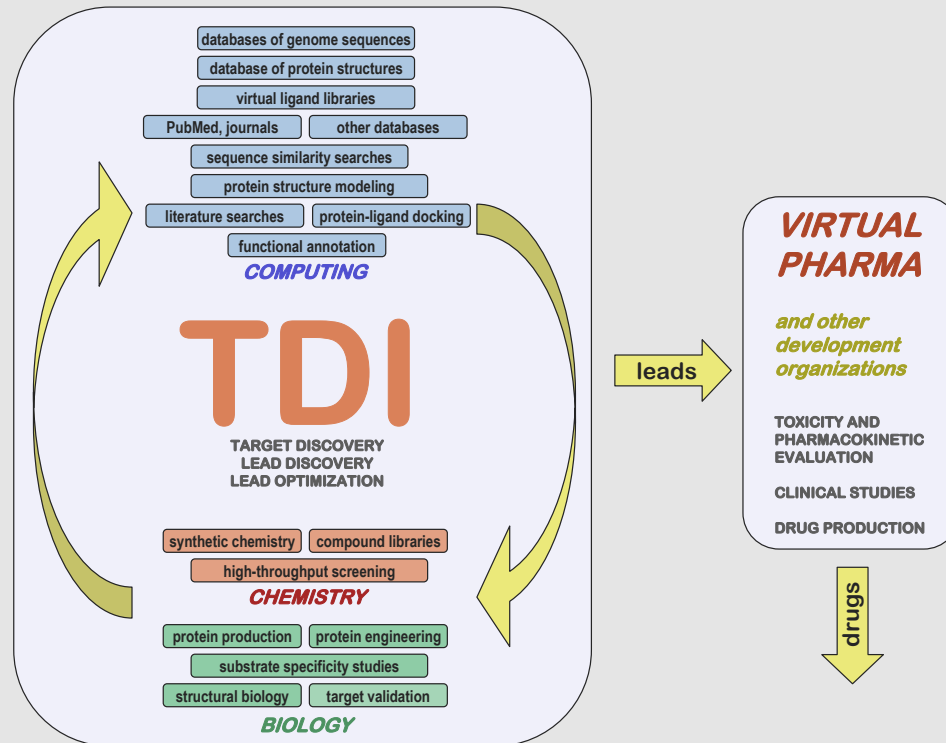


# The Tropical Disease Initiative

*Comparative docking on protein structure models from ten tropical disease genomes.*



Marc A. Marti-Renom

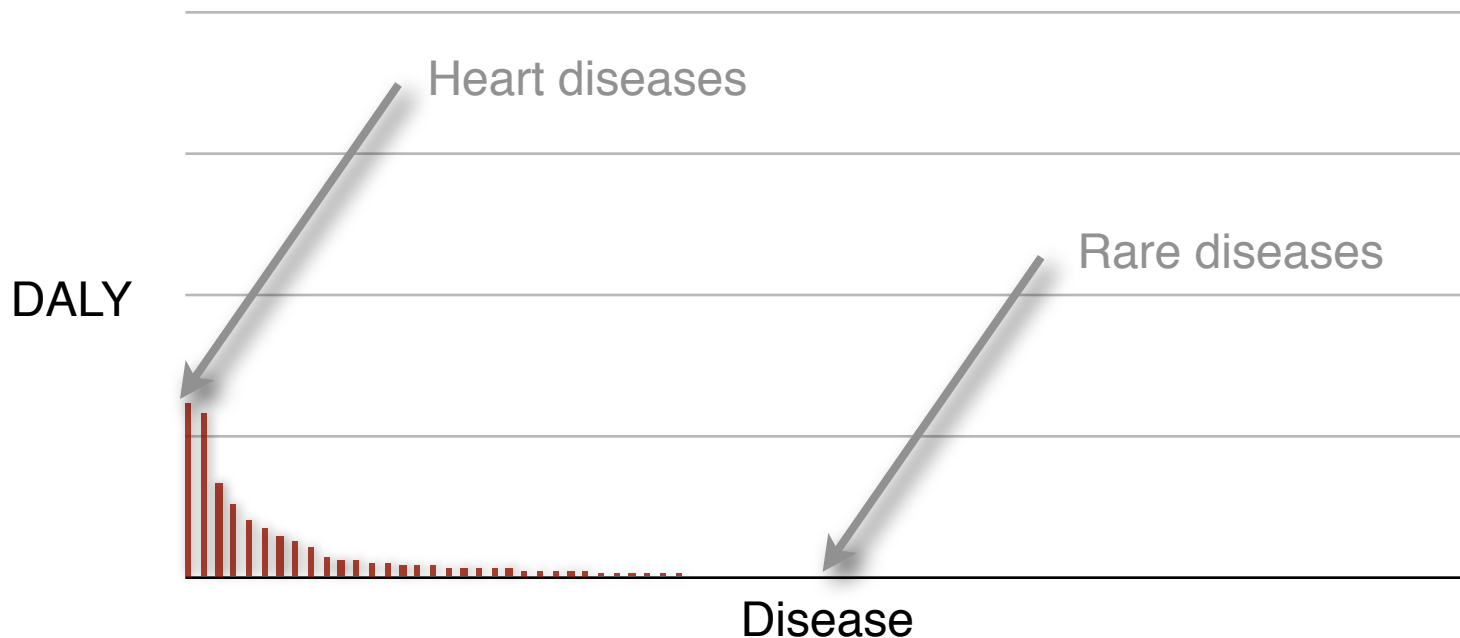
<http://bioinfo.cipf.es/sgu/>

Structural Genomics Unit  
Bioinformatics Department  
Prince Felipe Research Center (CIPF), Valencia, Spain



# Need is High in the Tail

- DALY Burden Per Disease in Developed Countries
- DALY Burden Per Disease in Developing Countries



Disease data taken from WHO, *World Health Report 2004*

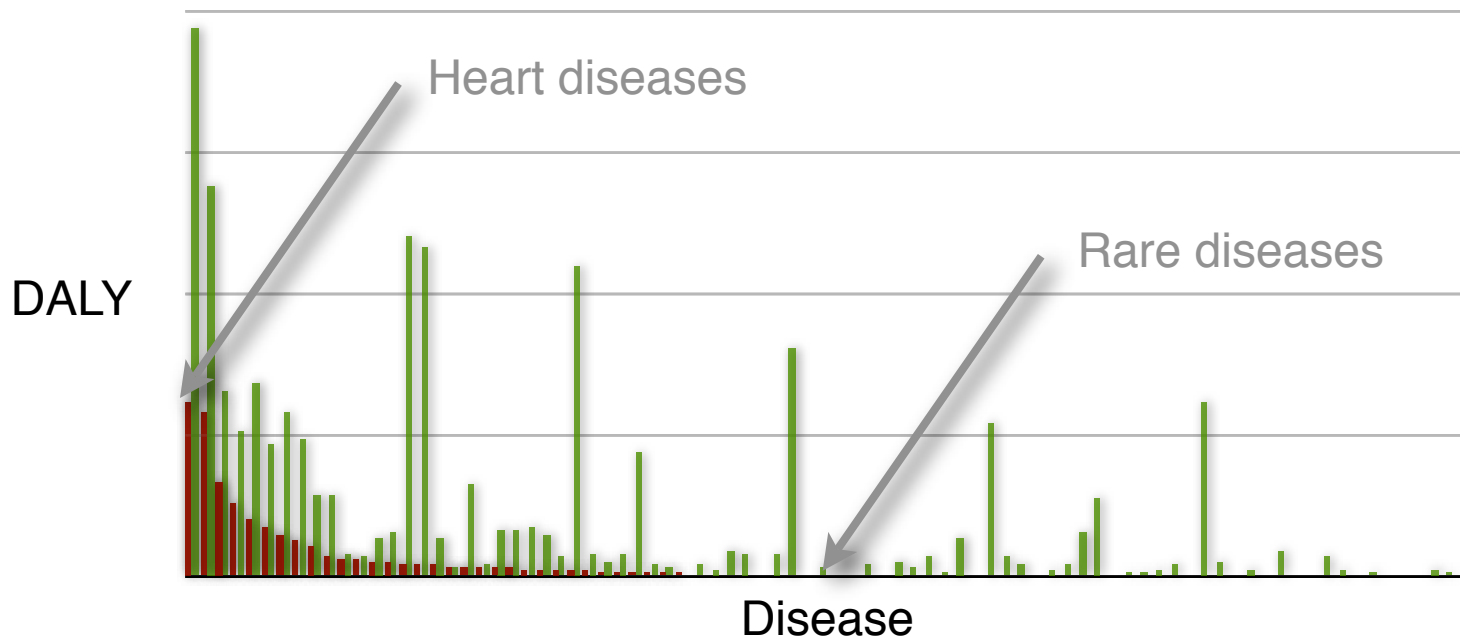
DALY - Disability adjusted life years

DALY is not a perfect measure of market size, but is certainly a good measure for importance.

*DALYs for a disease are the sum of the years of life lost due to premature mortality (YLL) in the population and the years lost due to disability (YLD) for incident cases of the health condition. The DALY is a health gap measure that extends the concept of potential years of life lost due to premature death (PYLL) to include equivalent years of 'healthy' life lost in states of less than full health, broadly termed disability. One DALY represents the loss of one year of equivalent full health.*

# Need is High in the Tail

- DALY Burden Per Disease in Developed Countries
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# “Unprofitable” Diseases and Global DALY (in 1000’s)

<b>Malaria*</b>	<b>46,486</b>
Tetanus	7,074
<b>Lymphatic filariasis*</b>	<b>5,777</b>
Syphilis	4,200
Trachoma	2,329
<b>Leishmaniasis*</b>	<b>2,090</b>
Ascariasis	1,817
<b>Schistosomiasis*</b>	<b>1,702</b>
<b>Trypanosomiasis*</b>	<b>1,525</b>

Trichuriasis	1,006
Japanese encephalitis	709
<b>Chagas Disease*</b>	<b>667</b>
<b>Dengue*</b>	<b>616</b>
<b>Onchocerciasis*</b>	<b>484</b>
<b>Leprosy*</b>	<b>199</b>
Diphtheria	185
Poliomyelitis	151
Hookworm disease	59

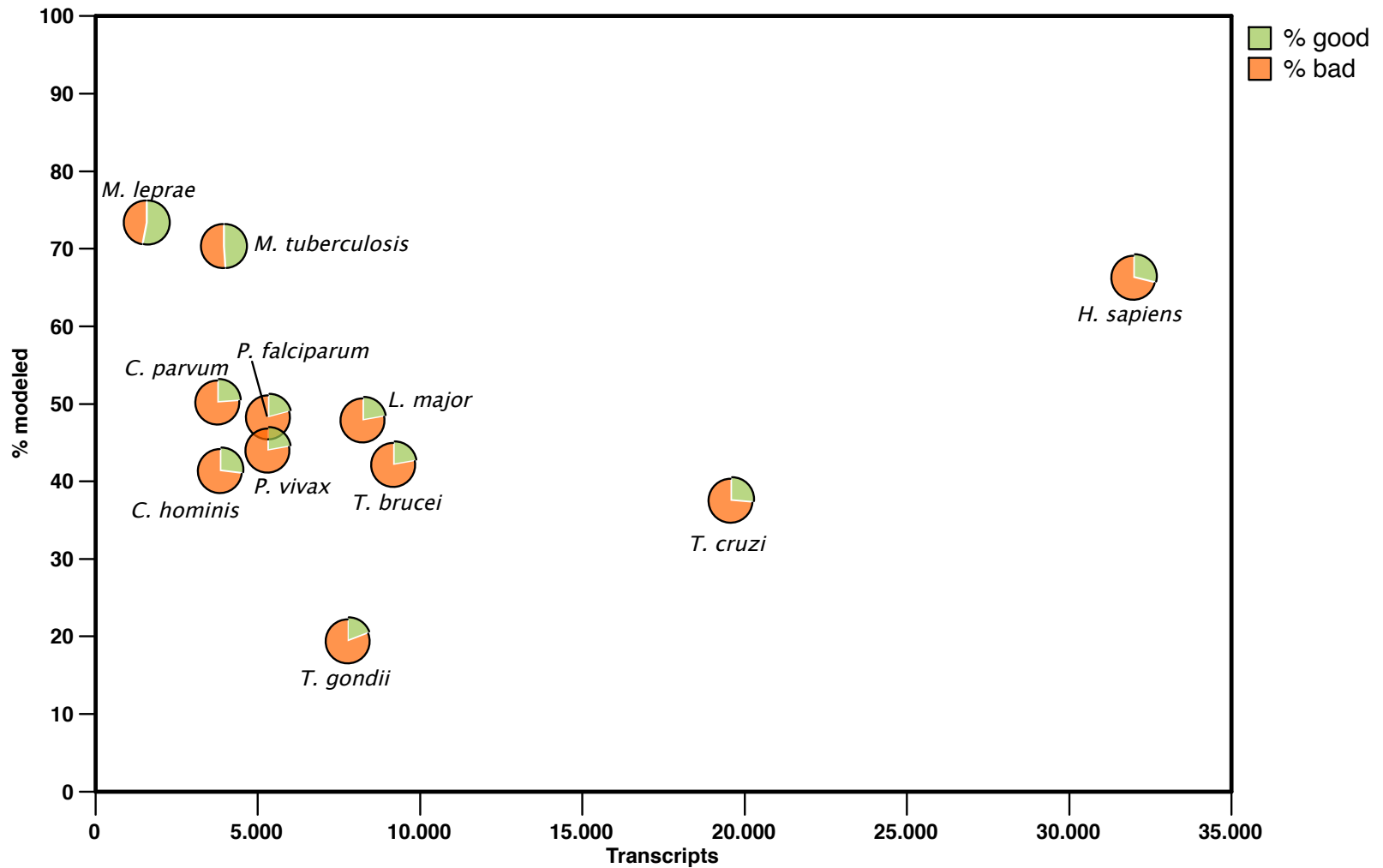
Disease data taken from WHO, *World Health Report 2004*

DALY - Disability adjusted life year in 1000’s.

\* Officially listed in the WHO Tropical Disease Research [disease portfolio](#).

# Modeling Genomes

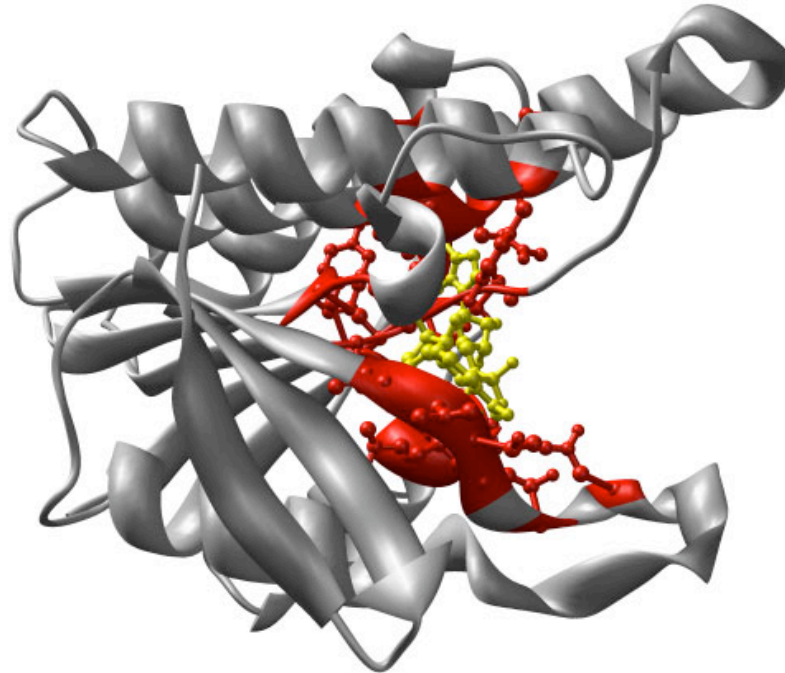
*data from models generated by ModPipe (Eswar, Pieper & Sali)*



*A good model has MPQS of 1.1 or higher*

# **AnnoLyze**

*Predicting binding sites in protein structure models.*



<http://bioinfo.cipf.es/sgu/services/DBAli/>  
<http://bioinfo.cipf.es/sgu/services/TDIModels/>

# DBAli<sub>v2.0</sub> database

<http://bioinfo.cipf.es/squ/services/DBAli/>

<http://www.salilab.org/DBAli/>

- ✓ Fully-automatic
- ✓ Data is kept up-to-date with PDB releases
- ✓ Tools for “on the fly” classification of families.
- ✓ Easy to navigate
- ✓ Provides tools for structure analysis

Does not provide a stable classification similar to that of CATH or SCOP

DBAli v2.0 home page

<http://salilab.org/DBAli/>

UCSF | Salilab | DBAli | MAMMOTH

## DBAli v2.0

Home

Search

Tools

Structural Genomics

Help

### DBAli. A Database of Structure Alignments.

[Marc A. Marti-Renom](#) and [Andrej Sali](#)

with the help of Madhusudhan M.S., E. Narayanan, B. Webb and A. Ortiz's [MAMMOTH](#) program.

This site contains an up-to-date all-against-all comparison of protein structures.

Currently, DBAli contains 1,210,877,705 pairwise structural alignments generated by [MAMMOTH](#) and family based multiple structure alignments for 28,184 non-redundant chains in PDB generated by [MODELLER](#). The database also includes several links to internal and external resources.

Use the links from the left frame to:

- Find structural relationships deposited in DBAli (Search)
- Analyze the data deposited in DBAli (Tools)
- Browse special pages for Structural Genomics
- Obtain help on how-to use DBAli (Help)

Site Map :: Reference :: Download :: Statistics :: Suggestions :: Report a problem Visitors: 6105 © 2000 - 2006 Marti-Renom

Pairwise structure alignments	
Last update:	July 5th, 2007
Number of chains:	93,307
Number of structure-structure comparisons:*	1,617,719,157
Multiple structure alignments	
Last update:	March 22nd, 2007
Number of representative chains:	31,848
Number of families:	11,900

Uses MAMMOTH for similarity detection

- ✓ VERY FAST!!!
- ✓ Good scoring system with significance

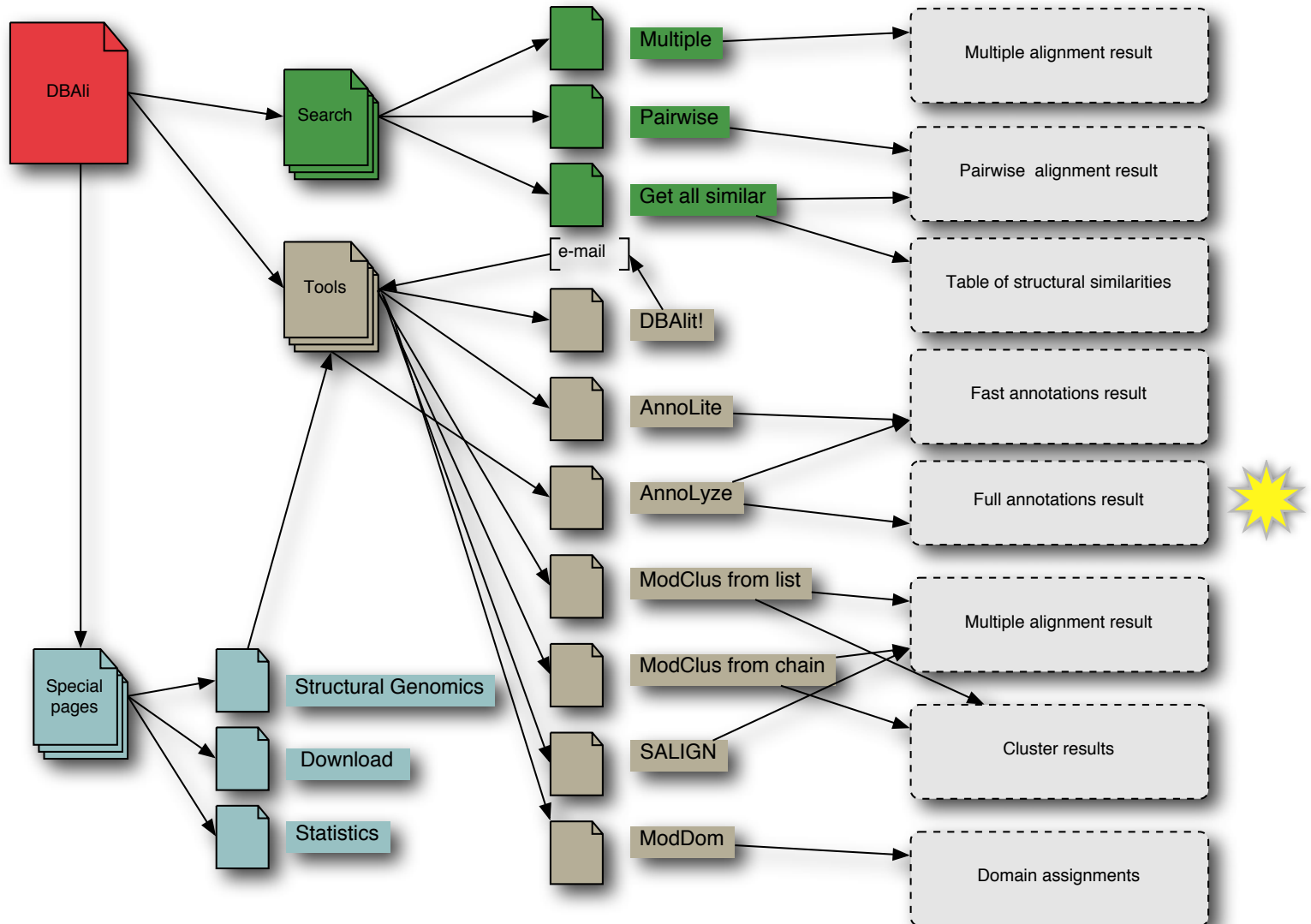
Ortiz AR, (2002) *Protein Sci.* 11 pp2606

Marti-Renom et al. 2001. *Bioinformatics.* 17, 746

# DBAli<sub>v2.0</sub> database

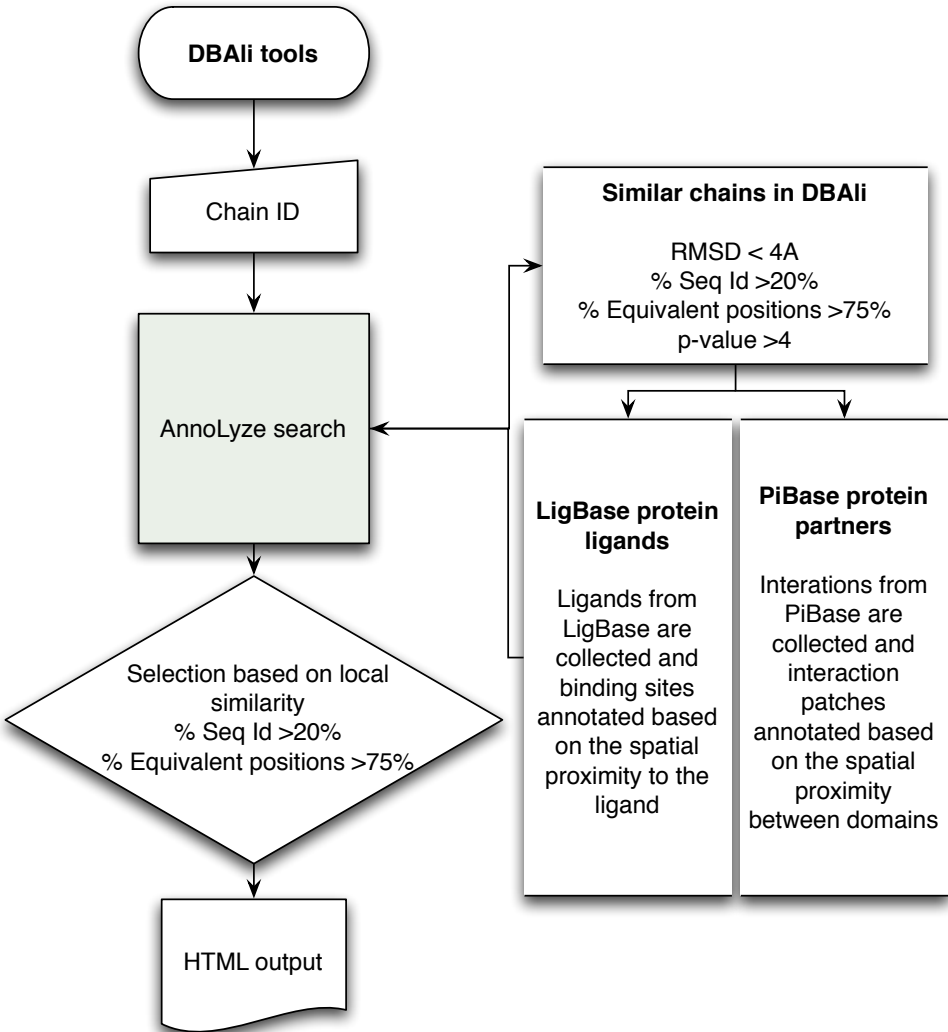
<http://bioinfo.cipf.es/squ/services/DBAli/>

<http://www.salilab.org/DBAli/>



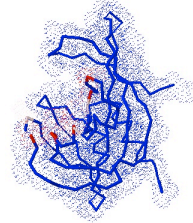


# Method



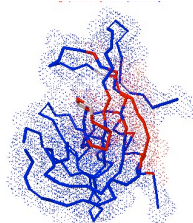
Inherited ligands: 4

Ligand	Av. binding site seq. id.	Av. residue conservation	Residues in predicted binding site (size proportional to the local conservation)
<a href="#">MO2</a>	59.03	<a href="#">0.185</a>	48 49 52 62 63 66 67 113 116
<a href="#">CRY</a>	20.00	<a href="#">0.111</a>	23 29 31 37 44 48 49 83 85 94 96 103 121
<a href="#">BOG</a>	20.00	<a href="#">0.111</a>	19 20 21 48 49 51 96 98 136
<a href="#">ACY</a>	15.87	<a href="#">0.163</a>	23 29 31 37 44 45 81 83 85 94 96 98 103 121 135



Inherited partners: 1

Partner	Av. binding site seq. id.	Av. residue conservation	Residues in predicted binding site (size proportional to the local conservation)
<a href="#">d.113.1.1</a>	23.68	<a href="#">0.948</a>	19 20 50 51 52 53 54 55 56 57 58 77 78 79 80 81 82 83 84 85 93 95 97 99 134 135 138 142 145



# Sensitivity .vs. Precision

	Optimal cut-off	Sensitivity (%) Recall or TPR	Precision (%)
Ligands	30%	71.9	13.7

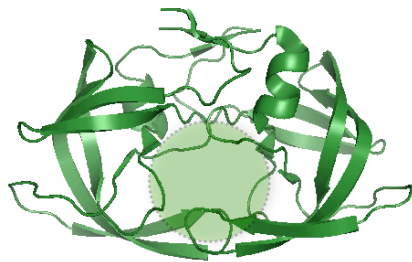
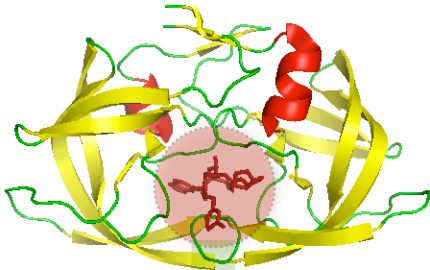
$$\text{Sensitivity} = \frac{TP}{TP + FN} \quad \text{Precision} = \frac{TP}{TP + FP}$$

**~90-95% of residues correctly predicted**

# Comparative docking

## 1. Expansion

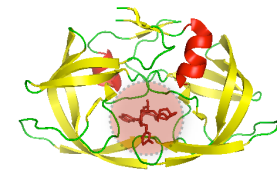
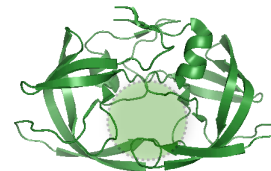
co-crystallized protein/ligand



crystallized protein

## 2. Inheritance

model



template

# Summary table

models with inherited ligands

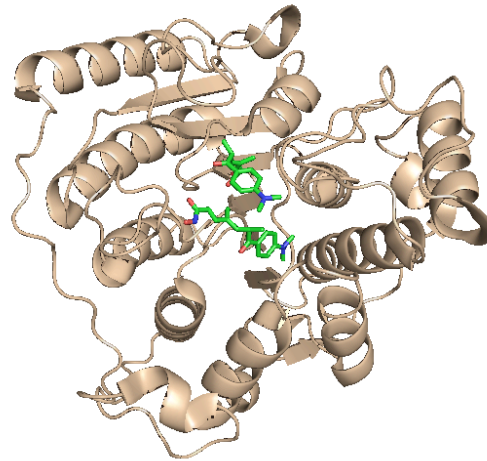
from 16,284 good models, 295 inherited a ligand/substance with at least one compound already approved by FDA and ready to be used from ZINC

	Transcripts	Good	Ligands	Lipinski	Lipinski+ZINC	FDA+ZINC
<i>C. hominis</i>	3,886	886	183	131	28	12 (10)
<i>C. parvum</i>	3,806	949	219	145	30	12 (10)
<i>L. major</i>	8,274	1,845	488	334	84	44 (34)
<i>M. leprae</i>	1,605	1,321	286	189	39	29 (25)
<i>M. tuberculosis</i>	3,991	2,887	404	285	71	44 (37)
<i>P. falciparum</i>	5,363	1,057	271	191	48	20 (16)
<i>P. vivax</i>	5,342	1,042	267	177	37	18 (15)
<i>T. brucei</i>	921	1,795	440	309	94	46 (36)
<i>T. cruzi</i>	19,607	3,915	730	493	127	62 (52)
<i>T. gondii</i>	7,793	587	174	124	28	8 (7)
<b>TOTAL</b>	<b>60,588</b>	<b>16,284</b>	<b>3,462</b>	<b>2,378</b>	<b>586</b>	<b>295 (242)</b>

# Example of inheritance (expansion)

*LmjF21.0680 from L. major “Histone deacetylase 2” (model 1)*

*Template 1t64A a human HDAC8 protein.*



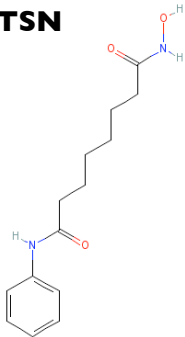
	Origen	Formula	Name	Cov.	Seq. Id. (%)
<b>ZN</b>	X-ray	Zn <sup>2+</sup>	Zinc ion	--	--
<b>NA</b>	X-ray	Na <sup>+</sup>	Sodium ion	--	--
<b>CA</b>	X-ray	Ca <sup>2+</sup>	Calcium ion	--	--
<b>TSN</b>	X-ray	C <sub>17</sub> H <sub>22</sub> N <sub>2</sub> O <sub>3</sub>	Trichostatin A	--	--
<b>SHH</b>	Expanded	C <sub>14</sub> H <sub>20</sub> N <sub>2</sub> O <sub>3</sub>	Octadenioic acid hydroxyamide phenylamide	100.00	83.8

# Example of inheritance (inheritance)

*LmjF21.0680 from L. major "Histone deacetylase 2" (model 1)*

	Formula	Name	Cov.	Seq. Id. (%)	Residues
<b>TSN</b>	C <sub>17</sub> H <sub>22</sub> N <sub>2</sub> O <sub>3</sub>	Trichostatin A	100.00	90.9	90 131 132 140 141 167
<b>SHH</b>	C <sub>14</sub> H <sub>20</sub> N <sub>2</sub> O <sub>3</sub>	Octadenioic acid hydroxyamide phenylamide	100.00	90.9	169 256 263 293 295

**TSN**



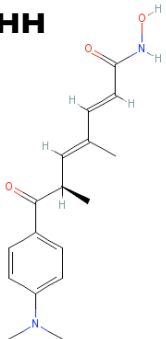
## suberoylanilide hydroxamic acid

**Pharmacological Action:**

[Anti-Inflammatory Agents, Non-Steroidal](#)  
[Antineoplastic Agents](#)  
[Enzyme Inhibitors](#)  
[Anticarcinogenic Agents](#)

Inhibits histone deacetylase 1 and 3

**SHH**



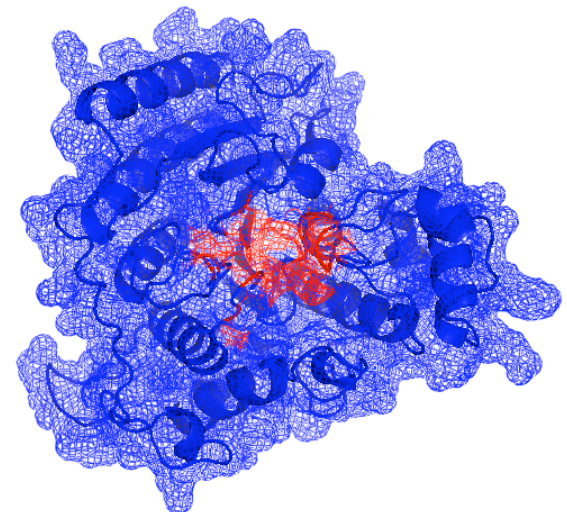
## trichostatin A

**Pharmacological Action:**

[Antibiotics, Antifungal](#)  
[Enzyme Inhibitors](#)  
[Protein Synthesis Inhibitors](#)

chelates zinc ion in the active site of histone deacetylases, resulting in preventing histone unpacking so DNA is less available for transcription

	LmjF21.0680.1.pdb
<b>Template</b>	1t64A
<b>Seq. Id (%)</b>	<b>38.00</b>
<b>MPQS</b>	1.47



# Example of inheritance (CDD-Roos-literature)

*LmjF21.0680 from L. major “Histone deacetylase 2” (model 1)*

*Proc. Natl. Acad. Sci. USA*  
Vol. 93, pp. 13143–13147, November 1996  
Medical Sciences

## **Apicidin: A novel antiprotozoal agent that inhibits parasite histone deacetylase**

(cyclic tetrapeptide/Apicomplexa/antiparasitic/malaria/coccidiosis)

SANDRA J. DARKIN-RATTRAY\*<sup>†</sup>, ANNE M. GURNETT\*, ROBERT W. MYERS\*, PAULA M. DULSKI\*,  
TAMI M. CRUMLEY\*, JOHN J. ALLOCCO\*, CHRISTINE CANNOVA\*, PETER T. MEINKE<sup>‡</sup>, STEVEN L. COLLETTI<sup>‡</sup>,  
MARIA A. BEDNAREK<sup>‡</sup>, SHEO B. SINGH<sup>§</sup>, MICHAEL A. GOETZ<sup>§</sup>, ANNE W. DOMBROWSKI<sup>§</sup>,  
JON D. POLISHOOK<sup>§</sup>, AND DENNIS M. SCHMATZ\*

Departments of \*Parasite Biochemistry and Cell Biology, <sup>‡</sup>Medicinal Chemistry, and <sup>§</sup>Natural Products Drug Discovery, Merck Research Laboratories, P.O. Box 2000, Rahway, NJ 07065

ANTIMICROBIAL AGENTS AND CHEMOTHERAPY, Apr. 2004, p. 1435–1436  
0066-4804/04/\$08.00+0 DOI: 10.1128/AAC.48.4.1435–1436.2004  
Copyright © 2004, American Society for Microbiology. All Rights Reserved.

Vol. 48, No. 4

## **Antimalarial and Antileishmanial Activities of Aroyl-Pyrrolyl-Hydroxyamides, a New Class of Histone Deacetylase Inhibitors**

# Models database

<http://bioinfo.cipf.es/sgu/services/TDIModels/>

The screenshot displays the TDIModels server interface. The browser address bar shows the URL <http://bioinfo.cipf.es/sgu/services/TDIModels/index.p>. The page title is "The TDIModels server".

**TDIModels**

Results for O96526 [O96526 Cdc2-related kinase (Cell division related protein )]  
Number of models: 2

**Model 1:**

JMOL   
This model has 1 predicted ligands.

Lipinski ZINC FDA Coverage	Seq. Id.
<input checked="" type="checkbox"/> NO3	100.00 100.00

SEQUENCE IDENTITY: 58.00  
MODPIPE QUALITY SCORE: 1.73  
TEMPLATE PDB: 1gz8  
TEMPLATE CHAIN: A  
TARGET LENGTH: 311  
TARGET BEGIN: 20  
TARGET END: 309  
[Download PDB file](#)

**Model 2:**

JMOL   
This model has 2 predicted ligands.

Lipinski ZINC FDA Coverage	Seq. Id.
<input type="checkbox"/> NO3	100.00 100.00
<input type="checkbox"/> KCX	100.00 93.75

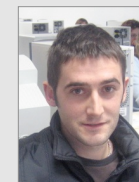
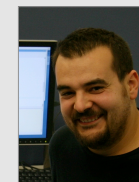
SEQUENCE IDENTITY: 29.00  
MODPIPE QUALITY SCORE: 1.13  
TEMPLATE PDB: 2cn5  
TEMPLATE CHAIN: A  
TARGET LENGTH: 311  
TARGET BEGIN: 1  
TARGET END: 311  
[Download PDB file](#)

[<- new search](#)

HELP:



# Acknowledgments



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NY-SGXRC

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Alfonso Valencia (CNB/UAM)

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