

Archiving spectra: use InChI and CML

Egon Willighagen 

Published June 22, 2007

Citation

Willighagen, E. (2007). Archiving spectra: use InChI and CML. In *chem-bla-ics*. chem-bla-ics. <https://doi.org/10.59350/4n5zx-19r45>

Keywords

Cml, Inchi, Nmr

Abstract

Ryan blogged in Archive This about some advices from ACD on how to store spectra in your electronic lab notebook.

Copyright

Copyright © Egon Willighagen 2007. Distributed under the terms of the [Creative Commons Attribution 4.0 International License](#), which permits unrestricted use, distribution, and reproduction in any medium, provided the original author and source are credited.

Ryan blogged in [Archive This](#) about some advices from ACD on how to store spectra in your electronic lab notebook.

Use InChI

This reminded me of a [discussion I had with with Colin](#) when he was at the CUBIC, which was about experimental sections. I proposed that the [InChI](#) should have a prominent place in the experimental section. An important argument for this is that it allows well-defined atom numbering to be used when writing down the NMR bits in that section: the InChI gives a unique numbering, so that the numbering used in the experimental section becomes author neutral. Because the InChI puts the carbons up front, the ^{13}C NMR details get numbers from 1-13, or whatever the carbon count is. For proton NMR it is not difficult either, they are simply numbered according to the heavy atom to which they are attached. For situations where two hydrogens attached to the same heavy atom have different shifts, then a and b can still be used. The numbers are easily added to 2D diagrams anyway.

If software vendors (e.g. [ACD](#) and [Bioclipse](#)) and publishers (e.g. ACS, [RSC](#), [Chemistry Central](#)) could adopt this proposal, then experimental sections immediately are better machine parseable and ready for automatic processing, such as discussed in my blog item [Chemical Archeology: OSCAR3 to NMRShiftDB.org](#) and by [Christoph at the ACS meeting](#), available as [PDF](#) and this 18MB [MP3](#).

Use CML

Even better is to use [CML](#) for this, or CMLSpect to be precise (paper is accepted, and should appear soon). This XML-based language allows the full semantic markup of all the experimental details and all the interesting assignments you want to archive. I would like to **challenge ACD** to follow Bioclipse's lead and provide export as CMLSpect for spectral assignments and markup of experimental details, in addition to the PDF in whatever format they prefer. Cheers for the work by Tobias and Stefan on spectrum support in Bioclipse!