ConSurf Job Information 10.03.2023 16:43

ConSurf Job Information

Go to Results of Job 1678459912 (until they are deleted from the server).

The information below is obtained from the header of PDB file AF vglut2NAT rank1 with segres consurf firstglance.pdb This PDB file can be saved to your computer.

Job Parameters

Job 1678459912 by ConSurf version 3.0 on 2023-03-10.

- PDB ID: AF VGLUT2NAT RANK1 WITH SEQRES
- Chain processed: A
- Chains in the deposited model identical in sequence to chain processed: A
- Multiple sequence alignment source: Generated by ConSurf
- Multiple sequence alignment method: MAFFT

SEQRES Length: 582 amino acids in

• Amino acids with coordinates (ATOM

• Unique homologous sequences found:

• Sequences used in the MSA: 150

experimental protein.

records): 294

• Phylogenetic tree source: Generated by ConSurf

- Homologous sequences obtained from: UNIREF90
- PSI Blast E value cutoff: 0.0001
- PSI Blast Iterations: 1
- Maximum homologous sequences to use: 150
- Conservation score method: **Bayes**
- Amino acid substitution model: LG

For an explanation of the Job Parameters and Results, please see the on-line documentation for the ConSurf Server.

Job Results

Number of amino acids at each conservation grade, 1-9: Black: excluding amino acids with insufficient data. Yellow: including amino acids with insufficient data,

shown only if different.

- 8 Variable 1.
- 2. 7
- 3. 22
- 4. 40
- 5. 32
- 6. 36
- 7. 60
- 8. 48
- 41 Conserved

294 Sum 294

- 0 (Amino acids with insufficient data)
- 294 (0% with insufficient data)

Close