# FOR DISCUSSION: InChl JNA/WASM Java+JavaScript Joint Development Proposal

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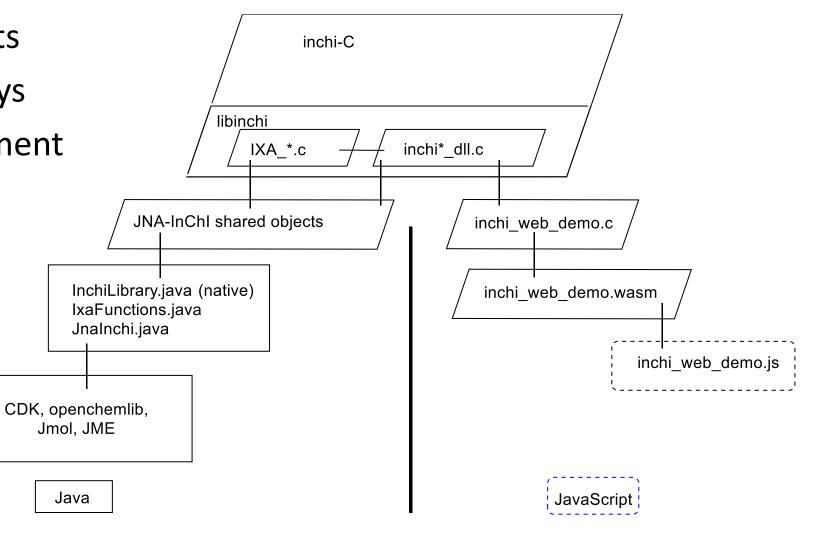
# Original configuration

Different starting points

Two divergent pathways

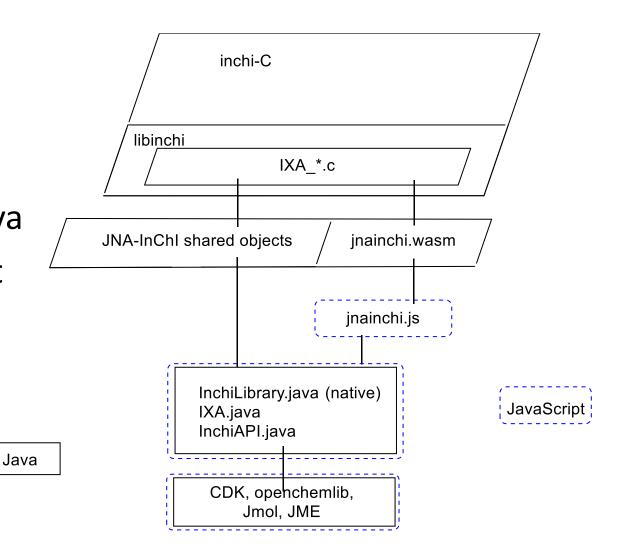
Independent development

Limited JavaScript dev



### Desired configuration

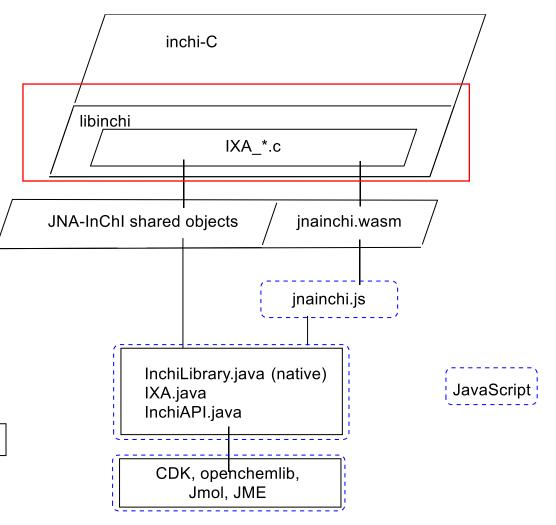
- Same starting point
- Common IXA-only interface
- Simple methods only
- No significant changes in Java
- Equivalent calls in JavaScript



No changes were required in inchi-C.

 Two methods were added to inchilib simply to round out the options and provide a built-in way to access the InChI version.
 These included IXA\_Builder\_GetInchiVersion and IXA\_MOL\_ReadAuxInfo

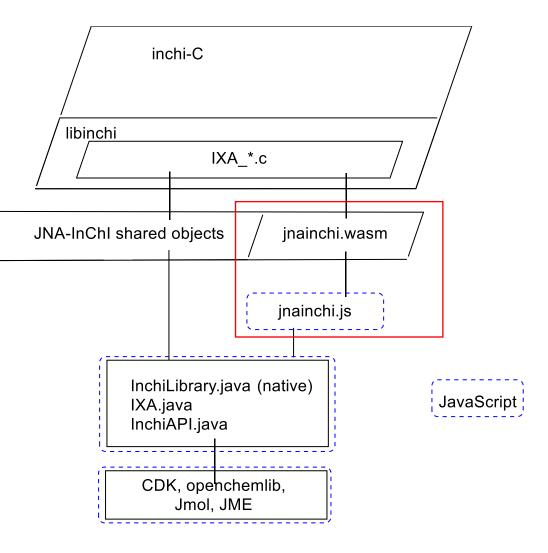
Java



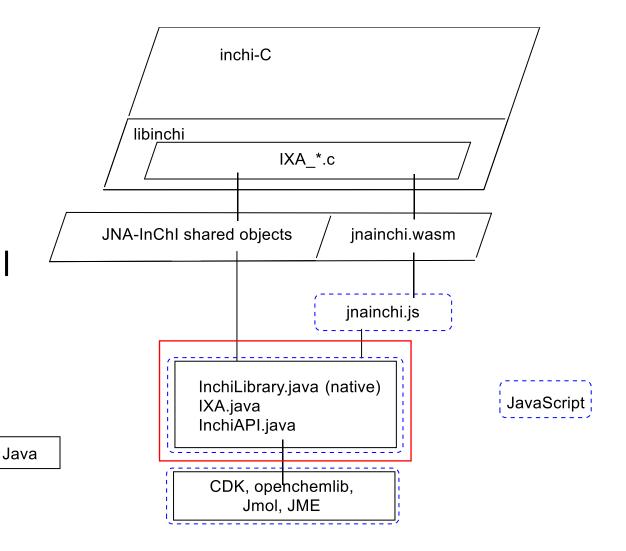
 WASM makefile adds IXA methods and renames output to match JNA native registration.

 Java2script Eclipse transpiler upgraded to implement "native"/ methods.

• SwingJS runtime adds Clazz.\_loadWasm, linking Java native methods with WASM exports when native methods\_are registered in JavaScript

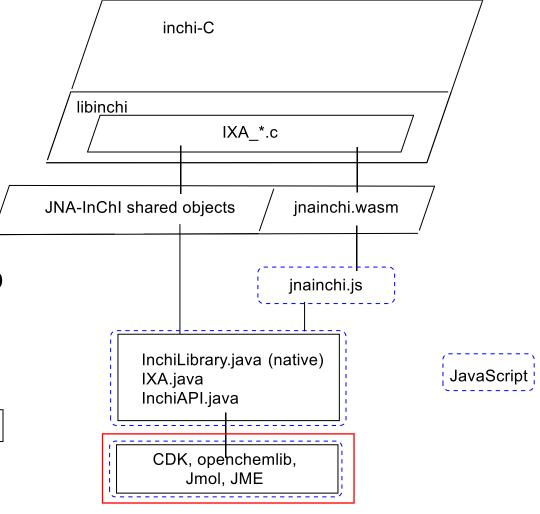


- JNA-InChI adds InchiAPI.java and IXA.java, avoiding all references to C structures.
- Adds additional InChI-to-Inchilnput flag "fixamide", which converts inchi-internal N=C-OH (iminol) to HN-C=O (amide) when creating an Inchilnput object from an InChI.



 CDK adds SwingJS-j2sExport class CDK.java as well as streamlined InchiGenerator.java and InchiToStructure.java classes functional in Java and JavaScript

 OCL configures inchiOCL class to handle all Java and JavaScript calls to InchiAPI

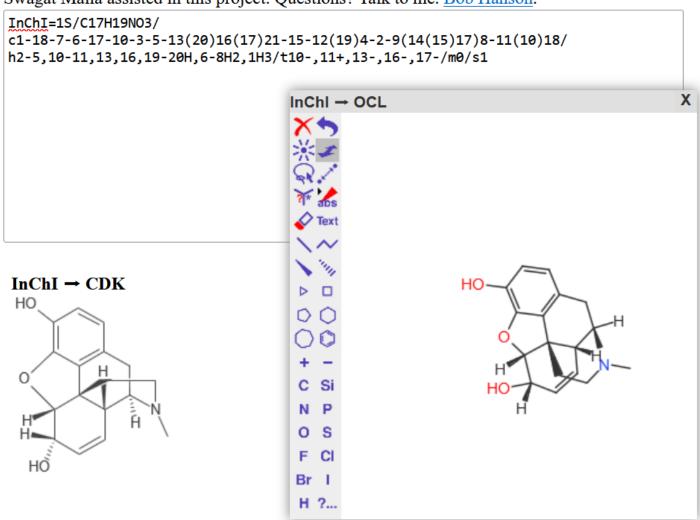


Java

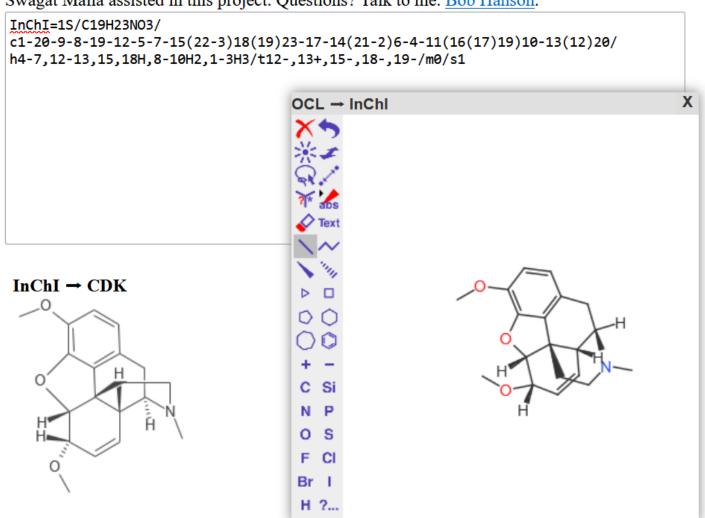
- Website demonstrating using both CDK and OpenChemLib in JavaScript for generating InChI from structures and structures from InChI.
- Includes InChI to structure editor, structure editor-to-InChI, CDX and CDXML file dropping, SMILES-to-InChI, MOL file-to-InChI, InChI to PNG, InChI-to-SVG, and OCL molecule-to-CDK molecule.

https://chemapps.stolaf.edu/inchi/cdk-web-demo/

To try it out, enter an InChI such as that for morphine or its enantiomer or this taxol MOL file (with amide fixing) or this AuxInfo or provide a SMILES. Go ahead and reverse the sign of one or more of the stereochemical indicators such as "13-" to "13+". Fun, huh? And, of course, you can drop your own MOL, CDX, or CDXML file into the OCL editor, or just modify the structure that is there. It should immediately update both the InChI and the CDK images. Many thanks to John Mayfield, Daniel Lowe, Jonathan Goodman, and others for assisting us in getting all these pieces working together. Interns Josh Charlton and Swagat Malla assisted in this project. Questions? Talk to me. Bob Hanson.



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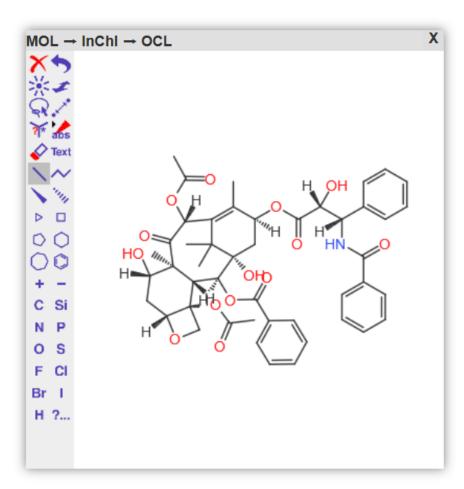


To try it out, enter an InChI such as that for <u>morphine</u> or <u>its enantiomer</u> or this <u>taxol MOL file (with amide fixing)</u> or this <u>AuxInfo</u> or provide a <u>SMILES</u>. Go ahead and reverse the sign of one or more of the stereochemical indicators such as "13-" to "13+". Fun, huh? And, of course, you can drop your own MOL, CDX, or CDXML file into the OCL editor, or just modify the structure that is there. It should immediately update both the InChI and the CDK images. Many thanks to John Mayfield, Daniel Lowe, Jonathan Goodman, and others for assisting us in getting all these pieces working together. Interns Josh Charlton and Swagat Malla assisted in this project. Questions? Talk to me. Bob Hanson.

InchI=1S/C47H51N014/
c1-25-31(60-43(56)36(52)35(28-16-10-7-11-17-28)48-41(54)29-18-12-8-13-19-29)23-47(5
7)40(61-42(55)30-20-14-9-15-21-30)38-45(6,32(51)22-33-46(38,24-58-33)62-27(3)50)39(
53)37(59-26(2)49)34(25)44(47,4)5/h7-21,31-33,35-38,40,51-52,57H,22-24H2,1-6H3,
(H,48,54)/t31-,32-,33+,35-,36+,37+,38-,40-,45+,46-,47+/m0/s1

#### MOL → InChI → MOL

#### $MOL \rightarrow InChI \rightarrow CDK$



# Coding Examples

The following examples use SwingJS. As such, the Java methods map directly to JavaScript functions. All of the functions **in bold** are written in Java. They are accessed in both Java and JavaScript. For JavaScript apps not using SwingJS, we would provide IXA calls using Emscripten methods *cwrap* and *ccall*. (And the apps would be responsible for anything else). But in SwingJS that is already taken care of by the JavaScript version of the native library registration process.

Classes swingjs.CDK and swingjs.OCL are imported in Java as CDK and OCL. In the JavaScript they are script-local references to var CDL = swingjs.CDK and var OCL = swingjs.OCL. These two classes provide equivalent calls in CDK and OCL in both Java and JavaScript for generally useful InChI-related methods. IXA in Java refers (currently) to io.github.dan2097.jnainchi.IXA, and in JavaScript refers to the same, but is also aliased to J2S.wasm.jnainchi.IXA.

# Coding examples (InChl-to-image, Java)

```
BufferedImage inchiToImage(String inchi) {
  try {
    Pointer hStatus = IXA.IXA STATUS Create();
    Pointer hMolecule = IXA.IXA MOL Create(hStatus);
    void IXA.IXA MOL ReadInChI(hStatus, hMolecule, inchi);
    InchiInput input = CDK.getInchiInputFromMoleculeHandle(hStatus, hMolecule, "fixamide");
    IAtomContainer mol = CDK.getCDKMoleculeFromInchiInput(input);
    BufferedImage image = (mol.getAtomCount() == 0 ? null : CDK.getImageFromCDKMolecule(mol));
    return image;
  } catch (e) {
    throw e;
  } finally {
                                                           Note that all of the IXA method parameters
    IXA.IXA STATUS Destroy(hStatus);
                                                           and returns are simple variable types or
    IXA.IXA_MOL_Destroy(null, hMolecule);
                                                           com.sun.jna.Pointer objects. These are
                                                           easily handled in JavaScript.
```

# Coding examples (InChl-to-image, JavaScript)

```
function inchiToImage(inchi, moreOptions) {
  try {
    var hStatus = IXA.IXA STATUS Create();
    var hMolecule = IXA.IXA_MOL_Create(hStatus);
    IXA.IXA MOL ReadInChI(hStatus, hMolecule, inchi);
    var input = CDK.getInchiInputFromMoleculeHandle(hStatus, hMolecule, "fixamide");
    var mol = CDK.getCDKMoleculeFromInchiInput(input);
    var image = (mol.getAtomCount$() == 0 ? null : CDK.getImageFromCDKMolecule(mol));
    return (image == null ? "" : CDK.getDataURIForImage(image);
} catch (e) {
    throw e;
  } finally {
                                                           For JavaScript, in the end we to turn the
    IXA.IXA STATUS Destroy(hStatus);
                                                           image into a DataURI in order to display it;
    IXA.IXA_MOL_Destroy(null, hMolecule);
                                                           this is not necessary in Java.
```

# Coding examples (SwingJS CDK to OCL molecule)

```
Java:
StereoMolecule cdkToOCLMolecule (IAtomContainer mol) {
    InchiInput input = CDK.getInchiInputFromCDKMolecule(mol);
    return OCL.getOCLMoleculeFromInchiInput(input);
}

JavaScript (SwingJS):

cdkToOCLMolecule = function(mol) {
    var input = CDK.getInchiInputFromCDKMolecule(mol);
    return OCL.getOCLMoleculeFromInchiInput(input);
}
These
```

These two versions are essentially identical. In both Java and JavaScript, converting simple structures from CDK to and from OCL is just a matter of passing them through Inchilnput. One need not actually create the InChI string. No string parsing is required, and no stereochemistry is lost.

# Moving Forward

- Obviously needs more discussion!
- Submit PR requests for InChI, JNA, OCL, and CDK
- Proposing two new InChl repositories:
  - Inchi-API-JNA
    - Focus on Java, with simple IXA methods only
  - Inchi-API-WASM
    - Focus on JavaScript, using simple IXA methods only
  - Coordinated development
- None of this requires the use of SwingJS; we only use that because it also gives us access to OpenChemLib and the CDK in JavaScript.