



CONNECTING RESEARCH AND RESEARCHERS

Getting Credit for Writing Scientific Software

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The Problem: Name Ambiguity

Reliable linking of researchers with their research is impossible using their names

The scholarly record is broken



The solution: unique author identifiers



Martin Fenner

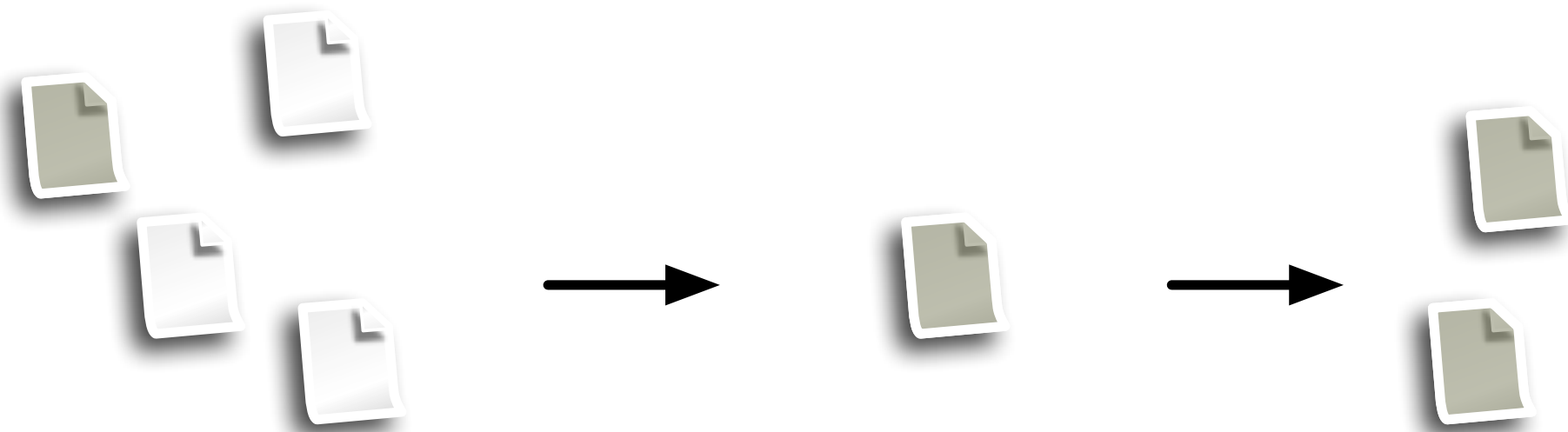
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Benefits: Discovery

Reliable linking of researchers with their research facilitates discovery

Easier to find research outputs by a particular researcher

Easier to find other researchers through co-authors



Benefits: Evaluation

Reliable linking of researchers with their research facilitates evaluation

Not just journal articles, but also posters, datasets, book chapters, software, etc.

Not just research outputs, but also funding, people, etc.



Why is ORCID a Solution?

Global	Not limited by discipline, institution or geography
Open	Inclusive and transparently governed not-for-profit Data and source code available under recognized open licenses
Integrated	Third-party seeding of profiles Part of institutional, publisher, and funding agency infrastructure

Who can create an ORCID identifier?

The Researcher



His University/Research Organization
with permission from the researcher





ORCID Members

ORCID launched in October, 2012 and had 27 member organizations as of January 6, 2013:

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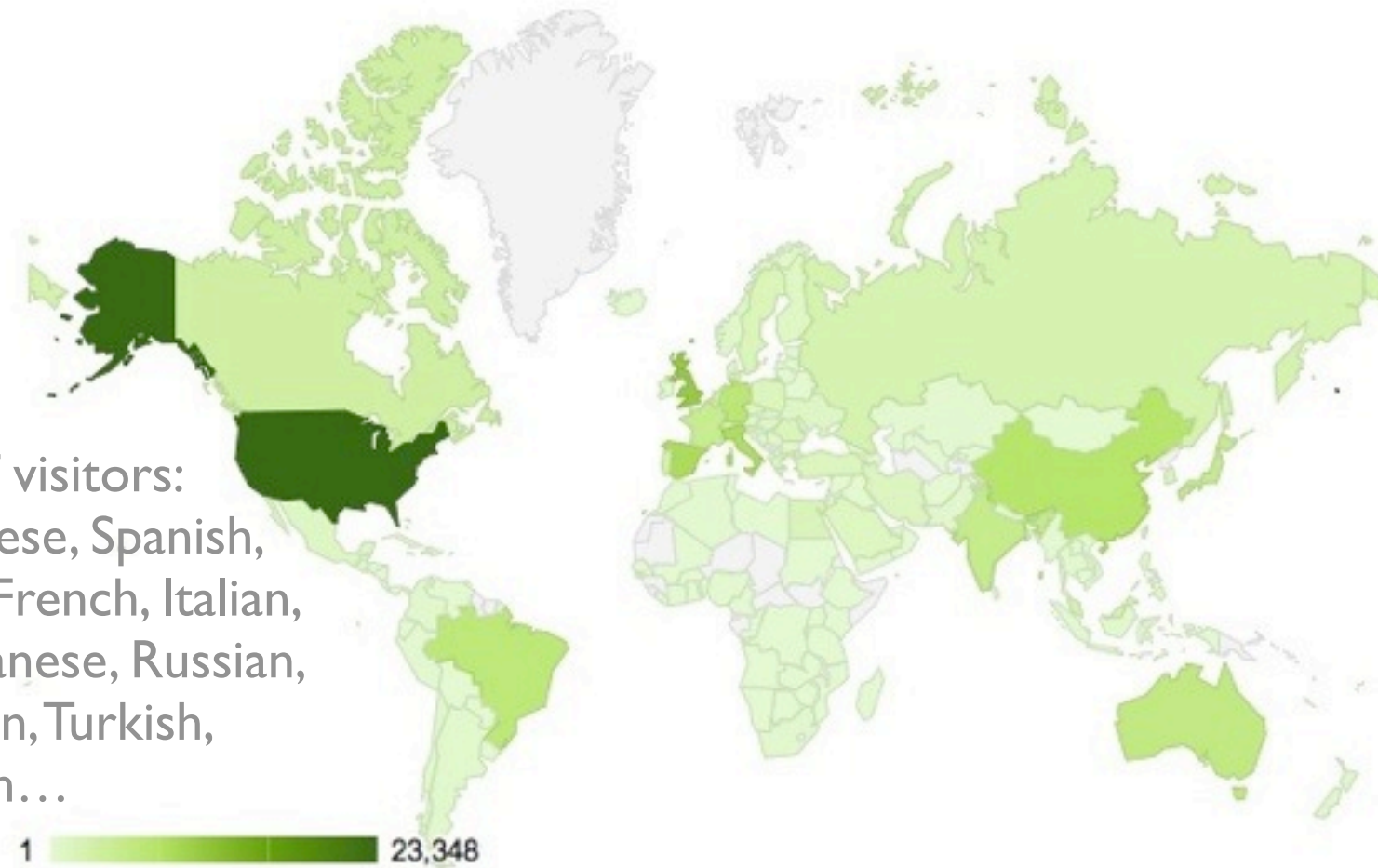
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27 countries with > 1000 visitors, 68 with >100.

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4. Italy	8,284	5.96%
5. China	7,435	5.35%
6. Germany	6,598	4.75%
7. Brazil	6,448	4.64%
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10. Japan	4,853	3.49%

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<http://orcid.org/register>

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Affiliations

Works UPDATE ↕

Blogs, Wikis und Podcasts im Unterricht Oct-2008

Public

DOI: [10.1002/biuz.200890080](https://doi.org/10.1002/biuz.200890080)

Fenner, M, 2008, 'Blogs, Wikis und Podcasts im Unterricht', *Biologie in unserer Zeit*, vol. 38, no. 5, pp. 284-286.

Carboplatin plus weekly docetaxel as salvage chemotherapy in docetaxel-resistant and castration-resistant prostate cancer:

Public

World Journal of Urology 2010

DOI: [10.1007/s00345-010-0527-5](https://doi.org/10.1007/s00345-010-0527-5)



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Thornton Group Software

The following software packages have been made publicly available by members of the Thornton Group.

Software	Description
ArchSchema	ArchSchema is a java program that generates dynamic plots of related Pfam domain architectures. Authors: Asif Tamuri and Roman Laskowski
LIGPLOT	LIGPLOT is a program which plots schematic diagrams of protein-ligand interactions. Authors: Andrew Wallace and Roman Laskowski .
PROCHECK	PROCHECK checks the stereochemical quality of protein structures. Authors: Roman Laskowski , Malcolm MacArthur, David Smith, David Jones, Gail Hutchinson, Louise Morris, David Moss and Janet Thornton.
PROCHECK-COMP	PROCOMP Compares residue-by-residue geometry of a set of closely-related protein structures. PROCOMP is part of PROCHECK (see above). Author: Roman Laskowski .
PROCHECK-NMR	PROCHECK-NMR is a suite of programs that have been derived from the PROCHECK programs to analyse ensembles of protein structures solved by NMR. Authors: Roman Laskowski and Malcolm MacArthur.
QUASIPROX	Quaternary Structure Inference of Proteins from their Crystals. Author: Hannes Posntingl.
SMSD	SMSD (small molecule subgraph detector) is a program to find similar molecules and reactions to a given one. Author: Syed Asad Rahman (asad@ebi.ac.uk).
SURFNET	Program to generate surfaces, and void regions between surfaces, from coordinate data supplied in a PDB file. Author: Roman Laskowski . N.B. This link is still hosted at UCL.

Contact

We would like to encourage laboratories wishing to discuss any collaborations to contact us. For information, comments and/or suggestions please [contact us](#).



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alm / 955 commits

ALM-158 fixed parsing for attributes in scienceseeker XML		
	mfenner authored 6 hours ago	latest commit 77af3e680e
app	6 hours ago	ALM-158 fixed parsing for attributes in scienceseeker XML [mfenner]
config	10 hours ago	ALM-193 increased timeout for jobs [mfenner]
db	2 days ago	ALM-153 tried to fix scienceseeker namespace errors [mfenner]
features	a day ago	ALM-154 display responses of past 31 days and not 30 days to be in sy... [mfenner]
lib	6 hours ago	ALM-158 fixed parsing for attributes in scienceseeker XML [mfenner]
log	11 months ago	initial checkin of a blank ruby on rails application [jsong-plos]
public	14 days ago	ALM-157 removed precompiled assets [mfenner]
script	8 months ago	Added Cucumber files [mfenner]
spec	6 hours ago	ALM-158 fixed parsing for attributes in scienceseeker XML [mfenner]
vendor	2 days ago	ALM-153 tried to fix scienceseeker namespace errors [mfenner]
.gitignore	7 days ago	ALM-178 don't include data_bags folder in version control [mfenner]



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#2	David Heinemeier Hansson	3595
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#5	Xavier Noria	1840
#6	Santiago Pastorino	1626
#7	Josh Peek	1230
#8	Rafael Mendonça França	1145
#9	Jon Leighton	1120
#10	Vijay Dev	1090
#11	Pratik Naik	945
#12	Yehuda Katz	895
#13	Carlos Antonio da Silva	888
#14	Michael Koziarski	757
#15	Rick Olson	595
#16	Piotr Sarnacki	537

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Brown, Dave • Brownrigg, Richard • Haley, Mary • Huang, Wei

[A case study for efficient management of high throughput primary lab datasource code](#) [version 1.1] # 2

[doi:10.5447/IPK/2011/0](#) Software : source code
Colmsee, Christian • Flemming, Steffen • Klapperstück, Matthias • Lange, Matthias • Scholz, Uwe

[Workspace](#) # 3

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Clark, S

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Ouyang, Jing • Wang, Xufeng • Qi, Minghao

[Polymer Modeler](#) [version 3.2] # 10

[doi:10.4231/D3PG1HN27](#) Software : Simulation Tool
Haley, Benjamin • Wilson, Nate • Li, Chunyu • Arguelles, Andrea • Jaramillo, Eugenio • (et. al.)

**Brief reports**[Open Access](#)

RMol: A Toolset for Transforming SD/Molfile structure information into R Objects

Martin Grabner, Kurt Varmuza and Matthias Dehmer

For all author emails, please [log on](#).

Source Code for Biology and Medicine 2012, **7**:12 doi:10.1186/1751-0473-7-12
Published: 14 November 2012

Abstract (provisional)

Background

The graph-theoretical analysis of molecular networks has a long tradition in chemoinformatics. As demonstrated frequently, a well designed format to encode chemical structures and structure-related information of organic compounds is the Molfile format. But when it comes to use modern programming languages for statistical data analysis in Bio- and Chemoinformatics, R as one of the most powerful free languages lacks tools to process R Molfile data collections and import molecular network data into R.

Results

We design an R object which allows a lossless information mapping of structural information from Molfiles into R objects. This provides the basis to use the RMol object as an anchor for connecting Molfile data collections with R libraries for analyzing graphs. Associated with the RMol objects, a set of R functions completes the toolset to organize, describe and manipulate the converted data sets. Further, we bypass R-typical limits for manipulating large data sets by storing R objects in bz-compressed serialized files instead of employing RData files.

Conclusions

By design, RMol is a R tool set without dependencies to other libraries or programming languages. It is useful to integrate into pipelines for serialized batch analysis by using network data and, therefore, helps to process sdf-data sets in R efficiently. It is freely available under the BSD licence. The script source can be downloaded from <http://sourceforge.net/p/rmol-toolset>.

Source Code for Biology and Medicine

Volume 7

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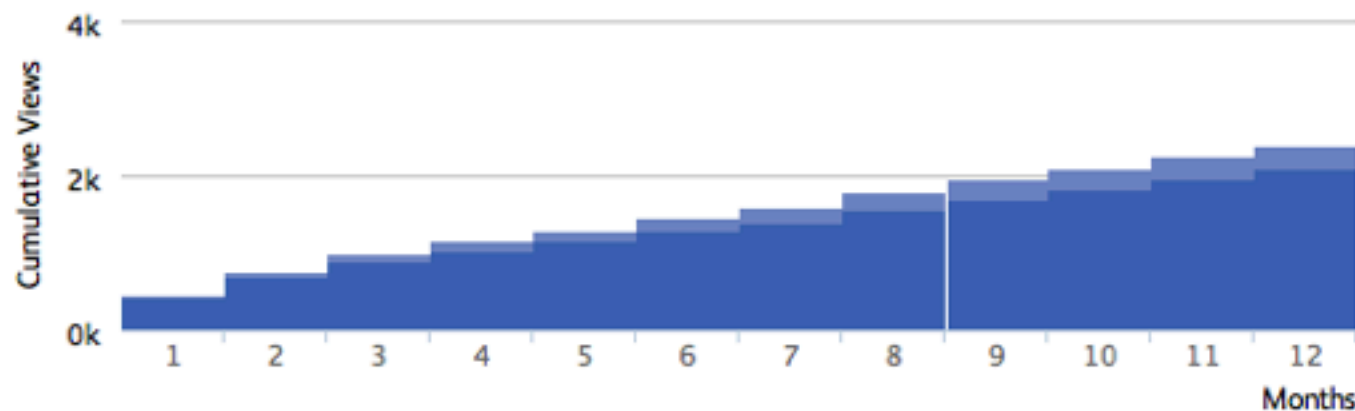
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- ChemInform (3,714)
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Published **6 Mar 2008** in **Nature** , volume **452** , issue **7183** , on pages **29** to **29**

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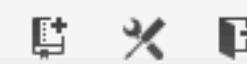
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kjw authored a day ago latest commit 334f0e0797

.bundle	6 months ago	Add bundler support. [Karl Ward]
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