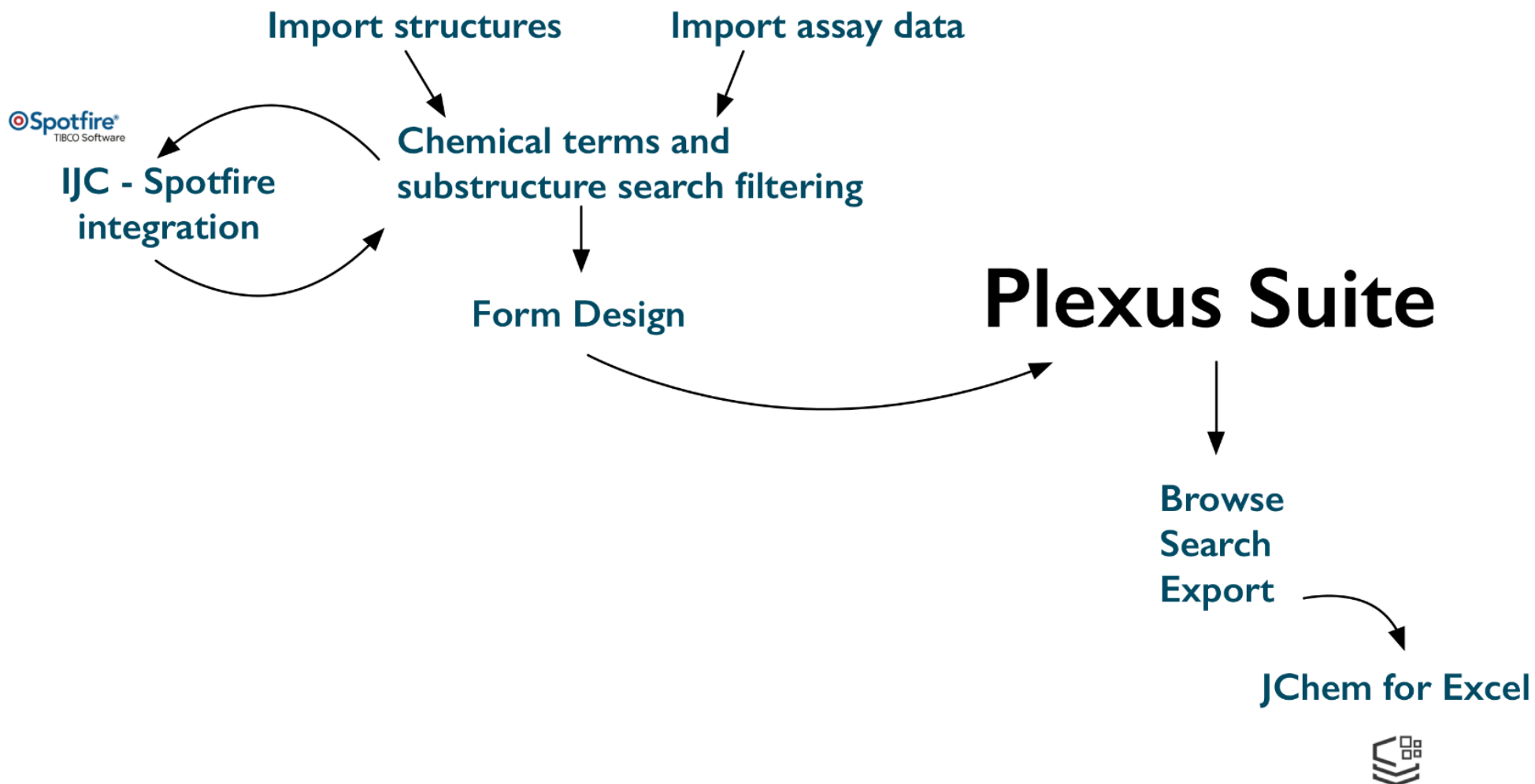


# INSTANT JCHEM (IJC)

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UGM Budapest 2015

# Typical Workflow Example





Projects [BI test] x Lists and queries

Dashboard

localdb [as admin]

- Structures
- Activities
- README.html

Query builder x

&lt;Not Queryable&gt;

## Forms

Search

Form	Data Tree	Schema	Project
Grid view for Structures	Structures	localdb	BI test
Grid view for Activities	Activities	localdb	BI test

## Quick Actions

Connect to Database...

Import File...

## Templates

Sample data

## Quick Start



## New features

[Export Wizard improvements](#)  
[Temporary Lists and queries improvements](#)  
[Improvements for In-list operator and calculated fields sorting](#)

New features in previous releases: [6.3](#), [6.2](#), [6.1](#)

## Documentation

[Release notes](#)  
[User's guide](#)  
[Developer's guide](#)  
[Admin's guide](#)  
[FAQ](#)

## Other links

[ChemAxon.com](#)  
[Query guide - search options](#)  
[Support forum](#)  
 Keyboard shortcuts for [Windows/Linux](#) and [Mac](#)  
[You should follow us on Twitter](#)  
[ChemAxon LinkedIn page](#)

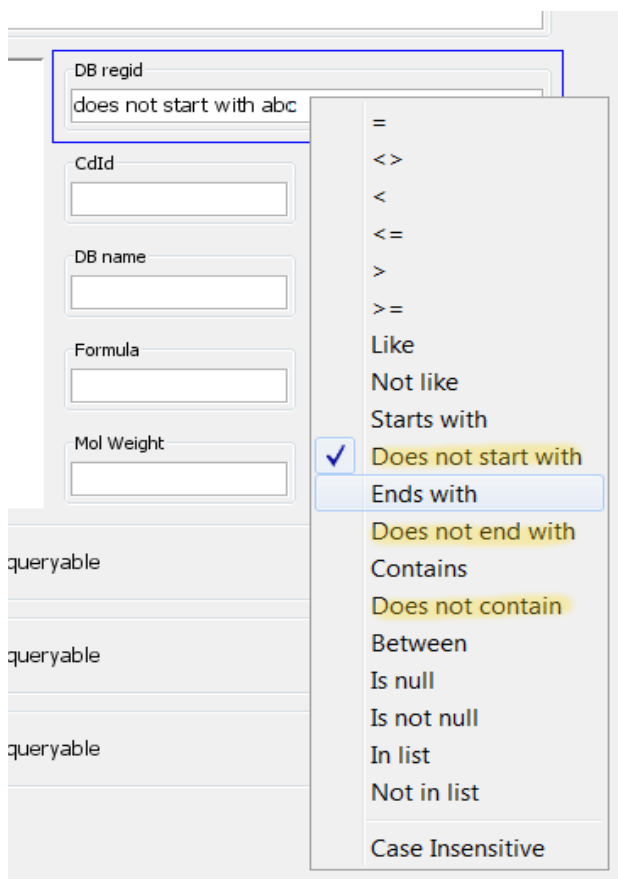
# Recent New Features



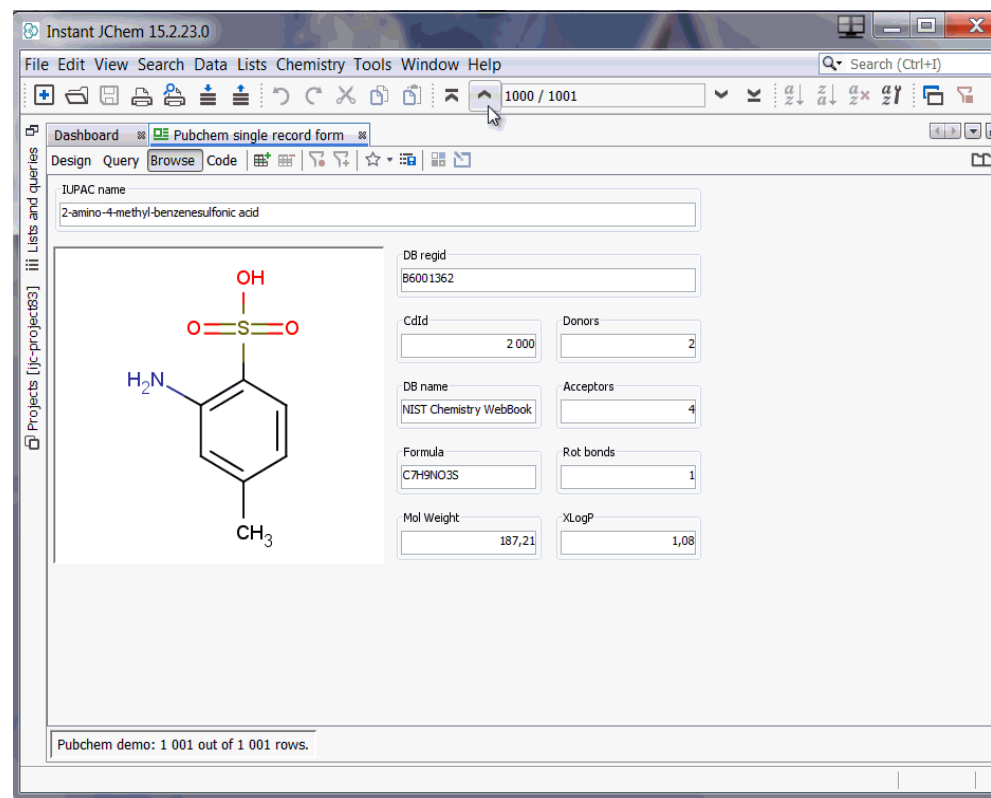
With the use of the newly developed Mass Spectra Display Widget, users can perform spectrum similarity searches based on a spectrum query, as well as search for spectra by structure or other chemical query terms.

# Recent new Features

## New Query Operators, aligned with Plexus Connect



## Improved "Add new row" dialog. It is possible to load values from existing rows.



# Recent New Features

The screenshot displays the Instant JChem 15.2.23.0 application window. The main content area shows a 'Pubchem single record form' for the compound (-)-acetylcarnitine. The interface includes a menu bar (File, Edit, View, Search, Data, Lists, Chemistry Tools, Window, Help) and a toolbar with various icons. The left sidebar shows a project list with 'Projects [ijc-project83]'. The main form contains the following fields:

- IUPAC name: (2-acetoxy-3-carboxylato-propyl)-trimethyl-ammonium
- DB regid: (-)-o-acetylcarnitine
- Cdid: 1 001
- Donors: 0
- DB name: BioCyc
- Acceptors: 4
- Formula: C9H17NO4
- Rot bonds: 5
- Mol Weight: 203,24
- XLogP: 3,41

Below the form, there are search links for Google, NCBI, and ChemSpider, all with the query '%28-%29o-acetylcarnitine'. At the bottom, a status bar indicates 'Pubchem demo: 1 000 out of 1 000 rows.' The background of the slide features a snowy landscape with evergreen trees.

Synchronize scripts between IJC schema and local disc folder.

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THANK YOU

