



# Avergen Pharmaceuticals GmbH

Founded in 2015 in Martinsried, Bavaria

# Challenge establishing a drug discovery platform



- ▶ We are establishing a new small molecule drug discovery platform for identifying inhibitors of protein-protein interaction (PPI)
- ▶ PPI are hard to drug targets. But PPI play pivotal roles in various diseases. There are dozens of validated targets for which no effective inhibitors exist.
- ▶ Our focus is in oncology.
- ▶ We are collaborating globally with research institutes and CROs.

# Challenge establishing a drug discovery platform

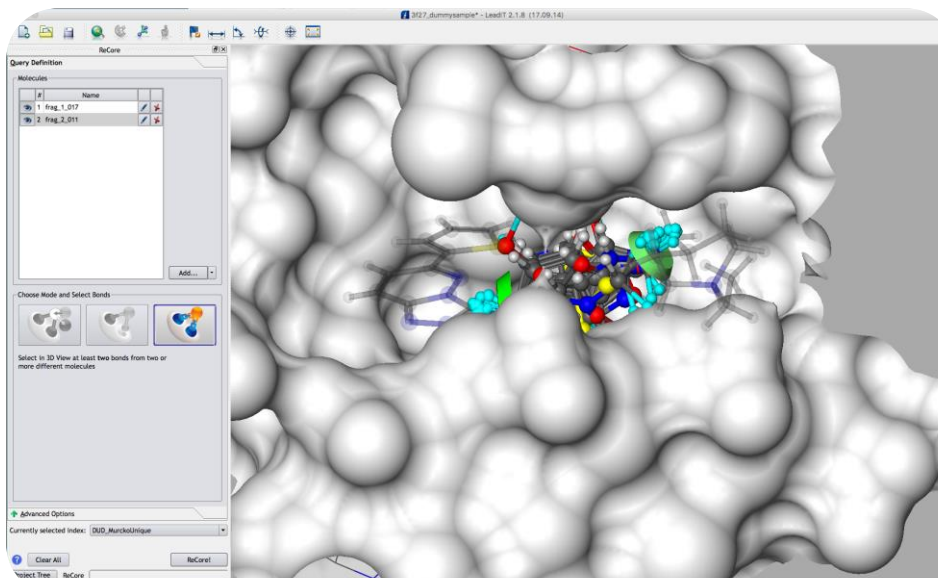
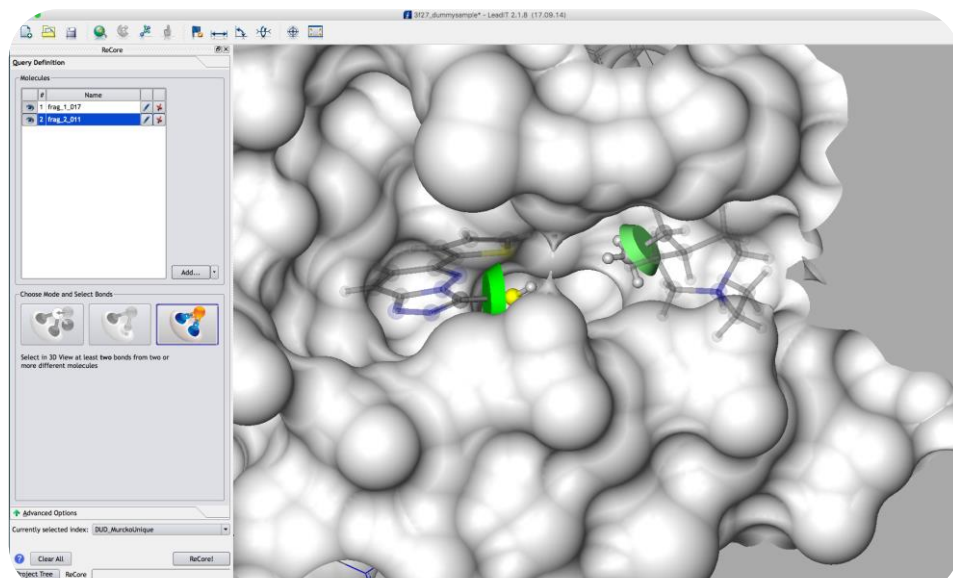


- ▶ Develop in a fast and economic way drug candidates for PPI drug targets which were until now thought undruggable.
- ▶ With the right software we could more efficiently select linkers for our fragments.
- ▶ This could reduce the number of linkers we have to synthesize and test in our screening assays.
- ▶ Lowering cost and speed up discovery time of new effective compounds.

# Automated linker selection with ReCore



- ▶ We link with ReCore fragments which bind the drug target.
- ▶ ReCore needs only two vectors to identify appropriate linkers.



# IT solution offered by BioSolveIT



- ▶ We created a new screening fragment data base according to specifications of our new platform technology.
- ▶ We started to create with BioSolveIT a new linker database:
  - ▶ We defined new shredding rules to generate linkers for ReCore.
  - ▶ The new linker database is optimized for our drug discovery platform. Thus the new linkers avoid solutions which do not fit into our platform specifications.
  - ▶ We filtered out unwanted chemistry moieties.

# Project progress and next steps



- ▶ We are redocking the new molecules in an iterative process and stepwise improve the shredding rules and fragment selection criteria. And thus improve our new linker and fragment data bases.
- ▶ The solutions are finally tested in our in vitro screening assays for their actual ability to bind the target structure and inhibit PPI.
- ▶ Finally, we want to use the new tools as part of our drug discovery platform to discover new PPI inhibitors for target structures which were until now regarded as “undruggable”.