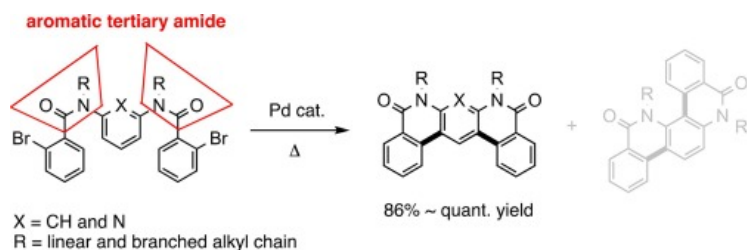
Palladium-Catalyzed intramolecular direct arylation of aromatic tertiary amide compounds revisited

Koji Takagi, Asuka Maeda, Ryotaro Tsunekawa

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Research article [Open access](#)

Pharmacological and quantum chemical studies of 2-aminobenzo[d]thiazol-3-ium 4-chlorobenzenesulphonate: Synthesis, spectral, thermal analysis and structural elucidation

C. Sudhakar, M. Saravanabhavan, K.S. Ramesh, V.N. Badavath, ... M. Sekar

Article 100442

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and grown as single crystals by slow solvent evaporation method at room temperature. Single crystal X-ray diffraction results reveals that 2ABT4CBS crystallized into a monoclinic crystal system with space group $P_{21/c}$. The molecular structure and functional groups were analyzed by spectral studies such as ^1H and ^{13}C NMR, UV-Visible and FT-IR. Thermal stability of the compound was determined by TG and DTA analysis. DFT was used in the calculation of the frontier molecular orbital (FMO), Mulliken load distribution and thermodynamic properties by employing B3LYP/6-31G (d,p) and HF/6-31G levels of theory. The results of the DNA binding demonstrate that the synthesized compound 2ABT4CBS interacts with CT-DNA in a considerable way. The antioxidant capabilities of 2ABT4CBS were also analyzed against DPPH radical and excellent radical scavenging abilities were identified. Molecular docking studies with PDB: 1RNA clearly suggested that synthesized compound

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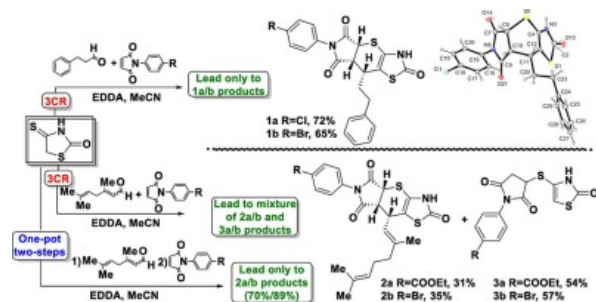
3-Phenylpropanal and citral in the multicomponent synthesis of novel thiopyrano[2,3-d]thiazoles

Andrii Lozynskiy, Andriy Karkhut, Svyatoslav Polovkovych, Olexandr Karpenko, ... Roman Lesyk

Article 100464

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Research article [Open access](#)

Stability of cannabidiol (CBD) in solvents and formulations: A GC-MS approach

Chiara Franco, Stefano Protti, Alessio Porta, Federica Pollastro, ... Daniele Merli

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A simple and effective Lewis acid assisted synthesis of indole-3-sulfonyl carbamates and sulfonamides using Burgess reagent

Sivalenka Vijayasaradhi, Ravula Srinivas, Uppara Ramesh, Koraboina Chandra Prakash, Kota Sathish

Article 100467

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Abstract

Abstract

Indole-3-sulfonyl carbamate served as a key intermediate for the synthesis of a HPK1 inhibitor, bearing a primary sulfonamide group. A simple and efficient synthesis of this key indole-3-sulfonyl carbamate was developed using the readily available Burgess reagent and Lewis acids. A systematic evaluation of a variety of Lewis acids with Burgess reagent to optimize the synthetic yield of indole-3-sulfonyl carbamate is reported in this paper.

Review article [Open access](#)

A minireview on diketopyrrolopyrrole chemistry: Historical perspective and recent developments

Salman A.L. Shaikh, Shailesh S. Birajdar, Sumit D. Ambore, Avinash L Puyad, ... Sheshanath V. Bhosale

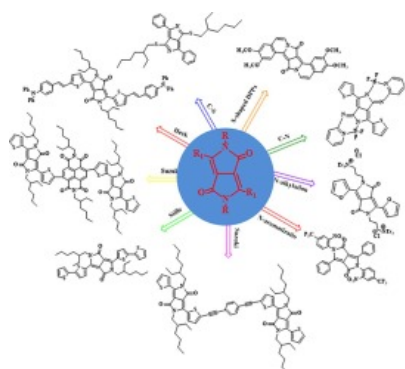
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Abstract

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Silica gel-assisted synthesis of benzo[*b*]thiophenes from *o*-(alkylsulfanyl)(ethynyl)benzenes

Kozo Toyota, Hiroki Tanaka, Taisei Hanagasaki

Article 100487

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under thermal conditions. Tertiary alkyl-substituted sulfanyl derivatives gave good results. Methylsulfanyl-substituted precursors also afforded benzo[*b*]thiophenes at higher temperature. Effect of modified silica gel was evaluated and the mechanism of this method was discussed. This method was also applied to preparation of 2-substituted thieno[3,2-*b*]thiophenes, which gave moderate yields.

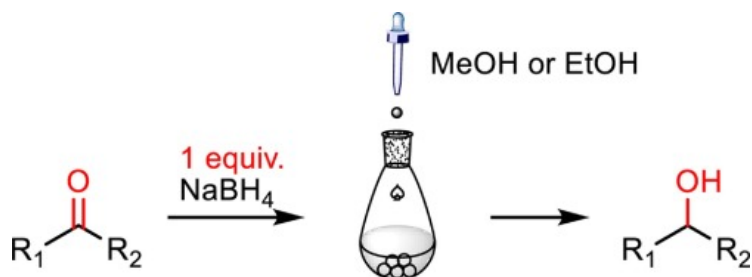
Research article [Open access](#)Reduction of aldehydes and ketones by a stoichiometric amount of NaBH₄ using a small amount of MeOH or EtOH

Kiyoshi Tanemura

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Research article [Open access](#)

Incredible colorimetric sensing behavior of pyrazole-based imine chemosensor towards copper (II) ion detection: synthesis, characterization and theoretical investigations

Malini Nelson, Harikrishnan Muniyasamy, Pavithra Ongi, Sankar Balakrishnan, ... Rajesh Jegathalaprathaban

Article 100501

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Abdolhamid Bamoniri, Nahid Yaghmaeiyan

Article 100505

[View PDF](#)[Article preview](#) **Abstract****Abstract**

Kaolin sulfonic acid nanoparticles (Nano-kaolin-SO₃H) as a recoverable solid catalyst was prepared by reaction of kaolin nanoparticles with chlorosulfonic acid. After characterization of kaolin-SO₃H nanoparticles by ATR, XRD, TEM, XRF, FESEM, EDS, EDS mapping, BET and TGA techniques, this solid acid used to synthesize highly substituted pyrazoles *via* condensation of hydrazine derivatives with various β -diketones. The reaction was proceeded in a short time, simple work-up, easy purification technique, high yields, solvent free conditions and above all environmentally benign.

Research article [Open access](#)KI-H₂O₂ promoted intramolecular oxidative C—H Functionalization: Synthesis of Benzo[*d*]thiazol-2-amines

Manish M. Katiya, Madhuri M. Sontakke, Madhukar G. Dhonde

Article 100512

[View PDF](#)[Article preview](#) **Abstract****Abstract**

This work reported a novel, efficient and green technique for the synthesis of *N*-substituted benzothiazoles using KI/H₂O₂ in DCM via intramolecular oxidative C—H functionalization was carried out. The simple isolation, mild reaction conditions, easy purification without column with good yields, rendering the methodology accessible herein highly eco-friendly compare to existing method.

Research article [Open access](#)

A synthetic methodology for the preparation of a water-soluble quaternized amino phosphine oxide ligand

Ioannis D. Kostas, Georgia Antonopoulou, Polydoros-Chrysovalantis Ioannou, Eleftherios Ferentinos, Panayotis Kyritsis

Article 100525

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Short communication *Open access*

Synthesis of advanced Maillard products: 2-formylpyrrole alkaloids

Brooks E. Maki, Xinyi Lily Feng, Maria R. Mosiyuchuk, Noah B. Davis

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Abstract

Abstract

A biologically-inspired method to access 2-formylpyrrole structures is reported using an Achmatowicz reaction and condensation sequence. This robust synthetic process has been used to access a variety of 2-formylpyrrole structures related to naturally isolated and bioactive compounds. Several families of pyrrole-based natural products have recently been isolated, and these structurally related compounds have been shown to demonstrate interesting bioactive properties. Synthesis targets include the natural products sinopyrrole C and pyrrolezanthine.

Research article *Open access*

Evaluation of an approach for the synthesis of C2 alkylated gramines

Khorshada Jahan, Tye Seideman, Shah Nawaz Ali, Mizzanoor Rahaman, M. Mahmum Hossain

Article 100535

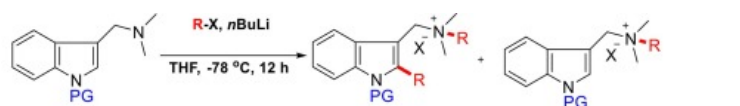
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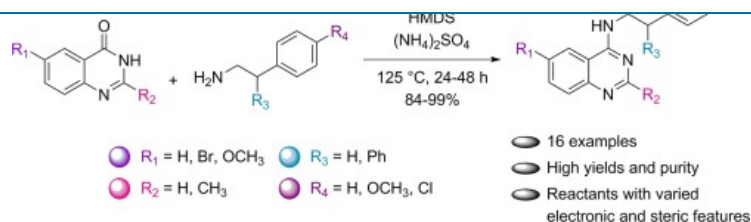
PG= Boc, Tosyl, AcO, Me, Bz etc

Research article *Open access*Synthesis of *N*-phenethylquinazolin-4-amines *via* silylation-amination mediated by hexamethyldisilazane

Guilherme Arraché Gonçalves, Flávio Castro do Nascimento, Sidnei Moura e Silva, Cristiano Valim Bizarro, ... Pablo Machado

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Synthesis and biological evaluation of mannosyl triazoles and varying the nature of substituents on the terminal phthalimido moiety in the aglycone backbone

Hussein Al-Mughaid, Younis Jaradat, Maha Khazaaleh

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Abstract

Abstract

In the present study, a facile synthesis of six monovalent α -D-mannoside ligands terminated with phthalimido moiety by Cu(I) catalyzed azide-alkyne cycloaddition reaction (CuAAC) has been achieved. All synthesized ligands were tested for their inhibitory activities against *E. coli* FimH adhesion using hemagglutination inhibition (HAI) assay and showed inhibitory activity in the range of HAI = 4.8–23.5 μ M comparing with 2-azidoethyl α -D-mannopyranoside **17** as the standard ligand (HAI = 135 μ M). Among them, ligand **21** (R = NO₂), displayed the best activity (HAI = 4.8 μ M) which was approximately 28 times more potent than the reference ligand **17**. Thus confirming the beneficial effect of lipophilic interactions between the aromatic aglycone and the tyrosine gate of FimH. We feel that our lead ligand will be a good platform for identification of more potent FimH inhibitors.

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“Click” synthesis of amphiphilic carbohydrate-alkyl triazole derivatives

Mario Daniel Contin, M. Leticia Bravi Costantino, Norma B. D'Accorso

Article 100558

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Abstract

Abstract

1,4-disubstituted 1*H*-1,2,3-triazole with a hydrophobic and hydrophilic moiety were synthesized by using Cu(I) catalyzed alkyne-azide 1,3-dipolar cycloaddition reaction. The hydrophilic moiety was D-galactopyranosyl or L-galactoyl while an alkyl chain of eight to sixteen carbons corresponds to the hydrophobic portion. All compounds were characterized by NMR and mass spectrometry. Partition coefficient was experimentally determined by HPLC ranging from 0.7 to 4.5. Compounds with the

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Fatemeh Sadegh, Hossein Tavakol

Article 100592



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Abstract**Abstract**

An efficient technique for producing the Ag/CoFe₂O₄ noble trimetallic aerogel is provided. The advantages of this approach are fast preparation, short gelation time, avoiding the use of surfactants, and preparation at room temperature. The Ag/CoFe₂O₄ aerogel was characterized by common analysis methods including FESEM, VSM, EDX mapping, TEM, BET, and XRD. The resulting aerogel exhibits outstanding catalytic activity, and durability in reducing different nitroarenes. The reactions were carried out using hydrazine hydrate as a reducing agent, a 0.03 g catalyst, and ethanol as the solvent at reflux conditions. Under these conditions, 13 derivatives of nitroarenes were successfully reduced to aniline derivatives in 4 h in 70–95 % yield. This reaction had some advantages such as no use of toxic organic solvents, mild reaction conditions, easy separation of the products without any tedious separation techniques, and a recyclable catalyst. The catalyst showed high recyclability in which only 8 % decrease was observed after five runs.

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A green protocol for one-pot synthesis of 3,4-dihydropyrano[*c*]chromenes and biscoumarins by employing WEB as an internal base

Vikrant Kumbhar, Sagar Gaiki, Rutik Raskar, Avinash Kumbhar, Bhushan Khairnar

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Abstract**Graphical abstract****Graphical abstract**Research article *Open access*

Selective silver (I)-catalyzed four-component gram-scale synthesis of novel 1,4-disubstituted 1,2,3-triazole-sulfonamides under heterogeneous catalysis and microwave irradiation in water

Ayoub El Mahmoudi, Hind El Masaoudi, Hamza Tachallait, Aicha Talha, ... Khalid Bougrin

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The Ag_3PO_4 -MEA catalyst was successfully prepared and proved to be an efficient and selective catalyst of the multicomponent synthesis of a new library of 1,4-disubstituted 1,2,3-triazole-sulfonamides **5a-n**. This new methodology cleanly proceeded in water under microwave activation using *N*-propargyl saccharin, NaN_3 or $\text{SO}_2(\text{N}_3)_2$, substituted benzyl chloride or protected sugar and non-conjugated primary amines, using Ag(I) in a method-comparison study to Cu(I) as conventional Sharpless catalyst. The proposed strategy allows a simple, rapid and scalable route to 1,2,3-triazoles sulfonamides derivatives **5** through *in-situ* ring-opening of saccharin system with primary amine. All compounds have been fully characterized by NMR, mass spectrometry and X-ray crystallography analysis. The catalyst has been prepared *via* a green Sol-Gel methodology and fully analyzed by X-ray Diffraction (XRD), Fourier Transform Infrared (FTIR) Spectroscopy and Scanning Electron Microscopy (SEM) coupled to Energy Dispersive X-ray Spectroscopy (EDX) and Brunauer-Emmett-Teller (BET) method. The catalyst is

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Cationic carbohydrate-based surfactants derived from renewable resources: Trends in synthetic methods

Abbas Abdulameer Salman

Article 100602

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Review article [Open access](#)

A review on biological and medicinal impact of heterocyclic compounds

Emranul Kabir, Monir Uzzaman

Article 100606

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Abstract

Heterocyclic compounds have gained a lot of attention because of their numerous significant medical and biological uses. Research interest on heterocyclic compounds is rapidly increasing due to the extensive synthetic study and functional utility. They are found in more than 90% of novel drugs, and span the gap between biology and chemistry, where so much scientific discover

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Synthesis of novel pyridazino[1,6-*a*]indole-2,4(1*H*,3*H*)-dione and pyridazino[1,6-*a*]indol-2(1*H*)-one via intramolecular electrophilic aromatic substitution

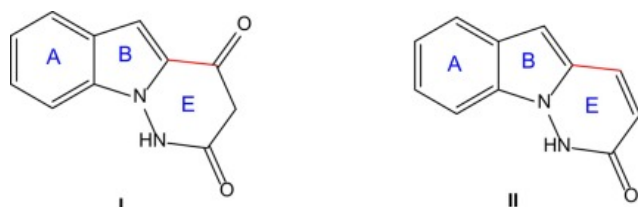
Katarzyna Ozdarska, Maud Petremant, Kai-Chen Wu, Erika Bourguet

Article 100612

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- Intramolecular electrophilic aromatic substitution
- Metal catalyst free
- Simple one-step synthesis

Research article Open access

Design, synthesis and biological evaluation of novel substituted indazole-1,2,3-triazolyl-1,3,4-oxadiazoles: Antimicrobial activity evaluation and docking study

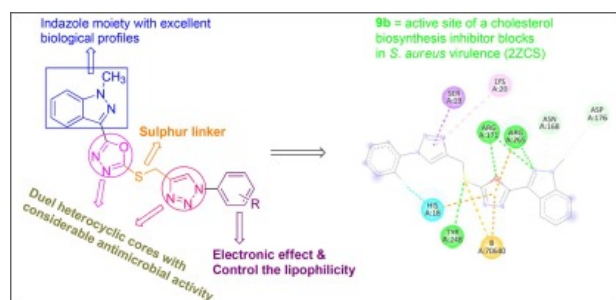
Ananda Kumar Dunga, Tejeswara Rao Allaka, Yugandhar Kethavarapu, Sunil Kumar Nechipadappu, ... Pilli V.V.N. Kishore

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Microwave-assisted synthesis of (6-((1-(4-aminophenyl)-1H-1,2,3-triazol-4-yl)methoxy)substituted benzofuran-2-yl(phenyl)methanones, evaluation of *in vitro* anticancer, antimicrobial activities and molecular docking on COVID-19

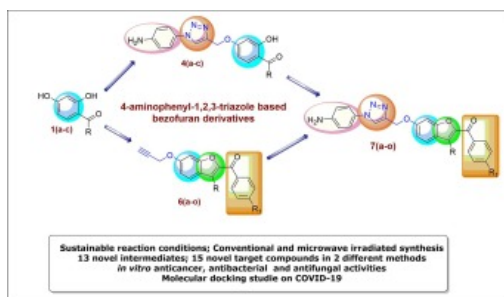
Ravinder Dharavath, M. Sarasija, K.N. Prathima, M. Ram Reddy, ... D. Ashok

Article 100628

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Sulfated polyborate as Bronsted acid catalyst for Knoevenagel condensation

Perna Ganwir, Priyanka Bandivadekar, Ganesh Chaturbhuji

Article 100632

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Research article [Open access](#)One-pot synthesis of novel 2-oxo(2*H*)-spiro[benzofuran-3,3'-pyrrolines] *via* 1,4-dipolar cycloaddition reaction

Mohammed M. Al-Mahadeen, Jalal A. Zahra, Mustafa M. El-Abadelah, Areej M. Jaber, Monther A. Khanfar

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Abstract

Abstract

An efficient synthesis of novel spiro[2-oxobenzofuran-3,3'-pyrrolines] is achieved *via* a three-component 1,4-dipolar cycloaddition reaction at rt involving benzofuran-2,3-diones, dimethyl acetylenedicarboxylate along with imidazo[1,5-*a*]pyridine, or imidazo[1,5-*a*]quinoline or imidazo[5,1-*a*]isoquinoline. The expected isomeric benzofuro[3,2-*f*]oxazepines (kinetic control cycloadducts) were, however, not detected. Structures of the title spiro products are evidenced from HRMS and NMR spectral data and further confirmed by single-crystal X-ray crystallography.

Research article [Open access](#)

Coumarin – benzimidazole hybrids: A review on diverse synthetic strategies

C.G. Arya, Munugala Chandrakanth, K. Fabitha, Neethu Mariam Thomas, ... Janardhan Banothu

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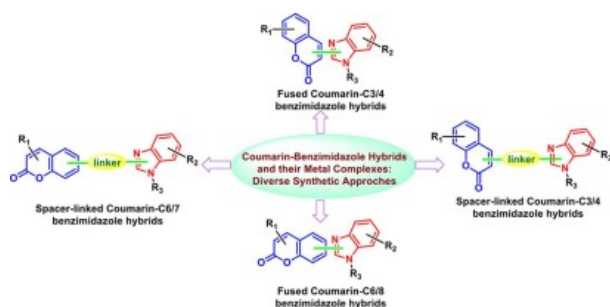
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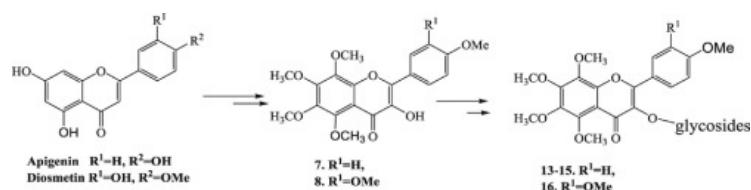
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The syntheses of polymethoxy flavonoid glycosides from apigenin and diosmetin and their cytotoxic activities on two human cancer cell lines

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Short communication *Open access*

Synthesis and biological evaluation of aryl 1,2,4-oxadiazole incorporated (2-(oxazol)-1H-imidazoles as anticancer agents

Suresh Bairi, Veerachamy Alagarsamy, Shyam Sunder Rachamalla, Laxminarayana Eppakayala

Article 100649

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Abstract

Abstract

A new series of 1,2,4-oxadiazole incorporated (2-(oxazol)-1H-imidazole derivatives **9a-j** have been developed and synthesized. Further, the newly prepared compounds (**9a-j**) compounds were evaluated for their preliminary anticancer properties against four human cancer cell lines like prostate cancer (PC3&DU-145), lung cancer (A549), and breast cancer (MCF-7) by using of MTT method and etoposide used as a positive reference. The obtained results indicated that most of the tested compounds displayed excellent to moderate activity. Among all derivatives, six compounds **9a**, **9b**, **9c**, **9d**, **9e**, and **9f** exhibited more potent anticancer activity than etoposide.

Research article *Open access*

Efficient synthesis of a novel euchromatic histone methyl transferase 2 (G9a) inhibitor

Davide Gornati, Roberta Sinisi, Stefania Bertuolo, Marilena De Matteo, Romano Di Fabio

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Abstract

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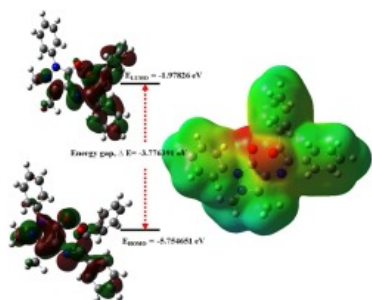
dihydro-3H-pyrazol-3-one

S. Anila Raj, V.G. Vidya, V. Preethi, V.G. Viju Kumar

Article 100665

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Graphical abstract

Research article [Open access](#)Colorimetric detection of Cu²⁺ by amino phenol based chemosensor

D.Y. Patil, N.B. Khadke, A.A. Patil, A.V. Borhade

Article 100658

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Abstract

A colorimetric chemosensor, based on 2-amino phenol Schiff base has been successfully synthesized by simple condensation reaction of 2-amino quinoline 3-carbaldehyde and 2-amino phenol. Synthesized Schiff base ligand 2-(2-aminoquinolin-3-yl methyleneamino) phenol (SBL) was characterised by various spectroscopic techniques. SBL displayed highly selective and sensitive Colorimetric enhancement to Cu²⁺ ion over 13 competing ions in ethanol–water (99:1/v:v) solution. UV titration experiment showed that SBL exclusively reacted with Cu²⁺ and the stoichiometric ratio is 1:1. The detection limit of SBL-Cu²⁺ was estimated to be 4.3 nM and 8.2×10^{-8} M, respectively.

Research article [Open access](#)Highly efficient magnetically separable Zn-Ag@L-arginine Fe₃O₄ catalyst for synthesis of 2-aryl-substituted benzimidazoles and multicomponent synthesis of pyrimidines

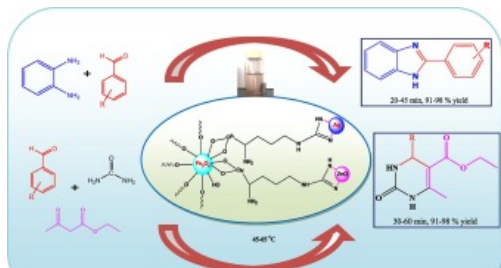
Padmakar A. Kulkarni, Sandeep S. Kahandal, Nitin A. Mirgane, Ashis Kumar Satpati, Suresh S. Shendage

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and multicomponent synthesis of pyrimidines.



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Synthesis, characterization, DFT studies and molecular docking investigation of 2-oxo-ethyl piperidine pentanamide-derived sulfonamides as anti-diabetic agents

Fredrick C. Asogwa, Ekoh C. Ogechi, H. Louis, Ugwu D. Izuchukwu, ... Okoro U. Chris

Article 100672

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Abstract

Abstract

The present study is focused on the synthesis, characterization, DFT Studies, and *in silico* biological evaluation of novel benzenesulfonamides as antidiabetic agents. The compounds were synthesized and characterized experimentally using the NMR, FT-IR and HRMS while the molecular electronic structure properties were investigated using DFT-based theoretical methods at the M06-2X/6-311++G (d, p) level of theory. The results of the calculated vibrational energies were in perfect agreement with those of the experimental frequencies for different functional groups. The natural bond orbital (NBO) analysis was also carried out to study the intramolecular charge transfer interactions and energies of stabilization. Molecular docking simulation showed that the compounds; 3-Methyl-2-(4-methylphenylsulfonamido)-N-(2-oxo-2(piperidin-1-yl) ethyl) pentanamide (MMOPEP), 3-Methyl-N-(2-oxo-2-(piperidin-1-yl)ethyl)-2-(phenylsulfonamido)pentanamide (MOPEPP) and, 2-Methyl-2-(4-nitrophenylsulfonamido)-N-(2-oxo-2-(piperidin-1-yl)ethyl)pentanamide (MNOPEP) fitted tightly to the amino

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Allylic arylation isomerisation of Morita-Baylis-Hillman adduct of isatin derivatives with arenes using FeCl₃ and K-10 clay

Sivaprakasam Sivaraman, Selvaraj Aruljothi, Vadivel Vaithyanathan

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Design, synthesis and evaluation of new alkylated pyrimidine derivatives as antibacterial agents

K.N. Hari, Boja Poojary, G. Chandrasehar

Article 100676

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Graphical abstract



1. Mild reaction conditions
2. Smoother workup
3. Easy access to Pyrimidine derivatives
4. Easy access to HPLC method

Research article *Open access*

2-aminophenol as a leading reactant for the one pot synthetic strategies towards benzoxazole derivatives: A systematic review

Jyoti Sharma, Praveena Mishra, Juli Bhadoria

Article 100670

[View PDF](#) [Article preview](#) **Abstract**

Abstract

Nitrogen containing heterocyclic compounds such as benzoxazole derivatives have been found to be pharmacologically active and are being employed as starting material for several drug discovery researches. Synthetic strategies invention to produce such beneficial moieties is on focus of synthetic organic chemists, as a result various efficient protocols have been developed since past decades to synthesize different medicinally useful benzoxazole derivatives. In this review, we attempt to organize recent synthetic strategies explored by chemists to produce benzoxazole derivatives by the reaction of 2-aminophenol with various reactants in different reaction conditions such as solvents, catalysts, temperature etc. Therefore, the updated synthetic information present in this review may be helpful to the researchers to develop better synthetic protocols for novel medicinally active derivatives.

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Synthesis of AuNPs decorated multi-valent Cu-Ni oxide nanoplates for electrochemical oxidation of methanol

Han Sun, Haiping Huang, Chao Hu, Yu Yan, ... Jing-Lin Chen

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Abstract

Abstract

Cu-Ni oxide nanocomposite was synthesized via solvothermal method with nickel (II) nitrate and copper nitrate as Ni and Cu resource respectively. AuNPs/oxide nanocomposite was obtained by dispersing the oxide into the pre-synthesized Au colloid. Transmission electron microscope (TEM) shows that the synthesized oxides are nanoplate morphology. Interplanar distance from the high resolution TEM (HRTEM) image shows that the Ni and Cu in the oxide are multivalent state. X-ray diffraction (XRD) and X-ray photoelectron spectroscopy (XPS) characterization further proves that the Ni and Cu are multivalent state, and the Au is Au⁰ form in the nanocomposite. EDS-mapping images confirm large numbers of AuNPs are uniformly absorbed onto the ultra-thin multivalent Cu-Ni oxide nanoplate. The electrochemical properties of multivalent oxide (m-v oxide) with or without AuNPs decoration are investigated in 0.1 M potassium chloride electrolyte with 2.0 mM Ferri/Ferro-Cyanide. When employed for the electrochemical catalysis in methanol oxidation reaction (MOR) the AuNPs/m-v oxide exhibits better

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Structures, spectral and photodynamic properties of two nitrosylruthenium (II) isomer complexes containing 8-quinolinolate and L-proline ligands

Chenyang Liu, Yu Wang, Ai Wang, Feng Su, Hongfei Wang

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a



b

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AIE or AIE(P)E-active transition metal complexes for highly sensitive detection of nitroaromatic explosives

Arumugam Ramdass, Veerasamy Sathish, Pounraj Thanasekaran

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Abstract

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This review focuses on the development of AIE or AIE(P)E-active transition metal complexes for the sensitive and selective detection of nitroexplosives.

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Investigation of phosphine donor properties to vanadium(V) nitrides

Linqing Mo, Hannah I. Barr, Aaron L. Odom

Article 100344

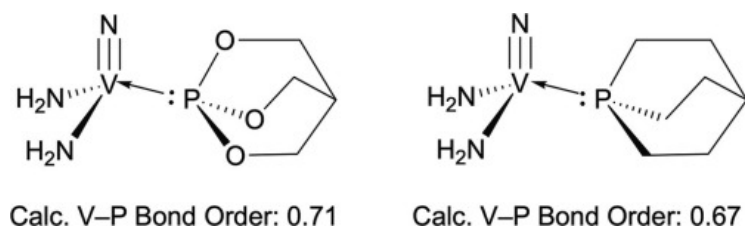
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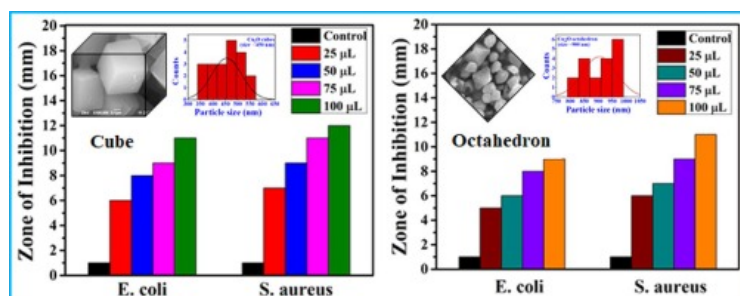
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Research article *Open access*Investigation on crystal facet-dependent antibacterial activity of Cu₂O crystals – A structural, morphological, and spectroscopy studies

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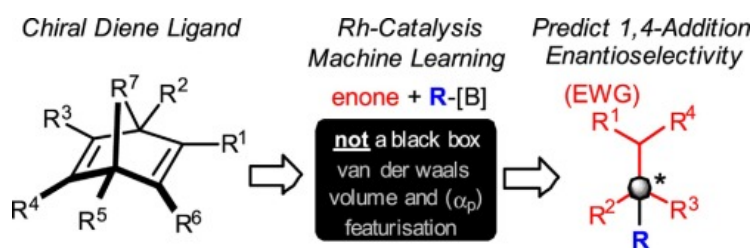
Machine learnt patterns in rhodium-catalysed asymmetric Michael addition using chiral diene ligands

Benjamin Owen, Katherine Wheelhouse, Graziela Figueredo, Ender Özcan, Simon Woodward

Article 100379

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Research article [Open access](#)

Hybrid polyphenolic Network/SPIONs aggregates with potential synergistic effects in MRI applications

A. Lazzarini, R. Colaiezzi, A. Galante, M. Passacantando, ... M. Crucianelli

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NIR-emission from Yb(III)- and Nd(III)-based complexes in the solid state sensitized by a ligand system absorbing in a broad UV and visible spectral window

Enrico Cavalli, Silvia Ruggieri, Silvia Mizzoni, Chiara Nardon, ... Fabio Piccinelli

Article 100388

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Abstract

Abstract

In this contribution, we present the synthesis, characterization and spectroscopic investigation of the heteroleptic (*R,R*)-YbL $\mathbf{1}$ (tta) and (*R,R*)-NdL $\mathbf{1}$ (tta) complexes (with tta = 2-thenoyltrifluoroacetate and L $\mathbf{1}$ = N,N'-bis(2-(8-hydroxyquinolate)methylidene)-1,2-(*R,R* or *S,S*)-cyclohexanediamine) in the solid state. The *f-f* metal-centered NIR luminescence emission of Nd(III) and Yb(III) is efficiently sensitized by both chromophoric ligands in a very broad range of wavelengths [from 250 to 600 nm, in the case of Nd(III) and from 250 to 650 nm, for Yb(III)]. A possible energy transfer mechanism is proposed: for (*R,R*)-NdL $\mathbf{1}$ (tta) complex a classical Ligand-to-Metal Energy Transfer (LMET) mechanism (*antenna effect*) is suggested, whilst in the case of the (*R,R*)-YbL $\mathbf{1}$ (tta) complex, the presence of a ligand-to-metal charge transfer (LMCT) state determines the sensitization of Yb(III) luminescence. We propose that this level is populated by the singlet and triplet excited states belonging to $\pi \rightarrow \pi^*$ and $n \rightarrow \pi^*$ transitions of both ligands and it can transfer the excitation

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