

Data-Producing Methods in CRC 985: Recommendations for Research Data Management in Large Interdisciplinary Projects

CRC 985: Functional Microgels and Microgel Systems

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Software availability:

No software was specifically developed for this project. The associated Jupyter Notebook can be found within the above-mentioned dataset.

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Abstract. Large, interdisciplinary projects produce various type of data underlying their published results. To gain a deeper understanding of the data produced, a survey was conducted in a project comprising the fields of chemistry, physics, engineering and life sciences, with the intention to improve the research data management.

Based on the collected information as well as feedback from researchers, we outline a holistic research data management approach, starting at the individual research group level. Here, we focus on data governance, documentation, and data exchange formats. We tie this together at the project level with a focus on data workflows for a collaborative data management and recommend data publication and archival solutions for this specific project. As a whole, this strives to provide researchers with the basic framework to efficiently work and manage their research data while producing understandable and reusable results in line with the FAIR principles.

1 Introduction

The collaborative research center (CRC)¹ 985 *Functional Microgels and Microgel Systems* has studied microgels, soft colloidal macromolecular compounds that find applications in many different fields, for over two funding periods, the current third funding period being its final. The project brings together research groups from numerous chemical institutes, chemical engineering, physics, biotechnology, and the life sciences, with RWTH Aachen University, DWI - Leibniz Institute for Interactive Materials, the RWTH Aachen University Hospital (UKA),

1. CRCs are long-term yet temporary research projects funded by the German Research Foundation (DFG). They can run a total of 12 years, with individual funding periods of 4 years.

8 and Forschungszentrum Jülich (FZJ) cooperating with each other. In total, roughly 40 groups,
9 currently involving approx. 90 principal investigators (PIs), post-doctoral researchers, or doctoral
10 researchers, have or are actively contributing to the project. Over 300 scientific publications
11 have been produced so far.

12 In the first funding period, which began in 2012, the research data management (RDM) struc-
13 ture included a Microsoft SharePoint, while Mattermost was introduced as an instant-message
14 communication system. On this basis, information could be shared and communicated across
15 research areas as well as internally in smaller groups. Furthermore, during the previous funding
16 periods, a sample management system was integrated into SharePoint to track sample history,
17 while implementing a universal naming system throughout the CRC and assigning persistent
18 identifiers (PIDs) [1]. Until the third funding period, the INF project largely focused on estab-
19 lishing collaborative digital systems in the first funding period and improving upon these to
20 increase acceptance in the second. At this point, consulting in terms of RDM also increased.

21 General guidelines for data publication were established, yet, most data was shared and stored
22 in a manner that did not follow any specific standards. The researchers' best practice has thus
23 been to document their work in the form of individually written texts, digital or analog, and
24 to save raw and/or processed measurement data in an individual project folder. Storing data
25 across projects with the same structure and making it accessible for future projects is challenging
26 with this approach. One reason for this is that different templates would have to be developed
27 individually for different tasks, or new software would have to be developed for this purpose
28 explicitly for this CRC. Similar statements regarding this problem description for projects of this
29 scale have been published in other CRCs [2], [3].

30 From today's perspective, proficient RDM requires much more, e.g., the sharing and archiving
31 of data according to the FAIR (findable, accessible, interoperable, reusable) principles that were
32 introduced in 2016 [4], coinciding with the second funding period as well as the establishment
33 of a central RDM team at RWTH Aachen University. At their core, these guiding principles
34 build upon one another to ultimately ensure a dataset's reusability. For research data, they carry
35 implications for both those producing the data, e.g., researchers, but also for those providing
36 infrastructure such as research data repositories [5]. Implementing practices and tools that enable
37 FAIR throughout each stage of a research project also facilitates FAIR in the long run. Large,
38 interdisciplinary projects can benefit from these practices as participants can efficiently find,
39 access, and (re)use data produced by their collaborating partners or predecessors, e.g., from
40 previous funding periods.

41 Fully functional RDM infrastructures and information standards are still a work in progress. The
42 German National Research Data Infrastructure (NFDI; German: Nationale Forschungsdatenin-
43 frastruktur) and its discipline-specific consortia aim to move this progress along [6]. In the area
44 of chemistry, NFDI4Chem strives to not only set up a system of repositories for data sharing and
45 archival, but also to establish minimum information and format standards to ensure data remains
46 reusable and interoperable [7]. These efforts should inform the research communities' RDM
47 practices, while the consortia also require researchers' input to best suit their needs.

48 As part of the CRC 985 Information and Infrastructure (INF) project, we present an overview of
49 the diversity in a research project of this magnitude in terms of the number of data-producing

50 methods and the variety of associated data. A survey to gather relevant information lays the
51 foundation of this work. Based on this information as well as on formal and informal exchange
52 with CRC project members, we discuss how to deal with such a variety of data in future projects
53 in terms of project preparation, recommended RDM practices regarding storage, publication,
54 archival and the accompanying data formats, and communication and awareness among participat-
55 ing researchers. Furthermore, as a project which includes many chemical and chemistry-related
56 disciplines, the information presented here can inform the efforts and goals within NFDI consortia
57 such as NFDI4Chem.

58 2 Methodology

59 Figure 1 shows the general approach taken for this work. Stage 1 focused on gathering information
60 within CRC 985. To this end, the INF project compiled a structured questionnaire [8] to survey
61 the data-producing methods and workflows throughout the CRC. It then acquired contacts for
62 RDM-related topics for the various research groups and subprojects by contacting the relevant
63 PI. The first version of the questionnaire was then distributed to the supplied contacts. In most
64 cases, the contacts named were PhD candidates working within CRC 985, yet, also included
65 more senior research staff in some cases.

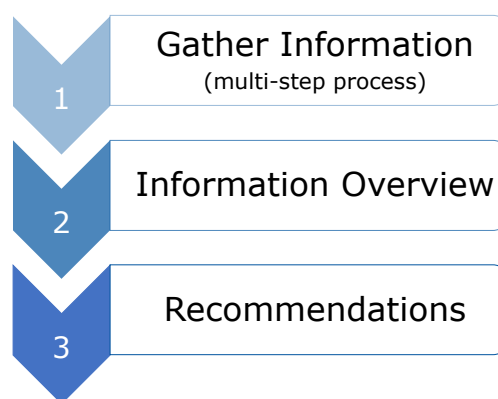


Figure 1: Targeted incremental approach to provide an overview of the project's data scope and set the basis for future RDM improvements.

66 The first version of the questionnaire focused on the methods themselves, aiming first and
67 foremost to understand technical aspects such as device specifications, output data formats and
68 volume, and frequency of use within for the CRC and within the respective research group. Two
69 issues soon became apparent: (1) The results lacked certain information that would be useful
70 to the INF project, especially regarding current RDM practices such as data workflows and
71 documentation, and (2) some terminology, such as metadata or controlled vocabulary (a term
72 added to the second version), or the questions themselves were unclear to the participants.

73 Thus, the questionnaire underwent two revisions. The third and final version split the question-
74 naire into two parts: one regarding each method used, gathering details as described above, and
75 a second regarding overall RDM practices such as the use of an electronic laboratory notebook
76 (ELN), the implementation of the CRC 985 policy on data, and the use of the sample management

77 system. Definitions of terminology were added as well. This granted participants the opportu-
78 nity to answer the questions independently and gather information in advance of face-to-face
79 exchanges. The first part now also included a question on data workflows, specifically, how data
80 are transferred from the device computer to other servers or data management systems, aiming
81 to determine if data workflows could benefit from automation.

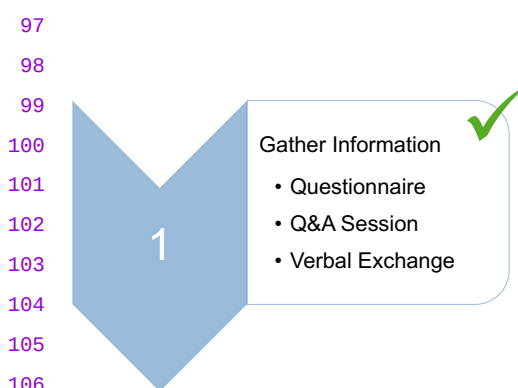
82 The questionnaire versions were maintained using the central CRC 985 SharePoint. These
83 surveys and exchanges took place starting in 2021 through 2023.

84 In the second stage, the INF project compiled an overview of the gathered information on data-
85 producing methods. This serves as a resource on available methods and contacts for CRC 985
86 and was therefore published on the project's SharePoint for easy reference.

87 The third stage, recommendations, employs the data collected and tabular overview created in the
88 previous stage as well as general information and feedback collected in a rather informal manner
89 in question and answer sessions as part of workshops or presentations. This informed the INF
90 project on the needs of the researchers. By drawing on knowledge provided by Fairsharing.org [9],
91 re3data.org [10], and NFDI4Chem [11] as well as central solutions offered by RWTH Aachen
92 University, recommendations for current and future projects on infrastructure options, e.g.,
93 working data storage, ELNs, and data publishing and archival services, are made. Furthermore,
94 areas that require additional work by infrastructure providers are pinpointed.

95 3 Results and Discussion

96 3.1 Stage 1: Gathering Information



107 **Figure 2:** Successful information gathering
108 through a questionnaire that was continuously
109 improved through question and answer sessions
110 and a close exchange with CRC 985 scientists.

The questionnaire created at the beginning of this study was used as a living document. Therefore, updates to the questions occurred throughout the first stage to better explain the questions and thus acquire more detailed information, as outlined in Section 2. The questionnaire successfully gathered information in a structured manner and allowed for a baseline to gain more detailed information. This required close face-to-face exchange between the research project members and members of the INF project. In total, 16 interviews were conducted, involving 13 research groups working within the project.

111 In addition, the INF project held seminars for
112 researchers to raise awareness with respect to RDM. Subsequent question and answer sessions
113 gave a further overview of the methodological diversity as well as other RDM-related concerns,
114 enabling the INF project to provide suggestions to facilitate RDM in the CRC 985. Therefore,
115 by combining a questionnaire as a living document with a close exchange between the data-
116 producing researchers, the first phase was successfully completed (Figure 2).

117 It should be noted that participation was voluntary and the knowledge of the participants regarding
118 RDM varied greatly. Thus, receiving a full and complete picture of RDM throughout the groups
119 involved in the CRC proved difficult, resulting in possibly incomplete information. To gain a
120 full and complete picture for a holistic RDM within such projects, INF projects must be better
121 integrated into the individual research groups, with responsibilities and points of contacts defined
122 from the onset, as further discussed in Section 3.3.

123 All versions of the questionnaire as well as the completed surveys can be found within the dataset
124 published on Radar4Chem [8]. The file naming convention includes the respective version for
125 each completed survey. Additional notes on verbal exchanges are included in the individual
126 documents.

127 3.2 Stage 2: Information Overview

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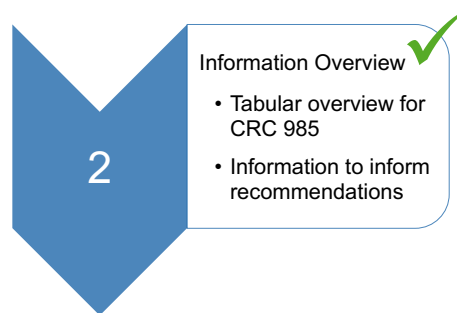


Figure 3: Successful information overview that tabulates all methods and resulting data volumes within CRC 985.

numerical methods, cover a broad context of disciplines. This rather coarse-grained depiction summarizes the methods into wider categories. It should be mentioned that the amount of devices and setups employed throughout the CRC gives rise to a large variety of data, including differences in the data output sizes and file types, even within a specific method. In total, 40 method categories were reported throughout the project. As this reporting was primarily voluntary and researchers may acquire, develop, or even switch methods as a project progresses, this number is approximate.

Figure 5 exhibits the resulting multitude of data output sizes. The majority of the methods produce data at or below the 1 GB mark, while five methods, namely high-resolution microscopy methods, such as superresolution fluorescence microscopy or tensiometry, and numerical methods, cross or go far beyond that mark. This must be taken into account for recommendations on storage, publication, and archival.

The survey results provide an overview of commonly used data formats for raw and exported data. This will be discussed in more detail in Section 3.3, with reported data formats provided in Table 1. During exchange with researchers and due to the responses presented below, it was clear that standard formats were not necessarily well-known, however, and therefore guidance

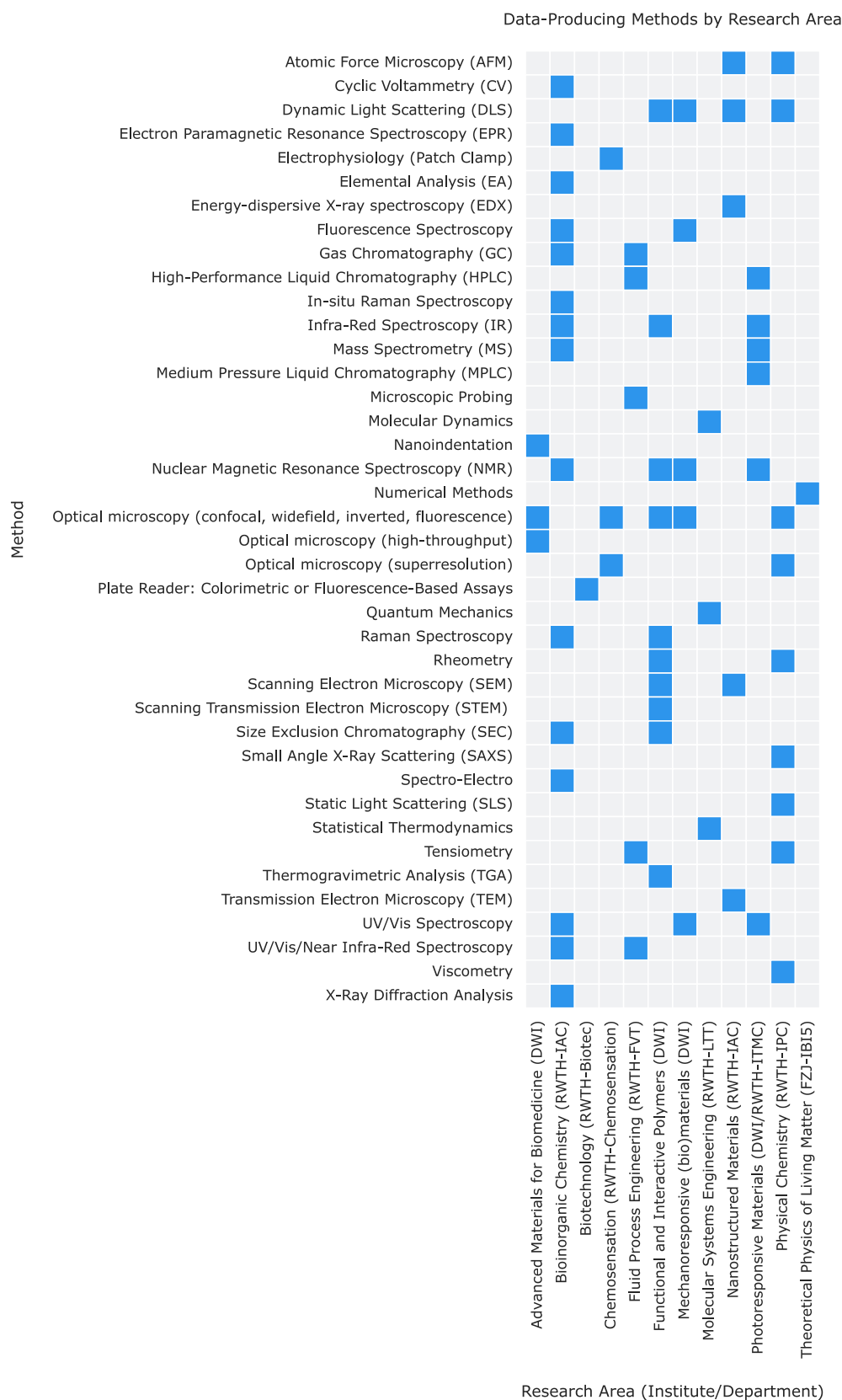


Figure 4: Methods reported according to their area of research in CRC 985. The employed or available methods range from spectroscopy, to microscopy, to numerical, representing the variety of disciplines involved in the project. Nevertheless, many methods are common to chemistry-related research. In total, 40 method categories were reported.

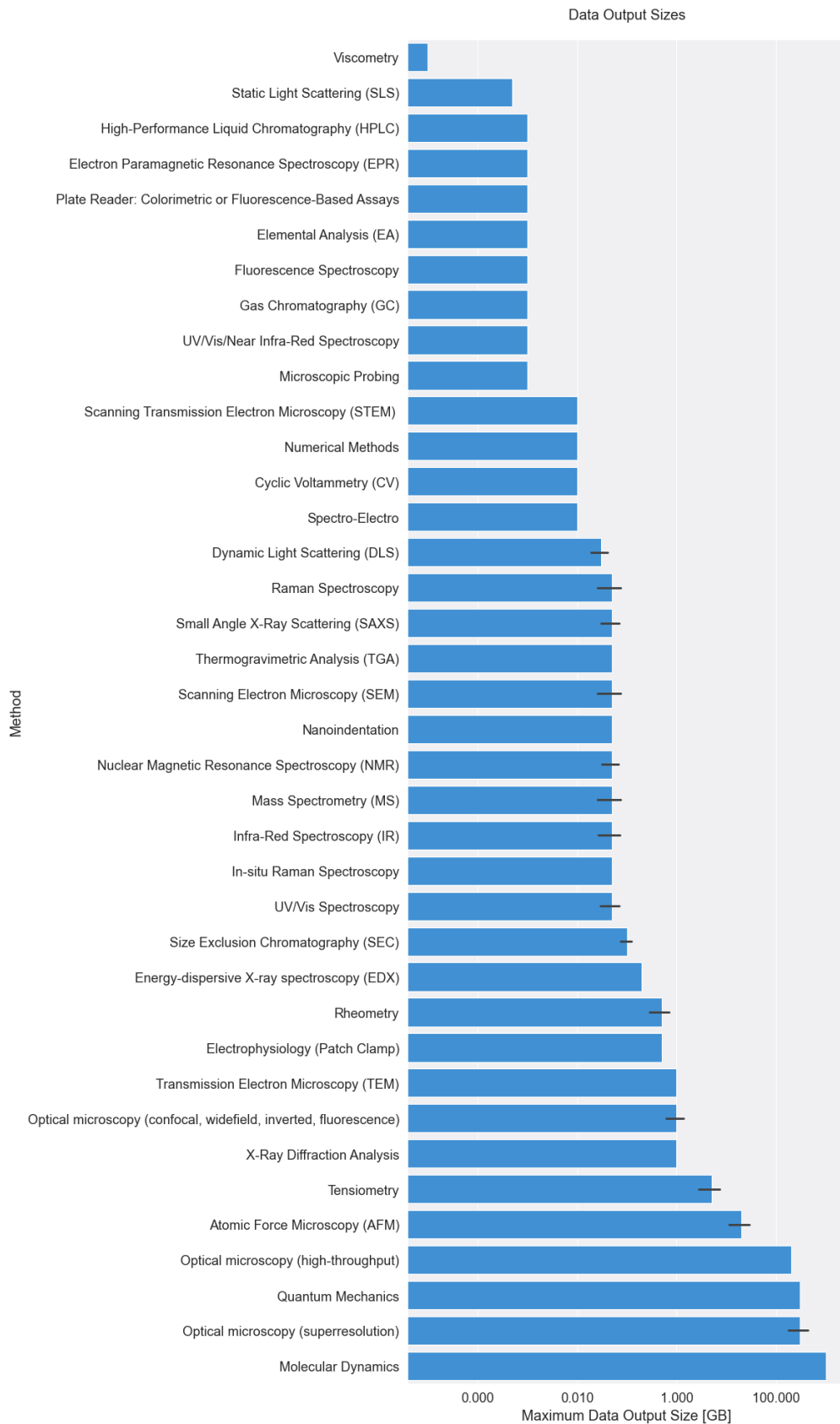


Figure 5: Methods and their output data sizes (logarithmic scale) reported in CRC 985. Most reported output sizes are smaller than 1 GB, with numeric and imaging methods far beyond that point and up to 1 TB. Where applicable, error bars indicate the standard deviation of the data output sizes reported for specific methods.

158 on data formats is required. This information was included on the shared overview table on the
159 SharePoint for project members to reference and to create general awareness. An anonymized
160 version of this table is also provided in the published dataset [8]. Furthermore, some information
161 was added to the table without specific surveys being carried out, rather, to add to the central
162 methods overview.

163 The questionnaire also addressed data documentation, especially regarding (uniform) metadata.
164 The responses reveal that, for most groups, very little uniform, machine-readable metadata are
165 recorded unless it is contained directly in the output data files. However, this information may
166 not always be contained in the exported version of the data, with which many members reported
167 working. Relevant information is often included directly in the file name, analog or ELNs, or
168 digitized in plain text, Microsoft Office, or Microsoft Excel files. Only one group mentioned
169 using controlled vocabularies.

170 It should be noted that, in some cases, project members, especially doctoral students, expressed
171 concerns in terms of data storage best practices, which data should be stored, published, and
172 archived at which stage (raw vs. exported or processed data), data organization, and data formats.
173 This was often expressed in informal conversations, workshop, or seminar settings.

174 Thus, the survey provided sufficient results to obtain an overview of the methodological diversity
175 and generated data that led to the successful completion of the second phase (Figure 3). In
176 addition to the data-producing methods, other foundational aspects and concerns regarding RDM
177 were collected and will be addressed in the following.

178 3.3 Stage 3: Recommendations

179 Based on the knowledge gained from the presented results, we derived the following recom-
180 mendations as outlined below. On the one hand, the data-producing method types and file sizes
181 influence aspects such as data publication platforms and recommended file types. On the other
182 hand, the project participants' accounts allow us to directly address the concerns and advise on
183 research data management best practices accordingly.

184 The main concerns reported were:

- 185 1. (Lack of) knowledge and implementation of data organization basics and best practices
186 regarding working data storage and structure
- 187 2. Internal data reuse, e.g., the ability to easily build upon a predecessor's work
- 188 3. Access to storage space for large amounts of (raw) data
- 189 4. Data exchange format standards
- 190 5. (Lack of) knowledge of data documentation best practices and minimum information
191 (metadata) standards
- 192 6. Publishing data underlying a journal article publication, e.g., which repository best suits
193 the research data and data access control (open access vs. closed access options)

194 These concerns were largely reported on a research group and not necessarily a project-specific
195 level. Many are interlinked and can thus be grouped together. Therefore, in the following, we will

196 discuss and make recommendations for data organization within a group, which involves working
197 data governance, data documentation, data formats, including minimum information (metadata)
198 standards as well as archival (covering points 1, 2, 3, 4, 5 above). Many of these aspects,
199 especially data governance, fall into the **planning** section of the research data lifecycle, depicted
200 in Figure 6. Here, RDM practices are planned and documented in data management plans (DMP)
201 or data policies. They are then carried out and updated throughout the data **production** and
202 **analysis** sections of the data lifecycle.

203 Together, these points ensure data can be reused by others within the group and also prepare data
204 for publication and reuse by those outside of an organization or project. We then recommended
205 repositories based on discipline and/or data acquisition methods employed, and how to reference
206 this data within a journal article (covering point 6 above). This allows others to **access** and **reuse**
207 the data, restarting the data lifecycle (Figure 6). Lastly, we outline how large, interdisciplinary
208 projects can tie the individual group RDM together in a collaborative data management.



Figure 6: The research data lifecycle depicts the typical stages of research data throughout a project. These include the planning of the project, which encompasses detailed planning on which research data will be generated or re-used as well as how it will be stored during and archived after the project. The active research phases include the data production and analysis phases, after which the data are preserved and access rights are determined, such as open-access in a public repository or closed access in an institutional archive. Those who have access to the data can then re-use it in the next project. At this point, the planning stage restarts the cycle [12].

209 For the further discussion of these points, we will use the following use cases to illustrate the
210 recommendation. These examples outline the status quo for specific methods within CRC 985 in
211 the third funding phase:

Case 1: Infrared Spectroscopy	Case 3: Superresolution Fluorescence Microscopy
<p>Status Quo</p> <ul style="list-style-type: none"> • Small data output (Table 5) • Data processing only possible on device computer • Limited metadata captured when exported to an open format • ELN available (Chemotion ELN) • Networked to institute server <p>Desired Outcome</p> <ul style="list-style-type: none"> • Enable data processing and analysis on computers other than the device computer • Automatically link data to the digital sample documentation 	<p>Status Quo</p> <ul style="list-style-type: none"> • Large data output (Table 5) • Limited uniform metadata automatically generated • Predecessors data not always understandable • ELN available (eLabFTW) <p>Desired Outcome</p> <ul style="list-style-type: none"> • Ensure complete data documentation/metadata record • Link data to digital documentation • Appropriate storage solution for large data volume

213 These examples represent typical cases. Infrared spectroscopy (IR) produces relatively small
 214 data output (just over 10 MB, see Figure 5), which is representative of a large portion of the
 215 methods reported and therefore storage space is of little concern. The issue lies rather in ensuring
 216 data and full metadata are exported and linked to the sample documentation, while enabling data
 217 processing from anywhere, not just through the device computer. This case is fairly representative
 218 for spectroscopy in general.

219 Superresolution Fluorescence Microscopy (SRFM) imaging reaches the 150 GB mark per mea-
 220 surement (see optical microscopy in Figure 5), which poses a challenge to the institutional storage
 221 solutions in the long run. Furthermore, the raw data does not include the full measurement
 222 parameters, such as which device setup and specific accessories that may have been used. An
 223 ELN, eLabFTW, is available to manually enter these parameters. The full dataset cannot be
 224 directly attached to this type of documentation due to the file size limitations of the standard
 225 database storage. Therefore, ensuring complete metadata and other documentation, automati-
 226 cally transferring the data to an appropriate storage solution, and linking the (meta)data and
 227 documentation to the measurement and analysis data is desirable. Due to the output data size
 228 and the need for improved documentation, this case represents not only other imaging methods.
 229 Certain RDM solutions may also be extended to computational chemistry, for example, where
 230 storage and uniform documentation of input parameters play an important role.

231 3.3.1 Data Governance

232 A general uncertainty regarding which data to store, e.g., raw vs. processed files, and how to
233 organize the stored data was reported, especially due to a lack of guidelines in this area. Thus,
234 doctoral researchers often establish their own individual directory structure, documentation
235 practices, software tools to use, file and sample naming conventions, and workflows. While
236 this works for the individual in the short term, establishing a holistic data governance within a
237 research group planning phase enables wider collaboration as it provides structure and guidance.
238 Proper data organization, first and foremost, ensures that those currently working with the data
239 can do so efficiently. Furthermore, it enables others to easily understand and therefore reuse or
240 build upon the data, from future doctoral students in the same group to external researchers with
241 whom the data may be shared.

242 Starting in the planning phase of research, it must be determined where to store data and how
243 this should be structured. A common practice, observed during exchange with researchers, is for
244 the individual to sort data in a folder bearing their name. However, creating common, structured
245 folder templates for each project and storing data accordingly—instead of associating it with the
246 person conducting the research—ensures the data can be correctly found in the years to come.
247 Central, shared storage options, such as institutional servers or rented server space from the
248 university's central service providers, are recommended, while access to individual folders is
249 controlled on an administrative level.

250 It must be clear to all group members at what stages research data should be saved. For example,
251 as with the cases outlined in Section 3.3, certain IR devices produce raw data in proprietary
252 formats, while exported data may be used to continue work on the researcher's computer. Raw
253 data may not be transferred as it cannot be opened without the device software. However, best
254 practice is to always store raw data, even if in proprietary format, in read-only folders within the
255 given directory structure.

256 These agreed upon practices and structures should be documented in a group-wide DMP as well
257 as plain-text README files contained within the directory structure for easy reference. Further
258 data policies and on- and off-boarding checklists ensure data are transferred smoothly from one
259 researcher to the next.

260 This planning and documentation does not stop with data organization and storage, but should
261 also include other aspects that will arise in data production and analysis, such as data exchange
262 formats for storage as well as preservation and reuse, documentation tools and standards, as
263 well as data archival and publication platforms to ensure preservation, access, and re-use, the
264 specifics of which are discussed in the following.

265 In this phase, clear documentation of the processes and data-producing methods also proves
266 useful to better understand where improvement may be required. For example, a group-level
267 project can fully assess the status quo to determine where data workflows may be improved and
268 where external help may be required,

269 These efforts not only aid in managing research and the corresponding as a group, but also
270 provide a reference for (external data) stewards or data managers, e.g., those involved in INF
271 projects, while providing contextual information for data publication.

272 3.3.2 Data Documentation

273 As noted, doctoral researchers often individually establish documentation practices. In turn, it
274 was often mentioned, that understanding a predecessors' data and work proved difficult. This
275 indicates that common, group-level documentation standards need to be established.

276 Using the above SRFM case as an example, the raw data obtained from the device does not
277 necessarily contain all relevant measurement parameters. For IR, raw data files cannot be opened
278 without the device software, while full etadata are not exported with all available data exports.
279 Thus, as a bare minimum, establishing templates and even metadata schema in text-based formats
280 such as YAML or JSON provides a simple, machine and human-readable format that may be
281 filled out for each dataset. Such files can then be stored directly alongside the data to give a
282 digital metadata record. This practice may be extended to digitally record and document research,
283 thereby documenting agreed-upon minimum information for an experiment, measurement, or
284 sample, and by following existing community standards, where available. These templates
285 should be established in the planning phase of the research data lifecycle and updated, when
286 necessary, throughout the data production and analysis phases (see Figure 6).

287 Up until here, this and the [previous section](#) cover very basic data storage and management that
288 does not employ any specialized tools or infrastructure, besides a well-managed central storage,
289 defined directory structure, and documentation using agreed-upon templates. This provides
290 group members, especially junior scientists, with the basic framework to operate in an efficient
291 and organized manner, while producing transparent results that are (re)usable by other current
292 and future research group members. However, sophisticated tools and platforms exist, and are
293 being continuously updated and improved, to further assist researchers in effective research data
294 management.

295 In many natural sciences, the laboratory journal stands as the staple of research documentation.
296 However, analog journals are not machine-readable and do not necessarily follow uniform
297 documentation standards. Digital counterparts, ELNs, offer a powerful solution to documenting
298 research in a digital and structured manner, while also managing and connecting the associated
299 research data. These platforms exist with a wide variety of styles and target user groups, from
300 the more synthetic chemistry focused Chemotion ELN [13]–[15] to the broadly customizable
301 eLabFTW [16], [17]. One group within the CRC transitioned to Chemotion ELN after the survey
302 had been conducted, while limited use of eLabFTW was reported, yet in a rather individualized
303 manner. Proprietary solutions such as FURTHRmind and mbook were also employed. Many
304 CRC members reported using analog journals or solutions such as MS Word and MS Excel files,
305 as noted above.

306 For ELNs, it is important to continue to follow data organization and documentation best practices.
307 While some ELNs, such as the Chemotion ELN, strive to adhere to minimum information
308 standards for supported methods, highly customizable instances or unsupported methods require
309 high-level organization from within the group or institute. As with the templates outlined
310 above, groups or institutes should agree on the information to record for their experiments and
311 create templates for the ELN. eLabFTW, for example, enables custom metadata and allows
312 for the creation of experiment templates. Chemotion has recently also expanded to include
313 LabIMotion [18] which enables custom modules for non-chemistry or not yet included methods.

314 Therefore, an ELN must be centrally managed and documented within the group, analogous
315 to the basic data organization and storage outlined above. This not only includes providing
316 templates and usage guidelines, but also training group members on ELN use.

317 For the examples, the IR use case involves a research group that employs the Chemotion ELN.
318 The ELN offers direct connections for many methods, including IR, which directly transfers
319 data and attaches it to an experiment [19]. It also offers ChemSpectra to edit the analytical
320 data [20]. These methods extract necessary metadata to complete the documentation, ensuring
321 documentation, research data as well as the analysis are bundled in one place.

322 For the SRFM use case, eLabFTW is available, which allows for structured metadata templates
323 to be established within experiment templates. Since not all relevant metadata are captured in
324 a given measurement, researchers can employ such templates to document their research and
325 manually enter any missing information. However, as opposed to IR, attaching SRFM data to
326 experiments within the ELN is not viable due to size limitations. Therefore, creating meaningful
327 links to the data within the documentation proves helpful.

328 For cases such as this, where increased storage is required while metadata management is at
329 the forefront, the RWTH Aachen IT Center has developed Coscine (short for Collaborative
330 Scientific Integration Environment) [21], [22]. This platform primarily aims to organize and
331 manage working research data in ongoing projects. On a group level, Coscine offers various
332 storage types, called resources, with a storage quota of up to 125 TB per project for participating
333 universities or groups involved in NFDI-related projects. Custom metadata application profiles
334 can be generated to fit group needs, which result in a fillable metadata form that includes metadata
335 validation for input values. Data within a project or subproject is organized into resources, each
336 of which has been assigned a specific application profile and a PID in the form of an ePIC [23],
337 which leads to a contact page. Therefore, groups can customize their data documentation and
338 storage structure to fit their needs and incorporate community-specific minimum information
339 standards. Details pertaining to the collaborative aspects of this platform will be discussed in
340 Section 3.3.4.

341 Both eLabFTW and Coscine offer a Representational State Transfer Application Programming
342 Interface (REST API). Such interfaces allow for information to be exchanged between the
343 platforms in an automated manner. Therefore, to maintain the local documentation using the
344 ELN while maintaining a connection to the associated raw and processed data, a Python script on
345 the device computer can transfer the measurement data to Coscine, while a link is added within
346 the ELN entry. Metadata from the ELN is then also mirrored in Coscine.

347 Similar templates workflows may be setup for different methods to ensure the datasets include
348 complete documentation for all methods employed within the group. Working from a basis
349 of well-structured and well-documented data organization, including governance and research
350 data documentation, established during the planning phase and implemented during the data
351 production and analysis phases of the research data lifecycle (Figure 6), provides the foundation
352 for RDM in collaborative projects. Maintenance of these practices and proper onboarding of
353 group members ensures adherence and avoids uncertainty.

354 3.3.3 Data Formats

355 Vendor software typically directs data formats for output data, which may be proprietary. Inter-
 356 operable data requires open and standardized data formats, which do not (yet) exist for every
 357 method [24]. For many methods, open export formats such as TEXT and comma-separated values
 358 (CSV) were reported, however, the associated metadata may be lost or incomplete upon export,
 359 as indicated for IR, for example. Furthermore, while these formats may be machine-readable to a
 360 certain extent, they are not necessarily machine-*understandable* as they lack a defined structure
 361 and semantic annotation.

362 As standard open data exchange formats exist for certain analysis methods within the CRC and
 363 since many of them were not mentioned in the survey responses, we gathered recommendations
 364 and summarized these in Table 1, sourcing information from FAIRsharing [9] and NFDI4Chem's
 365 Knowledge Base [11], as well as the Chemotion Repository documentation [25].

366 This information has also been shared on the CRC 985 SharePoint along with the method
 367 information outlined above. Gathering this information specifically arose from communication
 368 over the common misconception that data should always be stored and published as CSV or
 369 TEXT files. Other options exist, may even be supported by vendor software, and simply lack
 370 awareness.

371 **Table 1:** Data exchange formats recommended by FAIRsharing, NFDI4Chem, and the Chemotion
 Repository for selected methods reported within CRC 985 and common data formats or
 file extensions reported throughout the project. Formats sourced from FAIRsharing.org
 are cited accordingly, while those listed on NFDI4Chem's Knowledge Base and the
 Chemotion Repository Documentation are denoted accordingly. We recommend the
 adoption of formats printed in bold font.

method	data exchange format or file extension recommended by NFDI4Chm, FAIRsharing, and Chemotion Repository	data exchange formats within CRC 985
Chromatography	ANDI-MS [26], CSV ^a , TXT ^a	CSV, PDF, .vdt, .gcd
Colorimetric or Fluorescence-based Assays	-	.ruc (raw), ASCII (export including metadata)
Computational Chemistry	CHARMM Card File Format (CRD) [27]	ASCII, .log, .cosmo, .energy, .out, .gjf, .xyz, CSV (processed)
Cyclic Voltammetry (CV)	TXT ^a	.nox
Electrophysiology (patch clamp)	(patch -	.dat

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Electron Paramagnetic Resonance Spectroscopy (EPR)	TXT ^a		.spe, TXT (export)
Elemental Analysis (EA)	TXT ^a		TXT
Energy-dispersive X-ray spectroscopy (EDX)	-		TXT, JPEG (export), PNG (export)
Fluorescence spectroscopy	JCAMP-DX^a		OPJ, FDS, TXT (export), PDF (export)
IR Spectroscopy (IR)	JCAMP-DX [28] ^a , AniML [29] ^b	An-	.ispd, TXT (export), PDF (export)
Mass Spectrometry (MS)	JCAMP-DX [28], AniML [29] ^b , mzML [30] ^a	An-	.d, .bad, Xcalibur Raw file, TXT, .jws
Mechanical Surface Analysis (nanoindentation)	-	(standard data model: CWA 17552:2020 [31])	TXT
Microscopy	OME-TIFF [32]		.nid, .spm, .jpg-qi-image, .jpg-qi-data, TIFF^e , LIF, DM4 (TEM), JPEG (export), PNG (export), AVI (video), CSV, TXT
Nuclear Magnetic Resonance Spectroscopy (NMR)	NMR-STAR [33], CCPN [34], NMR-ML [35], NMRe-Data [36] (assignments) ^a , AniML [29] ^b , JCAMP-DX (raw) ^a		.mrnova, FID, PDF (export)
Raman Spectroscopy	JCAMP-DX^a , AniML [29] ^b		.icRaman, .sps, TXT (export), CSV (export), .spc (export), .xlsx (export)
Rheometry	-		.rdf, .tri, .iwp, CSV (export)
Dynamic Light Scattering	CSV ^b		ASC^d , .dts, .zmes^d , CSV (export), TXT (export)
Static Light Scattering	-		.d80, .txt (export, not all parameters included)
Small Angle X-Ray Scattering (SAXS)	-		.mpa, .info, .edf, .dat
Spectroelectrochemistry	-		.str8, TXT (export)

(Continued)

Tensiometry	PNG (contact angle measurements) ^a	.krs, .zip (export, contains all .krs and XML) , ^d XLSX (export or analysis results)
Thermal Analysis	-	.stad, .spp, TXT (export), CSV (export)
UV/Vis Spectroscopy	CSV, ^a JCAMP-DX ^c	.dsw, .bsk, .bkn, .str, .jws, .jwb, .ksd, .sre (ASCII), TXT (export), CSV (export)
X-Ray Diffraction Analysis (XRD)	CIF [37] (single crystal), ^a .xyd (powder) ^a	binary encoded frames (images), .p4p, .hkl, .res, CIF, .x

^a NFDI4Chem Knowledge Base^b under development according to FAIRsharing.org^c Chemotion Repository^d (meta)data accessible by common tools374 ^e preferably method-specific TIFF-formats that include extended metadata

375 The existing standard data exchange formats listed in Table 1 provide guidelines on formats
 376 to choose from, while recommended standards and common formats are highlighted in bold
 377 font. The exact format choice for each method will depend on available software and export or
 378 conversion tools and also the format data types specific repositories will accept for publication
 379 (see, for example, Chemotion Repository requirements in [25], [38]).

380 Notably, many methods do lack specific standards, for which the above-mentioned practice of
 381 documenting data appropriately and sharing data along with the associated metadata in open,
 382 text-based formats is advised. As the various efforts such as the NFDI consortia continue their
 383 work, more standards will become available. Furthermore, minimum information standards
 384 will continue to direct how data should be formatted and documented, further guiding format
 385 standards. Table 1 as well as the published overview [8] serve to inform the standards and
 386 infrastructure community on which formats researchers are employing in their day-to-day work
 387 and where standards are lacking.

388 For the example case IR, as the connection can be made to Chemotion ELN, the data should be
 389 exported to JCAMP-DX as advised by not only the Chemotion Repository as denoted in Table 1,
 390 but also the Chemotion ELN to allow for automatic data transfer. This format was not reported,
 391 yet it is supported by the vendor software. For SRFM, OME-TIFF may prove beneficial by
 392 adapting an instance of Omero on an institutional or university level [39]. Without this option,
 393 TIFF files are appropriate. Connecting the documentation and data management, as described,
 394 ensures full metadata annotation, especially since Coscine enables semantic metadata.

395 As with data organization and documentation, data exchange formats must be agreed upon as
 396 part of the planning stage of the data lifecycle (Figure 6), communicated within the group, and
 397 updated as more standards become available.

398 3.3.4 Collaboration

399 Up until now, the discussion has focused on the group level. Having a well-documented approach
400 to data organization, documentation, and the tools used helps in identifying how collaborative
401 projects such as CRCs and the contained subprojects can best manage data.

402 The CRC 985 INF project addressed sample tracking throughout a collaborative project involving
403 many different groups and institutes in previous funding periods [1], as described in Section 1.
404 This system aimed to solve a specific problem with sample traceability within the project, while
405 enabling project members to directly attach associated data to a (digital) sample. In this funding
406 period, the system was further improved. As such, metadata fields for better sample tracking
407 were added, enabling users to define who initially created the sample and who was currently
408 working with it. The main view was altered according to user feedback to only show the most
409 relevant information. This enabled researchers to better find relevant samples and data.

410 However, as shown in Figure 4, some research within the CRC may not involve physical samples,
411 for example, computational chemistry methods such as molecular dynamics. Furthermore,
412 SharePoint relies on database storage that cannot accommodate larger datasets. It is therefore
413 not suitable for methods with large (raw) data output, e.g., SRFM and numerical methods (see
414 Figure 5). For these cases, other systems can provide the necessary solutions. It should also be
415 noted that the metadata describe the sample rather than any attached data, and therefore would
416 still require external documentation to fully describe the dataset belonging to the sample if not
417 included directly within in the files.

418 A central ELN instance, that is used by all the members of the CRC, could provide one solution,
419 yet, this did not prove realistic in this CRC for multiple reason, from varying user and group needs
420 to the lack of a centralized solution offered by the university. As individual groups and institutes
421 have indeed implemented ELNs, exchange formats between these could assist in collaborations
422 in such projects. This is a central goal of the ELN Consortium [40], which currently involves ten
423 ELN providers, including Chemotion ELN and eLabFTW.

424 The RDM platform Coscine, described in Section 3.3.2, is intended for collaborative work- Roll
425 management occurs on a project level, therefore, members can be given access to their respective
426 subproject, with all data relevant to the project collected and documented in one place. As
427 described, a REST API allows for automated data workflows, e.g., between local servers or ELNs
428 and Coscine. As such, metadata, data, and identifiers may be mirrored between platforms, giving
429 members a working-group agnostic option. As outlined for SRFM, its large storage capacity
430 assists researchers where institutional servers or systems that rely on a database structure such as
431 SharePoint and some ELNs reach their limits. As such, it has been employed within CRC 985
432 not only for SRFM, but for computational chemistry data as well as tensiometry.

433 An example of such an automated workflow would be transferring measurement data from a
434 folder on an institutional server, such as a device computer or research group server, to a central
435 RDM platform such as Coscine. A script would, in a given time interval, check for new data,
436 parse the file for relevant metadata, and use the Coscine's API to transfer the individual files and
437 assign metadata in a structured manner. Thus, the data becomes available for project members
438 on one centralized system in an automated manner, while similar workflows can pull relevant

439 data from Coscine to their local storage and RDM solutions.

440 Implementing solutions that employ such interfacing options require scripts or programs, or even
441 software development for more complex tasks. These should be maintained on a system such as
442 RWTH Aachen University's GitLab instance to facilitate access and collaboration. It should be
443 clear what resources are available, aside from the API itself, such as networked computers and
444 other available hardware, and who is responsible for deploying and maintaining these systems
445 within the research group or institute. Staff with development skills may also be required,
446 depending on the complexity of the solution. Due to updates in a given software's API, updates
447 to technical implementations may be required.

448 3.3.5 Data Publication and Archival

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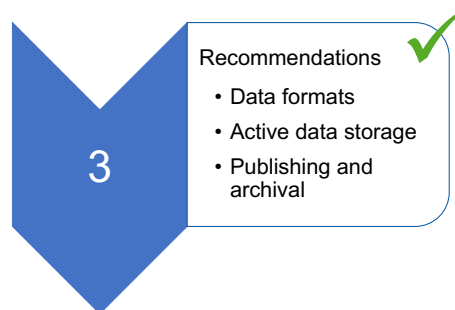
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Figure 7: Several recommendations could be made for active data storage, including data formats, documentation, and archival for a project on the scale of CRC 985.

repositories are preferred, as these are able to provide detailed metadata templates, enabling researchers to fully describe the published data. When using general or institutional repositories, adding as many (optional) metadata fields is best practice, while providing plain-text files for additional metadata and context. As institutional repositories may be used for reporting purposes, importing published datasets is also important, analogous to text publications.

Within these categories, we make the following recommendations for data sharing and archival in CRC 985 and similar projects, outlined in Table 2, which completes the final objective of this study (Figure 7). These were selected according to the methods reported within the conducted survey, the institutes involved in the CRC, while recommendations by NFDI4Chem [11] were preferred. Information on file sizes has been included to provide a reference as to which repository may accommodate larger data amounts for methods producing larger amounts of data.

Aside from facilitating research within groups as well as large projects, the aim to make data reusable according to FAIR also includes making the (meta)data available to others while describing how to access the data (Figure 6: Access and Re-Use). Therefore, a data policy was established during the second funding period [1], which stipulated that all data underlying a published journal article should be published as well.

Various options exist for such publications, with the three common categories being: (1) institutional repositories, (2) general repositories, and (3) community-specific repositories. Where possible, community-specific

475 **Table 2:** Repositories recommended for CRC 985 and projects with similar data types. Institutional repositories correspond to research institutes involved in the current project.

Repository (type)	Description [9]	Date Size Limits
Jülich DATA [41] (institutional)	A registry service to index all research data created at or in the context of Forschungszentrum Jülich, which may also be used to publish research data and software.	10 GB per file (depends on Dataverse installation); prefers links to larger datasets [42]
RWTH Publications Research Data [43] (institutional)	As part of the general RWTH Publications repository, data and software can be published by all RWTH Aachen University members and affiliates.	100 GB per file; 1 TB maximum over all files (gigamove) [44]
Chemotion Repository [45] (discipline-specific)	The repository supports the storage of data related to chemical samples or reactions, with a focus on data from synthetic and analytical work. While not a requirement, data may be submitted directly via the Chemotion ELN.	None; might limit it to 50 MB in future [46]

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Cambridge Structural Database (CSD) [47] (discipline-specific)	Established in 1965, the Cambridge Structural Database (CSD) is the a repository for small-molecule organic and metal-organic crystal 3D structures. Database records are automatically checked and manually curated by one of our expert in-house scientific editors. Every structure is enriched with chemical representations, as well as bibliographic, chemical and physical property information, adding further value to the raw structural data.	50 MB per file; 100 MB for the total size of files uploaded; exception for bigger files via email contact [48]
Inorganic Crystal Structure Database (ICSD) [49] (discipline-specific)	The world's largest database for fully determined inorganic crystal structures and contains the crystallographic data of published crystalline inorganic structures. Organometallic and theoretical structures have been added within the past years.	None; contact for file sizes > 10 TB [50]
ioChem-BD [51], [52] (discipline-specific)	IoChem-BD is a digital repository of Computational Chemistry and Materials results. A set of modules and tools aimed to manage large volumes of quantum chemistry results from a wide variety of broadly used simulation packages.	default 1 GB per upload; > 100 MB not to be uploaded by web interface [53]

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NOMAD Repository & Archive [54] (discipline-specific)	The NOMAD Repository and Archive stands for open access of scientific materials data. It enables the confirmatory analysis of materials data, their reuse, and repurposing. All data are available in their raw format as produced by the underlying code (Repository) and in a common, machine-processable, and well-defined data format (Archive).	32 GB per upload (maximum of 10 non-published uploads per user) [55]
RADAR4Chem [56], [57] (chemistry: general)	A low-threshold and easy-to-use service for sustainable publication and preservation of research data from all disciplines of chemistry. Currently, exclusive to publicly funded research institutions and universities in Germany.	10 GB per project [56]
Suprabank [58] (discipline-specific)	Curated, open resource for intermolecular interaction.	10 GB per user (can be adapted) [59]
zenodo [60] (general)	EU discipline-agnostic repository for data and other research results.	50 GB per data set [61]

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479 Certain repositories are also tied to ELNs, therefore providing direct data and metadata workflows.
 480 Going a step further, data may also be converted to standard open formats, as is the case with
 481 Chemotion ELN and Chemotion Repository, as mentioned in Section 3.2.

482 The published data should then be explicitly referenced via their DOI within the article using a
 483 data availability statement, which journals are increasingly requiring [62]. They may also be
 484 cited within the article itself. Especially in cases which involve multiple published datasets, this
 485 provides additional context for the reader.

486 As shown in Figure 5, much of the data volume falls into smaller sizes, with imaging and
 487 numerical methods requiring larger storage if all data were to be published. For these, the use
 488 of institutional repositories such as RWTH Publications Research data are the best option. For

489 some methods, such as Atomic Force Microscopy, not all extracted data must be published, yet
490 the scripts employed to do so could be. Hence, the data may be reproduced in the same manner
491 when needed, while the published data volume is held to a minimum in cases where repositories
492 limit quota. Otherwise, much of the produced can be published on subject-specific or general
493 chemistry repositories without too much concern for data volume. Furthermore, repositories
494 may offer more quota upon request.

495 In terms of data access control, most of the repositories mentioned offer embargo periods to
496 ensure the creators' first rights to the data. In addition, zenodo allows restricted access in cases
497 where data cannot be made public.

Research Data Repository Usage in CRC 985

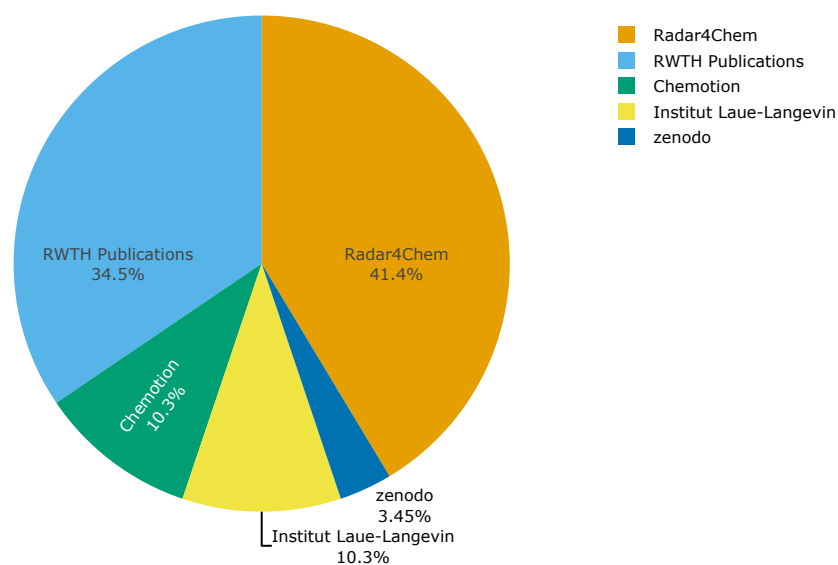


Figure 8: Research data repositories used to publish data underlying published articles in CRC 985. RADAR4Chem and RWTH Publications are widely used, followed by Chemotion and the institutional data repository for the Institut Laue-Langevin.

498 As shown in Figure 8, RADAR4Chem has proven itself as a readily-accepted data publication
499 platform, which may be attributed to its ease of use, the ability for data stewards to add standard
500 pre-filled metadata, as well as the recently-added notification system, allowing the INF project
501 to quickly respond to requests for dataset review. Institutional repositories found favor as well,
502 as RWTH Publications is used for 34.5% of data publications. Again, ease of use, but also a
503 certain trust in one's own services could be a strong factor here. For those using Chemotion
504 ELN, the direct publishing workflow to the Chemotion Repository considerably assists authors
505 in the publication process. In the example case for IR data, automated workflows from the
506 Chemotion ELN to the Chemotion Repository exist and enable simple data publication. Both
507 the Chemotion Repository and RADAR4Chem guarantee storage and accessibility for 10 years
508 or more, conforming with German Research Foundation (DFG) requirements; the data herein is
509 therefore successfully be deemed archived, while it can also be accessed and reused in accordance
510 with the research data lifecycle in Figure 6. RWTH Publications does not specifically list a

511 time span, but considers items published as archived as well. It should be noted that the Institut
512 Laue-Langevin carried out measurements for the CRC 985, the data for which is published on
513 the associated data repository, as indicated in Figure 8. This institutional data repository was
514 only omitted from Table 2 as only institutional repositories for direct participants were included.
515 , Typically, projects will amass more data than that, which has been published. This therefore
516 requires additional archive resources. For project members in CRC 985, the above-mentioned
517 Coscine also serves as an archiving space and may also be used where data access must be
518 controlled. It should be noted, however, that while the dataset PID may be used in a data
519 availability statement, the access restrictions should be stated. Furthermore, as the data has not
520 been published and received a DOI, it may not be cited.

521 The entire SharePoint, including the sample management system, will be archived under the
522 CRC's Coscine project, while members can gain access to the system to archive their data as
523 needed.

524 **3.4 Recommendations for Future CRCs and INF Projects**

525 The overarching role of INF projects within the CRC has largely been left out of the discussion
526 thus far. These central projects, however, can play a vital part in setting up and implementing
527 the above aspects.

528 Three aspects were identified within the CRC 985 INF project that should be considered for
529 future projects:

- 530 1. Support for project-wide data management plans and guidelines during project planning
531 stage
- 532 2. End-of-life plan for implemented infrastructure solutions
- 533 3. Sustainability of software solutions

534 To elaborate on 1., many workflows within research groups evolve naturally to fit the needs of
535 those carrying out much of the practical work, i.e., the individual doctoral researchers. However,
536 these tend to be highly individualistic and can be difficult to alter in order to streamline data
537 workflows. Therefore, providing clear guidelines on data organization and associated tools
538 is vital both within the group, but also across the project and should be established in the
539 planning phase. INF projects need to be involved at this stage and assist with infrastructure
540 planning and selection. Hence, overarching solutions can be available at the beginning of a
541 project to avoid implementing solutions and tools and altering workflows during ongoing work.
542 Individual workflows can then be developed within a given framework that facilitates data
543 storage, documentation, and exchange. This enables INF projects to focus on collaborative
544 workflows as opposed to improving individualized workflows, which proved difficult in CRC
545 985.

546 In terms of 2., the selected solutions require a detailed end-of-life management. It will not always
547 be possible to foresee which services and dependencies may become outdated over the lifetime
548 of a project. However, precautions and exit strategies to safeguard any and all data managed by
549 these services in a structured manner must exist.

550 As for 3., the software solutions developed by the INF project, e.g., data workflow scripts, should
551 be designed to outlive the project. The aspect of maintenance comes into play. Therefore, INF
552 projects should directly include individuals within the groups who are able to maintain these
553 solutions after the INF project is no longer available.

554 Overall, detailed, high-level planning for data management and the implementation of infrastruc-
555 ture solutions should involve INF projects at a very early stage of the project. Then, throughout
556 the project, members must be onboarded and continuously informed on common practices,
557 guidelines, and policies to ensure adherence.

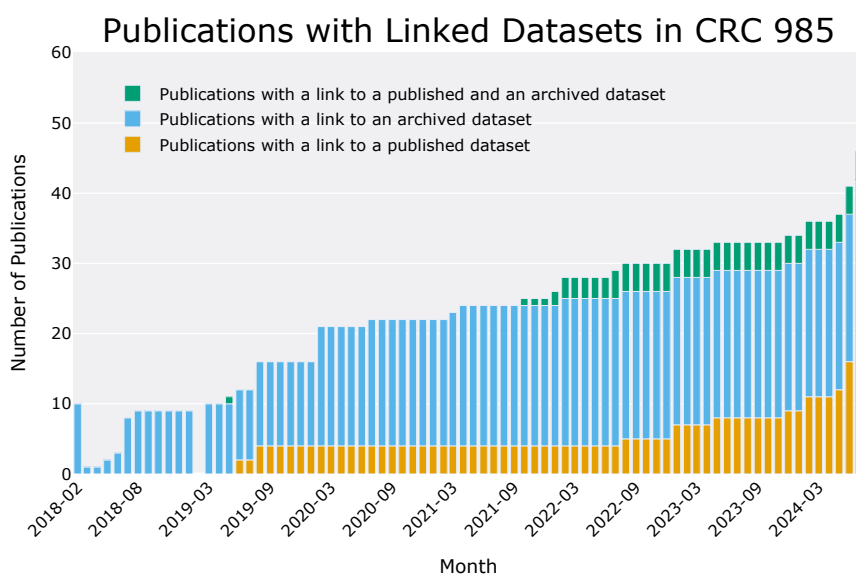


Figure 9: Publications with linked datasets according to RWTH Publications. Initially, linking archived (non-public) datasets was favorable in CRC 985, while publishing data becomes more common, especially in 2023 and 2024.

558 It should be noted that a readiness to publish data underlying published results generally exists
559 throughout CRC 985, especially in the third funding period. Figure 9 shows an increase in
560 (text) publications which are linked to a published dataset, especially in 2023 and 2024, while
561 archiving data in a non-public manner was preferred up until then. This data is recorded by
562 RWTH Publications, in which data as well as text publications within the CRC are recorded in
563 addition to its use as a data repository. This increase in text publications is likely due to general
564 changes in academic culture and awareness concerning data publication, but also the availability
565 of more platforms to easily do so. As noted in Section 3.3.5, RADAR4Chem, a service which
566 began in 2022, is greatly accepted. While its ease of use plays a role, the INF project also created
567 awareness of the repository.

568 For future INF projects, creating awareness of these platforms and workflows from the very
569 beginning should prove helpful, stressing their ease of use and how they conform to DFG
570 requirements on data publication and archival. INF members should be in exchange with

571 infrastructure providers to, on the one hand, stay up-to-date with developments, but also to
572 communicate researchers' requirements and expectations. This aids in increasing usability and
573 therefore acceptance, enabling researchers to make their data reusable.

574 4 Conclusion

575 Information on the data-producing methods and the associated data formats and data sizes in CRC
576 985 were collected in order to gain an overview of the diversity and derive RDM concepts and
577 structures for CRCs. The collected information is based on a structured survey, which collected
578 most of the details on the methods themselves, while formal as well as informal discussions in
579 various settings provided further feedback and deeper insight. Based on the information as a
580 whole, recommendations for this ongoing as well as future projects are made.

581 The gathered information paints a picture of the varied disciplines and the accordingly varied data
582 types and sizes. This underlines the need for standardized open exchange formats, as many of the
583 open export formats reported do not necessarily contain the required complete information in the
584 form of structured metadata to fully understand the acquired data. In order to assist in this, tools
585 from plain-text metadata templates to structured ELNs and data management platforms provide
586 essential machine-readable solutions for data documentation, assisting in data interoperability
587 and reuse.

588 The workflows and the RDM practices for each stage of the research data lifecycle (see Figure 6)
589 should be clearly defined and documented on a group level in advance. This information can
590 then feed into large projects such as CRCs, enabling informed decisions regarding RDM and
591 collaboration within the planning phase. In this way, data stewards within the INF project can
592 then establish policies, workflows, and infrastructures for collaboration within these institutional
593 frameworks while working closely with researchers.

594 For projects of the size of CRC 985, a one-size-fits-all solution, such as a uniform ELN and
595 repository where all (meta)data can be recorded in a well-structured manner, does not exist due
596 to the variety of analytical and experimental methods employed and the associated different
597 data formats and size requirements. Therefore, discipline-specific solutions found on a group
598 level require collaboration platforms that support RDM. Within CRC 985, Microsoft SharePoint
599 serves as collaboration platform, however, expectations regarding RDM evolved over the project
600 duration. FAIR data requires more structured and defined metadata on various levels. More
601 appropriate platforms for RDM have become available, including platforms such as the RWTH
602 Aachen University's Coscine as well as ELNs. This shows that, in addition to a minimum
603 standard which should be defined prior to the data production phase of the research data lifecycle
604 (see Figure 6), a certain flexibility should also be implemented to meet evolving requirements in
605 later funding periods.

606 With the requirement to publish all data underlying a text publication, ELNs and RDM platforms
607 can greatly assist researchers' workflows in FAIR data publication and archival in subject-specific
608 repositories by providing automated workflows. With much of this work still being in-progress
609 by infrastructure providers, future research projects will be able to greatly benefit, while current
610 work provides vital insight for these efforts.

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617 6 Roles and contributions

618 **Nicole A. Parks:** Conceptualization, Investigation, Writing, Visualization, Data Curation –
619 original draft

620 **Konstantin W. Kröckert:** Conceptualization, Investigation, Writing – original draft

621 **Fabian Claßen:** Conceptualization, Writing – original draft

622 **Walter Richtering:** Project Administration, Writing - review & editing

623 **Matthias Müller:** Project Administration, Writing - review & editing

624 **Sonja Herres-Pawlis:** Project Administration, Supervision, Writing – review & editing

625 References

- 626 [1] F. Claus, S. Kirchmeyer, M. S. Müller, and W. Richtering, “Das INF-Projekt im SFB 985
627 Funktionelle Mikrogele und Mikrogelsysteme,” *Bausteine Forschungsdatenmanagement*,
628 no. 2, pp. 104–111, Nov. 2019. DOI: [10.17192/bfdm.2019.2.8097](https://doi.org/10.17192/bfdm.2019.2.8097). (visited on
629 04/06/2023).
- 630 [2] M. Schröder, H. LeBlanc, S. Spors, and F. Krüger, “Intra-consortia data sharing platforms
631 for interdisciplinary collaborative research projects,” *it - Information Technology*, vol. 62,
632 no. 1, pp. 19–28, Feb. 2020, ISSN: 2196-7032, 1611-2776. DOI: [10.1515/itit-2019-
633 0039](https://doi.org/10.1515/itit-2019-0039). (visited on 02/03/2023).
- 634 [3] H.-J. Götze, M. Alten, H. Burger, *et al.*, “Data Management of the SFB 267 for the Andes
635 — from Ink and Paper to Digital Databases,” in *The Andes*, O. Oncken, G. Chong, G. Franz,
636 *et al.*, Eds., Springer Berlin Heidelberg, 2006, pp. 539–556, ISBN: 978-3-540-24329-8.
637 DOI: [10.1007/978-3-540-48684-8_26](https://doi.org/10.1007/978-3-540-48684-8_26). (visited on 03/08/2023).
- 638 [4] M. D. Wilkinson, M. Dumontier, I. J. Aalbersberg, *et al.*, “The FAIR Guiding Principles
639 for scientific data management and stewardship,” *Scientific Data*, vol. 3, no. 1, p. 160 018,
640 Mar. 2016, ISSN: 2052-4463. DOI: [10.1038/sdata.2016.18](https://doi.org/10.1038/sdata.2016.18). (visited on 01/27/2023).
- 641 [5] A. Kraft, *The FAIR Data Principles for Research Data*, Sep. 2017. (visited on 03/15/2023).
- 642 [6] N. Hartl, E. Wössner, and Y. Sure-Vetter, “Nationale Forschungsdateninfrastruktur (NFDI),”
643 *Informatik Spektrum*, vol. 44, no. 5, pp. 370–373, Oct. 2021, ISSN: 1432-122X. DOI:
644 [10.1007/s00287-021-01392-6](https://doi.org/10.1007/s00287-021-01392-6). (visited on 09/21/2023).

- 645 [7] C. Steinbeck, O. Koepler, F. Bach, *et al.*, “NFDI4Chem - Towards a National Research
646 Data Infrastructure for Chemistry in Germany,” *Research Ideas and Outcomes*, vol. 6,
647 e55852, Jun. 2020, ISSN: 2367-7163. DOI: [10.3897/rio.6.e55852](https://doi.org/10.3897/rio.6.e55852). (visited on
648 04/22/2022).
- 649 [8] N. A. Parks, K. Kröckert, F. Claßen, M. Müller, S. Herres-Pawlis, and W. Richterig, *Dataset*
650 *belonging to the publication "Data-Producing Methods in CRC 985: Recommendations*
651 *for Research Data Management in Large Interdisciplinary Projects"*, 2023. DOI: [10.22000/1793](https://doi.org/10.22000/1793).
652
- 653 [9] FAIRsharing | Home. [Online]. Available: <https://fairsharing.org/> (visited on
654 04/04/2023).
- 655 [10] Re3data - Registry of Research Data Repositories. DOI: [10.17616/R3D](https://doi.org/10.17616/R3D). [Online].
656 Available: <https://www.re3data.org/> (visited on 12/21/2022).
- 657 [11] NFDI4Chem Knowledge Base | NFDI4Chem Knowledge Base. [Online]. Available: <https://knowledgebase.nfdi4chem.de> (visited on 04/04/2023).
658
- 659 [12] FDM-Team RWTH Aachen University, *Research Data Life Cycle*, 2022. (visited on
660 04/27/2022).
- 661 [13] P. Tremouilhac, A. Nguyen, Y.-C. Huang, *et al.*, “Chemotion ELN: An Open Source
662 electronic lab notebook for chemists in academia,” *Journal of Cheminformatics*, vol. 9,
663 no. 1, p. 54, Sep. 2017, ISSN: 1758-2946. DOI: [10.1186/s13321-017-0240-0](https://doi.org/10.1186/s13321-017-0240-0). (visited
664 on 04/25/2022).
- 665 [14] PiTrem, *ComPlat/chemotion_ELN: Chemotion ELN 0.9.1*, Jun. 2021. DOI: [10.5281/ze
666 nodo.4899080](https://doi.org/10.5281/zenodo.4899080). [Online]. Available: <https://zenodo.org/record/4899080> (visited
667 on 01/26/2023).
- 668 [15] *Chemotion*. [Online]. Available: <https://eln.chemotion.net/home> (visited on
669 04/25/2022).
- 670 [16] N. Carpi, A. Minges, and M. Piel, “eLabFTW: An open source laboratory notebook for
671 research labs,” *The Journal of Open Source Software*, vol. 2, no. 12, p. 146, Apr. 2017,
672 ISSN: 2475-9066. DOI: [10.21105/joss.00146](https://doi.org/10.21105/joss.00146). (visited on 04/05/2023).
- 673 [17] *eLabFTW - Open Source Laboratory Notebook*. [Online]. Available: [https://www.ela
674 bftw.net](https://www.elabftw.net) (visited on 04/05/2023).
- 675 [18] *The LabIMotion Extension | Chemotion*, Mar. 2023. [Online]. Available: [https://chem
676 otion.net/docs/labimotion](https://chemotion.net/docs/labimotion) (visited on 04/17/2023).
- 677 [19] J. Potthoff, P. Tremouilhac, P. Hodapp, B. Neumair, S. Bräse, and N. Jung, “Procedures for
678 systematic capture and management of analytical data in academia,” *Analytica Chimica*
679 *Acta: X*, vol. 1, p. 100 007, Mar. 2019, ISSN: 2590-1346. DOI: [10.1016/j.acax.2019
680 .100007](https://doi.org/10.1016/j.acax.2019.100007). (visited on 05/02/2024).
- 681 [20] Y.-C. Huang, P. Tremouilhac, A. Nguyen, N. Jung, and S. Bräse, “ChemSpectra: A web-
682 based spectra editor for analytical data,” *Journal of Cheminformatics*, vol. 13, no. 1,
683 p. 8, Dec. 2021, ISSN: 1758-2946. DOI: [10.1186/s13321-020-00481-0](https://doi.org/10.1186/s13321-020-00481-0). (visited on
684 08/16/2024).

- 685 [21] M. Politze, F. Claus, B. D. Brenger, M. A. Yazdi, B. P. A. Heinrichs, and A. Schwarz,
686 “How to Manage IT Resources in Research Projects? Towards a Collaborative Scientific
687 Integration Environment,” *European journal of higher education IT*, 2020. DOI: [10.181](https://doi.org/10.18154/RWTH-2020-11948)
688 [54/RWTH-2020-11948](https://doi.org/10.18154/RWTH-2020-11948).
- 689 [22] *Coscine | The research data management platform*. [Online]. Available: <https://coscine.de/> (visited on 03/17/2023).
- 691 [23] *Persistent Identifiers for eResearch*. [Online]. Available: <https://www.pidconsortium.net/> (visited on 03/16/2023).
- 693 [24] D. Rauh, C. Blankenburg, T. G. Fischer, *et al.*, “Data format standards in analytical
694 chemistry,” *Pure and Applied Chemistry*, vol. 94, no. 6, pp. 725–736, Jun. 2022, ISSN:
695 1365-3075. DOI: [10.1515/pac-2021-3101](https://doi.org/10.1515/pac-2021-3101). (visited on 03/16/2023).
- 696 [25] *For Data Files | Chemotion*, https://chemotion.net/docs/repo/details_standards/files, Oct.
697 2023. (visited on 01/16/2024).
- 698 [26] FAIRsharing Team, *FAIRsharing record for: Analytical Data Interchange Protocol for*
699 *Chromatographic Data*. DOI: [10.25504/FAIRSHARING.D7795C](https://doi.org/10.25504/FAIRSHARING.D7795C). (visited on 03/29/2023).
- 700 [27] FAIRsharing Team, *FAIRsharing record for: CHARMM Card File Format*, 2015. DOI:
701 [10.25504/FAIRSHARING.7HP91K](https://doi.org/10.25504/FAIRSHARING.7HP91K). (visited on 03/29/2023).
- 702 [28] FAIRsharing Team, *FAIRsharing record for: Joint Committee on Atomic and Molecular*
703 *Physical data - working group on Data eXchange*, 2018. DOI: [10.25504/FAIRSHARING.V8NVE2](https://doi.org/10.25504/FAIRSHARING.V8NVE2). (visited on 03/29/2023).
- 705 [29] FAIRsharing Team, *FAIRsharing record for: Analytical Information Markup Language*,
706 2015. DOI: [10.25504/FAIRSHARING.6CS4BF](https://doi.org/10.25504/FAIRSHARING.6CS4BF). (visited on 03/29/2023).
- 707 [30] FAIRsharing Team, *FAIRsharing record for: Mz Markup Language*, 2015. DOI: [10.2550](https://doi.org/10.25504/FAIRSHARING.26DMBA)
708 [4/FAIRSHARING.26DMBA](https://doi.org/10.25504/FAIRSHARING.26DMBA). (visited on 03/29/2023).
- 709 [31] FAIRsharing Team, *FAIRsharing record for: CWA 17552:2020 Engineering materials*
710 *- Electronic data interchange - Instrumented indentation test data*. DOI: [10.25504](https://doi.org/10.25504/FAIRSHARING.5C379F)
711 [/FAIRSHARING.5C379F](https://doi.org/10.25504/FAIRSHARING.5C379F). (visited on 03/28/2023).
- 712 [32] FAIRsharing Team, *FAIRsharing record for: Open Microscopy Environment - Tagged Im-*
713 *age File Format*, 2015. DOI: [10.25504/FAIRSHARING.CQ8TG2](https://doi.org/10.25504/FAIRSHARING.CQ8TG2). (visited on 03/28/2023).
- 714 [33] FAIRsharing Team, *FAIRsharing record for: NMR Self-defining Text Archive and Retrieval*
715 *format*, 2015. DOI: [10.25504/FAIRSHARING.2CHXXC](https://doi.org/10.25504/FAIRSHARING.2CHXXC). (visited on 03/29/2023).
- 716 [34] FAIRsharing Team, *FAIRsharing record for: Collaborative Computing Project for NMR*,
717 2015. DOI: [10.25504/FAIRSHARING.AVW5Q](https://doi.org/10.25504/FAIRSHARING.AVW5Q). (visited on 03/29/2023).
- 718 [35] FAIRsharing Team, *FAIRsharing record for: Nuclear Magnetic Resonance Markup Lan-*
719 *guage*, 2015. DOI: [10.25504/FAIRSHARING.ES03FK](https://doi.org/10.25504/FAIRSHARING.ES03FK). (visited on 03/29/2023).
- 720 [36] FAIRsharing Team, *FAIRsharing record for: Nuclear Magnetic Resonance Extracted Data*
721 *Format*. DOI: [10.25504/FAIRSHARING.8AE3D0](https://doi.org/10.25504/FAIRSHARING.8AE3D0). (visited on 03/29/2023).
- 722 [37] FAIRsharing Team, *FAIRsharing record for: Crystallographic Information Framework*,
723 2015. DOI: [10.25504/FAIRSHARING.ZR52G5](https://doi.org/10.25504/FAIRSHARING.ZR52G5). (visited on 03/29/2023).

- 724 [38] P. Tremouilhac, P.-C. Huang, C.-L. Lin, *et al.*, “Chemotion Repository, a Curated Repository for Reaction Information and Analytical Data,” *Chemistry–Methods*, vol. 1, no. 1, pp. 8–11, 2021, ISSN: 2628-9725. DOI: [10.1002/cmtd.202000034](https://doi.org/10.1002/cmtd.202000034). (visited on 04/18/2023).
- 725
726
727
- 728 [39] C. Allan, J.-M. Burel, J. Moore, *et al.*, “OMERO: Flexible, model-driven data management for experimental biology,” *Nature Methods*, vol. 9, no. 3, pp. 245–253, Mar. 2012, ISSN: 1548-7091, 1548-7105. DOI: [10.1038/nmeth.1896](https://doi.org/10.1038/nmeth.1896). (visited on 08/20/2024).
- 729
730
- 731 [40] *The ELN Consortium*. [Online]. Available: <https://github.com/TheELNConsortium> (visited on 04/06/2023).
- 732
- 733 [41] Re3data.Org, “Jülich DATA,” 23 dataverses, 29 datasets, 1170 files, 2021. DOI: [10.17616/R31NJMYC](https://doi.org/10.17616/R31NJMYC). (visited on 04/06/2023).
- 734
- 735 [42] *Dataset + File Management — Jülich DATA documentation*. [Online]. Available: <https://apps.fz-juelich.de/fdm/staging/mode-of-access/user/dataset-management.html%5C#file-upload> (visited on 10/13/2023).
- 736
737
- 738 [43] Re3data.Org, “RWTH Publications Research Data,” 319 research datasets, 2018. DOI: [10.17616/R33N6J](https://doi.org/10.17616/R33N6J). (visited on 04/06/2023).
- 739
- 740 [44] *GigaMove - RWTH AACHEN UNIVERSITY IT Center - English*. [Online]. Available: <https://www.itc.rwth-aachen.de/cms/it-center/Services/Kollaboration/~smiiti/GigaMove/?lidx=1> (visited on 10/13/2023).
- 741
742
- 743 [45] FAIRsharing Team, *FAIRsharing record for: Chemotion repository*, 2018. DOI: [10.25504/FAIRSHARING.IAGXCR](https://doi.org/10.25504/FAIRSHARING.IAGXCR). (visited on 04/18/2023).
- 744
- 745 [46] *Frequently Asked Questions (FAQ) | Chemotion*, Oct. 2023. [Online]. Available: <https://chemotion.net/docs/repo/faq> (visited on 10/13/2023).
- 746
- 747 [47] FAIRsharing Team, *FAIRsharing record for: The Cambridge Structural Database*, 2015. DOI: [10.25504/FAIRSHARING.VS7865](https://doi.org/10.25504/FAIRSHARING.VS7865). (visited on 04/18/2023).
- 748
- 749 [48] *Deposit - The Cambridge Crystallographic Data Centre (CCDC)*. [Online]. Available: <https://www.ccdc.cam.ac.uk/deposit/upload> (visited on 10/13/2023).
- 750
- 751 [49] FAIRsharing Team, *FAIRsharing record for: Inorganic Crystal Structure Database*. DOI: [10.25504/FAIRSHARING.A95199](https://doi.org/10.25504/FAIRSHARING.A95199). (visited on 04/18/2023).
- 752
- 753 [50] D. A. Steudel, D. S. Rühl, D. R. Hinek, and S. Rehme, “Scientific Manual ICSD Database,”
- 754 [51] C. Bo, M. Alvarez, N. Lopez, F. Maseras, J. M. Poblet, and C. De Graaf, *ioChem-BD Find central service*, Nov. 2017. DOI: [10.19061/iochem-bd-find](https://doi.org/10.19061/iochem-bd-find). (visited on 04/18/2023).
- 755
- 756 [52] FAIRsharing Team, *FAIRsharing record for: ioChem-BD*, 2018. DOI: [10.25504/FAIRSHARING.LW6A1](https://doi.org/10.25504/FAIRSHARING.LW6A1). (visited on 09/21/2023).
- 757
- 758 [53] *Set upload limits — ioChem-BD documentation*. [Online]. Available: <https://docs.iochem-bd.org/en/latest/faqs/admin/setup-upload-limits.html> (visited on 10/13/2023).
- 759
760
- 761 [54] FAIRsharing Team, *FAIRsharing record for: NoMaD Repository*, 2018. DOI: [10.25504/FAIRSHARING.AQ20QN](https://doi.org/10.25504/FAIRSHARING.AQ20QN). (visited on 04/18/2023).
- 762

- 763 [55] *How to upload data — NOMAD Repository and Archive documentation*. [Online]. Avail-
764 able: <https://nomad-lab.eu/prod/rae/docs/upload.html> (visited on 10/13/2023).
- 765 [56] *RADAR4Chem | RADAR*. [Online]. Available: <https://radar.products.fiz-karlsruhe.de/de/radarabout/radar4chem> (visited on 04/18/2023).
- 767 [57] FAIRsharing Team, *FAIRsharing record for: RADAR*. DOI: [10.25504/FAIRSHARING.601A27](https://doi.org/10.25504/FAIRSHARING.601A27). (visited on 04/18/2023).
- 769 [58] FAIRsharing Team, *FAIRsharing record for: SupraBank*, 2018. DOI: [10.25504/FAIRSHARING.VJWUT7](https://doi.org/10.25504/FAIRSHARING.VJWUT7). (visited on 04/18/2023).
- 771 [59] *SupraBank*. [Online]. Available: https://suprabank.org/terms%5C_of%5C_service (visited on 10/13/2023).
- 773 [60] FAIRsharing Team, *FAIRsharing record for: Zenodo*, 2018. DOI: [10.25504/FAIRSHARING.WY4EGF](https://doi.org/10.25504/FAIRSHARING.WY4EGF). (visited on 04/18/2023).
- 775 [61] *Zenodo - Research. Shared*. [Online]. Available: <https://help.zenodo.org/faq/>
776 (visited on 10/13/2023).
- 777 [62] N. A. Parks, T. G. Fischer, C. Blankenburg, *et al.*, “The current landscape of author
778 guidelines in chemistry through the lens of research data sharing,” *Pure and Applied*
779 *Chemistry*, Feb. 2023, ISSN: 1365-3075. DOI: [10.1515/pac-2022-1001](https://doi.org/10.1515/pac-2022-1001). (visited on
780 04/06/2023).