

pdb format:

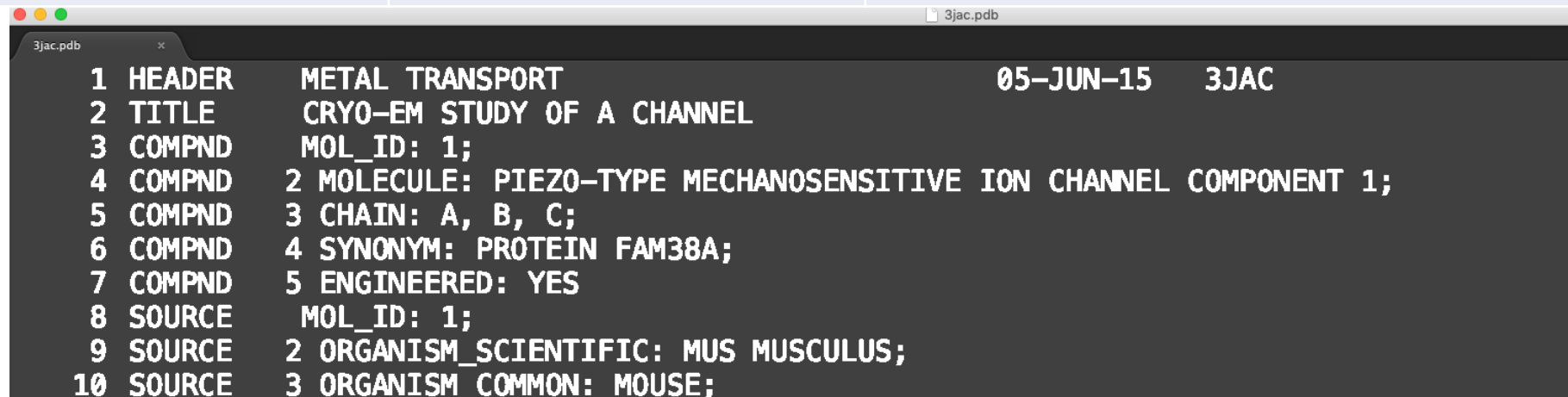
PDB files contain information about the 3D structure of proteins.

One record per line

The first six characters of the record determine the record type

HEADER : short description & PDB identifier

Columns	Example	Description
1-6	HEADER	Fixed string identifying each line.
63-66	3JAC	PDB ID



```
1  HEADER      METAL TRANSPORT                      05-JUN-15   3JAC
2  TITLE      CRYO-EM STUDY OF A CHANNEL
3  COMPND     MOL_ID: 1;
4  COMPND     2 MOLECULE: PIEZO-TYPE MECHANOSENSITIVE ION CHANNEL COMPONENT 1;
5  COMPND     3 CHAIN: A, B, C;
6  COMPND     4 SYNONYM: PROTEIN FAM38A;
7  COMPND     5 ENGINEERED: YES
8  SOURCE      MOL_ID: 1;
9  SOURCE      2 ORGANISM_SCIENTIFIC: MUS MUSCULUS;
10 SOURCE      3 ORGANISM_COMMON: MOUSE;
```

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PDB files contain information about the 3D structure of proteins.

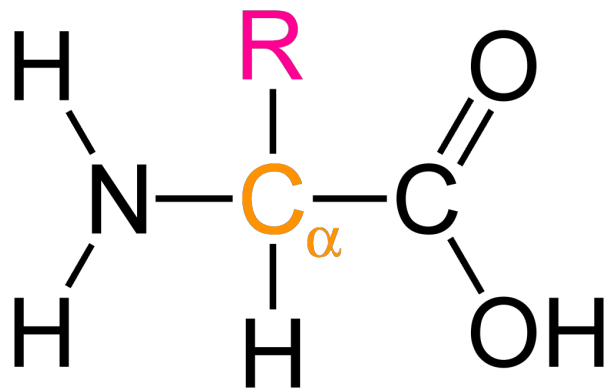
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HEADER : short description & PDB identifier

ATOM: Specify the chain, residue and atom information, including the coordinates, for one atom

ATOM	3554	N	VAL	B	567	-19.159	3.100	-6.625	1.00	30.00	N
ATOM	3555	CA	VAL	B	567	-18.816	4.270	-5.821	1.00	30.00	C
ATOM	3556	C	VAL	B	567	-17.366	4.679	-6.056	1.00	30.00	C
ATOM	3557	O	VAL	B	567	-17.044	5.865	-6.150	1.00	30.00	O
ATOM	3558	CB	VAL	B	567	-19.032	4.008	-4.323	1.00	20.00	C
ATOM	3559	CG1	VAL	B	567	-18.868	5.296	-3.542	1.00	20.00	C
ATOM	3560	CG2	VAL	B	567	-20.406	3.412	-4.080	1.00	20.00	C



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ATOM	3558	CB	VAL	B	567	-19.032	4.008	-4.323	1.00	20.00	C
ATOM	3559	CG1	VAL	B	567	-18.868	5.296	-3.542	1.00	20.00	C
ATOM	3560	CG2	VAL	B	567	-20.406	3.412	-4.080	1.00	20.00	C

Description of the fields in an **ATOM** record.

Columns	Example	Description
1-6	ATOM	Fixed string identifying each line.
7-11	7507	Atom serial number.
13-16	CA	Atom name.
18-20	ILE	Residue name.
22	B	Chain identifier.
23 - 26	255	Residue sequence number.
31 - 38	35.578	X coordinate of atom.
39 - 46	9.357	Y coordinate of atom.
47 - 54	5.792	Z coordinate of atom.

ATOM 3560 CG2 VAL B 567 -20.406 3.412 -4.080 1.00 20.00 C