

CAS
SciFinderⁿ
Quick Reference Guide

Contents

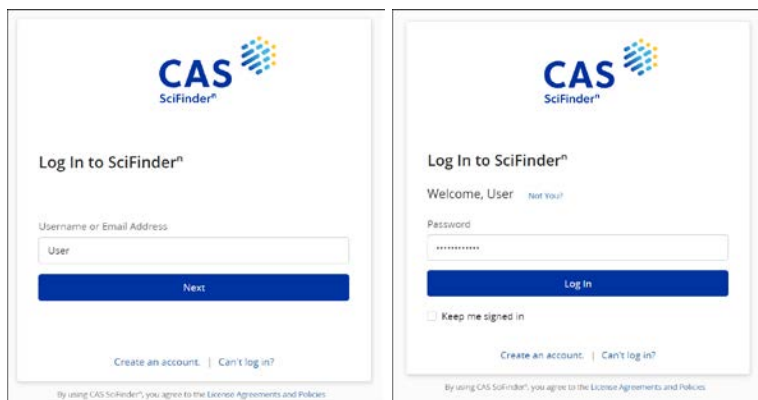
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Welcome to CAS SciFinder[®]

This Quick Reference Guide will show you how to start using CAS SciFinder[®], the industry's most trusted and comprehensive chemistry relevance engine.

First, open the CAS SciFinder[®] Login page: <https://scifinder-n.cas.org>.

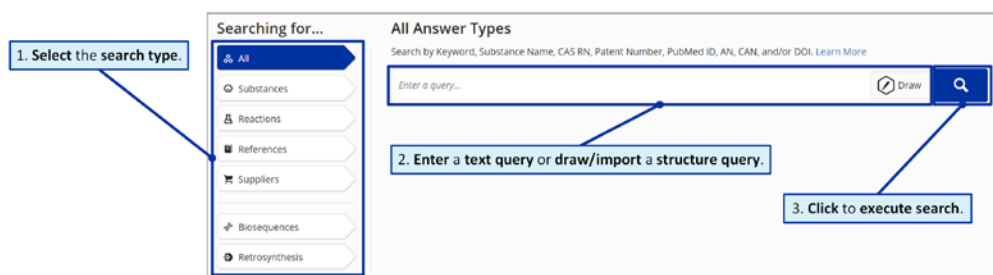
Log in using your CAS SciFinder[®] **Username** and **Password**.



Search

Search for the result type you need using a keyword, substance name, CAS Registry Number[®], patent number, or structure.

Note: You may enter a document object identifier (DOI) in the **All** and **References** searches.



Using **Advanced Search** for **References** and **Substances**, you may search by specific information type (e.g., author name or substance property).

Patent Markush Search: To conduct a patent markush search, select **Substances**, draw/import the query using the Structure Editor, and then check the box for **Search Patent Markush**.



Substance Results

Filter by structure match.

Retrieve related data for all results.

Download results.

Save results/search, create alert.

Email results.

Click the X to remove a filter or Clear All Filters. Multiple filters in dropdown menu.

Clear All Filters

Structure Match

- As Drawn (13)
- Substructure (36)
- Similarity (123K)
- Analyze Structure Precision
- Chemscapc Analysis

Filtering: Reference Role: 2 Selected X Commercial Availability: Available X

2 Selected 5 Results

Sort: Relevance **View: Full**

Keep or remove selected results.

Change result display.

Sort results by relevance, CAS RN, Molecular Formula or Weight, and Number of References or Suppliers.

Click to display data visualization.

Create Chemscapc Analysis

Filter Behavior

Filter by Exclude

- Reaction Role
- Reference Role
- Commercial Availability
- Number of Components
- Molecular Weight
- Stereochemistry
- Substance Class
- Isotopes
- Metals
- Experimental Property
- Experimental Spectrum
- Bioactivity Indicator
- Target Indicator
- Regulatory Data by List
- Search Within Results

Select filters to focus results.

Retrieve related data for a specific result.

Click property name to view more information on Substance Detail.

Click to view substance information.

Click to select result.

Click to open Substance Detail.

Click to download the contents (name and number value) for all or applied filters as an .xlsx file.

Download filter data from this result set.

51234-28-7

70280-67-0

66934-19-8

70062-36-1

Key Physical Properties

Key Physical Properties	Value	Condition
Molecular Weight	301.72	-
Boiling Point (Predicted)	446.2±30.0 °C	Press: 760 Torr
Density (Predicted)	1.362±0.06 g/cm ³	Temp: 20 °C; Press: 760 Torr
pKa (Predicted)	4.36±0.30	Most Acidic Temp: 25 °C

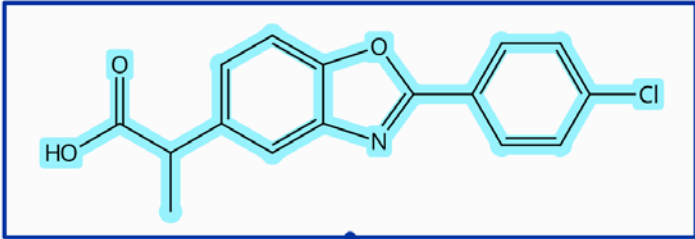
Chemical Structure: C1=CC=C(C=C1)C2=CC=C(C=C2)C3=CC=C(C=C3)C4=CC=C(C=C4)C5=CC=C(C=C5)C6=CC=C(C=C6)C7=CC=C(C=C7)C8=CC=C(C=C8)C9=CC=C(C=C9)C10=CC=C(C=C10)C11=CC=C(C=C11)C12=CC=C(C=C12)C13=CC=C(C=C13)C14=CC=C(C=C14)C15=CC=C(C=C15)C16=CC=C(C=C16)C17=CC=C(C=C17)C18=CC=C(C=C18)C19=CC=C(C=C19)C20=CC=C(C=C20)C21=CC=C(C=C21)C22=CC=C(C=C22)C23=CC=C(C=C23)C24=CC=C(C=C24)C25=CC=C(C=C25)C26=CC=C(C=C26)C27=CC=C(C=C27)C28=CC=C(C=C28)C29=CC=C(C=C29)C30=CC=C(C=C30)C31=CC=C(C=C31)C32=CC=C(C=C32)C33=CC=C(C=C33)C34=CC=C(C=C34)C35=CC=C(C=C35)C36=CC=C(C=C36)C37=CC=C(C=C37)C38=CC=C(C=C38)C39=CC=C(C=C39)C40=CC=C(C=C40)C41=CC=C(C=C41)C42=CC=C(C=C42)C43=CC=C(C=C43)C44=CC=C(C=C44)C45=CC=C(C=C45)C46=CC=C(C=C46)C47=CC=C(C=C47)C48=CC=C(C=C48)C49=CC=C(C=C49)C50=CC=C(C=C50)C51=CC=C(C=C51)C52=CC=C(C=C52)C53=CC=C(C=C53)C54=CC=C(C=C54)C55=CC=C(C=C55)C56=CC=C(C=C56)C57=CC=C(C=C57)C58=CC=C(C=C58)C59=CC=C(C=C59)C60=CC=C(C=C60)C61=CC=C(C=C61)C62=CC=C(C=C62)C63=CC=C(C=C63)C64=CC=C(C=C64)C65=CC=C(C=C65)C66=CC=C(C=C66)C67=CC=C(C=C67)C68=CC=C(C=C68)C69=CC=C(C=C69)C70=CC=C(C=C70)C71=CC=C(C=C71)C72=CC=C(C=C72)C73=CC=C(C=C73)C74=CC=C(C=C74)C75=CC=C(C=C75)C76=CC=C(C=C76)C77=CC=C(C=C77)C78=CC=C(C=C78)C79=CC=C(C=C79)C80=CC=C(C=C80)C81=CC=C(C=C81)C82=CC=C(C=C82)C83=CC=C(C=C83)C84=CC=C(C=C84)C85=CC=C(C=C85)C86=CC=C(C=C86)C87=CC=C(C=C87)C88=CC=C(C=C88)C89=CC=C(C=C89)C90=CC=C(C=C90)C91=CC=C(C=C91)C92=CC=C(C=C92)C93=CC=C(C=C93)C94=CC=C(C=C94)C95=CC=C(C=C95)C96=CC=C(C=C96)C97=CC=C(C=C97)C98=CC=C(C=C98)C99=CC=C(C=C99)C100=CC=C(C=C100)

Substance Detail

Substance Detail

References (944)
Reactions (36)
Suppliers (35)

CAS Registry Number
51234-28-7



C₁₆H₁₂ClNO₃
5-Benzoxazoleacetic acid, 2-(4-chlorophenyl)-α-methyl- (94)

Key Physical Properties	Value	Condition
Molecular Weight	301.72	-
Melting Point (Experimental)	189-190 °C	-
Boiling Point (Predicted)	446.2±30.0 °C	Press: 760 Torr
Density (Predicted)	1.362±0.06 g/cm ³	Temp: 20 °C; Press: 760 Torr
pKa (Predicted)	4.36±0.30	Most Acidic Temp: 25 °C

Save detail.

Download detail.

Email detail.

Expand All | Collapse All

- Other Names and Identifiers
- Experimental Properties
- Experimental Spectra
- Predicted Properties
- Predicted Spectra
- Bioactivity Indicators
- Target Indicators
- Regulatory Information

Regulatory List: EINECS, REACH, VNECI
 Confidential Business Information: Public

Expand All | Collapse All

 - Regulatory Synonyms (8)
 - Details by Country/International & Other Lists
- Additional Details

Reference Results

The screenshot displays the 'References' section of the CAS SciFinder interface. It features a left-hand sidebar for filtering, a main results area with a list of search results, and a top navigation bar. Blue callout boxes with arrows point to specific elements, providing instructions on how to use them.

Callouts and their locations:

- Retrieve related data for all results.** Points to the 'Substances', 'Reactions', and 'Citing' dropdown menus at the top.
- Download results.** Points to the download icon in the top right.
- Save results/search, create alert.** Points to the 'Save and Alert' icon in the top right.
- Email results.** Points to the email icon in the top right.
- Click the X to remove a filter or Clear All Filters. Multiple filters in dropdown menu.** Points to the 'X' icon on a filter dropdown and the 'Clear All Filters' button.
- Sort: Relevance** and **View: Partial Abstract** are also labeled.
- Select filters to focus results.** Points to the 'Filter Behavior' sidebar.
- Keep or remove selected results.** Points to the checkbox next to a result number.
- Change result display.** Points to the 'View: Partial Abstract' dropdown.
- Sort results by relevance, times cited, accession number, or publication date.** Points to the 'Sort: Relevance' dropdown.
- Retrieve related data for a specific result.** Points to the 'Substances (6)', 'Reactions (0)', and 'Citing (63)' dropdowns below a result.
- Click to select result.** Points to a checkbox next to a result number.
- Click to open Reference Detail.** Points to the 'Full Text' dropdown for a specific result.
- Click to download the contents (name and number value) for all or applied filters as an .xlsx file.** Points to the 'Download filter data from this result set.' button.
- Click to access full-text viewing options.** Points to the 'PatentPak' dropdown.
- Click to access patent information viewing options.** Points to the 'PatentPak' dropdown.

Reference Detail

Reference Detail Retrieve data related to reference. Save detail. Download detail. Email detail.

Substances (3) Reactions (0) Citing (6) Citation Map

PATENT

Patent Number: WO9834612
 Publication Date: 1998-08-13
 Application Number: WO1998-EP648
 Application Date: 1998-02-03
 Kind Code: A1
 Assignee: The Boots Company PLC, United Kingdom

Source: World Intellectual Property Organization
 CODEN: PIXXD2

Database Information: AN: 1998:548531, CAN: 129:180142, CPlus

Language: English

Pharmaceutical compositions containing **ibuprofen** and domperidone for the treatment of migraine
 By: Pankhania, Manendra Govind; Yurdakul, Saruhan
 Migraine, ibuprofen, or a salt thereof and a domperidone, or a salt thereof, containing a therapeutically effective amount of ibuprofen 60, domperidone 1.5, and stearic acid 0.6%.

Keywords: pharmaceutical tablet, **ibuprofen**, domperidone, migraine treatment

Patent Family

Patent	Language	Kind Code	PatentPak Options	Publication Date	Application Number	Application Date
WO9834612	English	A1	PDF PDF+ Viewer	1998-08-13	WO1998-EP648	1998-02-03
CA2279184	Undetermined	A1		1998-08-13	CA1998-2279184	1998-02-03
EP1017300	Undetermined	A1		2000-07-12	EP1998-906926	1998-02-03
HU225043	Hungarian	B1		2006-05-29	HU2000-1104	1998-02-03

View All

Priority Application

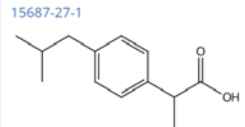
Priority Application Number	Application Date
GB1997-2392	1997-02-06
WO1998-EP648	

Expand All | Collapse All

IPC Data
 Concepts
 Substances

Substances (3)

15687-27-1

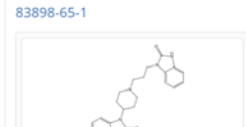


C₁₃H₁₈O₂
Ibuprofen

PatentPak

Role: Biological Activity or Effector, Except Adverse, Biological Study, Unclassified, Therapeutic Use, Biological Study, Uses

83898-65-1

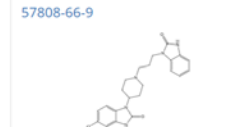


C₂₂H₂₄ClN₅O₂·C₄H₄O₄
Domperidone maleate

PatentPak

Role: Biological Activity or Effector, Except Adverse, Biological Study, Unclassified, Therapeutic Use, Biological Study, Uses

57808-66-9



C₂₂H₂₄ClN₅O₂
Domperidone

PatentPak

Role: Biological Activity or Effector, Except Adverse, Biological Study, Unclassified, Therapeutic Use, Biological Study, Uses

Formulations
 Cited Documents

View interactive version of patent that highlights specific locations of indexed substances.
 Set citing alert.
 View map of references this document cites and references that cite this document.
 Click to access full-text viewing options.
 Request a prior art analysis that contains both patent and non-patent literature.
 Click a PatentPak option view the patent source document.
 View IPC codes for basic patent and patent family members.
 View concepts that characterize the reference's general subject matter.
 View the reference's indexed substances.
 View the reference's formulation data.
 View the reference's citations.

Reaction Results

Retrieve related data for all results. (References dropdown)

Download results. (Download icon)

Save results/search, create alert. (Save and Alert icon)

Email results. (Email icon)

Click the X to remove a filter or Clear All Filters. Multiple filters in dropdown menu. (Filtering: Number of Steps: 2 X Experimental Protocols: Synthetic Methods X)

Clear All Filters. (Clear All Filters button)

1 Selected 29 Results

Keep or remove selected results. (1 Selected dropdown)

Group reactions by scheme or document. (Group: By Scheme dropdown)

Change result display. (View: Expanded dropdown)

Select filters to focus results. (Filter Behavior sidebar)

View substance vendors. (Suppliers (98) button)

Click for options to view details, generate retrosynthesis plan, and edit/download structure file. (Expand Scheme dropdown)

Click to open reaction reference's detail page. (Reaction Summary section)

Process for preparing varenicline, varenicline intermediates, and pharmaceutically acceptable salts thereof (Reaction Summary text)

Click to access full-text viewing options. (PatentPak Full Text dropdown)

Click to access patent information viewing options. (PatentPak dropdown)

View reaction's detail page. (View Reaction Detail button)

View experimental procedure for reaction. (Experimental Protocols button)

Click to download the contents (name and number value) for all or applied filters as an .xlsx file. (Download filter data icon)

Reaction Detail

Reaction Detail

Suppliers (19) Suppliers (78) Suppliers (40)

Download detail. Save detail. Email detail. View substance vendors. Click to open reaction reference's detail page.

Steps: 2

Suppliers (63)

View alternative reactions for the same product. Alternative Steps (7)

Click tabs to view steps in multi-step reactions.

Stage	Reagents	Catalysts	Solvents	Conditions
1	Sodium periodate	Benzyltriethylammonium chloride	Dichloromethane Water	10 °C; 1 h
2	-	-	-	1 min, 25 - 30 °C
3	Sodium cyanoborohydride	-	Dichloromethane	0 °C; 0 °C → 30 °C; 0.5 - 1 h
4	Sodium carbonate	-	Water	1 h, pH 9

CAS Reaction Number: 31-032-CAS-650575

Experimental Protocols

Synthetic Methods Experimental Procedure

Products: 2,3,4,5-Tetrahydro-3-(phenylmethyl)-1,5-methano-1H-3-benzazepine

Reactants: 1,2,3,4-Tetrahydro-1,4-methanonaphthalene-2,3-diol, Benzylamine

Reagents: Sodium periodate, Sodium cyanoborohydride, Sodium carbonate

Catalysts: Benzyltriethylammonium chloride

Solvents: Dichloromethane, Water

Procedure: 1. Stir a 1,2,3,4-tetrahydro-1,4-methanonaphthalene-2,3-diol (100 g) in a mixture of water (2600 ml) and methylene chloride (1040 ml) under nitrogen at 10 °C. 2. Add sodium periodate (127.6 g) and triethylbenzyl ammonium chloride (10 g) and stir the resulting mixture for 1 hour.

PATENT

Process for preparing varenicline, varenicline intermediates, and pharmaceutically acceptable salts thereof

By: Shekhawat, Kundan Singh; et al
View All

World Intellectual Property Organization

PatentPak Full Text

Click to access patent information viewing options.

Click to access full-text viewing options.

Patent Number: WO2010023561

Publication Date: 2010-03-04

Application Number: WO2009-1B7081

Application Date: 2009-08-31

Kind Code: A1

Assignee: Actavis Group PTC ehf, Iceland

Supplier Results

Suppliers

Click the X to remove a filter or Clear All Filters. Multiple filters in dropdown menu.

Download results. Email results.

Clear All Filters

Sort results by relevance, supplier name, shipping speed, or purity.

Sort: Relevance

Select filters to focus results.

Filter Behavior

Filter by Exclude

Preferred Suppliers

Supplier

Purity

Quantity

Ships Within

Stock Status

Order From Supplier

Country/Region

Filter Content Report

Download filter data from this result set.

Filtering: Purity: 95-98% X

1 Selected 21 Results

Supplier Substance Purity Purchasing Details Availability

1 ASTATECH 51234-28-7 BENOXAPROFEN 95-98% 0.1 g, USD 3500 Synthesis on demand

Click to open Substance Detail.

AstaTech Product List

United States

Last Updated: 31 Mar 2022

View details, generate retrosynthesis plan, and edit/download structure file.

Open product ordering page on supplier's website.

Order From Supplier

2 Arspichem Product List 51234-28-7 2-(2-(4-chlorophenyl)benzo[d]oxazol-5-yl)propanoic acid 95-98% Typically in stock

Open product information page on supplier's website.

Product Information

3 Alchem Pharmtech, Inc. 51234-28-7 2-(2-(4-chlorophenyl)benzo[d]oxazol-5-yl)propanoic acid 95-98% Maintained in stock

Click to select result.

Click thumbs up/down to set supplier preference.

Click to download the contents (name and number value) for all or applied filters as an .xlsx file.

Supplier Detail

Supplier Detail

Click thumbs up to set supplier as Preferred or thumbs down for Non-Preferred.

Download detail. Email detail.

Click to open Substance Detail.

Web <https://www.AstaTechInc.com>

Email sales@astatechinc.com

Phone 215-785-3197

Substance Information

CAS Registry Number 51234-28-7

CAS Name Benoxaprofen

Chemical Name BENOXAPROFEN

Order Number C90147

Purity 95%

Quantity, Price 0.1 g, USD 3500
0.25 g, USD 6900

Stock Status Synthesis on demand

Ships Within 8 weeks

Pricing Information 31 Mar 2022

Last Updated

Open product ordering page on supplier's website.

Order From Supplier

Additional Contact Information

AstaTech, Inc.
Keystone Business Park
2525 Pearl Buck Road
Bristol, PA, 19007
United States

Fax 215-785-2656

Click to view details, generate retrosynthesis plan, and edit/download structure file.

Biosequence Results

Biosequences

Retrieve related data for all results.

Download results.

Email results.

References

Click to select another sequence.

Sort results by Alignment Identity, E-Value, Query Coverage, or Subject Coverage.

Sort: Alignment Identity View: Expanded

Change result display.

Click to display data visualization.

Click to view references indexing the sequence.

Select filters to focus results.

Click to view CAS RN, NCBI Identifier, sequence length, NCBI Organism, and tabular sequence.

Click to patent/journal results containing a matching sequence.

BLAST Search Details

Sequence Type: Protein
Search Within: Nucleotides
BLAST Algorithm: TBLASTn-fast
NCBI Included: No
Alignment Identity: -
Query Coverage: 90%
E-Value: 10
Match with Gaps?: Yes
Gap Costs: Existence 11
Extension 1

Bioscape Analysis

Visually explore sequence similarity with a new tool. Learn more about Bioscape.

Create Bioscape Analysis

Filter by

E-Value
0 to 10⁶

Query Coverage %
0 to 100

Subject Coverage %
0 to 100

Alignment Identity %
60.5 to 100

Apply Reset Filters

Query Details View More

> pdb|1E57|C Chain C, Physalis Mottle Virus

83 Results

Alignment Identity: 100%

Matches: 188
Mismatches: 0

Query 1 188

Subject 1 1,255

View Less

Alignment Subject References

Alignment Data
BLAST Score: 889
E-Value: 3.1542e-118

Q 1 MDSSEVVK QASIPAPGSI LSPNTEQSP AIVLPQFEA ttfgtaetaa C
S 540 MDSSEVVK QASIPAPGSI LSPNTEQSP AIVLPQFEA TTFGTASTAA QVSLQTADPI TKLTAPYRHA 609
SPATPTQILR VYGGQSFVLG GATSAAKTIE VPLNLDSVNR 140
SPATP
SIQIS
S 680 MLKDSVITYD TPKLLAYSRA PTNPSKIPTA SIQISGRIRL SKPMLIAN 727

Alignment Identity: 100%

Matches: 188
Mismatches: 0

Query 1 188

Subject 1 6,673

View Less

Alignment Subject References

Alignment Data
BLAST Score: 889
E-Value: 1.71882e-108

Bioscape

Bioscape visualizes the similarity and patent landscape for a set of sequence results. The location of the sequence bar in the visualization corresponds to the similarity of the sequence to the query, and the height of the sequence bar corresponds to the number of patents in which the sequence has been published.

The screenshot shows the Bioscape interface with several callout boxes:

- Refine sequence result bars by similarity.** Points to the 'Sequence Similarity' slider, which ranges from 97% to 100%.
- Refine sequence result bars by patent keyword and simple legal status.** Points to the search filters on the left side of the interface.
- Change how sequence result bars display.** Points to a button with a bar chart icon at the top center.
- Multiple CAS RNs Associated with this Sequence** (with a close button 'X'). This panel includes:
 - If present, click to view substances.** Points to a 'View Substances' button.
 - Click to view relevant patents.** Points to a 'View Patents' button.
- Queried sequence.** Points to a specific vertical bar in the main visualization.
- Click a bar to view its patent count and sequence length.** Points to a bar in the visualization.
- Click the Select Sequence button and then click-and-drag to select multiple sequence results for viewing.** Points to a 'Select Sequence' button at the bottom.

At the bottom of the interface, it says 'Powered by patsnap' and 'Copyright © 2022 American Chemical Society. All Rights Reserved. | 京ICP备13047075号-3'. There are also links for 'Help', 'Contact Us', and 'Legal'.

Chemscape

Chemscape visualizes the similarity and patent landscape for a set of substance results. The location of the substance bar in the visualization corresponds to the similarity of the substance to the query and the height of the substance in the visualization corresponds to the number of patents in which the substance has been published.

The screenshot displays the Chemscape visualization in CAS SciFinder. The interface includes a top navigation bar with the CAS SciFinder logo and a search bar. Below the search bar, there are several callout boxes providing instructions:

- View and manage your saved Chemscape.** (Points to the top-left navigation icons)
- Group and refine Chemscape structures to show key information.** (Points to the top-left navigation icons)
- Add new structures to your Chemscape and indicate their position.** (Points to the top-left navigation icons)
- Refine your Chemscape by keyword or an exact match to a chemical structure.** (Points to the top-left navigation icons)
- Change how structure result bars display.** (Points to the 'Structural Similarity' slider)
- Click to open Substance Detail page.** (Points to the 'Olanarib' substance bar)
- Click to view relevant patents.** (Points to the 'Patent Count' field in the detail panel)
- Click a bar to view its structure and number of associated patents.** (Points to a bar in the visualization)
- Click to save your Chemscape for later access in MyChemscape.** (Points to the 'Save' button in the bottom right)
- Click the Select Structure button and then click-and-drag to select multiple structure results for viewing or a new Chemscape.** (Points to the 'Select Structure' button)

The visualization itself shows a 3D bar chart where the x-axis represents similarity to the queried sequence and the z-axis represents the number of patents. A 'Queried sequence' is highlighted in the center. The detail panel for Olanarib (763113-22-0) shows its chemical structure, molecular formula (C₂₄H₂₉FN₄O₃), SMILES string, and a list of synonyms including AZD 2281 and Lynparza.

Saved Searches and Results

Filter by

- Result Type
 - Reactions (4)
 - References (4)
 - Substances (2)
- Alerts
 - Unviewed (1)
 - Alerts Set (2)
 - No Alerts (8)
- Tags
 - Citing Alert (1)
 - For Review (1)
 - NSAID (1)

Combine Saved Results

Combine

Set/edit an alert on saved query or view and manage alert results.

Filter saved items.

Click to select item.

Delete selected item.

Email selected item.

Click to open Reference Detail.

View references citing the saved item.

Edit name.

Run a saved query.

Combine previously saved queries or selected results.

Create/add tags.

Click to remove tag.

Display saved result set.

View Saved

Rerun Search

View Citing

Alerts (10)

1 Selected

Ibuprofen pharma compositions

January 19, 2022, 1:13 PM

References

Citing Alert on: Dihydrocodeine/ibuprofen pharmaceutical compositions

Alerts Add Tags Citing Alert

Ibuprofen References

October 30, 2018, 10:53 AM

References ibuprofen

Alerts Add Tags For Review NSAID

Teflon Subtracted Intersected 306A and 8510273

October 1, 2018, 4:20 PM

Combined From Reaction Saved Items

Search History

Filter history by selected search types.

Filter by

- Result Type
 - All (4)
 - Biosequences (19)
 - Patent Markush (1)
 - Reactions (13)
 - References (64)
 - Retrosynthesis (16)
 - Substances (36)
 - Suppliers (8)
- Date

Start Date End Date

mm/dd/yyyy to mm/dd/yyyy

May, 2022

1 2 3 4 5 6 7

8 9 10 11 12 13 14

15 16 17 18 19

22 23 24 25 26

29 30 31

Display search history for a specified date range.

Your Search History

53 Searches

1 Selected

Delete searches.

May 2, 2022

Substances

1:47 PM

As Drawn (7)
Substructure (1,724)
Similarity (608K)

Rerun search to retrieve latest results.

Edit search and then rerun.

Rerun Search

Edit Search

March 28, 2022

Retrosynthesis

3:12 PM

Synthetic Depth: 3
Predicted Rules: Common
Break & Protect Bonds: No
Starting Material Cost Limit:
\$1,000.00/mol

Open retrosynthesis plan or edit options and rerun.

Open Plan

Edit Search

Complete

CAS SciFinder[®] Support

To access CAS SciFinder[®] in-application support, click the **Help** link at the bottom of any page or select **Help** from the **Account** menu.



For additional assistance using CAS SciFinder[®], please contact the **CAS Customer Center**:

- **Hours:** 8:00 a.m. to 6:00 p.m. EST Monday – Friday.
- **Phone:**
 - 1-800-753-4227 (North America)
 - +1-614-447-3700 (outside North America)
 - **Option 2:** General information or account-related questions
 - **Option 3:** Assistance with search strategies, database content, or using a product
 - **Option 4:** Technical assistance with software set up, installation, and configuration
- **Email:** help@cas.org
- **Web:** <https://www.cas.org/contact>

If desired, ask for a CAS SciFinder[®] Familiarization Training Session visit or online session.