# FAIR Service Descriptions: enriching life science SPARQL endpoints

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#### Abstract

SPARQL service descriptions allow for rich information schemas describing the data inside SPARQL endpoints. Rewriting information schema (re)-discovery queries to queries using an existing one can give major performance benefits. Rich service descriptions have many use cases beyond query rewriting.

#### Keywords

SPARQL, RDF, Information schema, Query rewriting

A significant challenge for users of SPARQL endpoints is discovering the shape and quantity of the data exposed inside them. The W3C standards for SPARQL allow for a Service Description (SD), enumerating the capabilities and capacities of SPARQL endpoint. The Swiss-Prot group provides extensive service descriptions for it's SPARQL endpoints: (https://hamap.expasy.org/sparql, https://beta.swisslipids.org/sparql,https://sparql.rheadb.org/sparql and https://sparql.uniprot.org/sparql).

A SD contain metadata about a SPARQL endpoint, such as when it was updated and which ontologies it uses. Such a SD can be seen as an information schema for a SPARQL endpoint. Using the Service Description [1], VoID [2] and VoID-Ext [3] vocabularies. We store these in in-dependent named graphs, which we always name as address of the SPARQL endpoint + ./well-known/void. e.g. https://sparql.rhea-db.org/.well-known/void.

FAIR SDs have many use cases, such as:

- Query optimization and dataset visualizations. The tool SPEX which generates entity relationship diagrams uses these in part if they are available.
- Generating ShACL files describing the shape of the data in a SPARQL endpoint.
- Generate APIs in languages such as R or Python to access the data in the SPARQL endpoint. To be demonstrated in the CHIST-ERA: Open Research Data - TRIPLE project.
- License and last updated information for FAIR data monitors.

As an example: a common SPARQL query people are thought to use is to discover how many distinct classes there are in a SPARQL endpoint shown in listing:1. For large datasets like UniProt this is a non-trivial. Imagine running it as a classical unix pipeline like listing:2. Then be surprised that this takes a few days to run if you have enough disk space and memory that

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is. This is because there are more than 140 billion distinct triples in UniProt. Of course having such a SD is not enough as the people who are used to using such queries won't change to use a different query on an "information schema" by default. This means we need to rewrite the query (listing:1) to a query in the form of (listing:3). Query rewriting needs to take into account variations in prefix, white-space and variable naming. We solve this by using a SPARQL parser from the RDF4j project use the abstract SPARQL algebra for the query matching and rewrite. The original query with is redirected to a new location with a new query (http 301).

```
Listing 1: "Count distinct classes used in a SPARQL endpoint."

SELECT (COUNT(DISTINCT ? class) AS ? classes)

WHERE { ? subject a ? class . }
```

```
Listing 2: "Simple pipeline to count the unique classes in an ntriples file."
sort -u all_triples_in_uniprot.nt | grep rdf:type | sort -u |
wc -1
```

Listing 3: "Rewritten SPARQL query to retrieve the count of the distinct classes in the endpoint." **SELECT** (COUNT(**DISTINCT** ?classesRaw) AS ?classes) **FROM** <http://sparql.uniprot.org/.well-known/void> **WHERE** { [] <http://rdfs.org/ns/void#class> ?classesRaw . }

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