# Fragalysis - #651 Fix data download

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Date	Who	Update
23/09	Duncan	Added estimate
05/10	Duncan	Document approved in meeting 05/10/2021

# Introduction

#### Issue #651

Hard to get the coordinates that you want.

#### Minimal:

- Currently visible structures
- Structures with tag

### Optional:

- Download builder dialog probably generate temporary (or hidden) tags
  - o (Maybe redundant, because the search dialog would stand in -
- Fix search tools (epic) #654

Always by specific endpoint (fully restful), with toggles to configure (like for the direct view URL)

- (might include setting up a restful endpoint to create tags)
- needs backend work for assembling zipfile on FE request

### **Assumption**

In this analysis, it is assumed that the user wants a subset of the information in the current download file (and not yet stuff from other tables). Once the basic process has been designed it could then be extended to pull, for example, other information from the Fragalysis database.

### As-Is (Current)

When a target is uploaded, it is in the form of a zip file. This is unzipped and processed in the backend to load the Fragalysis tables. Some files (pdb/map/bound) are copied into common folders and direct links to these files are stored in the *Protein* table - but not all files included are directly referenced. After the file is processed an archive is created and it is this that is downloaded in its entirety when the "download structures" link is pressed in Fragalysis. The archive is also available under the target name in the media directory.

The available download options will be provided by the direct links and information in *Protein* as well as linked information provided in the *Target* and *Molecule* tables .

#### *Protein* direct links to files:

The Protein table has direct links to files in the media directory - so from a list of Fragalysis protein codes/required types for a target, a file containing the files that these links reference could be assembled and returned in a zip file.

pdb\_info: varchar-255 bound\_info: varchar-255 cif\_info: varchar-255 mtz\_info: varchar-255 map\_info: varchar-255 diff\_info: varchar-255 event\_info: varchar-255 sigmaa\_info: varchar-255

#### Metadata:

In addition to the individual file links there is also a link to the metadata.csv file that was uploaded with the target set in table *Target*.

If *metadata* is requested, this file should also be included in the download, but as is, in its entirety. No attempt at this stage will be made to extract information for individual proteins - but the code will be factored so that this can be added later.

Note that the *target id* is available on the protein table.

### SDF, Smiles and Transformation matrix files:

When a protein is uploaded with the target loader, a related *Molecule* record is created if ligand information exists (Molecule can be read using it's foreign key: *prot\_id*) with contents extracted from an SDF file. Furthermore, the Smiles is also extracted from the sdf mol

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information and added to Molecule. The SDF and Smiles are desired to be part of the download.

The Transformation Matrix is not yet loaded into Fragalysis - but this is also desired to be in the download.

To simplify the solution, it has been agreed that links to the SDF file and Transformation Matrix (a CSV to be added to the Target upload) will be explicitly added to *viewer\_protein* so that they can be requested in the same way as the other file types.

If <smiles> is requested, then the *smiles* information in the, potentially multiple following #652, molecule/component records attached to the protein will be put in a text file created for each molecule.

### **Tags**

A tag can currently be created attached to a session project or a molecule. These have an additional\_info field that is simply json. If we can work out a structure that is understandable to the backend, then maybe an extract template could be stored within the additional info and referred to by the tag.

### **Proposed Solution**

### **Database Changes**

*Viewer\_protein* will be extended with the following fields:

sdf info: varchar-255

trans\_matrix\_info : varchar-255

### Target Uploader

The sdf file will now be explicitly saved in a new media subvolume 'sdfs' (as well as with the target as it is now) and have a link created in the *sdf\_info* field in the protein record.

If a transformation matrix file is supplied in the zip file, it will be explicitly saved in a new media subvolume (as well as with the target) and have a link created in the *trans\_matrix\_info* field in the protein record.

### Download API

A new API will be developed for this download - as a rough model I have based this on the existing *api/dicttocsv* API that was developed recently for Fragalysis.

#### The API will consist of:

- a POST form containing the requested fields/files. This will compile the zip file on the media server and return a link to it.
- A GET request that can be used to download the compiled zip file from the media server using the link.

### Structure of the zip file

The zip file will be structured as follows:

- Each type of file will have it's own folder in the structure and
- Each filename will generally start with the protein code. File names should be unique
  as they are made unique on the upload. I suggest the smiles text file have a suffix of
  .smi.
- The metadata.csv file (if requested) will be placed on the root folder (named metadata.csv).

### api/download\_structures

POST ../api/download structures/

 Construct a zip file based on the requested parameters and return a temporary download link for the file.

#### Params:

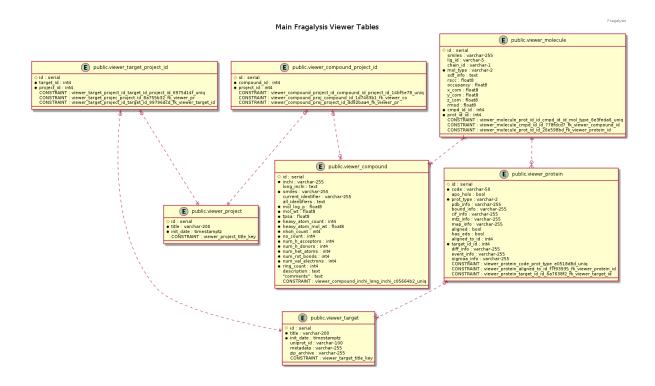
- Target name (string)
- List of Protein.codes (only the first part up to the colon will be used)
- pdb\_info: boolean when true, adds the file linked to in the *Protein.pdb\_info* field to the zip file in the relevant folder (if the link works and file is present)
- bound info : boolean ditto for Protein.bound info
- cif\_info : boolean ditto for Protein.cif\_info
- mtz info: boolean ditto for *Protein.mtz info*
- map info: boolean ditto for *Protein.map info*
- diff info : boolean ditto for Protein.diff info
- event\_info : boolean ditto for *Protein.event\_info*
- sigmaa\_info : boolean ditto for *Protein.sigmaa\_info*
- sdf\_info : boolean ditto for Protein.sdf\_info
- transformation matrix info : boolean ditto for Protein.trans matrix info
- metadata info: boolean add the metadata.csv file to the root folder.

If any dead links are found (i.e. links that are not spaces, but where the file does not exist), an *error.csv* file will be created with a list of Protein code/Dead link, so that these can be investigated.

### GET ../api/download\_structures/<download link>

- Returns the file from the download location.

# **Data Model**



Two new fields will be added to the *viewer\_protein* table.

### **Estimates**

Estimate includes API tests and documentation updates where appropriate.

I estimate the work described in this document to take around 5 days to complete.