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Education

Doctorate, Erciyes University, Fen Fakültesi, Kimya, Turkey 1987 - 1990

Postgraduate, Erciyes University, Fen Fakültesi, Kimya, Turkey 1985 - 1987

Undergraduate, Inonu University, Temel Bilimler Fakültesi, Kimya Mühendisliği, Turkey 1979 - 1983

Dissertations

Doctorate, 4-BENZOIL-5-FENİL-FURAN-2,3-DİON'UN ÇEŞİTLİ SEMIKARBAZON'LAR, aminoguanidinler ve amidlerle reaksiyonları, Erciyes Üniversitesi, Fen Bilimleri Enstitüsü, Kimya, 1990

Research Areas

Simulation and Modelling, Bioinformatics, Chemistry, Organic Chemistry, Chemistry of Heterocyclic Compounds, Natural Sciences

Academic Positions

Professor, Erciyes University, Fen Fakültesi, Kimya, 2003 - Continues

Academic and Administrative Experience

Head of Department, Erciyes University, Fen Fakültesi, Kimya, 2016 - Continues

Supervised Theses

Sarıpinar E., N-(aminoamidino)tiyoüre türevlerinin pfa ortamında sıklızasyon reaksiyonlarının araştırılması,
Postgraduate, H.DOĞAN(Student), 2019

Sarıpinar E., Benzotiyenopirimidin bileşik serisinin elektron konformasyon-genetik algoritma (EC-GA) yöntemiyle 4D-QSAR incelenmesi, Postgraduate, H.KESKİN(Student), 2019

Sarıpinar E., Pirimidin ve imidazol çekirdeği içeren bileşiklerin sentezi ve in vitro sitotoksik aktivite çalışmaları, Doctorate, S.ÇAĞLAR(Student), 2019

Sarıpinar E., Fenilpirazin, alkinilfenoksiasetik asit, ftalazin türevlerinin elektron konformasyonal-genetik algoritma (EC-GA) metodu ile 4D-QSAR analizi, Doctorate, S.KÖPRÜ(Student), 2018

Patat Ş., SARIPINAR E., REACTION OF N-(AMINOAMIDINO)THIOUREA DERIVATIVES WITH DIMETHYLACETYLENE DICARBOXYLATE, Postgraduate, İ.ADNAN(Student), 2018

Sarıpinar E., Elektron konformasyonel-genetik algoritma 4D-QSAR metodu ile piridin karboksilik asit, oksadiazol, pirimidin ve oksazoltürevlerine ait farmakofor gruplarının belirlenmesi ve biyoaktivite hesabı, Doctorate, B.TÜZÜN(Student), 2018

Sarıpinar E., Çok fonksiyonlu 4-oksotiyaoazolidin sentezi ve kuantum hesaplamaları, Postgraduate, N.DOĞAN(Student), 2017

Sarıpinar E., Elektron konformasyonel-genetik algoritma metodu ile fenilpirazolglutamik asit piperazin bileşiklerinde farmakofor belirlenmesi ve 4D-QSAR analizi, Postgraduate, R.BAHAR(Student), 2016

Sarıpinar E., Dipeptidil bor bileşiklerinin elektron konformasyonel-genetik algoritma (EC-GA) metodu ile 4D-QSAR analizi, Postgraduate, S.ÇATALKAYA(Student), 2014

Sarıpinar E., Rutenyum(II) aren komplekslerinin elektron konformasyonel-genetik algoritma (EC-GA) metodu ile farmakofor modellemesi ve 4D QSAR analizi, Postgraduate, S.ÇAĞLAR(Student), 2013

Sarıpinar E., Anti-HIV etkili HEPT, TIBO, tiyazolidin ve DABO türevlerinin elektron konformasyonel-genetik algoritma metodu kullanılarak yeni bir 4D-QSAR yöntemi ile aktif gruplarının belirlenmesi, Doctorate, L.AKYÜZ(Student), 2010

Sarıpinar E., 3-(1,1-diokso-2h-(1,2,4)-benzotiyodiazin-3-il)-4-hidroksi-2(1h)-kinolin bileşik serisinin hepatit c inhibitörü olarak elektron konformasyon-genetik algoritma (EC-GA) yöntemi ile QSAR incelenmesi, Postgraduate, S.AYDIN(Student), 2010

Sarıpinar E., Elektron konformasyonel-genetik algoritma qsar metodu ile penisilin türevlerine ait farmakofor grupların belirlenmesi ve biyoaktivite hesabı, Postgraduate, E.YANMAZ(Student), 2010

Sarıpinar E., Elektron konformasyonel genetik algoritma 4d-QSAR metodu ile pirazol, benzotriazin, dibenzazosin ve kinazolin serilerinde farmakofor belirlenmesi ve biyoaktivite hesabı, Doctorate, K.ŞAHİN(Student), 2010

Sarıpinar E., 1,4-dihidropiridin, artemisinin ve triazolpiridinoksazol türevlerinin nicel yapı aktivite ilişkilerinin incelenmesi ve biyoaktif yapının belirlenmesi, Doctorate, N.GEÇEN(Student), 2010

Sarıpinar E., ddd, Doctorate, M.SAŞMACI(Student), 2009

Sarıpinar E., HEPT bileşik serisinin HIV-1 inhibitörü olarak elektron konformasyon-genetik algoritma (EC-GA) yöntemi ile QSAR incelenmesi, Postgraduate, E.KAYA(Student), 2008

Sarıpinar E., Adenosin bileşik serisinin anti-tümör aktivitesinin elektron konformasyon-genetik algoritma (EC-GA) yöntemi ile QSAR incelenmesi, Postgraduate, Y.YAZAR(Student), 2008

Sarıpinar E., F2-NH-DABOs ve S-DABOs bileşik serilerinin HIV-1 inhibitörü olarak EC yöntemi (elektron konformasyonel) ile QSAR incelenmesi, Postgraduate, S.GÖKTAŞ(Student), 2007

Sarıpinar E., Benzodiazepinon (TIBO) bileşiklerinin HIV-1 inhibitörü olarak CoMCET yöntemi ile 3D-QSAR incelenmesi, Postgraduate, T.HİZCAN(Student), 2005

Sarıpinar E., 2-amino-6-arilsülfonil benzonitril bileşiklerinin HIV-1 inhibitörü olarak CoMCET metodu ile QSAR incelenmesi, Postgraduate, L.KULAK(Student), 2005

Sarıpinar E., Katı süper asit katalizörü kullanılarak aromatik hidrokarbonların uzun zincirli olefinlerle alkilasyonu, Postgraduate, E.BULUT(Student), 2005

Sarıpinar E., Keten ve oksoketenlerin çeşitli 1,3 dipolar bileşiklerle reaksiyonlarının kuantum kimyasal incelenmesi, Postgraduate, Ş.ŞAHİNDAŞ(Student), 2004

Sarıpinar E., Konjuge doymamış karbonil bileşiklerine organometal bileşiklerinin katılma mekanizmasının kuantum kimyasal incelenmesi, Postgraduate, A.YENER(Student), 2003

Sarıpinar E., Furan-2,3-dion'un semikarbanon ve fenilhidrazin ile reaksiyon mekanizmasının semiempirik (PM3) yönteme araştırılması, Postgraduate, N.YAZIR(Student), 2002

Sarıpinar E., Fonksiyonel yeni pirimidin, piridazin ve imidazol türevlerinin sentezi: Deneysel verileri ve semiempirik (AM1) hesaplamalar, Postgraduate, Ç.YILMAZ(Student), 2000

Sarıpinar E., Visinal furan ve pirrol-2,3-dion sistemlerinin hidrazin ve o-fenilendiamin türevleriyle reaksiyonlarının deneysel ve semiempirik (PM3) verileri, Postgraduate, İ.ÖNCEL(Student), 1999

Sarıpinar E., N-arylpirrol-2,3-dion türevlerinin sentezi, termolizi, hidrzin hidrat ve o-difenilendiaminle reaksiyonları deneysel veriler ve semiempirik hesaplamalar, Postgraduate, S.KARATAŞ(Student), 1997
Sarıpinar E., 4-p Metoksibenzoil -5-p metokspipeni (-2,3-) furandionun çeşitli dienofil ve aminlerle reaksiyonları, Postgraduate, L.GÖKTAŞ(Student), 1995

Journal articles indexed in SCI, SSCI, and AHCI

- I. **Imidazole-based hydrazones as potent anti-colon cancer agents: Design, synthesis, biological evaluation and computational studies**
Tapera M., Doğan E., Şahin K., Gözkamane G. A., Kekeçmuhammed H., Sandal S., GÜRKAN A., BORA R. E., Anber A., Durdagi S., et al.
Journal of Molecular Structure, vol.1318, 2024 (SCI-Expanded)
- II. **Novel 1,2,4-triazole-maleamic acid derivatives: synthesis and evaluation as anticancer agents with carbonic anhydrase inhibitory activity**
Tapera M., Kekeçmuhammed H., TÜZÜN B., DURNA DAŞTAN S., Safa Çelik M., Taslimi P., DAŞTAN T., Topcu K. S. B., Cacan E., Şahin O., et al.
Journal of Molecular Structure, vol.1313, 2024 (SCI-Expanded)
- III. **The Suppressive Effect of Novel Hydrazones-Tethered Imidazoles in HCT-116 and HT-29 Colorectal Cancer Cells: Synthesis, Biological Activity and Molecular Modeling Studies**
Kekeçmuhammed H., Tapera M., Sahin K., Sever B. O., Anber A. M., Bora R. E., AVŞAR T., Kılıç T., Güner E., Sarıpinar E., et al.
CHEMISTRYSELECT, vol.9, no.19, 2024 (SCI-Expanded)
- IV. **Molecular hybrids integrated with imidazole and hydrazone structural motifs: Design, synthesis, biological evaluation, and molecular docking studies**
Tapera M., Kekeçmuhammed H., SARIPINAR E., DOĞAN M., TÜZÜN B., KOÇYİĞİT Ü. M., Çetin F. N.
Journal of Molecular Liquids, vol.391, 2023 (SCI-Expanded)
- V. **Design, synthesis, molecular docking and biological evaluation of 1,2,4-triazole derivatives possessing a hydrazone moiety as anti-breast cancer agents**
Tapera M., Kekeçmuhammed H., Tunc C. U., Kutlu A. U., ÇELİK İ., Zorlu Y., AYDIN Ö., SARIPINAR E.
New Journal of Chemistry, vol.47, no.24, pp.11602-11614, 2023 (SCI-Expanded)
- VI. **Synthesis, Biological Activity Evaluation and Molecular Docking of Imidazole Derivatives Possessing Hydrazone Moiety**
Kekeçmuhammed H., Tapera M., Aydoğdu E., SARIPINAR E., Aydin Karatas E., Mehtap Uc E., Akyuz M., TÜZÜN B., GÜLÇİN İ., Emin Bora R. E., et al.
Chemistry and Biodiversity, vol.20, no.6, 2023 (SCI-Expanded)
- VII. **Experimental and theoretical studies on 3-(4-chlorophenyl)-5-(4-ethoxyphenyl)-4,5-dihydropyrazole-1-carbonitrile: DFT quantum mechanical calculation, vibrational band analysis, prediction of activity spectra, and molecular docking**
Zeyrek C. T., Akman S., İLHAN İ. Ö., KÖKBUDAK Z., SARIPINAR E., AKKOÇ S.
Journal of Molecular Structure, vol.1276, 2023 (SCI-Expanded)
- VIII. **Synthesis, carbonic anhydrase inhibitory activity, anticancer activity and molecular docking studies of new imidazolyl hydrazone derivatives**
Tapera M., Kekeçmuhammed H., TÜZÜN B., SARIPINAR E., KOÇYİĞİT Ü. M., Yıldırım E., DOĞAN M., Zorlu Y.
Journal of Molecular Structure, vol.1269, 2022 (SCI-Expanded)
- IX. **Synthesis, characterization, anti-tuberculosis activity and molecular modeling studies of thiourea derivatives bearing aminoguanidine moiety**
Tapera M., Kekeçmuhammed H., Sahin K., Krishna V. S., Lherbet C., Homerset H., Chebaiki M., Tonjum T., Mourey L., Zorlu Y., et al.
JOURNAL OF MOLECULAR STRUCTURE, vol.1270, 2022 (SCI-Expanded)
- X. **Synthesis, Characterization, Biological Activity and Molecular Modeling Studies of Novel**

Aminoguanidine Derivatives

Doğan N., Çağlar Yavuz S., Sahin K., Orhan M. D., Muhammed H. K., Calis S., Öztürk Küp F., Avsar T., Akkoç S., Tapera M., et al.

CHEMISTRYSELECT, vol.7, no.45, pp.1-20, 2022 (SCI-Expanded)

- XI. **Antiproliferative activity and molecular docking studies of new 4-oxothiazolidin-5-ylidene acetate derivatives containing guanylhydrazone moiety**

Al-Janabi I. A. S., Yavuz S. Ç., KÖPRÜ S., Tapera M., Kekeçmuhammed H., AKKOÇ S., TÜZÜN B., PATAT Ş., SARIPINAR E. Journal of Molecular Structure, vol.1258, 2022 (SCI-Expanded)

- XII. **Synthesis, Molecular Docking and Antiproliferative Activity Studies of a Thiazole-Based Compound Linked to Hydrazone Moiety**

Kekeçmuhammed H., Tapera M., Tüzün B., Akkoç S., Zorlu Y., Sarıpinar E.
CHEMISTRYSELECT, vol.7, no.26, pp.1-10, 2022 (SCI-Expanded)

- XIII. **Combined 4D-QSAR and target-based approaches for the determination of bioactive Isatin derivatives**

Şahin K., Sarıpinar E., Durdağı S.

Sar And Qsar In Environmental Research, vol.1, no.1, pp.1-25, 2021 (SCI-Expanded)

- XIV. **Efficient synthesis and molecular docking studies of new pyrimidine-chromeno hybrid derivatives as potential antiproliferative agents**

Yavuz S. Ç., Akkoç S., Tüzün B., Şahin O., Sarıpinar E.

Synthetic Communications, vol.51, pp.2135-2159, 2021 (SCI-Expanded)

- XV. **Synthesis of novel heterocyclic compounds containing pyrimidine nucleus using the Biginelli reaction: Antiproliferative activity and docking studies**

Yavuz S. C., Akkoç S., Türkmenoğlu B., Sarıpinar E.

JOURNAL OF HETEROCYCLIC CHEMISTRY, vol.57, no.6, pp.2615-2627, 2020 (SCI-Expanded)

- XVI. **A novel hybrid method named electron conformational genetic algorithm as a 4D QSAR investigation to calculate the biological activity of the tetrahydrodibenzazosines**

Sahin K., Sarıpinar E.

JOURNAL OF COMPUTATIONAL CHEMISTRY, vol.41, no.11, pp.1091-1104, 2020 (SCI-Expanded)

- XVII. **Discovery of hydrazone containing thiadiazoles as Mycobacterium tuberculosis growth and enoyl acyl carrier protein reductase (InhA) inhibitors**

Dogan H., DOĞAN Ş. D., GÜNDÜZ M. G., Krishna V. S., Lherbet C., Sriram D., Sahin O., Sarıpinar E.

EUROPEAN JOURNAL OF MEDICINAL CHEMISTRY, vol.188, 2020 (SCI-Expanded)

- XVIII. **Discovery of hydrazone containing thiadiazoles as Mycobacterium tuberculosis growth and enoyl acyl carrier protein reductase (InhA) inhibitors**

Doğan H., Doğan D., Gözde Gündüz, Lherbet C., Sriram D., Şahin O., Sarıpinar E.

European Journal Of Medicinal Chemistry, vol.1, no.1, pp.1-10, 2020 (SCI-Expanded)

- XIX. **The Effect of Stereoisomerism on the 4D-QSAR Study of Some Dipeptidyl Boron Derivatives**

Çatalkaya S., Sabancı N., Çağlar Yavuz S., Sarıpinar E.

Computational Biology And Chemistry, vol.1, no.1, pp.1-16, 2019 (SCI-Expanded)

- XX. **The cytotoxic activities of imidazole derivatives prepared from various guanylhydrazone and phenylglyoxal monohydrate**

Yavuz S. C., AKKOÇ S., SARIPINAR E.

SYNTHETIC COMMUNICATIONS, vol.49, no.22, pp.3198-3209, 2019 (SCI-Expanded)

- XXI. **Molecular docking and 4D-QSAR model of methanone derivatives by electron conformational-genetic algorithm method**

Tüzün B., Sarıpinar E.

Journal Of The Iranian Chemical Society, vol.1, no.1, pp.1-16, 2019 (SCI-Expanded)

- XXII. **DNA Cleavage Properties, Antimicrobial and Cytotoxic Activity and 4D-QSAR Analysis of Some Pyrazole Derivatives**

Köprü S., Kup F. O., Sabancı N., Çadır M., Bulut D. C., Duman F., İlhan İ. Ö., Sarıpinar E.

LETTERS IN DRUG DESIGN & DISCOVERY, vol.16, pp.904-918, 2019 (SCI-Expanded)

- XXIII. **4D-QSAR analysis and pharmacophore modeling for alkynylphenoxyacetic acids as CRTh2 (DP2) receptor antagonists**
Köprü S., Sarıpinar E.
TURKISH JOURNAL OF CHEMISTRY, vol.42, pp.1577-1597, 2018 (SCI-Expanded)
- XXIV. **Synthesis and Single Crystal X-Ray Structure of 3,5-Diphenyl-1-(phenylacetyl)-4,5-dihydro-1H-pyrazol**
Kansız S., Tinmaz F., İLHAN I. Ö., SARIPINAR E., Dege N.
CRYSTALLOGRAPHY REPORTS, vol.63, no.6, pp.937-941, 2018 (SCI-Expanded)
- XXV. **Pharmacophore Modelling and 4D-QSAR Study of Ruthenium(II) Arene Complexes as Anticancer Agents (Inhibitors) by Electron Conformational-Genetic Algorithm Method**
Yavuz S., Sabancı N., SARIPINAR E.
CURRENT COMPUTER-AIDED DRUG DESIGN, vol.14, no.1, pp.79-94, 2018 (SCI-Expanded)
- XXVI. **4D-QSAR Study of Some Pyrazole Pyridine Carboxylic Acid Derivatives By Electron Conformational-Genetic Algorithm Method**
Tuzun B., YAVUZ S. C., Sabancı N., SARIPINAR E.
CURRENT COMPUTER-AIDED DRUG DESIGN, vol.14, no.4, pp.370-384, 2018 (SCI-Expanded)
- XXVII. **Crystal and Molecular Structure of 3,5-Diphenyl-4,5-dihydro-2-phenylcarboxamide-1H-pyrazole**
Tanak H., İLHAN I. Ö., Dege N., Akin N., SARIPINAR E.
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- XXVIII. **4D-QSAR investigation and pharmacophore identification of pyrrolo[2,1-c][1,4]benzodiazepines using electron conformational-genetic algorithm method**
Ozalp A., Yavuz S., Sabancı N., Copur F., Kökbudak Z., Sarıpinar E.
SAR AND QSAR IN ENVIRONMENTAL RESEARCH, vol.27, pp.317-342, 2016 (SCI-Expanded)
- XXIX. **Crystal structure of 1-acetyl-3-(4-methylphenyl)-5-phenyl-4,5-dihydro-1H-pyrazole**
Bulbul H., TINMAZ F., Dege N., İLHAN I. Ö., SARIPINAR E.
CRYSTALLOGRAPHY REPORTS, vol.59, no.7, pp.1071-1073, 2014 (SCI-Expanded)
- XXX. **Synthesis, spectroscopic (FT-IR/NMR) characterization, X-ray structure and DFT studies on (E)-2-(1-phenylethylidene) hydrazinecarboximidamide nitrate hemimethanol**
Ozdemir N., Inkaya E., SARIPINAR E., Akyuz L., İLHAN I. Ö., Aydin S., Dincer M., Buyukgungor O.
SPECTROCHIMICA ACTA PART A-MOLECULAR AND BIOMOLECULAR SPECTROSCOPY, vol.114, pp.175-182, 2013 (SCI-Expanded)
- XXXI. **Conformation depends on 4D-QSAR analysis using EC-GA method: pharmacophore identification and bioactivity prediction of TIBOs as non-nucleoside reverse transcriptase inhibitors**
Akyuz L., SARIPINAR E.
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- XXXII. **4D-QSAR study of HEPT derivatives by electron conformational-genetic algorithm method**
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- XXXIII. **Application of electron conformational-genetic algorithm approach to 1,4-dihydropyridines as calcium channel antagonists: pharmacophore identification and bioactivity prediction**
Gecen N., SARIPINAR E., Yanmaz E., Sahin K.
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- XXXIV. **The Reactions of Some Pyrazole-3-carboxylic Acid and Carboxylic Acid Chlorides with Various Alcohols**
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- XXXV. **Reactions of Some Pyrazole-3-Carboxylic Acid and Carboxylic Acid Chlorides with Various Alcohols**
İLHAN I. Ö., Bahadir O., Onal Z., SARIPINAR E.
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- XXXVI. **4D-QSAR analysis and pharmacophore modeling: Electron conformational-genetic algorithm approach for penicillins**

- Yanmaz E, SARIPINAR E, Sahin K, Gecen N, Copur F.
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- XXXVII. Quantitative bioactivity prediction and pharmacophore identification for benzotriazine derivatives using the electron conformational-genetic algorithm in QSAR
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- XXXVIII. Pharmacophore identification and bioactivity prediction for triaminotriazine derivatives by electron conformational-genetic algorithm QSAR method
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- XXXIX. Synthesis and Reactions of Some 1H-Pyrazole-3 Carboxylic Acid Chloride"
İlhan İ. Ö., Zuhal S., Kökbudak Z., Sarıpinar E.
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- XL. SYNTHESIS AND REACTIONS OF SOME 1H-PYRAZOLE-3 CARBOXYLIC ACID CHLORIDE
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- XLI. New beta-tricarbonyl compounds: Synthesis, reactions with urea and some thioureas
Sacmacı M., Alkan A., SAÇMACI Ş., SARIPINAR E., ŞAHİN E.
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- XLII. A convenient synthesis of functionalized 1H-pyrimidine-2-ones/thiones, pyridazine and imidazole; Experimental data and PM3 calculations
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HETEROCYCLES, vol.68, no.10, pp.2045-2062, 2006 (SCI-Expanded)
- XLIII. 2-(1-phenylethylideneamino) guanidine
DEMİR S., Dincer M., Sarıpinar E.
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- XLIV. A new method for the preparation of pyridazine systems: Experimental data and semiempirical PM3 calculations
Ünal D., Sarıpinar E., Akcamur Y.
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DINCER M., OZDEMIR N., Sarıpinar E., KULAK L., BUYUKGUNGOR O.
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- XLVI. Reaction of some furan-2,3-diones with various 1,2-phenylenediamines
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- XLVII. Synthesis and thermolysis of the 2,3-dihydro-1H-pyrole-2,3-diones, pseudopericyclic reactions of formyl(N-phenylimidoyl)ketene: Experimental data and PM3 calculations.
Sarıpinar E., KARATAS S.
JOURNAL OF HETEROCYCLIC CHEMISTRY, vol.42, no.5, pp.787-796, 2005 (SCI-Expanded)
- XLVIII. 5-(4-methoxybenzoyl)-4-(4-methoxyphenyl)-1-[1-(4-methoxyphenyl)ethylidenenamino]pyrimidin-2(1H)-one
OZCELIK S., DINCER M., Sarıpinar E., YILMAZ C.
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- XLIX. Synthesis of some pyrazole-3-carboxylic acid-hydrazide and pyrazolopyridazine compounds
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JOURNAL OF HETEROCYCLIC CHEMISTRY, vol.42, no.1, pp.117-120, 2005 (SCI-Expanded)
- L. Synthesis of novel proponohydrazides and their hydrolysis reactions
SACMACI M., Sarıpinar E., AKCAMUR Y.
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- LI. Reactions of cyclic oxaryl compounds, synthesis of functionalized hydrazono-2H-imidazol-4,5(1H,

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Saripinar E., AKCAMUR Y., KOLLENZ G.
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- LII. Double Mitsunobu reactions of cis-cycloalk-2-ene-1,4-diols and 3,4-epoxycycloalkenes: Rearrangements of allylic diazides**
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- LIII. Reactions of cyclic oxaryl compounds, reaction of some furan-2,3-diones with various hydrazine derivatives**
ILHAN I., AKCAMUR Y., Saripinar E., ASLAN E.
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