10.03.2023 16:38 ConSurf Job Information

ConSurf Job Information

Go to Results of Job 1678461301 (until they are deleted from the server).

The information below is obtained from the header of PDB file AF_vglut3NAT_rank1_with_seqres_consurf_firstglance.pdb

This PDB file can be saved to your computer.

Job Parameters

Job 1678461301 by ConSurf version 3.0 on 2023-03-10.

- PDB ID: AF_VGLUT3NAT_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: 589 amino acids in

• Amino acids with coordinates (ATOM

• Unique homologous sequences found:

• Sequences used in the MSA: 150

experimental protein.

records): 287

• 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 <u>ConSurf</u> 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.

- 12 Variable
- 2. **7**
- 3. **23**
- 4. 31
- 5. **39**
- *6*. **25**
- *7*. **62**
- 8. **51**
- 9. 37 Conserved

Sum 287 287

- + 0 (Amino acids with insufficient data)
- = 287 (0% with insufficient data)

Close