



Full wwPDB/EMDatabank EM Map/Model Validation Report ⓘ

Jan 25, 2017 – 11:45 AM EST

PDB ID : 5SYG
EMDB ID: : EMD-8323
Title : Cryo-EM reconstruction of zampanolide-bound microtubule
Authors : Kellogg, E.H.; Hejab, N.M.A.; Nogales, E.
Deposited on : 2016-08-11
Resolution : 3.50 Å(reported)

This is a Full wwPDB/EMDatabank EM Map/Model Validation Report
for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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) : rb-20028442

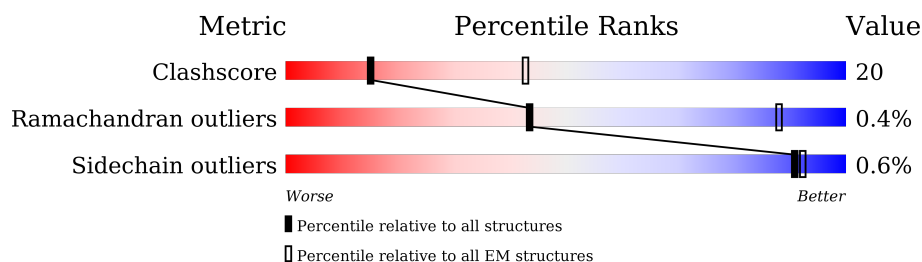
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	
2	B	426	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	ZPN	B	502	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6792 atoms, of which 0 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.

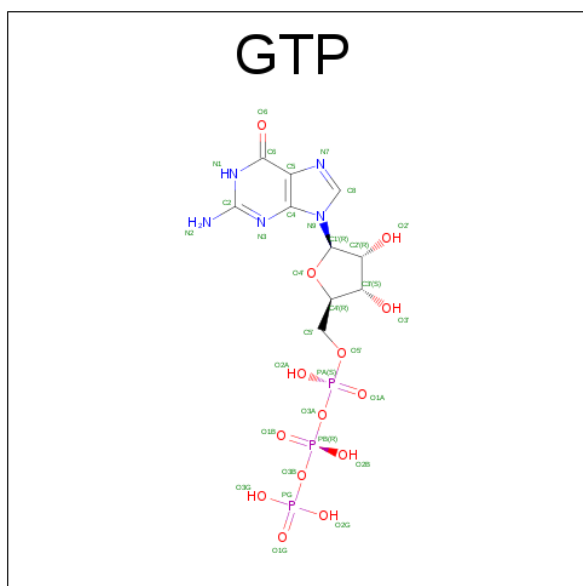
- Molecule 1 is a protein called Tubulin alpha chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	426	3343	2121	569	631	22	0	0

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	426	3352	2106	575	645	26	0	0

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

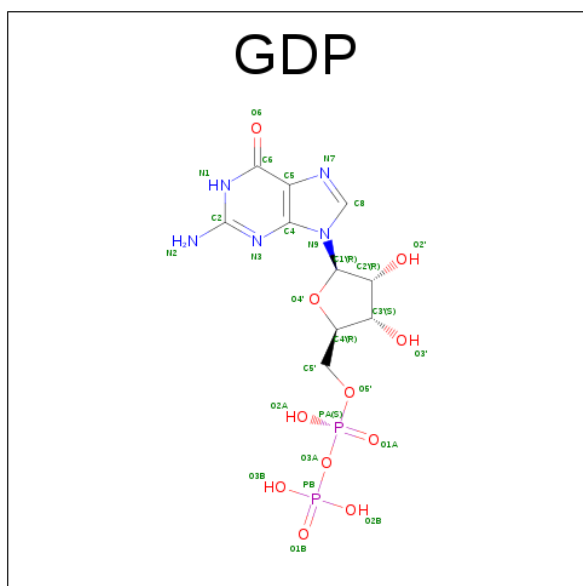


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	32	10	5	14	3	0

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

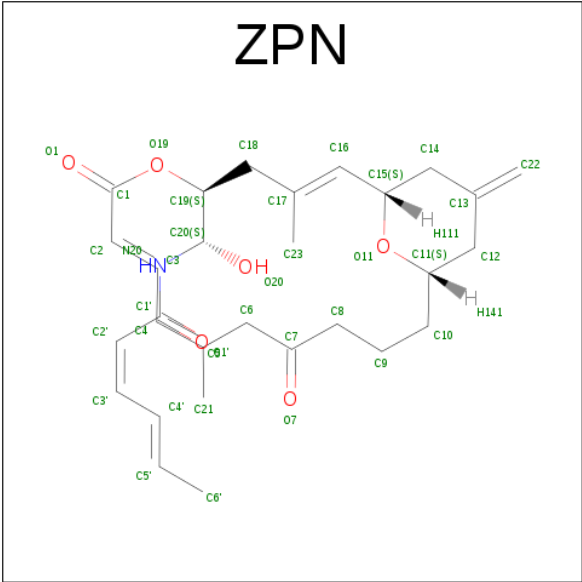
Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Mg	0
			1	1	

- 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	

- Molecule 6 is (2Z,4E)-N-[(S)-[(1S,2E,5S,8E,10Z,17S)-3,11-dimethyl-19-methylidene-7,13-dioxo-6,21-dioxabicyclo[15.3.1]henicosa-2,8,10-trien-5-yl](hydroxy)methyl]hexa-2,4-dienamide (three-letter code: ZPN) (formula: $C_{29}H_{39}NO_6$).

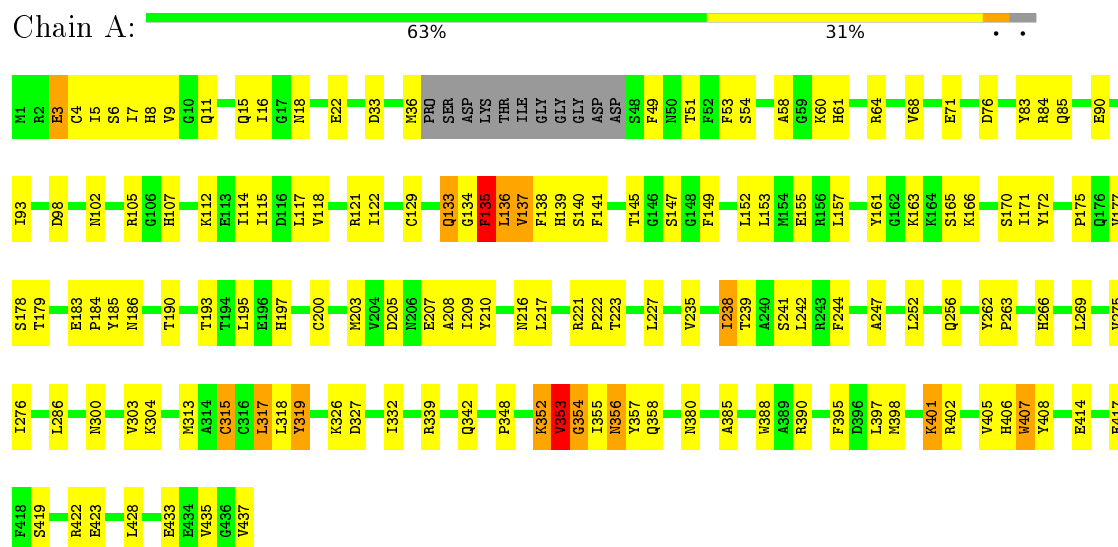


Mol	Chain	Residues	Atoms				AltConf
6	B	1	Total	C	N	O	0
			36	29	1	6	

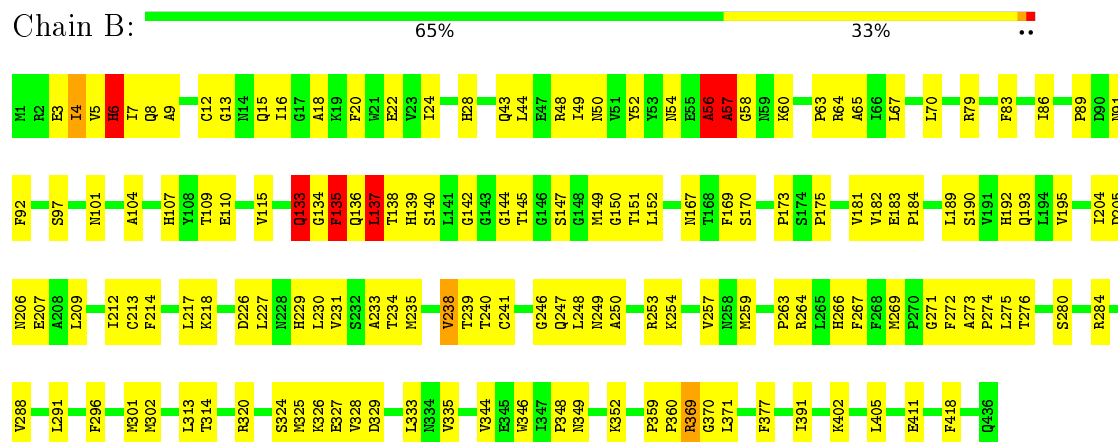
3 Residue-property plots [i](#)

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 chain



• Molecule 2: Tubulin beta chain



4 Experimental information

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of segments used	17069	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)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	27500	Depositor
Image detector	Not provided	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.65	7/3419 (0.2%)	0.85	21/4639 (0.5%)
2	B	0.52	1/3427 (0.0%)	0.76	8/4641 (0.2%)
All	All	0.59	8/6846 (0.1%)	0.81	29/9280 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	B	0	5
All	All	0	8

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	315	CYS	C-O	7.56	1.37	1.23
1	A	356	ASN	N-CA	6.79	1.59	1.46
1	A	317	LEU	C-O	6.70	1.36	1.23
1	A	354	GLY	N-CA	6.05	1.55	1.46
1	A	352	LYS	C-O	5.92	1.34	1.23
2	B	238	VAL	C-O	5.84	1.34	1.23
1	A	407	TRP	CB-CG	-5.28	1.40	1.50
1	A	238	ILE	C-O	5.05	1.32	1.23

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	353	VAL	C-N-CA	-8.08	105.33	122.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	317	LEU	CA-C-O	8.07	137.06	120.10
2	B	56	ALA	O-C-N	-8.00	109.91	122.70
1	A	319	TYR	N-CA-C	-7.64	90.37	111.00
1	A	315	CYS	CA-C-O	7.63	136.12	120.10
2	B	133	GLN	O-C-N	-7.58	110.31	123.20
2	B	6	HIS	C-N-CA	-7.47	103.02	121.70
1	A	354	GLY	CA-C-O	-7.14	107.74	120.60
1	A	352	LYS	N-CA-C	-7.09	91.86	111.00
1	A	6	SER	C-N-CA	-7.05	104.06	121.70
1	A	8	HIS	C-N-CA	-6.77	104.77	121.70
1	A	133	GLN	O-C-N	-6.76	111.70	123.20
2	B	8	GLN	C-N-CA	-6.75	104.83	121.70
1	A	315	CYS	O-C-N	-6.66	112.05	122.70
1	A	315	CYS	C-N-CA	-6.31	105.92	121.70
1	A	352	LYS	O-C-N	6.08	132.44	122.70
1	A	137	VAL	O-C-N	-6.07	112.98	122.70
1	A	317	LEU	O-C-N	-5.89	113.27	122.70
1	A	317	LEU	C-N-CA	-5.79	107.22	121.70
1	A	355	ILE	C-N-CA	-5.71	107.42	121.70
2	B	135	PHE	O-C-N	-5.67	113.63	122.70
1	A	4	CYS	C-N-CA	-5.54	107.86	121.70
1	A	135	PHE	O-C-N	-5.46	113.97	122.70
2	B	137	LEU	O-C-N	-5.46	113.97	122.70
1	A	315	CYS	N-CA-C	5.32	125.35	111.00
2	B	4	ILE	C-N-CA	-5.27	108.52	121.70
1	A	356	ASN	O-C-N	5.26	131.11	122.70
1	A	356	ASN	C-N-CA	5.05	134.32	121.70
2	B	57	ALA	N-CA-C	-5.03	97.42	111.00

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	133	GLN	Mainchain
1	A	135	PHE	Mainchain
1	A	353	VAL	Peptide
2	B	133	GLN	Mainchain
2	B	135	PHE	Mainchain
2	B	137	LEU	Mainchain
2	B	56	ALA	Mainchain
2	B	6	HIS	Peptide

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	3343	0	3263	132	0
2	B	3352	0	3237	164	0
3	A	32	0	12	0	0
4	A	1	0	0	0	0
5	B	28	0	12	0	0
6	B	36	0	39	36	0
All	All	6792	0	6563	261	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 20.

All (261) 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:229:HIS:NE2	6:B:502:ZPN:H92	1.39	1.34
2:B:229:HIS:CD2	6:B:502:ZPN:H92	1.64	1.31
2:B:272:PHE:CZ	6:B:502:ZPN:H232	1.71	1.25
2:B:276:THR:N	6:B:502:ZPN:O1'	1.70	1.25
2:B:276:THR:OG1	6:B:502:ZPN:H9	1.50	1.12
2:B:272:PHE:HZ	6:B:502:ZPN:H232	1.03	1.04
2:B:233:ALA:HA	6:B:502:ZPN:C22	1.89	1.02
2:B:229:HIS:NE2	6:B:502:ZPN:C9	2.26	0.99
2:B:229:HIS:CD2	6:B:502:ZPN:C9	2.46	0.98
1:A:7:ILE:N	1:A:136:LEU:O	1.95	0.97
1:A:5:ILE:O	1:A:135:PHE:HA	1.66	0.96
2:B:5:VAL:O	2:B:135:PHE:HA	1.65	0.94
2:B:7:ILE:O	2:B:137:LEU:HA	1.70	0.92
2:B:276:THR:HB	6:B:502:ZPN:C4'	2.03	0.89
1:A:9:VAL:N	1:A:138:PHE:O	2.05	0.88
1:A:315:CYS:O	1:A:352:LYS:O	1.92	0.86
2:B:233:ALA:HA	6:B:502:ZPN:H223	1.57	0.86
2:B:272:PHE:CZ	6:B:502:ZPN:C23	2.58	0.84
2:B:276:THR:CB	6:B:502:ZPN:H9	2.08	0.83
1:A:222:PRO:O	2:B:324:SER:OG	1.97	0.82
1:A:317:LEU:O	1:A:354:GLY:O	1.99	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:LYS:HD3	2:B:346:TRP:CH2	2.16	0.80
1:A:93:ILE:HD11	1:A:121:ARG:HG3	1.62	0.80
1:A:216:ASN:HB3	1:A:275:VAL:O	1.85	0.77
1:A:177:VAL:HG22	2:B:333:LEU:CD1	2.15	0.77
1:A:401:LYS:HG2	2:B:346:TRP:CE3	2.19	0.77
1:A:136:LEU:HD12	1:A:136:LEU:H	1.51	0.75
2:B:275:LEU:HA	6:B:502:ZPN:O1'	1.89	0.73
2:B:233:ALA:CA	6:B:502:ZPN:C22	2.66	0.73
2:B:230:LEU:O	2:B:233:ALA:HB3	1.89	0.72
1:A:5:ILE:N	1:A:134:GLY:O	2.22	0.72
2:B:271:GLY:HA3	2:B:377:PHE:HB3	1.70	0.71
2:B:9:ALA:N	2:B:138:THR:O	2.23	0.71
2:B:276:THR:HB	6:B:502:ZPN:C5'	2.21	0.71
1:A:177:VAL:HG22	2:B:333:LEU:HD11	1.74	0.70
1:A:114:ILE:O	1:A:117:LEU:N	2.25	0.69
2:B:313:LEU:HD23	2:B:344:VAL:HG21	1.73	0.69
1:A:172:TYR:HB2	1:A:203:MET:HB3	1.73	0.69
2:B:276:THR:OG1	6:B:502:ZPN:C6'	2.37	0.69
2:B:5:VAL:N	2:B:134:GLY:O	2.16	0.69
2:B:115:VAL:HG21	2:B:152:LEU:HD23	1.75	0.68
1:A:7:ILE:O	1:A:137:VAL:HA	1.92	0.68
1:A:5:ILE:HB	1:A:135:PHE:HB3	1.75	0.68
2:B:89:PRO:HA	2:B:92:PHE:HD1	1.58	0.68
1:A:22:GLU:HG2	1:A:83:TYR:CE1	2.29	0.67
2:B:272:PHE:HZ	6:B:502:ZPN:C23	1.95	0.67
2:B:271:GLY:N	2:B:377:PHE:O	2.23	0.67
1:A:407:TRP:CZ2	2:B:257:VAL:HG22	2.30	0.67
1:A:11:GLN:NE2	2:B:247:GLN:O	2.29	0.66
2:B:226:ASP:O	2:B:229:HIS:HB3	1.95	0.66
1:A:145:THR:O	1:A:149:PHE:HB3	1.96	0.65
1:A:7:ILE:O	1:A:138:PHE:N	2.30	0.65
1:A:406:HIS:CG	2:B:263:PRO:HG3	2.32	0.65
1:A:332:ILE:HD11	1:A:353:VAL:HG11	1.78	0.65
1:A:222:PRO:HD2	2:B:326:LYS:HD3	1.78	0.65
2:B:101:ASN:HA	2:B:144:GLY:HA3	1.79	0.64
2:B:275:LEU:CA	6:B:502:ZPN:O1'	2.46	0.64
2:B:57:ALA:O	2:B:60:LYS:N	2.26	0.64
1:A:401:LYS:HD3	2:B:346:TRP:CZ3	2.32	0.63
1:A:33:ASP:HA	1:A:85:GLN:HB2	1.80	0.63
2:B:6:HIS:CE1	2:B:136:GLN:HB2	2.33	0.63
1:A:112:LYS:O	1:A:115:ILE:HG22	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:LEU:HD22	1:A:428:LEU:HD11	1.81	0.62
1:A:105:ARG:HH12	2:B:253:ARG:HD3	1.65	0.62
1:A:216:ASN:ND2	1:A:300:ASN:OD1	2.33	0.62
1:A:178:SER:HB3	2:B:349:ASN:ND2	2.15	0.61
1:A:11:GLN:HE22	2:B:248:LEU:HA	1.65	0.61
1:A:179:THR:O	2:B:352:LYS:HG2	2.00	0.60
2:B:56:ALA:O	2:B:60:LYS:HB2	2.01	0.60
1:A:313:MET:N	1:A:380:ASN:O	2.33	0.60
1:A:71:GLU:HB2	1:A:98:ASP:HB3	1.83	0.60
2:B:284:ARG:HD3	6:B:502:ZPN:H7	1.84	0.59
2:B:22:GLU:HB2	2:B:83:PHE:HE2	1.66	0.59
1:A:317:LEU:HB3	1:A:319:TYR:CE1	2.38	0.59
2:B:250:ALA:HA	2:B:254:LYS:HE2	1.83	0.58
1:A:107:HIS:O	1:A:107:HIS:CD2	2.56	0.58
1:A:177:VAL:HG13	2:B:333:LEU:HD11	1.85	0.58
2:B:20:PHE:CZ	2:B:24:ILE:HD11	2.38	0.58
1:A:178:SER:HB3	2:B:349:ASN:CG	2.24	0.57
2:B:145:THR:O	2:B:149:MET:HB3	2.02	0.57
2:B:274:PRO:HB2	6:B:502:ZPN:C4'	2.35	0.57
2:B:169:PHE:CE2	2:B:235:MET:HG2	2.39	0.57
1:A:397:LEU:HD23	2:B:348:PRO:HD3	1.86	0.57
1:A:5:ILE:O	1:A:135:PHE:CA	2.48	0.57
1:A:223:THR:HB	2:B:324:SER:HA	1.87	0.57
1:A:407:TRP:CE2	2:B:257:VAL:HG22	2.40	0.57
1:A:177:VAL:HG22	2:B:333:LEU:HD13	1.85	0.56
1:A:147:SER:HG	1:A:190:THR:HG1	1.49	0.56
2:B:65:ALA:O	2:B:91:ASN:ND2	2.39	0.56
1:A:84:ARG:HH21	1:A:85:GLN:HE22	1.54	0.56
2:B:195:VAL:HG22	2:B:267:PHE:HZ	1.70	0.56
2:B:240:THR:HG21	2:B:320:ARG:CZ	2.37	0.55
2:B:275:LEU:C	6:B:502:ZPN:O1'	2.41	0.55
1:A:339:ARG:O	1:A:342:GLN:NE2	2.39	0.55
2:B:276:THR:N	6:B:502:ZPN:H4'1	2.20	0.55
2:B:136:GLN:HG2	2:B:167:ASN:OD1	2.07	0.55
1:A:105:ARG:NH1	2:B:253:ARG:HD3	2.22	0.55
2:B:229:HIS:CE1	6:B:502:ZPN:H92	2.31	0.55
1:A:147:SER:OG	1:A:190:THR:OG1	2.10	0.54
2:B:214:PHE:O	2:B:218:LYS:HA	2.07	0.54
1:A:58:ALA:HB3	1:A:60:LYS:HG2	1.89	0.54
2:B:192:HIS:CD2	2:B:193:GLN:HE21	2.25	0.54
2:B:205:ASP:OD2	2:B:207:GLU:HB3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:63:PRO:HD3	2:B:86:ILE:HG22	1.90	0.54
1:A:139:HIS:ND1	1:A:140:SER:O	2.41	0.53
1:A:3:GLU:OE2	1:A:129:CYS:HB3	2.09	0.53
1:A:186:ASN:OD1	1:A:408:TYR:OH	2.26	0.53
1:A:177:VAL:HG11	2:B:329:ASP:HB3	1.90	0.53
2:B:4:ILE:HG12	2:B:133:GLN:O	2.08	0.53
1:A:326:LYS:HG3	1:A:327:ASP:N	2.23	0.53
2:B:173:PRO:HD2	2:B:391:ILE:HD11	1.90	0.52
2:B:204:ILE:HD13	2:B:231:VAL:HG22	1.92	0.52
2:B:288:VAL:HG11	2:B:328:VAL:HA	1.90	0.52
1:A:36:MET:HB2	1:A:61:HIS:CD2	2.44	0.52
1:A:114:ILE:HG22	1:A:118:VAL:HG23	1.91	0.52
1:A:139:HIS:NE2	1:A:170:SER:OG	2.38	0.52
6:B:502:ZPN:C4	6:B:502:ZPN:H81	2.40	0.52
2:B:217:LEU:HD22	6:B:502:ZPN:O20	2.09	0.52
1:A:84:ARG:NH2	1:A:85:GLN:NE2	2.59	0.51
2:B:263:PRO:O	2:B:264:ARG:HB2	2.10	0.51
1:A:221:ARG:HD3	2:B:327:GLU:OE1	2.11	0.51
2:B:233:ALA:CB	6:B:502:ZPN:C22	2.89	0.51
2:B:276:THR:CG2	2:B:280:SER:HB2	2.41	0.50
2:B:229:HIS:NE2	6:B:502:ZPN:C8	2.73	0.50
2:B:54:ASN:HB2	2:B:64:ARG:CZ	2.42	0.50
1:A:53:PHE:HB3	1:A:61:HIS:HB3	1.92	0.49
1:A:185:TYR:HA	1:A:395:PHE:CD1	2.48	0.49
1:A:402:ARG:HB3	1:A:405:VAL:CG1	2.41	0.49
2:B:107:HIS:HD2	2:B:151:THR:HB	1.77	0.49
1:A:276:ILE:HD11	1:A:286:LEU:HD11	1.95	0.49
1:A:210:TYR:HB3	2:B:326:LYS:HD2	1.95	0.49
2:B:212:ILE:O	2:B:217:LEU:HG	2.12	0.49
2:B:248:LEU:HD11	2:B:325:MET:HE1	1.94	0.49
2:B:181:VAL:HG23	2:B:182:VAL:HG13	1.94	0.49
2:B:234:THR:HG21	2:B:302:MET:SD	2.53	0.49
2:B:273:ALA:HB1	2:B:291:LEU:HD12	1.94	0.48
1:A:16:ILE:HD11	1:A:171:ILE:HD11	1.95	0.48
1:A:433:GLU:O	1:A:437:VAL:HG23	2.13	0.48
2:B:276:THR:CB	6:B:502:ZPN:C6'	2.87	0.48
1:A:175:PRO:HB3	1:A:390:ARG:CZ	2.43	0.48
2:B:205:ASP:OD1	2:B:206:ASN:N	2.46	0.48
2:B:89:PRO:HA	2:B:92:PHE:CD1	2.45	0.48
1:A:155:GLU:HG3	1:A:197:HIS:CE1	2.49	0.48
2:B:209:LEU:HD22	2:B:230:LEU:HD13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:272:PHE:HB3	2:B:275:LEU:HD11	1.96	0.48
2:B:79:ARG:HD3	2:B:92:PHE:CE2	2.49	0.48
1:A:5:ILE:HB	1:A:135:PHE:CB	2.43	0.47
1:A:401:LYS:CG	2:B:346:TRP:CE3	2.92	0.47
1:A:54:SER:HB2	1:A:64:ARG:HH12	1.80	0.47
2:B:52:TYR:OH	2:B:239:THR:HG21	2.14	0.47
2:B:70:LEU:HB3	2:B:97:SER:O	2.14	0.47
1:A:141:PHE:CZ	1:A:170:SER:HB2	2.49	0.47
1:A:18:ASN:O	1:A:22:GLU:HG3	2.15	0.47
1:A:239:THR:O	1:A:242:LEU:N	2.47	0.47
2:B:189:LEU:HD11	2:B:418:PHE:CZ	2.50	0.47
1:A:102:ASN:HD21	2:B:257:VAL:HG21	1.79	0.47
1:A:165:SER:OG	1:A:256:GLN:NE2	2.48	0.47
2:B:250:ALA:HA	2:B:254:LYS:CE	2.45	0.47
1:A:221:ARG:HB2	2:B:324:SER:CB	2.45	0.47
2:B:229:HIS:CD2	6:B:502:ZPN:C7	2.98	0.47
1:A:313:MET:SD	1:A:435:VAL:HG11	2.55	0.46
2:B:57:ALA:O	2:B:58:GLY:C	2.53	0.46
2:B:360:PRO:HB2	2:B:369:ARG:HG2	1.97	0.46
1:A:49:PHE:CE1	1:A:61:HIS:CE1	3.03	0.46
2:B:147:SER:OG	2:B:190:SER:HB2	2.16	0.46
2:B:296:PHE:CE1	2:B:335:VAL:HG11	2.51	0.46
1:A:76:ASP:OD2	2:B:48:ARG:NH2	2.49	0.46
2:B:9:ALA:O	2:B:13:GLY:HA3	2.16	0.46
2:B:22:GLU:HB2	2:B:83:PHE:CE2	2.50	0.46
2:B:246:GLY:N	2:B:249:ASN:OD1	2.49	0.46
1:A:205:ASP:OD1	1:A:207:GLU:N	2.48	0.46
1:A:208:ALA:HB2	1:A:304:LYS:HB2	1.98	0.46
2:B:253:ARG:O	2:B:257:VAL:HG23	2.15	0.45
1:A:135:PHE:O	1:A:135:PHE:CD1	2.68	0.45
2:B:230:LEU:HD23	6:B:502:ZPN:C23	2.46	0.45
2:B:344:VAL:HG12	2:B:346:TRP:HB2	1.97	0.45
1:A:319:TYR:H	1:A:356:ASN:H	1.63	0.45
1:A:401:LYS:HE2	2:B:346:TRP:CE2	2.52	0.45
2:B:9:ALA:HB3	2:B:139:HIS:ND1	2.32	0.45
2:B:266:HIS:ND1	2:B:266:HIS:O	2.47	0.45
1:A:414:GLU:O	1:A:417:GLU:HB3	2.16	0.45
1:A:114:ILE:CG2	1:A:118:VAL:HG23	2.47	0.45
2:B:16:ILE:HB	2:B:138:THR:HG21	1.98	0.45
2:B:269:MET:HB3	2:B:301:MET:SD	2.57	0.45
2:B:9:ALA:HB3	2:B:139:HIS:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:HIS:CE1	1:A:193:THR:HG21	2.52	0.45
1:A:222:PRO:HD2	2:B:326:LYS:CD	2.44	0.45
2:B:274:PRO:HB2	6:B:502:ZPN:C3'	2.46	0.45
2:B:56:ALA:O	2:B:57:ALA:CB	2.63	0.45
1:A:406:HIS:CB	2:B:263:PRO:HG3	2.46	0.44
1:A:107:HIS:CD2	1:A:152:LEU:HB2	2.53	0.44
1:A:161:TYR:O	1:A:163:LYS:N	2.39	0.44
2:B:151:THR:HG21	2:B:193:GLN:HB2	2.00	0.44
2:B:369:ARG:HG3	6:B:502:ZPN:H151	1.98	0.44
2:B:276:THR:HB	6:B:502:ZPN:H9	1.97	0.44
1:A:15:GLN:O	1:A:18:ASN:HB2	2.18	0.44
1:A:90:GLU:HB3	1:A:121:ARG:NE	2.33	0.44
1:A:135:PHE:CZ	1:A:166:LYS:HG2	2.52	0.44
1:A:422:ARG:NH2	1:A:423:GLU:OE2	2.51	0.44
2:B:107:HIS:O	2:B:152:LEU:HD22	2.17	0.44
2:B:238:VAL:O	2:B:241:CYS:HB2	2.17	0.44
1:A:419:SER:O	1:A:423:GLU:HG2	2.18	0.44
2:B:67:LEU:HD12	2:B:92:PHE:CD2	2.53	0.44
1:A:137:VAL:O	1:A:137:VAL:HG12	2.17	0.43
2:B:139:HIS:CE1	2:B:150:GLY:HA3	2.53	0.43
2:B:15:GLN:O	2:B:18:ALA:HB3	2.18	0.43
2:B:142:GLY:HA3	2:B:183:GLU:HG3	1.99	0.43
1:A:185:TYR:HA	1:A:395:PHE:HD1	1.82	0.43
2:B:359:PRO:HB2	2:B:371:LEU:O	2.17	0.43
2:B:7:ILE:O	2:B:138:THR:N	2.51	0.43
2:B:107:HIS:CD2	2:B:151:THR:HB	2.52	0.43
2:B:5:VAL:O	2:B:136:GLN:N	2.51	0.43
1:A:107:HIS:HE1	1:A:193:THR:HG21	1.84	0.43
1:A:217:LEU:HG	1:A:275:VAL:HG12	2.01	0.43
1:A:247:ALA:HB2	1:A:357:TYR:CE1	2.53	0.43
1:A:3:GLU:HB3	1:A:51:THR:HA	2.00	0.43
1:A:84:ARG:NH2	1:A:85:GLN:HE22	2.13	0.43
2:B:104:ALA:HB1	2:B:411:GLU:HB2	2.01	0.43
1:A:68:VAL:HG11	1:A:149:PHE:CZ	2.53	0.43
2:B:360:PRO:O	2:B:370:GLY:N	2.51	0.43
1:A:242:LEU:HD11	1:A:252:LEU:HG	2.00	0.43
2:B:183:GLU:HB3	2:B:184:PRO:HD3	2.00	0.43
2:B:213:CYS:SG	2:B:227:LEU:HD23	2.59	0.43
2:B:28:HIS:ND1	2:B:43:GLN:O	2.45	0.43
2:B:189:LEU:O	2:B:192:HIS:HB3	2.18	0.42
1:A:269:LEU:HD22	1:A:303:VAL:HG21	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:MET:SD	2:B:348:PRO:HD2	2.60	0.42
1:A:401:LYS:HA	1:A:401:LYS:HD2	1.12	0.42
1:A:222:PRO:HG2	2:B:326:LYS:HD2	2.02	0.42
1:A:177:VAL:CG1	2:B:329:ASP:HB3	2.50	0.42
1:A:183:GLU:HB3	1:A:184:PRO:HD3	2.00	0.42
2:B:12:CYS:SG	2:B:140:SER:HB3	2.60	0.42
2:B:52:TYR:OH	2:B:136:GLN:OE1	2.37	0.42
1:A:244:PHE:CD1	1:A:358:GLN:HG2	2.55	0.42
1:A:7:ILE:HD13	1:A:153:LEU:HD21	2.01	0.42
1:A:319:TYR:O	1:A:356:ASN:N	2.53	0.41
1:A:200:CYS:HA	1:A:266:HIS:CD2	2.54	0.41
1:A:262:TYR:CD2	1:A:263:PRO:HD2	2.55	0.41
1:A:385:ALA:O	1:A:388:TRP:HB2	2.20	0.41
1:A:177:VAL:CG2	2:B:333:LEU:HD11	2.47	0.41
1:A:209:ILE:HG21	1:A:227:LEU:HG	2.02	0.41
1:A:319:TYR:O	1:A:356:ASN:O	2.38	0.41
2:B:109:THR:OG1	2:B:110:GLU:N	2.53	0.41
2:B:276:THR:HB	6:B:502:ZPN:C6'	2.48	0.41
2:B:5:VAL:O	2:B:135:PHE:CA	2.51	0.41
1:A:241:SER:HA	1:A:356:ASN:HD22	1.85	0.41
1:A:401:LYS:O	1:A:402:ARG:HB2	2.21	0.41
1:A:402:ARG:HB3	1:A:405:VAL:HG11	2.02	0.41
6:B:502:ZPN:H31	6:B:502:ZPN:H61	1.84	0.41
1:A:122:ILE:HD13	1:A:157:LEU:HD21	2.03	0.41
1:A:401:LYS:CG	2:B:346:TRP:CD2	3.04	0.41
1:A:223:THR:HA	2:B:325:MET:HB2	2.03	0.41
1:A:318:LEU:HD23	1:A:354:GLY:HA3	2.01	0.41
2:B:402:LYS:HB3	2:B:405:LEU:HD22	2.03	0.41
1:A:235:VAL:HA	1:A:238:ILE:HD12	2.02	0.41
2:B:175:PRO:HD2	2:B:205:ASP:OD2	2.21	0.41
2:B:259:MET:O	2:B:314:THR:HG21	2.21	0.41
2:B:151:THR:CG2	2:B:193:GLN:HB2	2.51	0.40
2:B:3:GLU:HG3	2:B:50:ASN:O	2.22	0.40
2:B:44:LEU:HA	2:B:49:ILE:HB	2.03	0.40
2:B:344:VAL:HG12	2:B:346:TRP:H	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	422/437 (97%)	402 (95%)	19 (4%)	1 (0%)	52	88
2	B	424/426 (100%)	401 (95%)	21 (5%)	2 (0%)	34	78
All	All	846/863 (98%)	803 (95%)	40 (5%)	3 (0%)	43	81

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	57	ALA
2	B	369	ARG
1	A	348	PRO

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	359/368 (98%)	356 (99%)	3 (1%)	86	95
2	B	367/367 (100%)	366 (100%)	1 (0%)	94	99
All	All	726/735 (99%)	722 (99%)	4 (1%)	91	97

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

Mol	Chain	Res	Type
1	A	3	GLU
1	A	136	LEU
1	A	401	LYS

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Mol	Chain	Res	Type
2	B	170	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	61	HIS
1	A	133	GLN
1	A	256	GLN
2	B	102	ASN
2	B	107	HIS
2	B	193	GLN
2	B	258	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	0.92	1 (3%)	29,54,54	1.61	3 (10%)
5	GDP	B	501	-	24,30,30	1.12	2 (8%)	26,47,47	1.95	5 (19%)
6	ZPN	B	502	-	35,37,37	0.86	1 (2%)	36,48,48	1.93	11 (30%)

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
6	ZPN	B	502	-	-	0/40/53/53	0/0/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	GTP	C6-N1	2.97	1.38	1.33
5	B	501	GDP	C5-C4	2.99	1.47	1.40
5	B	501	GDP	C6-C5	3.36	1.48	1.41
6	B	502	ZPN	C20-N20	3.50	1.47	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	GTP	N3-C2-N1	-5.14	120.56	127.56
6	B	502	ZPN	C3-C4-C5	-4.69	121.93	127.69
5	B	501	GDP	C5-C6-N1	-4.61	117.50	123.52
3	A	501	GTP	C5-C6-N1	-3.69	118.69	123.52
5	B	501	GDP	N3-C2-N1	-3.57	122.71	127.56
5	B	501	GDP	C6-C5-C4	-3.56	116.79	120.86
6	B	502	ZPN	C10-C11-C12	-2.79	105.90	113.04
6	B	502	ZPN	C3'-C4'-C5'	-2.70	119.87	125.07
6	B	502	ZPN	O1'-C1'-N20	-2.60	118.25	122.34
6	B	502	ZPN	C23-C17-C18	-2.53	112.12	115.58
6	B	502	ZPN	C19-C20-N20	-2.41	101.80	110.39
5	B	501	GDP	C1'-N9-C4	-2.33	124.20	126.81
6	B	502	ZPN	O19-C1-O1	-2.26	119.92	123.39
6	B	502	ZPN	C9-C8-C7	-2.04	110.11	115.04
6	B	502	ZPN	O1'-C1'-C2'	2.58	127.44	122.92
6	B	502	ZPN	C19-O19-C1	2.76	121.69	117.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	GTP	C6-N1-C2	3.97	120.54	115.88
5	B	501	GDP	C6-N1-C2	6.06	122.98	115.88
6	B	502	ZPN	O19-C19-C18	6.10	122.16	106.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	502	ZPN	36	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.