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OPEN LETTER

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# The *bio.tools* registry of software tools and data resources for the life sciences

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## Abstract

Bioinformaticians and biologists rely increasingly upon workflows for the flexible utilization of the many life science tools that are needed to optimally convert data into knowledge. We outline a pan-European enterprise to provide a catalogue (<https://bio.tools>) of tools and databases that can be used in these workflows. *bio.tools* not only lists where to find resources, but also provides a wide variety of practical information.

A myriad of providers - from individual scientists to large service organizations - have created thousands of databases and tools, serving a dynamic domain spanning biology, biotechnology and medicine. Scholars must contend with intrinsically complex biological data, integrated into hundreds of data formats for analysis by a vast array of methods and diverse types of software, deployments and interfaces. Developments are often ad hoc, and in the absence of a source of unified information, it is not easy to assess the scope and compatibility of new resources in context of global offerings. For example, software may lack a formalized description of its scientific and technical function, and the absence of persistent, unique tool identifiers confounds reliable citation and reproducibility of analyses. There are significant barriers to find and connect the right tools among a multitude of possibilities, making the work of the bioinformatician - developing practical workflows for scientific discovery - far from trivial.

Since the 1980s various initiatives, at a local or more global level, have catalogued bioinformatics resources to advertise their wares and guide scientists in their choices. Early single investigator initiatives include the famous Pedro's List of weblinks and Gunnar von Heijne's 1987 book 'Sequence Analysis in Molecular Biology: Treasure Trove or Trivial Pursuit' [1]. Contemporary examples include international service providers [2], laboratories (<https://www.rostlab.org/>), software suites (<https://www.bioconductor.org/>), deployment solutions [3, 4], scientific publishers [5], WIKIs such as [msutils.org](https://msutils.org), and online catalogues [6] including commercial offerings such as from omicX (<https://omictools.com/>) and open lists such as from the BIG Data Center initiative (<https://bigd.big.ac.cn/tools>) based in Beijing. Such collections serve their communities well, but when taken as a corpus of information about tools in general, present a fragmented information landscape, with much redundancy. Owing to a lack of commonly adopted information standards, it can be difficult to understand what is available and compare different approaches to the same problem. Web search engines like Google provide the entry point for searches, but yield results reflecting mostly historic prevalences, insufficiently structured to allow ready comparison. Thus, in an era of highly efficient Web searches generally, barriers remain to the efficient utilization of bioinformatics resources, with continued use of suboptimal offerings, slow uptake of new tools and reinvention of existing functions.

A practical first step [7] towards a sustainable and unified resource registry engaged enthusiastic individuals from the spectrum of European bioinformatics, to share and maintain information about resources within their scope. This effort is now joined by the 22 nodes of

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ELIXIR (<https://www.elixir-europe.org/>), the European Infrastructure for Biological Information. Our aims include:

1. scientists can find, understand and compare tools for computational experiments, and access the wealth of data resources
2. bioinformaticians have clues about compatibility of tools with various data types and formats, thus, what might readily be chained into functional workflows
3. developers can find and assess implementations of desired functionality, encouraging reuse and repurposing over reinvention
4. end-users can easily find supplementary information, such as benchmarking results or training courses
5. facility managers can see the status (emerging, mature or legacy) of a resource, including licensing, and assess its applications and technical performance during service design
6. funders and reviewers have an overview of productions at various hierarchical levels such as individual, institutional or even national
7. tool developers and service providers can contribute to the registry in simple but effective ways
8. information about the legacy of resource developments does not get lost

Fulfillment of these aims requires upkeep of a high quality, non-redundant corpus of information, that is integrated with deployment solutions, scientific literature and pertinent activities including benchmarking, monitoring and training, and which can adapt to the bioinformatics landscape of tomorrow. The burden is therefore onerous. Developers and providers are best placed - and motivated - to document their own productions, but given the complex landscape they require sustained coordination and support. They have been left alone in this critical activity, and it is no surprise that a unified and enduring catalogue has remained an elusive goal. A major community-driven effort is required, sustained by long term institutional commitments. ELIXIR, as the linchpin of a network of diverse research infrastructures, is ideally positioned to promote a common strategy and deliver a portal that is broadly relevant across a range of disciplines and user groups.

Our portal (<https://bio.tools>), which has developed steadily over 5 years, now includes over 250,000 annotations on some 12,000 resources. All types of application software are within scope, across all life science domains globally. This includes everything from simple command-line tools and Web applications, to databases, workflows and integrated workbenches. Most entries describe open source or freely

accessible tools with straightforward functions, which are therefore readily combinable into functional workflows. Accessions are assigned a unique tool identifier: a manually verified, URL-safe version of the supplied tool name. When used in combination with a version label assigned by a developer, the tool IDs provide a pragmatic means to cite and trace software, especially in the absence of a traditional publication. The IDs are used in persistent *bio.tools* URLs, resolving to Tool Cards of essential information. *bio.tools* mandates only bare-bones information (name, short description and homepage), whilst supporting rich description of 50 salient scientific, technical and administrative attributes. Resource descriptions must conform to rigorous semantics and syntax, defined in a formalized schema, *biotoolsSchema* (<https://github.com/bio-tools/biotoolsSchema>). Controlled vocabularies are used extensively, and provide concise, consistent and therefore comparable information, for the convenience of the user. For example, tools may be annotated with specific topics, operations, input and output data types and supported formats from the EDAM ontology [8]. Standard identifiers are used where possible, e.g. DOIs for publications, and verbose information, such as documentation or citation instructions, are referenced by URL. Hence, the dizzying complexity of bioinformatics software is reduced to collections of readily understandable functional units, put in scientific and technical context, including information to enable access and use. The aggregation and standardization of data under the portal can help end-users in very practical ways. Consider for example a biologist who is surveying recently published tools in a general scientific area, or for a specific computational task, and wants to identify those which are freely available for use. They can search *bio.tools* using specific EDAM topics and operations to quickly make a list of candidate tools and compare alternatives, drilling down to tools available under open license and with a recent publication. Without *bio.tools* they would need to manually search and browse a large number of web pages, ranging from software repositories (e.g. GitHub) to scientific literature resources (e.g. PubMed), which can be a time-consuming and difficult process.

The initiative upholds open science principles [9], and thus far has benefited from 1127 contributors from 422 domains. Contributions to date are mostly from Europe and the USA, which simply reflects *bio.tools'* European foundation and the high volume of American tools. There are, of course, vibrant bioinformatics communities all over the world, and we warmly welcome and encourage their participation. Direct curation assistance is available from the core *bio.tools* team, through collaboration with ELIXIR partners and at community-led workshops. The effort expected from providers is thus reduced to a relatively small and maintainable level, and we hope to attract and retain many new contributors and collaborators. Direct participation in the project and

re-use of the registry is strongly encouraged. Practical information describing how scientific communities and individual software developers can contribute are available online [10, 11]. Access to the portal is unrestricted and both the registry content and portal source code (<https://github.com/bio-tools/biotoolsRegistry/>) are freely available under open license (CC BY 4.0 and GPL-3.0 respectively).

We have summarized our vision and progress towards a solution of a global and major challenge: a uniform means by which to describe, publish, discover and cite bioinformatics resources. *bio.tools* is a step towards a central point of unified information, to avoid the rewriting of resource descriptions in so many different contexts. The current implementation upholds the FAIR data principles [12] and, with progressive development, will help make bioinformatics resources more findable and accessible, and somewhat more interoperable and reusable. To fully realize our vision, however, involves much ongoing work:

1. inclusion of information about online services, deployment solutions and supported data formats, to provide users with information about availability and uptime, and enable tool use and applications such as automated workflow composition [13].
2. ease the curation process, e.g. by curation tools [14], and utilities [15] to pull tool information from workbench environments such as Galaxy, or, where applicable, directly from code repositories such as GitHub, and by new linting utilities (e.g. <https://github.com/bio-tools/biotoolslint>) to identify and fix inconsistencies in annotations.
3. leverage specialized community efforts (<https://www.elixir-europe.org/communities>) and biomedical science research infrastructures internationally, to expand coverage and improve quality in areas such as proteomics, metabolomics and bioimaging.
4. stable metadata sharing mechanisms for institutional collections such as IFB tools (<https://www.france-bioinformatique.fr/en/services/tools>) and specialized registries such as BioContainers [16].
5. inclusion of Web APIs and services for accessing the multitude of biological databases, e.g. by developing systems [17] that leverage community standards such as OpenAPI (<https://www.openapis.org>).
6. expose quality metrics to provide a trustworthy and rational means for tool assessment, including scientific benchmarking of analytical tools and monitoring of service technical robustness, from platforms such as ELIXIR openEBench (<https://openebench.bsc.es>).
7. services [18] to combine and export *bio.tools* data with execution-layer information in specific workflow configuration formats such as used by Galaxy [19] or a generic one such as the Common Workflow Language (<https://www.commonwl.org/>).
8. more convenient and powerful interfaces and features for query formulation, searching and browsing.
9. enhancing the management of user profiles and crediting of contributions, e.g. using ELIXIR AAI [20] federated user identity management, which incorporates researcher identities such as ORCID (<https://orcid.org/>).
10. crosslink with portals such as ELIXIR TeSS [21] (training resources) and FAIRSharing [22] (data standards), in order to make navigation of the broader bioinformatics resource landscape more coherent and convenient

With community support, *bio.tools* can become a standard way to disseminate publicly-funded software development. The primary long-term challenge is to nurture the community around it and ensure the portal matches end-user requirements. Here, the anchoring within ELIXIR allows us to draw upon a coordinated, European-wide community of experts, including national service managers. Long-term support from these partners, and synergistic relationships with community projects and other major international initiatives, will sustain the portal in the long term, allowing for secure planning and investment. We welcome collaborations with all scholars on common goals, and encourage life scientists worldwide to join forces in a task that can greatly benefit the whole community.

## Additional file

**Additional file 1:** Review history. (DOCX 20 kb)

### Review history

The review history is available as Additional file 1.

### Authors' contributions

Jl led the work described and prepared the manuscript. HI, ER and PC developed the *bio.tools* website. HI, HM, MK and VS contributed to the technical development of registry, its content and the EDAM ontology. SD, TR and AS contributed to the content. BG, NB, RJ, KR and GV contributed to the technical development of registry. RL, HS, BP, RSV, JV, HP, IJ, RH, TN, AV, SC, JG, FZ, BS, BL, CB, AO, OC, JvH and PL coordinated the European institutional contributions to the registry content and manuscript. *bio.tools* is coordinated on behalf of ELIXIR by the Danish ELIXIR Node under the leadership of SB. All authors read and approved the final manuscript.

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### Availability of data and materials

The *bio.tools* content is freely available to all via the *bio.tools* API under the Creative Commons Attribution license (CC BY 4.0). The source code of the registry is available under standard GPL 3.0 license from GitHub (<https://github.com/bio-tools/biotoolsRegistry/>) [23]. EDAM and biotoolsSchema are licensed under the Creative Commons Attribution-ShareAlike 4.0 International License (CC BY-SA 4.0).

### Ethics approval and consent to participate

Not applicable.

### Competing interests

The authors declare that they have no competing interests.

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