



جامعة الكرخ للعلوم
كلية علوم الطاقة والبيئة
السيرة الذاتية العلمية

المعلومات الشخصية

الإسم الثلاثي واللقب	خطاب عدنان عبد الخفاجي
المواليد	١٩٨٧١٠٧٦١١
الحالة الاجتماعية	متزوج
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التخصص العام	علوم كيمياء
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سنة الحصول على الشهادة	القسم / الكلية	الجامعة المانحة للشهادة	الشهادة الحاصل عليها
٢٠٢١	كلية العلوم اقسام الكيمياء	جامعة غازي عينتاب	دكتوراه

السلك الوظيفي

سنة اشغال المكان الوظيفي	المكان الوظيفي	العنوان الوظيفي
٢٠٢٣	جامعة الكرخ للعلوم	مدرس دكتور

المشروعات البحثية التخصصية لخدمة البيئة والمجتمع أو تطوير التعليم

اسم البحث	كلمات مفتاحية عن البحث	تاريخ النشر

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المهام التدريسية

العام الدراسي	المرحلة الدراسية	التخصص العام	اسم المادة
٢٠٢٤-٢٠٢٣	المستوى الثاني	كيمياء	كيمياء حياتية
٢٠٢٣-٢٠٢٢	المستوى الثالث	علوم الحياة	تصنيف حيوي
٢٠٢٤-٢٠٢٣	المستوى الثاني	لغة انكليزي	لغة انكليزية ٢

النشاطات والمهام العلمية

البحوث المنشورة

Yagiz G., Noma S.A.A., Altundas A., Al-khafaji K., Taskin-Tok T., Ates B.
 AUTHOR FULL NAMES: Yagiz, Güler (57458840600); Noma, Samir Abbas Ali (57202381427); Altundas, Aliye (6508047760); Al-khafaji, Khattab (57215680494); Taskin-Tok, Tugba (56268532500); Ates, Burhan(٦٥٠٧٥٠٤١٢٣)
 57458840600٦٥٠٧٥٠٤١٢٣ ; ٥٦٢٦٨٥٣٢٥٠٠ ; ٥٧٢١٥٦٨٠٤٩٤ ; ٦٥٠٨٠٤٧٧٦٠ ; ٥٧٢٠٢٣٨١٤٢٧ ;
 Synthesis, inhibition properties against xanthine oxidase and molecular docking studies of dimethyl N-benzyl-1H-1,2,3-triazole-4,5-dicarboxylate and (N-benzyl-1H-1,2,3-triazole-4,5-diyl)dimethanol derivatives



, ١٠٨ , (٢٠٢١) art. no. 104654

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Salman H.A., Hussin A.M., Hassan A.H., Al Sabbah H., Al-Khafaji K.

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56711021200٥٧٢١٥٦٨٠٤٩٤ ; ١٧١٣٤١٩٣٣٠٠ ; ٥٨٦٤٠٣٤٦٦٠٠ ; ٥٧٣٧٣٥٢٠٩٠٠ ;

The prevalence of potential side effects of COVID-19 vaccines among vaccinated Iraqi people:
a prospective cross-sectional study

(٢٠٢٢)

DOI: 10.1108/AGJSR-04-2023-0144

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Combination of QSAR, molecular docking, molecular dynamic simulation and MM-PBSA: analogues of lopinavir and favipiravir as potential drug candidates against COVID-19

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DOI: 10.1080/07391102.2020.1850355

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Saleh, Mohanad Yakdhan (57220102602); Abdelhakem, Adel Mohamed (57220106211); Alanazi, Amer M.

(54942111300); Islam, Md Ataul(٧٤٠٣٦٩٣٢٧٩)

٧٤٠٣٦٩٣٢٧٩ ; ٥٤٩٤٢١١١٣٠٠ ; ٥٧٢٢٠١٠٦٢١١ ; ٥٧٢٢٠١٠٢٦٠٢ ; ٥٧٢١٥٦٨٠٤٩٤ ; ٥٧٢١٥٠٦١٠٩٤

Identification of naphthyridine and quinoline derivatives as potential Nsp16-Nsp10 inhibitors: a pharmacoinformatics study

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٥٨٢٦٥٨٧٦٤٠٠ ; ٥٧٢٠١٠٦٨٠٧١ ; ٥٧٢٢١٧١٨٥٤٩ ; ١٦١٧٥٢٦٩٢٠٠ ; ٥٧٢١٥٦٨٠٤٩٤ ; ٥٧٢٢٠١١٥٢٧١



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,(١) ١٢ ,(٢٠٢٣)art. no. 21

DOI: 10.1007/s13721-023-00416-3

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Understanding the mechanism of amygdalin's multifunctional anti-cancer action using computational approach

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٥٦٢٦٨٥٣٢٥٠٠ ;٥٧٢١٥٦٨٠٤٩٤

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٥٥٩٩١٧٩٢١٠٠

Antiemetic effects of sclareol, possibly through 5-HT3 and D2 receptor interaction pathways: In-vivo and in-silico studies

,١٨١ ,(٢٠٢٣)art. no. 114068

DOI: 10.1016/j.fct.2023.114068

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٤٢٠٦٠٩٥٦٤٠٠ ; ٥٧٢١٥٦٨٠٤٩٤ ; ٥٧٢٢٠٢٠٧٧٦١ ; ٥٧١٨٨٨٢٧٤٢٧

Synthesis, anti-inflammatory effects, molecular docking and molecular dynamics studies of 4-hydroxy coumarin derivatives as inhibitors of COX-II enzyme

, ١٢٤٧ , (٢٠٢٢) art. no. 131377

DOI: 10.1016/j.molstruc.2021.131377

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٥٦٢٦٨٥٣٢٥٠٠ ; ٥٧٢١٥٦٨٠٤٩٤

Computational techniques for studying protein-protein interactions

, (٢٠٢٢) pp. 125 - 135

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Investigation of berberine and its derivatives in Sars Cov-2 main protease structure by molecular docking, PROTOX-II and ADMET methods: in machine learning and in silico study

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DOI: 10.1080/07391102.2022.2142848

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Bioinformatic and computational analysis for predominant mutations of the Nrf2/Keap1 complex in pediatric leukemia

, (١٢) ٣٩ , (٢٠٢١) pp. 4290 - 4303

DOI: 10.1080/07391102.2020.1775702

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Anxiolytic-like Effect of Quercetin Possibly through GABA Receptor Interaction Pathway: In Vivo and In Silico Studies

, (٢١) ٢٧ , (٢٠٢٢) art. no. 7149

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Al-Khafaji K., AL-Duhaidahawi D., Taskin Tok T.

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٥٦٢٦٨٥٣٢٥٠٠ ; ٥٧١٨٨٨٢٧٤٢٧ ; ٥٧٢١٥٦٨٠٤٩٤

Using integrated computational approaches to identify safe and rapid treatment for SARS-CoV-2

, (٩) ٣٩ , (٢٠٢١) pp. 3387 - 3395

DOI: 10.1080/07391102.2020.1764392

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57323075000 ; ٥٧٣١٥٤٥٦١٠٠ ; ٥٧٤٦٨٢٤٠٧٠٠ ; ٥٧٩٧٠٤٨٧٧٠٠ ; ٥٧٤٦٧٨٧٧٤٠٠ ; ٥٧٢١٥٦٨٠٤٩٤ ;
٥٧٣١٨٥٨٠٦٠٠ ; ٥٧٢٢٣٢١٠١٧٠ ; ٥٧٨١٤٤٠٣٦٠٠ ; ٥٥٣٩٩٠٢٩٨٠٠ ; ٥٨٠٢٣٥١٤١٠٠ ; ٥٨٠٧٤٣٨٩٢٠٠
٥٧٢١٨٢٢٧٣٩٦ ; ٥٧٢٢٣٩٦٤١٨٩ ; ٥٨٥٩٧١٨٧١٠٠ ; ٥٧٢٢٥٣٧٥٨٦ ; ٥٥٨٤٠٧٩٨٦٠٠ ; ٥٧١٩٤٤٥٤٣٦٣
٥٨٤٨٣٥١٦٩٠٠ ; ٥٥٤٥٧٩٦١٠٠ ; ١٦٤٤٤٠١٦٧٠٠ ; ٢٤٧٦٤٧٨٣٦٠٠

Recognizing novel drugs against Keap1 in Alzheimer's disease using machine learning grounded computational studies

, ١٥ , (٢٠٢٢) art. no. 1036552

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Hossain R., Al-Khafaji K., Khan R.A., Sarkar C., Shahazul Islam M.D., Dey D., Jain D., Faria F., Akbor R., Atolani O., Oliveira S.M.R., Siyadatpanah A., De Lourdes Pereira M., Islam M.T.

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; ٥٧٢٢٦٥٦٧٠٥٨ ; ٥٧٢٢٦٥٦٩٢٢٤ ; ٥٧٢١٦٣٥٠١٢٠ ; ٥٧٢٢٦٥٦٤٦٨ ; ٥٧٢١٥٦٨٠٤٩٤ ; ٥٧٢١٧٣٣٥٩٨٠
؛ ٥٥٦٤٦١٢٢١٠٠ ; ٥٧٢٢١٦٨٥٨٢٠ ; ٣٥٧٤٨٦٦٦٤٠٠ ; ٥٧٢٢٦٥٦٢٨٦٠ ; ٥٧٢١٩٨٥٠٩٦٢ ; ٥٧٢١٣٠٦٢٥٢٩
٥٥٩٩١٧٩٢١٠٠ ; ٣٥٥١٧٩٨٨٩٠٠

Quercetin and/or ascorbic acid modulatory effect on phenobarbital-induced sleeping mice possibly through gabaa and gabab receptor interaction pathway

, (٨) ١٤ , (٢٠٢١) art. no. 721

DOI: 10.3390/ph14080721

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GC-MS analysis, and evaluation of protective effect of Piper chaba stem bark against paracetamol-induced liver damage in Sprague-Dawley rats: Possible defensive mechanism by targeting CYP2E1 enzyme through in silico study



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Computer-based identification of potential compounds from *Salviae miltiorrhizae* against Neirisaral adhesion A regulatory protein

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Superior photo-induced antibacterial/antibiofilm activities of ZnPcs/TiO₂ and computational simulation studies

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Phytochemical constituents of *Inula britannica* as potential inhibitors of dihydrofolate reductase: A strategic approach against shigellosis

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DOI: 10.1016/j cmpb.2020.105660

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Target prediction, computational identification, and network-based pharmacology of most potential phytoconstituent in medicinal leaves of *Justicia adhatoda* against SARS-CoV-2

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New tamoxifen analogs for breast cancer therapy: synthesis, aromatase inhibition and in silico studies

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Diterpenes/diterpenoids and their derivatives as potential bioactive leads against dengue virus: A computational and network pharmacology study

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The (NHC)PdBr₂(2-aminopyridine) complexes: synthesis, characterization, molecular docking study, and inhibitor effects on the human serum carbonic anhydrase and serum bovine xanthine oxidase

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Nature-Derived Compounds as Potential Bioactive Leads against CDK9-Induced Cancer: Computational and Network Pharmacology Approaches

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DOI: 10.3390/pr10122512

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A detailed understanding of the COL10A1 and SOX9 genes interaction based on potentially damaging mutations in gastric cancer using computational techniques

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الجان والتکلیفات

اللجنة / التکلیف	امر اللجنة / التکلیف	الرقم الأمر	السنة
العمل في شعبة التسجيل	تكليف	٦٤٢/٩ م/ع	٢٠٢٣
لجنة امتحانية في قسم علوم البيئة	لجنة		٢٠٢٣
لجنة ضمان الجودة ومقاييس المختبرات في قسم علوم البيئة	لجنة		٢٠٢٣
لجنة علمية في قسم علوم البيئة	لجنة		٢٠٢٣

المؤتمرات والدورات العلمية

اساسيات في المعلوماتي الحيوية- دورة

براءات الاختراع

العدد - التاريخ	طلب تسجيل الاختراع	الادعاء	الموضوع/ الفكرة

تقييم براءات الاختراع

الموضوع/ الفكرة	الادعاء	طلب تسجيل الاختراع	التاريخ

تقييم البحث

عنوان البحث	المجلة	العدد	التاريخ
Antitumor activity against human promyelocytic leukemia and in silico studies of some benzoxazines	Journal of Biomolecular Structure and Dynamics		٣١-٠٨-٢٠٢٢
Target SARS-CoV-2: Theoretical	Journal of		٢٧-١١-٢٠٢٠

		Biomolecular Structure and Dynamics	Exploration On Clinical Suitability of Certain Drugs

كتب الشكر والتقدير

التعهد الإلكتروني

اني عضو الهيئة التدريسية (خلطاب عدنان عبد) أتعهد بصحة المعلومات المدرجة اعلاه
واتحمل كافة التبعات القانونية في حالة مخالفة المعلومات التي ستدرج في الموقع الالكتروني الخاص بجامعتنا.

نعم , اوافق على التعهد اعلاه

ملا ، لا اوافق على التعهد اعلاه

مصادقة السيد رئيس الجامعة



ملاحظة : في حالة الموافقة على التعهد الالكتروني اعلاه , يرجى اختيار اللون الابيض في الخانة المرافقة ادنى التعهد " كلا , لا اوافق على التعهد اعلاه" والعكس صحيح.