index.php?tab=DOCKING&page=RESULTS&jobId=6wzu_RoccustyrnaTM_gs_convs_1XX2XXpdb_62b306a...

Table

Rank	File ID	Compound	Affinity	Total Energy	vdW Energy	Elec. Energy
1	c359152ed8	ligand 1	-4.830	-86.676	-0.000	0.113
		run 4	-4.830	-86.676	-0.000	0.113
		run 4	-4.830	-86.675	-0.000	0.114
		run 4	-4.830	-86.675	-0.000	0.114
2	6f9df35049	ligand 1	-4.657	-28.226	-0.000	0.036
		run 12	-4.657	-28.226	-0.000	0.036
		run 12	-4.657	-28.226	-0.000	0.036
		run 12	-4.657	-28.226	-0.000	0.036

« 1 »

3D View

Use your mouse to drag, rotate, and zoom in and out of the structure:

`r` *auto view* `shift + left_mouse + drag` *zoom* `ctrl + right_mouse + drag` *rotate*

index.php?tab=DOCKING&page=RESULTS&jobId=6wzu_RoccustyrnaTM_gs_convs_1XX2XXpdb_62b306a...

« 1 »

3D View

Use your mouse to drag, rotate, and zoom in and out of the structure.

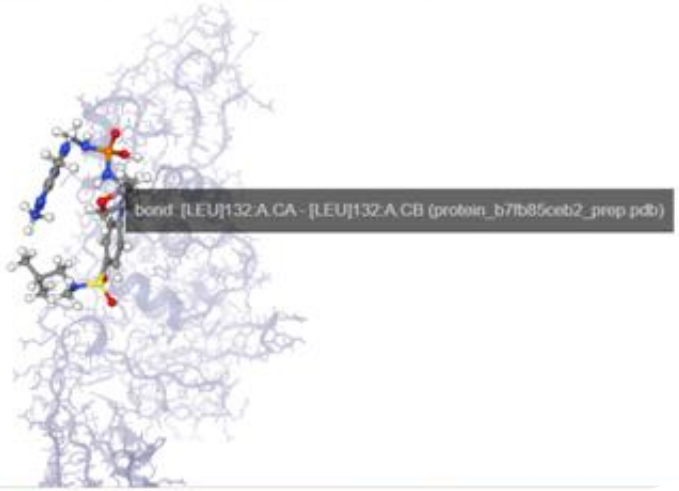
`r` *auto view* `shift + left_mouse + drag` *zoom* `ctrl + right_mouse + drag` *rotate*

index.php?tab=DOCKING&page=RESULTS&jobId=6wzu_RoccustyrnaTM_gs_convs_1XX2XXpdb_62b306a...

run 12	-4.657	-28.226	-0.000	0.036
run 12	-4.657	-28.226	-0.000	0.036

« 1 »

3D View

Use your mouse to drag, rotate, and zoom in and out of the structure.

`r` *auto view* `shift + left_mouse + drag` *zoom* `ctrl + right_mouse + drag` *rotate*