



wwPDB EM Map/Model Validation Report ⓘ

Apr 26, 2016 – 06:26 PM BST

PDB ID : 5IJ0
EMDB ID: : EMD-8094
Title : Cryo EM density of microtubule assembled from human TUBB3
Authors : Ti, S.-C.; Pamula, M.C.; Howes, S.C.; Duellberg, C.; Nicholas, C.I.; Kleiner, R.E.; Forth, S.; Surrey, T.; Nogales, E.; Kapoor, T.M.
Deposited on : 2016-03-01
Resolution : 3.80 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk27461

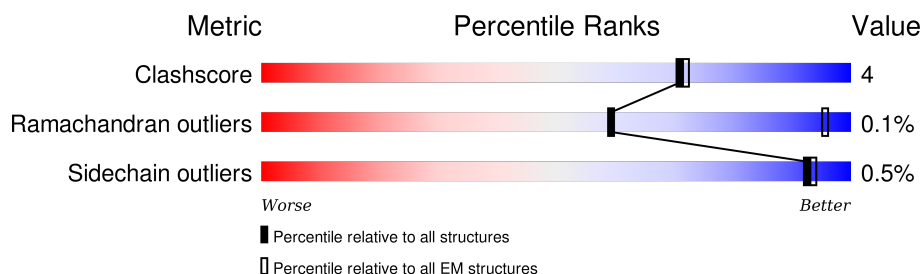
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	437	 85% 13% .
2	B	426	 90% 10%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 13304 atoms, of which 6523 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

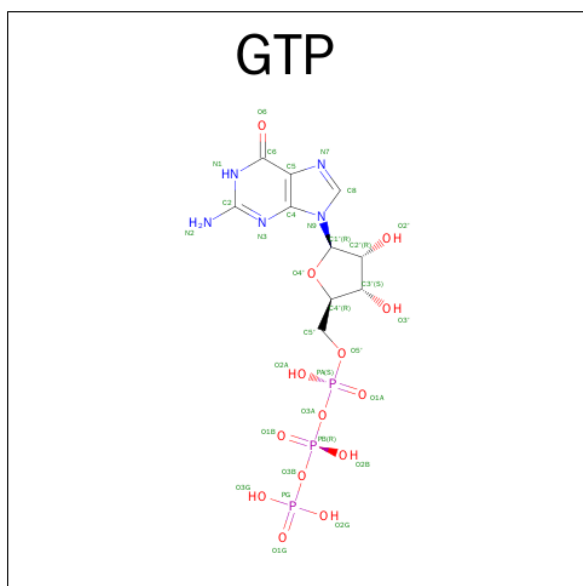
- Molecule 1 is a protein called Tubulin alpha-1B chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	A	430	6653	2136	3283	573	639	22	0	0

- Molecule 2 is a protein called Tubulin beta-3 chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	B	426	6590	2108	3240	574	642	26	0	0

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).

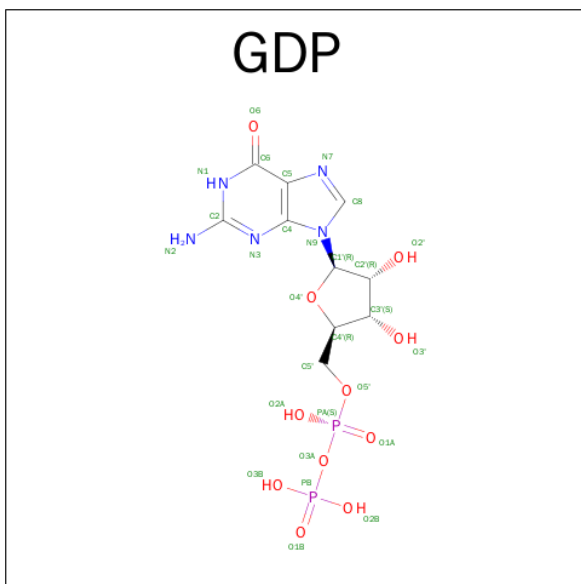


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	32	10	5	14	3	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Mg	0
			1	1	

- Molecule 5 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).

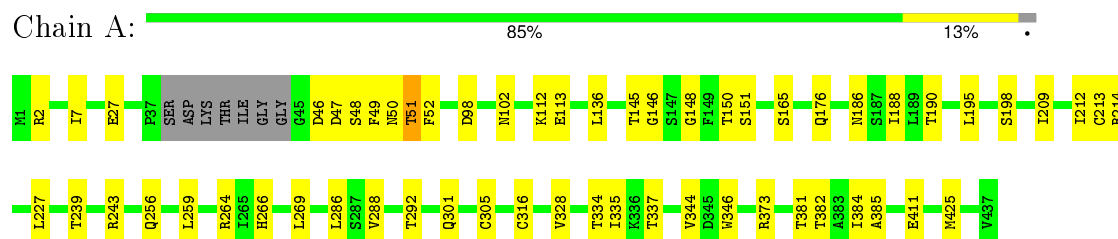


Mol	Chain	Residues	Atoms					AltConf
5	B	1	Total	C	N	O	P	0
			28	10	5	11	2	

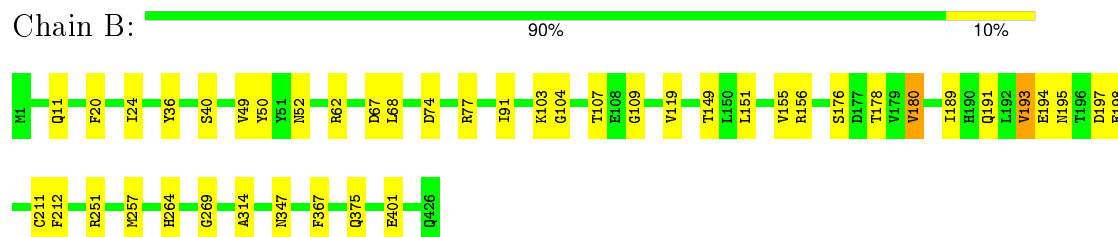
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Tubulin alpha-1B chain



- Molecule 2: Tubulin beta-3 chain



4 Experimental information

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	24268	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	Not provided	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GDP, GTP, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.24	0/3447	0.39	0/4679
2	B	0.23	0/3425	0.38	0/4638
All	All	0.24	0/6872	0.39	0/9317

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3370	3283	3284	33	0
2	B	3350	3240	3242	28	0
3	A	32	0	12	1	0
4	A	1	0	0	0	0
5	B	28	0	12	1	0
All	All	6781	6523	6550	59	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (59) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:11:GLN:NE2	5:B:501:GDP:O3B	2.27	0.67
2:B:52:ASN:OD1	2:B:62:ARG:NH2	2.29	0.66
1:A:186:ASN:O	1:A:190:THR:OG1	2.11	0.64
2:B:49:VAL:O	2:B:62:ARG:NH1	2.31	0.63
1:A:98:ASP:O	2:B:251:ARG:NH2	2.32	0.62
2:B:74:ASP:OD1	2:B:77:ARG:NH2	2.33	0.61
2:B:156:ARG:NE	2:B:195:ASN:O	2.31	0.60
2:B:91:ILE:HD11	2:B:119:VAL:HG21	1.83	0.59
1:A:301:GLN:NE2	1:A:305:CYS:O	2.35	0.59
1:A:292:THR:HB	1:A:335:ILE:HD11	1.87	0.55
1:A:212:ILE:HD11	1:A:301:GLN:O	2.08	0.54
1:A:239:THR:O	1:A:243:ARG:N	2.41	0.54
2:B:103:LYS:O	2:B:107:THR:OG1	2.09	0.54
1:A:334:THR:O	1:A:337:THR:OG1	2.26	0.52
1:A:2:ARG:O	1:A:51:THR:OG1	2.14	0.52
2:B:49:VAL:HG13	2:B:50:TYR:N	2.25	0.51
2:B:176:SER:OG	2:B:178:THR:O	2.29	0.51
2:B:191:GLN:NE2	2:B:194:GLU:OE2	2.41	0.51
2:B:189:ILE:O	2:B:193:VAL:N	2.44	0.51
1:A:48:SER:OG	1:A:243:ARG:O	2.18	0.51
1:A:176:GLN:O	2:B:347:ASN:ND2	2.41	0.51
1:A:209:ILE:CG2	1:A:227:LEU:HD22	2.42	0.50
1:A:47:ASP:O	1:A:50:ASN:N	2.46	0.49
1:A:46:ASP:O	1:A:49:PHE:N	2.35	0.48
1:A:102:ASN:ND2	1:A:411:GLU:OE1	2.46	0.48
2:B:20:PHE:CZ	2:B:24:ILE:HD11	2.48	0.48
2:B:20:PHE:CE2	2:B:24:ILE:HD11	2.49	0.48
1:A:269:LEU:HD21	1:A:384:ILE:HD13	1.95	0.48
1:A:382:THR:O	1:A:385:ALA:HB3	2.15	0.47
1:A:148:GLY:O	1:A:151:SER:OG	2.27	0.47
2:B:67:ASP:OD1	2:B:68:LEU:N	2.48	0.46
2:B:198:GLU:HA	2:B:264:HIS:HB2	1.98	0.46
2:B:49:VAL:HG13	2:B:50:TYR:H	1.80	0.46
1:A:381:THR:O	1:A:384:ILE:HG22	2.16	0.45
2:B:180:VAL:HG12	2:B:180:VAL:O	2.17	0.45
1:A:195:LEU:HD11	1:A:264:ARG:O	2.17	0.44
2:B:156:ARG:NH2	2:B:197:ASP:OD1	2.50	0.44
1:A:112:LYS:NZ	1:A:113:GLU:OE2	2.49	0.44
1:A:146:GLY:O	1:A:150:THR:OG1	2.09	0.44
2:B:257:MET:SD	2:B:314:ALA:HB2	2.58	0.43
1:A:7:ILE:N	1:A:136:LEU:O	2.48	0.43
2:B:211:CYS:SG	2:B:212:PHE:N	2.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:LEU:HD21	1:A:316:CYS:HB2	2.02	0.42
1:A:209:ILE:HG22	1:A:227:LEU:HD22	2.01	0.42
1:A:145:THR:OG1	3:A:501:GTP:O2G	2.24	0.42
2:B:149:THR:HG23	2:B:191:GLN:CB	2.50	0.42
1:A:344:VAL:HG11	1:A:346:TRP:NE1	2.35	0.42
1:A:213:CYS:SG	1:A:214:ARG:N	2.92	0.42
1:A:286:LEU:O	1:A:373:ARG:NH1	2.53	0.41
2:B:103:LYS:NZ	2:B:401:GLU:OE2	2.41	0.41
1:A:188:ILE:HD12	1:A:425:MET:HG3	2.01	0.41
2:B:151:LEU:O	2:B:155:VAL:HG23	2.21	0.41
2:B:104:GLY:O	2:B:109:GLY:N	2.49	0.41
1:A:165:SER:HA	1:A:256:GLN:HE22	1.86	0.40
2:B:269:GLY:N	2:B:367:PHE:O	2.53	0.40
1:A:198:SER:O	1:A:266:HIS:NE2	2.54	0.40
1:A:27:GLU:OE1	1:A:243:ARG:NH2	2.53	0.40
1:A:288:VAL:HG11	1:A:328:VAL:HG22	2.03	0.40
2:B:36:TYR:OH	2:B:40:SER:O	2.39	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	426/437 (98%)	405 (95%)	21 (5%)	0	100	100
2	B	424/426 (100%)	411 (97%)	12 (3%)	1 (0%)	52	86
All	All	850/863 (98%)	816 (96%)	33 (4%)	1 (0%)	59	90

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	180	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	363/368 (99%)	361 (99%)	2 (1%)	90	96
2	B	367/367 (100%)	365 (100%)	2 (0%)	92	97
All	All	730/735 (99%)	726 (100%)	4 (0%)	92	97

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	THR
1	A	52	PHE
2	B	193	VAL
2	B	375	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	256	GLN
2	B	256	ASN
2	B	264	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	GTP	A	501	4	26,34,34	0.94	1 (3%)	29,54,54	1.55	4 (13%)
5	GDP	B	501	-	24,30,30	1.15	2 (8%)	26,47,47	1.82	4 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GTP	A	501	4	-	0/18/38/38	0/3/3/3
5	GDP	B	501	-	-	0/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	GTP	C6-N1	2.99	1.38	1.33
5	B	501	GDP	C5-C4	3.08	1.47	1.40
5	B	501	GDP	C6-C5	3.61	1.48	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	GTP	N3-C2-N1	-5.40	120.21	127.56
5	B	501	GDP	C5-C6-N1	-4.24	117.98	123.52
5	B	501	GDP	N3-C2-N1	-3.48	122.82	127.56
5	B	501	GDP	C6-C5-C4	-3.18	117.22	120.86
3	A	501	GTP	C5-C6-N1	-2.93	119.69	123.52
3	A	501	GTP	C4'-O4'-C1'	2.05	111.81	109.64
3	A	501	GTP	C6-N1-C2	3.30	119.75	115.88
5	B	501	GDP	C6-N1-C2	5.53	122.37	115.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	GTP	1	0
5	B	501	GDP	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.