



Full wwPDB EM Validation Report ⓘ

Mar 29, 2026 – 10:30 AM UTC

PDB ID : 6VPO / pdb_00006vpo
EMDB ID : EMD-21314
Title : Cryo-EM structure of microtubule-bound KLP61F motor domain in the AMPPNP state
Authors : Bodrug, T.; Wilson-Kubalek, E.M.; Nithianantham, S.; Debs, G.; Sindelar, C.V.; Milligan, R.; Al-Bassam, J.
Deposited on : 2020-02-04
Resolution : 4.40 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

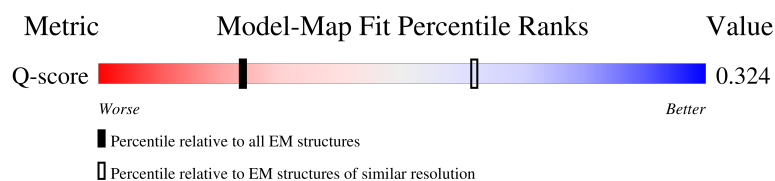
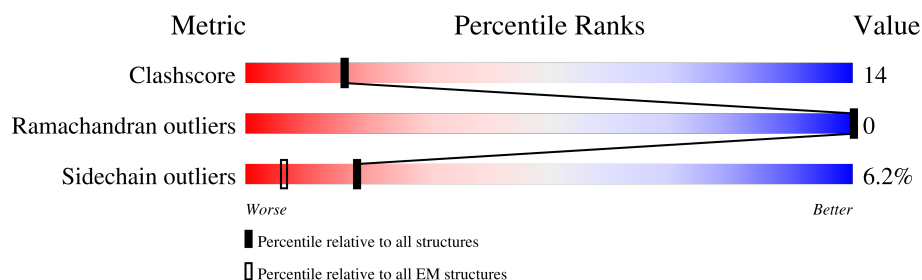
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.40 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	3132 (3.91 - 4.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	451	
2	B	445	
3	C	377	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9266 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-1A chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	422	Total	C	N	O	S	0	0
			3295	2087	561	626	21		

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	421	Total	C	N	O	S	0	0
			3304	2077	565	637	25		

- Molecule 3 is a protein called Kinesin-like protein Klp61F.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	330	Total	C	N	O	S	0	0
			2575	1604	456	505	10		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	370	LEU	-	expression tag	UNP P46863
C	371	GLU	-	expression tag	UNP P46863
C	372	HIS	-	expression tag	UNP P46863
C	373	HIS	-	expression tag	UNP P46863
C	374	HIS	-	expression tag	UNP P46863
C	375	HIS	-	expression tag	UNP P46863
C	376	HIS	-	expression tag	UNP P46863
C	377	HIS	-	expression tag	UNP P46863

- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).

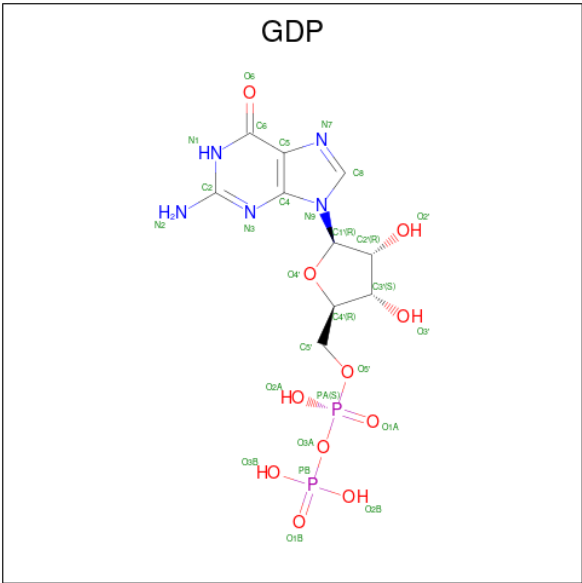


Mol	Chain	Residues	Atoms					AltConf
4	A	1	Total	C	N	O	P	0
			32	10	5	14	3	

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

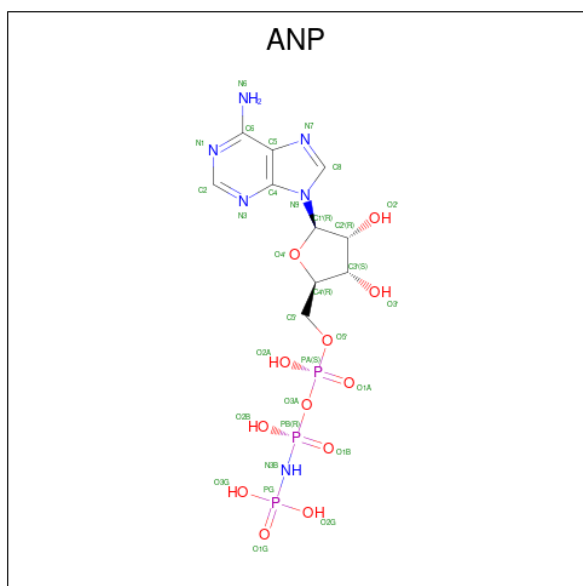
Mol	Chain	Residues	Atoms		AltConf
5	B	1	Total	Mg	0
			1	1	

- Molecule 6 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

- Molecule 7 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$) (labeled as "Ligand of Interest" by depositor).

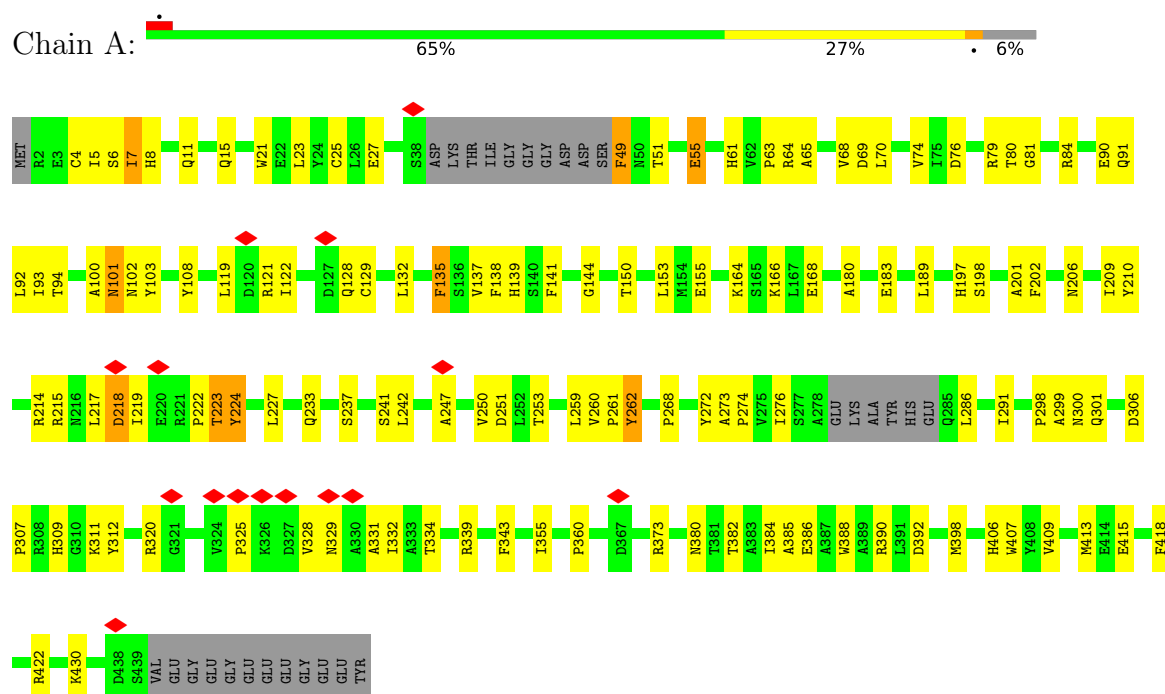


Mol	Chain	Residues	Atoms					AltConf
7	C	1	Total	C	N	O	P	0
			31	10	6	12	3	

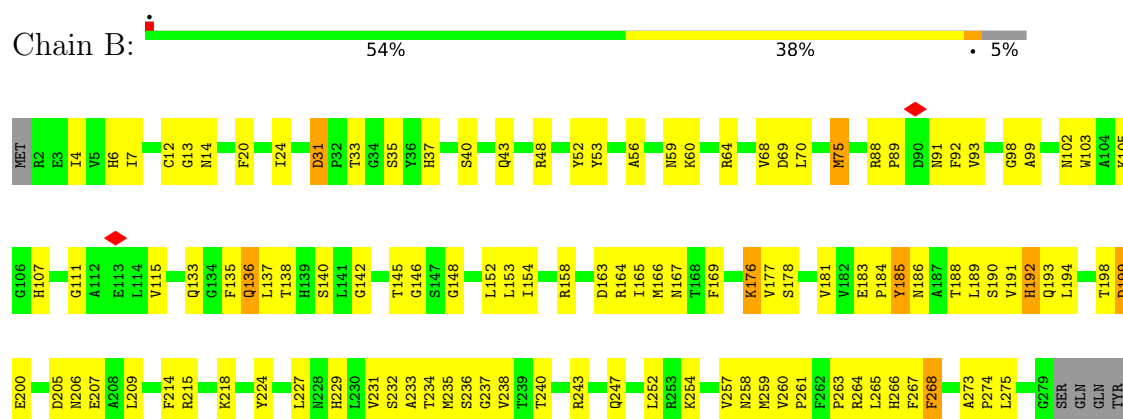
3 Residue-property plots

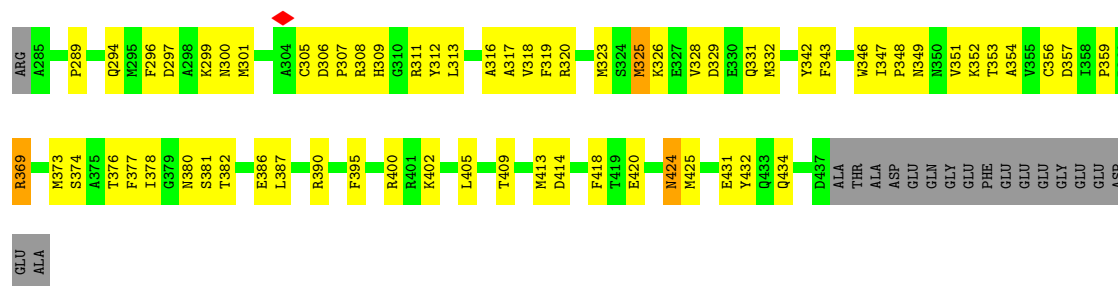
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. 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. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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-1A chain

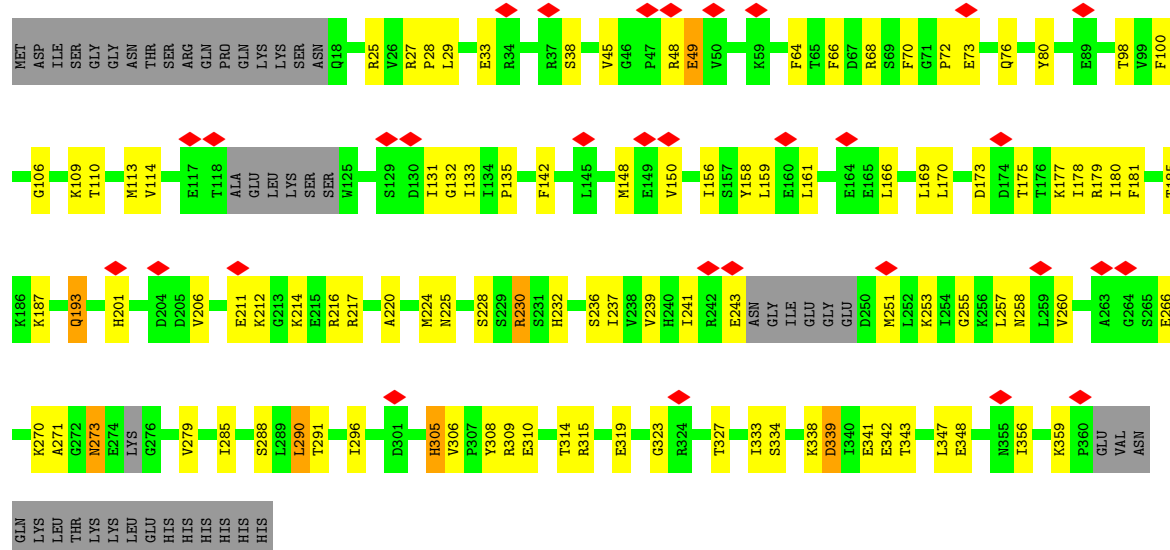


• Molecule 2: Tubulin beta chain





• Molecule 3: Kinesin-like protein Klp61F



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-34.61°, rise=17.4 Å, axial sym=C1	Depositor
Number of segments used	39220	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{Å}^2$)	38	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.110	Depositor
Minimum map value	-0.057	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.025	Depositor
Map size (Å)	83.84, 175.54, 103.49	wwPDB
Map dimensions	64, 134, 79	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.31, 1.31, 1.31	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: GTP, ANP, MG, GDP

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	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.16	0/3369	0.34	0/4573
2	B	0.16	0/3377	0.31	0/4575
3	C	0.13	0/2606	0.31	0/3515
All	All	0.15	0/9352	0.32	0/12663

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 [i](#)

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	3295	0	3205	86	0
2	B	3304	0	3185	130	0
3	C	2575	0	2589	58	0
4	A	32	0	12	1	0
5	B	1	0	0	0	0
6	B	28	0	12	2	0
7	C	31	0	13	5	0
All	All	9266	0	9016	260	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 14.

All (260) 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:115:VAL:HG21	2:B:152:LEU:HD23	1.67	0.77
2:B:320:ARG:HB2	2:B:374:SER:HB3	1.66	0.77
3:C:179:ARG:H	3:C:193:GLN:HE22	1.32	0.76
1:A:398:MET:HE3	2:B:348:PRO:HD2	1.70	0.74
3:C:100:PHE:HB3	3:C:260:VAL:HB	1.69	0.74
2:B:136:GLN:HG2	2:B:167:ASN:HD21	1.53	0.73
2:B:267:PHE:O	2:B:380:ASN:ND2	2.22	0.73
2:B:238:VAL:HG22	2:B:378:ILE:HD11	1.74	0.69
2:B:400:ARG:O	2:B:402:LYS:NZ	2.25	0.69
2:B:352:LYS:NZ	2:B:353:THR:O	2.26	0.69
1:A:5:ILE:HG13	1:A:135:PHE:HB3	1.74	0.69
2:B:154:ILE:HG23	2:B:166:MET:HE2	1.74	0.68
2:B:176:LYS:HG2	2:B:177:VAL:HG23	1.74	0.68
1:A:11:GLN:NE2	2:B:247:GLN:O	2.27	0.67
1:A:68:VAL:HG22	1:A:93:ILE:HB	1.77	0.67
2:B:14:ASN:ND2	2:B:75:MET:SD	2.69	0.66
3:C:64:PHE:HB3	3:C:66:PHE:HE1	1.60	0.66
2:B:297:ASP:OD2	2:B:299:LYS:NZ	2.28	0.66
2:B:35:SER:OG	2:B:60:LYS:NZ	2.28	0.65
2:B:332:MET:HG3	2:B:351:VAL:HG11	1.79	0.64
1:A:215:ARG:HH12	1:A:299:ALA:HB1	1.63	0.64
2:B:254:LYS:HD3	2:B:258:ASN:HD21	1.62	0.64
2:B:99:ALA:HB3	2:B:145:THR:HG22	1.79	0.64
1:A:233:GLN:NE2	1:A:272:TYR:OH	2.31	0.64
2:B:268:PHE:HD2	2:B:380:ASN:HB2	1.63	0.64
1:A:307:PRO:HB3	1:A:312:TYR:HE1	1.61	0.63
3:C:333:ILE:HD13	3:C:347:LEU:HD21	1.80	0.63
1:A:81:GLY:O	1:A:84:ARG:NH1	2.32	0.63
1:A:100:ALA:O	1:A:102:ASN:ND2	2.32	0.63
1:A:132:LEU:O	1:A:164:LYS:NZ	2.22	0.62
3:C:25:ARG:HH12	3:C:76:GLN:HE22	1.46	0.62
1:A:108:TYR:HB2	1:A:413:MET:HE2	1.82	0.62
1:A:180:ALA:N	1:A:183:GLU:OE2	2.31	0.62
1:A:392:ASP:OD2	1:A:422:ARG:NH2	2.32	0.62
2:B:420:GLU:OE2	3:C:315:ARG:NH2	2.32	0.62
2:B:325:MET:HA	2:B:328:VAL:HB	1.82	0.61
2:B:205:ASP:OD1	2:B:207:GLU:N	2.33	0.61
1:A:320:ARG:HB3	1:A:360:PRO:HD3	1.81	0.61
2:B:206:ASN:ND2	6:B:502:GDP:O2'	2.34	0.60
3:C:49:GLU:HG2	3:C:68:ARG:HA	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:215:ARG:O	2:B:218:LYS:NZ	2.32	0.60
3:C:109:LYS:NZ	7:C:501:ANP:O1G	2.34	0.60
1:A:311:LYS:H	1:A:382:THR:HG1	1.50	0.60
1:A:385:ALA:HA	1:A:388:TRP:HD1	1.67	0.59
2:B:294:GLN:O	2:B:300:ASN:ND2	2.34	0.59
3:C:150:VAL:HG23	3:C:243:GLU:HB3	1.83	0.59
1:A:286:LEU:HB3	1:A:291:ILE:HD11	1.85	0.59
1:A:409:VAL:HG21	3:C:290:LEU:HD13	1.85	0.59
1:A:49:PHE:HD1	1:A:51:THR:HG23	1.68	0.59
2:B:247:GLN:NE2	2:B:357:ASP:OD1	2.36	0.58
2:B:209:LEU:HD22	2:B:227:LEU:HB3	1.85	0.58
1:A:8:HIS:HA	1:A:138:PHE:HB2	1.85	0.58
2:B:405:LEU:O	2:B:409:THR:N	2.33	0.58
3:C:27:ARG:NH1	3:C:28:PRO:O	2.37	0.58
2:B:99:ALA:O	2:B:102:ASN:ND2	2.37	0.57
2:B:264:ARG:HG3	3:C:310:GLU:HG3	1.86	0.57
1:A:329:ASN:HD22	1:A:332:ILE:HD12	1.70	0.57
2:B:274:PRO:HG3	2:B:374:SER:HA	1.87	0.57
1:A:63:PRO:O	1:A:91:GLN:NE2	2.33	0.56
7:C:501:ANP:O2G	7:C:501:ANP:O2A	2.23	0.56
2:B:329:ASP:HA	2:B:332:MET:HE3	1.88	0.56
1:A:64:ARG:NH1	1:A:129:CYS:SG	2.79	0.56
1:A:209:ILE:HG22	1:A:227:LEU:HD22	1.87	0.56
2:B:261:PRO:HG3	2:B:313:LEU:HD13	1.85	0.56
7:C:501:ANP:O1G	7:C:501:ANP:O2B	2.22	0.56
2:B:145:THR:OG1	6:B:502:GDP:O2B	2.23	0.56
1:A:7:ILE:HD13	1:A:153:LEU:HD21	1.88	0.55
2:B:320:ARG:HB3	2:B:359:PRO:HA	1.89	0.55
1:A:166:LYS:HE3	1:A:198:SER:HA	1.89	0.55
1:A:223:THR:OG1	1:A:224:TYR:N	2.40	0.55
2:B:166:MET:HB2	2:B:199:ASP:H	1.71	0.55
2:B:4:ILE:HG12	2:B:133:GLN:HB3	1.89	0.55
3:C:80:TYR:HD2	3:C:133:ILE:HG23	1.71	0.54
1:A:90:GLU:HG3	1:A:121:ARG:HH21	1.73	0.54
1:A:218:ASP:OD1	1:A:218:ASP:N	2.40	0.54
2:B:328:VAL:O	2:B:332:MET:N	2.37	0.53
2:B:386:GLU:HB3	2:B:390:ARG:HH12	1.72	0.53
1:A:6:SER:HB2	1:A:65:ALA:HA	1.90	0.53
2:B:311:ARG:H	2:B:382:THR:HG1	1.57	0.53
2:B:31:ASP:OD2	2:B:33:THR:OG1	2.26	0.53
1:A:268:PRO:HG3	1:A:380:ASN:HD21	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:ASN:HA	2:B:200:GLU:HG3	1.91	0.53
1:A:237:SER:O	1:A:241:SER:OG	2.28	0.52
1:A:331:ALA:O	1:A:334:THR:OG1	2.24	0.52
3:C:98:THR:HB	3:C:327:THR:HG23	1.90	0.52
2:B:142:GLY:O	2:B:186:ASN:ND2	2.43	0.52
2:B:349:ASN:ND2	2:B:351:VAL:O	2.42	0.52
3:C:230:ARG:NH2	3:C:266:GLU:OE1	2.42	0.52
3:C:333:ILE:HB	3:C:343:THR:HG23	1.91	0.52
1:A:5:ILE:HG22	1:A:64:ARG:HB3	1.92	0.51
1:A:320:ARG:HH21	1:A:360:PRO:HA	1.74	0.51
2:B:89:PRO:HA	2:B:92:PHE:HD1	1.75	0.51
2:B:191:VAL:HB	2:B:425:MET:HE1	1.92	0.51
2:B:316:ALA:HB3	2:B:378:ILE:HB	1.93	0.51
3:C:113:MET:HA	3:C:133:ILE:HD12	1.93	0.51
2:B:301:MET:HG3	2:B:307:PRO:HG3	1.93	0.51
2:B:318:VAL:HB	2:B:376:THR:HB	1.93	0.51
2:B:237:GLY:HA3	2:B:376:THR:HG21	1.92	0.51
3:C:237:ILE:N	3:C:257:LEU:O	2.39	0.51
2:B:7:ILE:HB	2:B:137:LEU:HG	1.92	0.51
2:B:236:SER:O	2:B:240:THR:OG1	2.24	0.51
2:B:69:ASP:OD1	2:B:70:LEU:N	2.41	0.50
3:C:25:ARG:HA	3:C:70:PHE:HB2	1.93	0.50
3:C:239:VAL:O	3:C:255:GLY:N	2.43	0.50
1:A:21:TRP:CZ2	1:A:65:ALA:HB2	2.47	0.50
1:A:407:TRP:CD1	2:B:257:VAL:HG13	2.47	0.50
1:A:242:LEU:HD22	1:A:250:VAL:H	1.77	0.49
3:C:161:LEU:HB3	3:C:232:HIS:HB2	1.94	0.49
3:C:251:MET:O	3:C:253:LYS:NZ	2.44	0.49
3:C:296:ILE:HG23	3:C:356:ILE:HD11	1.94	0.49
1:A:251:ASP:OD2	1:A:253:THR:OG1	2.23	0.49
3:C:225:ASN:HB3	3:C:228:SER:HB2	1.93	0.49
1:A:286:LEU:O	1:A:373:ARG:NH2	2.43	0.49
2:B:107:HIS:HB2	2:B:148:GLY:HA2	1.95	0.49
1:A:119:LEU:HD23	1:A:122:ILE:HD12	1.95	0.49
1:A:306:ASP:OD2	1:A:309:HIS:N	2.46	0.49
3:C:106:GLY:H	7:C:501:ANP:HNB1	1.60	0.49
3:C:179:ARG:N	3:C:193:GLN:HE22	2.06	0.49
3:C:241:ILE:N	3:C:253:LYS:O	2.41	0.48
2:B:158:ARG:NH1	2:B:199:ASP:OD1	2.47	0.48
2:B:260:VAL:HG22	2:B:266:HIS:HB3	1.95	0.48
2:B:166:MET:N	2:B:199:ASP:OD2	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:312:TYR:CD2	2:B:381:SER:HB2	2.49	0.48
3:C:64:PHE:HB3	3:C:66:PHE:CE1	2.45	0.48
2:B:192:HIS:HB3	2:B:193:GLN:HE21	1.79	0.48
2:B:186:ASN:O	2:B:190:SER:N	2.47	0.47
2:B:24:ILE:HD12	2:B:243:ARG:HH22	1.80	0.47
1:A:103:TYR:CD2	1:A:189:LEU:HD13	2.49	0.47
1:A:222:PRO:HD2	2:B:326:LYS:HD2	1.96	0.47
1:A:155:GLU:HG3	1:A:197:HIS:CE1	2.50	0.47
2:B:68:VAL:HG22	2:B:93:VAL:HB	1.95	0.47
3:C:131:ILE:HG12	3:C:135:PRO:HG2	1.97	0.47
1:A:210:TYR:CE1	2:B:326:LYS:HB2	2.49	0.47
1:A:325:PRO:HG3	1:A:355:ILE:HD13	1.97	0.47
2:B:240:THR:HG22	2:B:356:CYS:SG	2.55	0.47
2:B:178:SER:HB2	2:B:183:GLU:OE1	2.14	0.47
2:B:189:LEU:HD11	2:B:418:PHE:CE1	2.51	0.46
3:C:185:THR:O	3:C:187:LYS:NZ	2.39	0.46
3:C:236:SER:HA	3:C:258:ASN:HA	1.97	0.46
1:A:101:ASN:HA	1:A:144:GLY:HA3	1.98	0.46
1:A:206:ASN:OD1	1:A:206:ASN:N	2.48	0.46
3:C:273:ASN:OD1	3:C:273:ASN:N	2.47	0.46
1:A:217:LEU:HB3	1:A:219:ILE:HG12	1.97	0.46
1:A:260:VAL:HG12	1:A:262:TYR:O	2.15	0.46
2:B:346:TRP:CD1	2:B:347:ILE:HG13	2.51	0.46
2:B:53:TYR:O	2:B:64:ARG:NH2	2.49	0.46
2:B:188:THR:HG21	2:B:395:PHE:CG	2.51	0.46
3:C:158:TYR:O	3:C:169:LEU:N	2.47	0.46
1:A:55:GLU:OE2	1:A:61:HIS:NE2	2.49	0.45
3:C:173:ASP:OD1	3:C:175:THR:OG1	2.30	0.45
2:B:305:CYS:SG	2:B:386:GLU:HB2	2.57	0.45
1:A:79:ARG:HD3	1:A:92:LEU:HD13	1.99	0.45
2:B:103:TRP:HZ3	2:B:413:MET:HE3	1.81	0.45
2:B:318:VAL:HG22	2:B:354:ALA:HB3	1.99	0.45
3:C:80:TYR:CD2	3:C:133:ILE:HG23	2.51	0.45
2:B:183:GLU:N	2:B:184:PRO:HD2	2.32	0.45
2:B:231:VAL:O	2:B:234:THR:OG1	2.28	0.45
2:B:312:TYR:HB2	2:B:343:PHE:CD2	2.52	0.45
2:B:176:LYS:HD2	2:B:207:GLU:HG3	1.99	0.45
2:B:296:PHE:HZ	2:B:342:TYR:HD1	1.65	0.45
3:C:214:LYS:HA	3:C:217:ARG:HB2	1.99	0.45
1:A:23:LEU:O	1:A:27:GLU:N	2.41	0.45
2:B:209:LEU:HB3	2:B:227:LEU:HD13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:273:ALA:O	2:B:294:GLN:NE2	2.50	0.45
2:B:329:ASP:HA	2:B:332:MET:HB3	1.99	0.45
1:A:210:TYR:O	1:A:214:ARG:HG2	2.17	0.44
2:B:75:MET:HE3	2:B:92:PHE:HD2	1.81	0.44
3:C:224:MET:HG2	3:C:271:ALA:HA	1.98	0.44
2:B:102:ASN:HD22	2:B:105:LYS:HD3	1.81	0.44
3:C:339:ASP:OD1	3:C:339:ASP:N	2.49	0.44
2:B:308:ARG:HG2	2:B:342:TYR:CE1	2.52	0.44
2:B:312:TYR:HB2	2:B:343:PHE:HD2	1.83	0.44
1:A:259:LEU:HG	1:A:380:ASN:HD22	1.81	0.44
1:A:261:PRO:HG2	1:A:262:TYR:CZ	2.53	0.44
1:A:298:PRO:HA	1:A:301:GLN:HG3	1.98	0.44
2:B:163:ASP:OD2	2:B:164:ARG:NH2	2.50	0.44
3:C:220:ALA:HB3	3:C:230:ARG:HD2	1.99	0.44
1:A:21:TRP:CZ3	1:A:63:PRO:HB3	2.53	0.44
2:B:40:SER:HB2	2:B:43:GLN:HG3	2.00	0.44
1:A:76:ASP:OD2	2:B:48:ARG:NH2	2.49	0.44
1:A:139:HIS:CD2	1:A:150:THR:HG21	2.52	0.44
1:A:406:HIS:CG	2:B:263:PRO:HG3	2.53	0.43
2:B:289:PRO:HB3	2:B:331:GLN:NE2	2.33	0.43
1:A:21:TRP:O	1:A:25:CYS:N	2.40	0.43
2:B:194:LEU:O	2:B:198:THR:N	2.51	0.43
2:B:88:ARG:HG3	2:B:91:ASN:OD1	2.19	0.43
1:A:210:TYR:CE1	1:A:227:LEU:HD11	2.54	0.43
1:A:224:TYR:HD1	1:A:224:TYR:HA	1.74	0.43
2:B:12:CYS:SG	2:B:140:SER:HB2	2.59	0.43
2:B:89:PRO:HA	2:B:92:PHE:CD1	2.54	0.43
1:A:273:ALA:N	1:A:274:PRO:HD3	2.34	0.43
2:B:56:ALA:HB3	2:B:60:LYS:HB2	2.01	0.43
3:C:156:ILE:HG13	3:C:170:LEU:HD22	2.01	0.43
3:C:288:SER:O	3:C:291:THR:OG1	2.34	0.43
3:C:28:PRO:HG3	3:C:72:PRO:HB2	2.01	0.43
1:A:76:ASP:O	1:A:80:THR:OG1	2.22	0.42
2:B:145:THR:OG1	2:B:146:GLY:N	2.52	0.42
3:C:212:LYS:O	3:C:216:ARG:N	2.50	0.42
1:A:233:GLN:HG2	1:A:272:TYR:CE2	2.53	0.42
2:B:296:PHE:CZ	2:B:342:TYR:HD1	2.37	0.42
1:A:69:ASP:OD1	1:A:70:LEU:N	2.52	0.42
2:B:20:PHE:O	2:B:24:ILE:HG12	2.20	0.42
2:B:352:LYS:HA	2:B:352:LYS:HD2	1.88	0.42
3:C:323:GLY:O	3:C:359:LYS:N	2.42	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:285:ILE:O	3:C:288:SER:OG	2.35	0.42
3:C:308:TYR:O	3:C:314:THR:OG1	2.24	0.42
1:A:233:GLN:HG2	1:A:272:TYR:CZ	2.55	0.42
2:B:20:PHE:HE1	2:B:235:MET:HB2	1.84	0.42
2:B:369:ARG:HH11	2:B:369:ARG:HB3	1.85	0.42
1:A:168:GLU:HB2	1:A:201:ALA:HA	2.01	0.42
1:A:247:ALA:HB3	1:A:355:ILE:HG21	2.02	0.42
2:B:265:LEU:HD21	2:B:431:GLU:HB3	2.02	0.42
2:B:135:PHE:N	2:B:165:ILE:O	2.50	0.42
3:C:110:THR:OG1	7:C:501:ANP:O1G	2.38	0.42
2:B:52:TYR:OH	2:B:136:GLN:OE1	2.36	0.42
2:B:200:GLU:HB2	2:B:268:PHE:CE1	2.54	0.42
2:B:382:THR:HG22	2:B:432:TYR:HD2	1.84	0.42
2:B:12:CYS:SG	2:B:13:GLY:N	2.92	0.42
2:B:111:GLY:O	2:B:115:VAL:HG23	2.20	0.42
3:C:27:ARG:NH2	3:C:334:SER:OG	2.53	0.41
3:C:114:VAL:O	3:C:132:GLY:HA3	2.20	0.41
3:C:166:LEU:HB2	3:C:178:ILE:HB	2.01	0.41
2:B:431:GLU:HA	3:C:305:HIS:NE2	2.34	0.41
2:B:273:ALA:O	2:B:275:LEU:N	2.52	0.41
2:B:306:ASP:HB3	2:B:309:HIS:CG	2.55	0.41
2:B:70:LEU:HB3	2:B:98:GLY:HA2	2.01	0.41
1:A:386:GLU:HB3	1:A:390:ARG:HH12	1.85	0.41
1:A:406:HIS:CD2	2:B:263:PRO:HG3	2.56	0.41
3:C:211:GLU:HA	3:C:214:LYS:HE2	2.02	0.41
1:A:384:ILE:HD12	1:A:384:ILE:HA	1.91	0.41
2:B:420:GLU:CD	3:C:181:PHE:HA	2.45	0.41
1:A:180:ALA:HA	2:B:352:LYS:HD2	2.03	0.41
1:A:137:VAL:HG12	1:A:139:HIS:HD2	1.86	0.41
1:A:168:GLU:O	1:A:202:PHE:N	2.44	0.41
2:B:206:ASN:HA	2:B:209:LEU:HD12	2.01	0.41
2:B:317:ALA:HB1	2:B:319:PHE:CE1	2.56	0.41
1:A:11:GLN:HE21	1:A:15:GLN:NE2	2.19	0.41
1:A:101:ASN:ND2	4:A:501:GTP:O3G	2.54	0.41
1:A:135:PHE:HE1	1:A:166:LYS:HG2	1.85	0.41
1:A:189:LEU:HD11	1:A:418:PHE:HE2	1.86	0.41
2:B:138:THR:HA	2:B:169:PHE:O	2.21	0.41
2:B:189:LEU:HD21	2:B:418:PHE:CE1	2.56	0.41
2:B:229:HIS:O	2:B:232:SER:OG	2.39	0.41
2:B:233:ALA:O	2:B:236:SER:OG	2.25	0.41
2:B:259:MET:HE2	2:B:378:ILE:HG22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:424:ASN:OD1	3:C:309:ARG:NH2	2.40	0.41
2:B:185:TYR:HD1	2:B:185:TYR:HA	1.70	0.40
2:B:301:MET:HE1	2:B:377:PHE:CZ	2.56	0.40
1:A:141:PHE:HD1	1:A:141:PHE:HA	1.77	0.40
1:A:276:ILE:HD13	1:A:276:ILE:HA	1.95	0.40
1:A:329:ASN:HD22	1:A:329:ASN:HA	1.71	0.40
2:B:323:MET:HE2	2:B:373:MET:HB3	2.03	0.40
3:C:45:VAL:HG11	3:C:48:ARG:HD3	2.02	0.40
1:A:385:ALA:HA	1:A:388:TRP:CD1	2.51	0.40
2:B:434:GLN:OE1	3:C:305:HIS:HD2	2.05	0.40
2:B:52:TYR:HD1	2:B:52:TYR:HA	1.78	0.40
2:B:107:HIS:ND1	2:B:152:LEU:HB2	2.36	0.40
2:B:275:LEU:HD23	2:B:275:LEU:HA	1.78	0.40
3:C:33:GLU:HG3	3:C:38:SER:HB3	2.04	0.40
3:C:306:VAL:HB	3:C:308:TYR:CZ	2.56	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	416/451 (92%)	398 (96%)	18 (4%)	0	100	100
2	B	417/445 (94%)	403 (97%)	14 (3%)	0	100	100
3	C	322/377 (85%)	310 (96%)	12 (4%)	0	100	100
All	All	1155/1273 (91%)	1111 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	355/377 (94%)	336 (95%)	19 (5%)	20	42
2	B	362/381 (95%)	341 (94%)	21 (6%)	18	40
3	C	292/341 (86%)	269 (92%)	23 (8%)	11	32
All	All	1009/1099 (92%)	946 (94%)	63 (6%)	18	38

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	CYS
1	A	7	ILE
1	A	49	PHE
1	A	55	GLU
1	A	74	VAL
1	A	94	THR
1	A	101	ASN
1	A	128	GLN
1	A	135	PHE
1	A	218	ASP
1	A	223	THR
1	A	224	TYR
1	A	262	TYR
1	A	300	ASN
1	A	328	VAL
1	A	339	ARG
1	A	343	PHE
1	A	415	GLU
1	A	430	LYS
2	B	6	HIS
2	B	31	ASP
2	B	37	HIS
2	B	59	ASN
2	B	75	MET
2	B	136	GLN
2	B	153	LEU

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Mol	Chain	Res	Type
2	B	176	LYS
2	B	181	VAL
2	B	185	TYR
2	B	192	HIS
2	B	199	ASP
2	B	214	PHE
2	B	224	TYR
2	B	252	LEU
2	B	268	PHE
2	B	325	MET
2	B	369	ARG
2	B	387	LEU
2	B	414	ASP
2	B	424	ASN
3	C	29	LEU
3	C	49	GLU
3	C	73	GLU
3	C	142	PHE
3	C	148	MET
3	C	159	LEU
3	C	177	LYS
3	C	180	ILE
3	C	193	GLN
3	C	201	HIS
3	C	206	VAL
3	C	230	ARG
3	C	270	LYS
3	C	273	ASN
3	C	279	VAL
3	C	290	LEU
3	C	305	HIS
3	C	319	GLU
3	C	338	LYS
3	C	339	ASP
3	C	341	GLU
3	C	342	GLU
3	C	348	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN

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Mol	Chain	Res	Type
1	A	101	ASN
1	A	128	GLN
1	A	133	GLN
1	A	197	HIS
1	A	258	ASN
1	A	329	ASN
2	B	59	ASN
2	B	96	GLN
2	B	102	ASN
2	B	193	GLN
2	B	258	ASN
2	B	293	GLN
2	B	294	GLN
2	B	336	GLN
2	B	339	ASN
2	B	385	GLN
3	C	193	GLN
3	C	284	ASN
3	C	305	HIS
3	C	337	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
4	GTP	A	501	5	33,34,34	0.90	1 (3%)	50,54,54	1.55	9 (18%)
7	ANP	C	501	-	33,33,33	0.95	4 (12%)	45,52,52	0.58	0
6	GDP	B	502	-	29,30,30	1.16	3 (10%)	45,47,47	1.73	7 (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
4	GTP	A	501	5	-	5/22/38/38	0/3/3/3
7	ANP	C	501	-	-	9/18/38/38	0/3/3/3
6	GDP	B	502	-	-	4/16/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	502	GDP	C5-C4	3.01	1.47	1.38
6	B	502	GDP	C6-N1	-2.66	1.33	1.38
7	C	501	ANP	PB-O3A	-2.62	1.55	1.59
7	C	501	ANP	PG-O1G	2.48	1.49	1.46
7	C	501	ANP	PG-N3B	2.38	1.69	1.63
7	C	501	ANP	PB-O1B	2.26	1.49	1.46
6	B	502	GDP	C5-N7	-2.23	1.34	1.39
4	A	501	GTP	C2-N3	2.10	1.38	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	502	GDP	C5-C4-N3	-5.98	118.88	128.39
4	A	501	GTP	C5-C4-N3	-5.09	120.30	128.39
6	B	502	GDP	C2-N3-C4	4.84	120.64	112.30
4	A	501	GTP	C2-N3-C4	4.49	120.03	112.30
6	B	502	GDP	N9-C4-N3	4.45	134.84	125.95
4	A	501	GTP	N9-C4-N3	3.21	132.38	125.95
6	B	502	GDP	C6-C5-N7	3.15	136.03	130.29
4	A	501	GTP	C2-N1-C6	-2.92	119.81	125.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	501	GTP	N9-C8-N7	-2.53	108.71	113.40
6	B	502	GDP	C4-C5-N7	-2.50	106.71	110.67
4	A	501	GTP	O6-C6-C5	-2.48	119.97	126.53
4	A	501	GTP	C8-N7-C5	2.44	108.60	104.26
4	A	501	GTP	C5-C6-N1	2.43	119.45	113.25
4	A	501	GTP	C3'-C2'-C1'	2.20	105.63	101.46
6	B	502	GDP	C3'-C2'-C1'	2.14	105.51	101.46
6	B	502	GDP	O6-C6-C5	-2.11	120.97	126.53

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	501	GTP	C5'-O5'-PA-O1A
4	A	501	GTP	C5'-O5'-PA-O2A
6	B	502	GDP	C5'-O5'-PA-O3A
6	B	502	GDP	C5'-O5'-PA-O1A
6	B	502	GDP	O4'-C4'-C5'-O5'
7	C	501	ANP	PB-N3B-PG-O1G
7	C	501	ANP	PG-N3B-PB-O1B
7	C	501	ANP	C5'-O5'-PA-O1A
7	C	501	ANP	C5'-O5'-PA-O2A
7	C	501	ANP	C5'-O5'-PA-O3A
6	B	502	GDP	C3'-C4'-C5'-O5'
7	C	501	ANP	O4'-C4'-C5'-O5'
7	C	501	ANP	C3'-C4'-C5'-O5'
4	A	501	GTP	C3'-C4'-C5'-O5'
4	A	501	GTP	O4'-C4'-C5'-O5'
4	A	501	GTP	C5'-O5'-PA-O3A
7	C	501	ANP	PA-O3A-PB-O2B
7	C	501	ANP	PA-O3A-PB-O1B

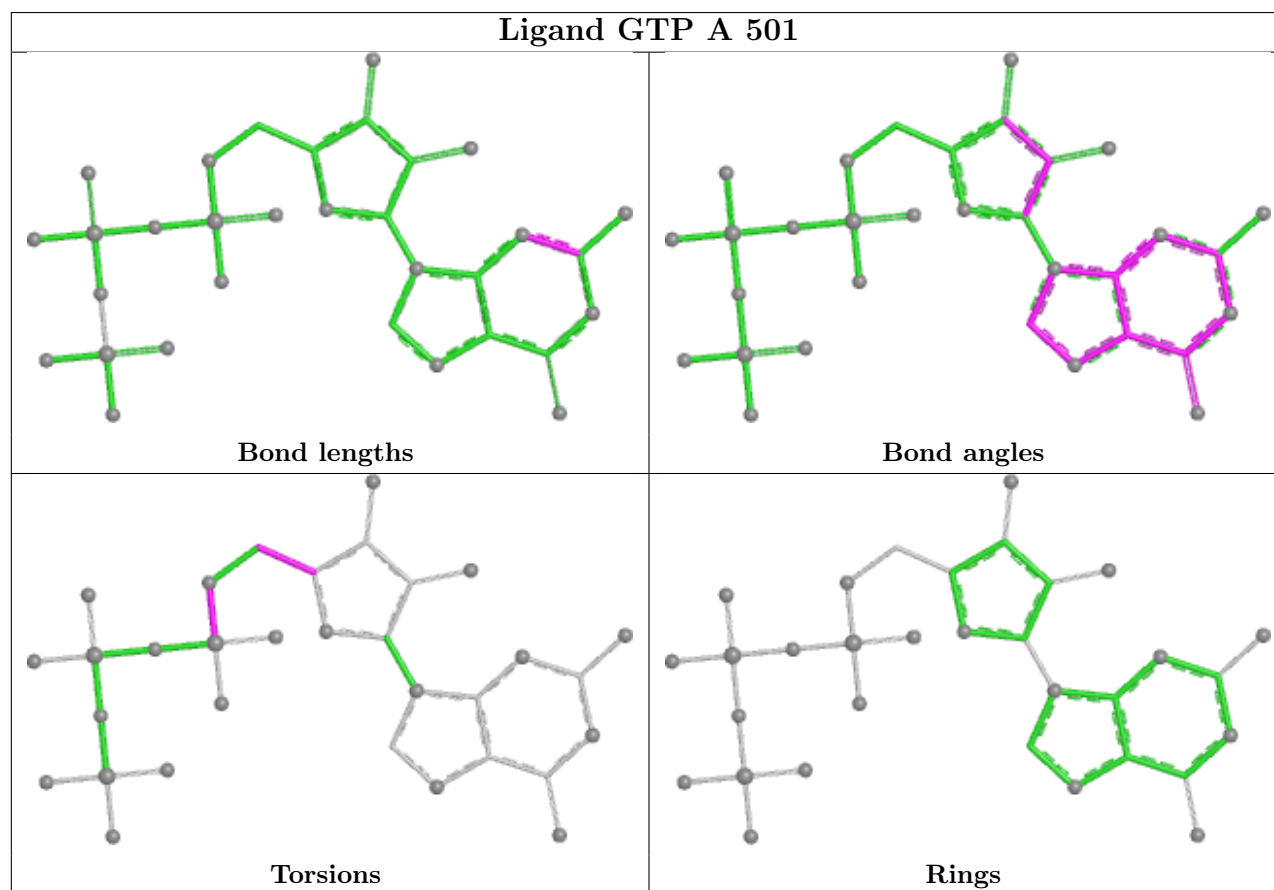
There are no ring outliers.

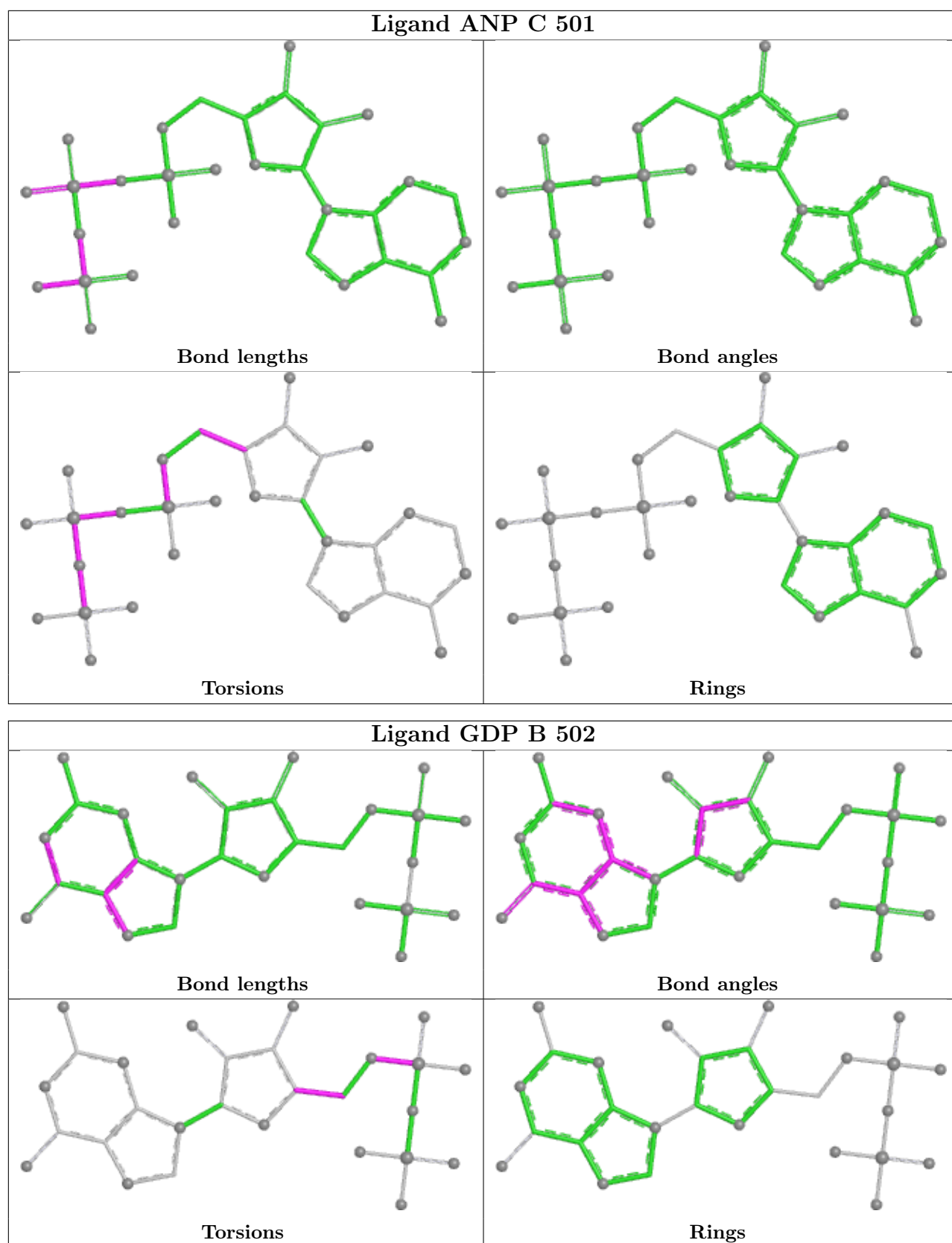
3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	501	GTP	1	0
7	C	501	ANP	5	0
6	B	502	GDP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

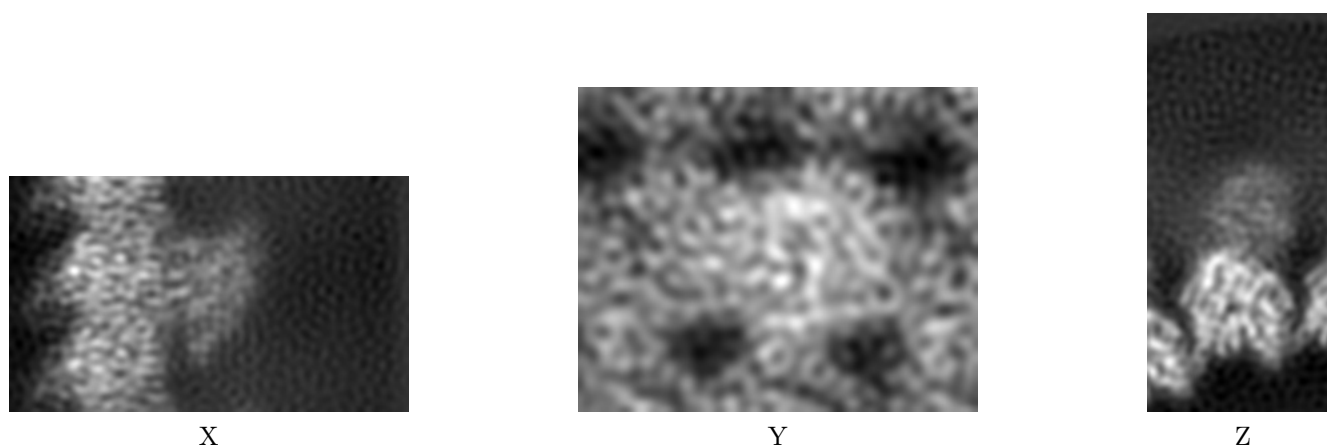
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21314. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

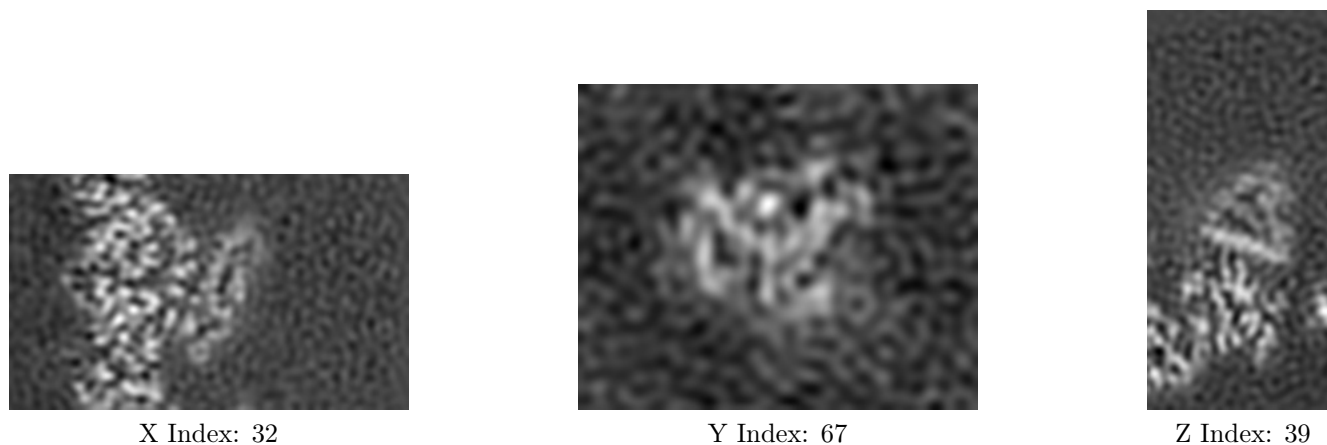
6.1.1 Primary map



The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

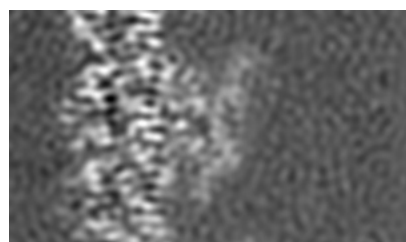
6.2.1 Primary map



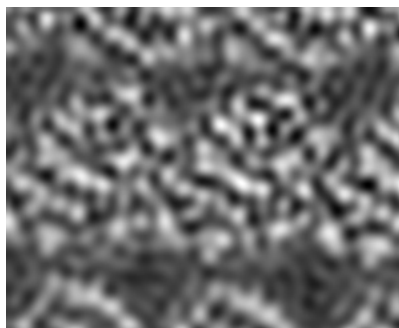
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

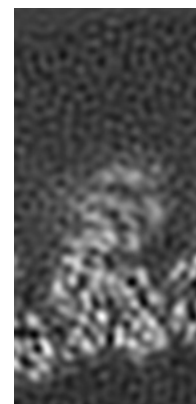
6.3.1 Primary map



X Index: 27



Y Index: 30

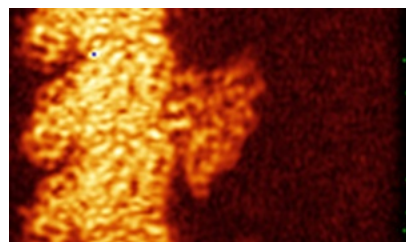


Z Index: 46

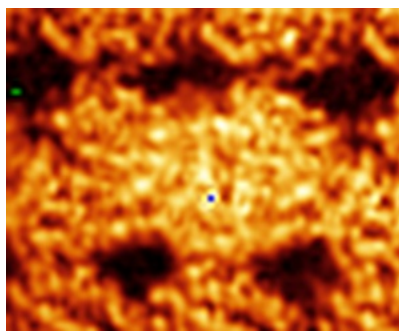
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

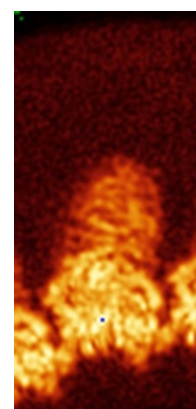
6.4.1 Primary map



X



Y

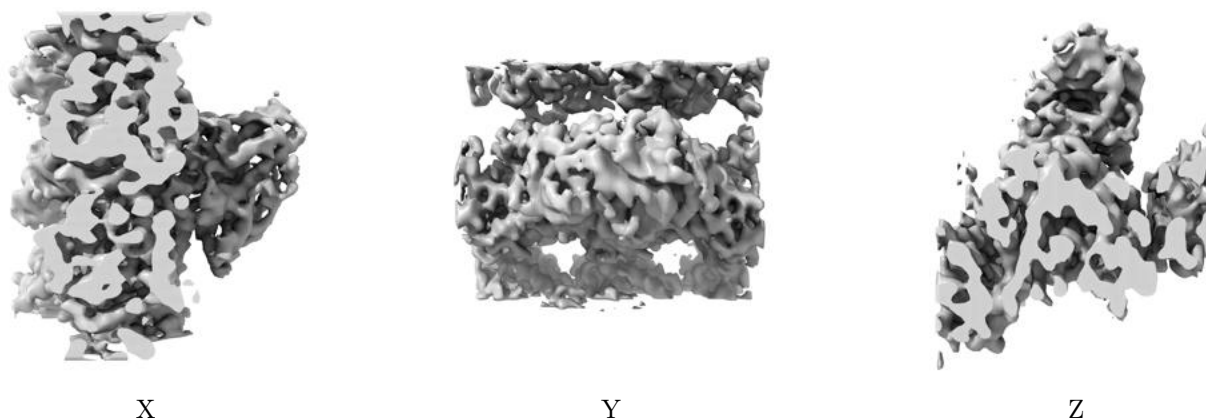


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.025. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

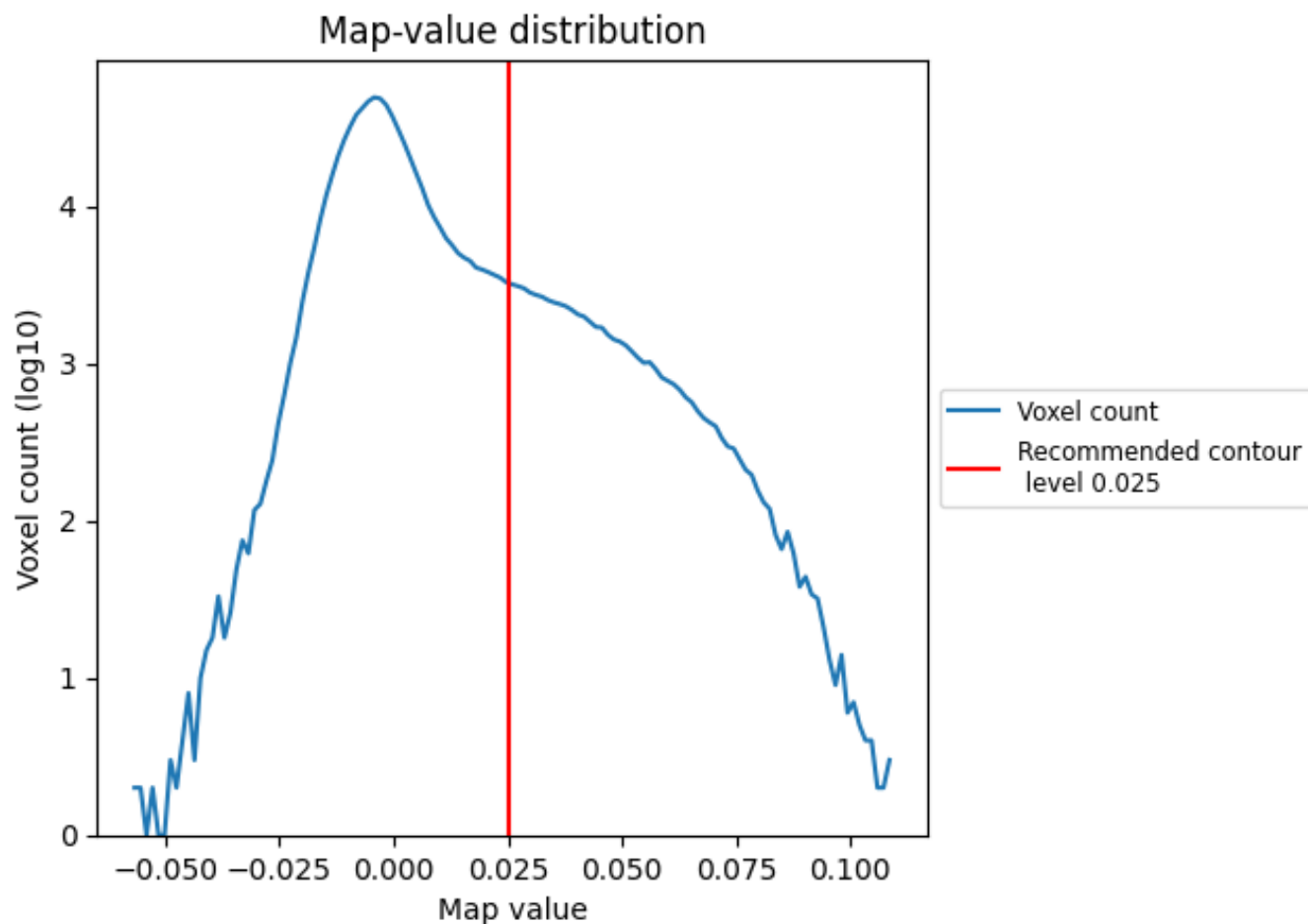
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

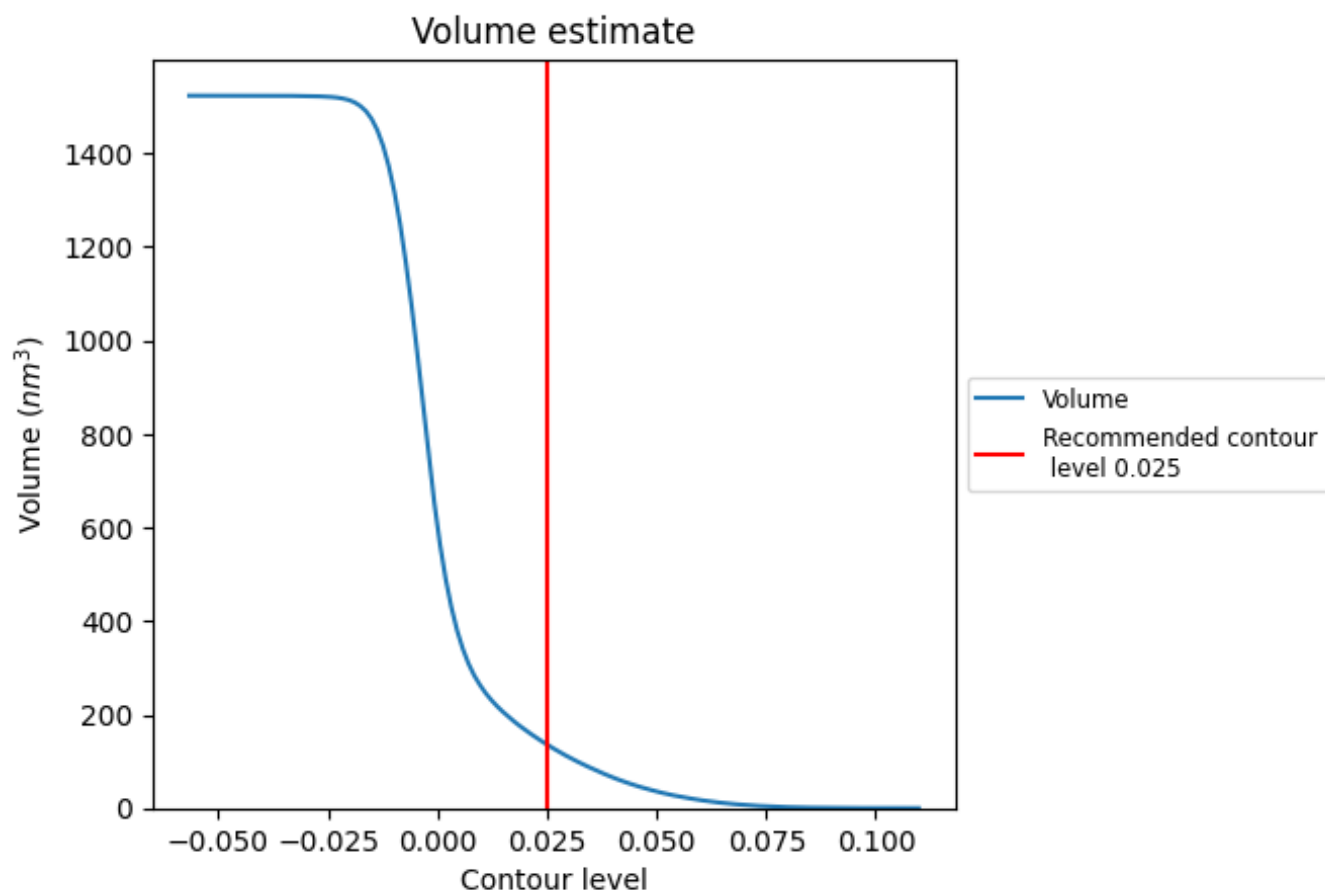
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

7.2 Volume estimate [i](#)



The volume at the recommended contour level is 136 nm³; this corresponds to an approximate mass of 123 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum [i](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

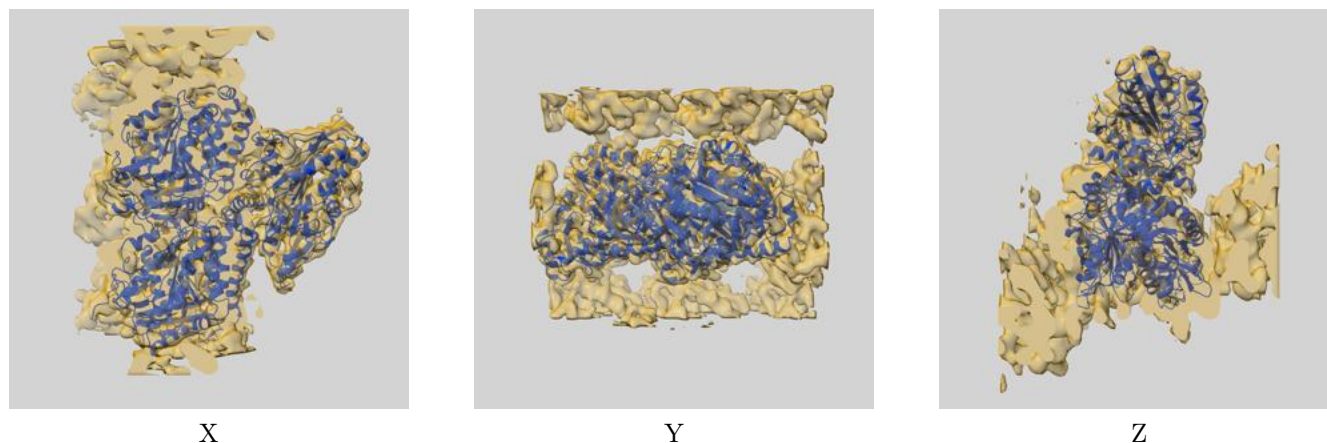
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

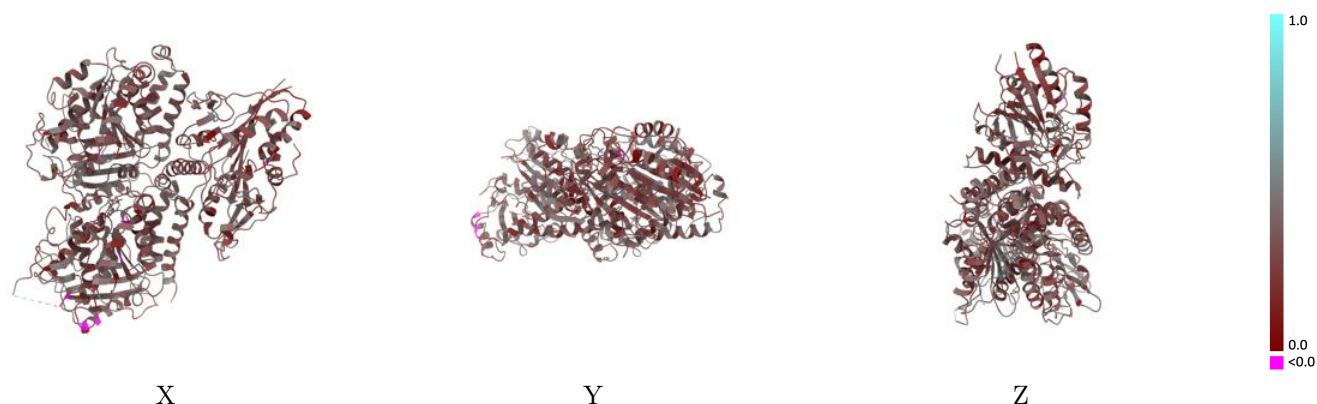
This section contains information regarding the fit between EMDB map EMD-21314 and PDB model 6VPO. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



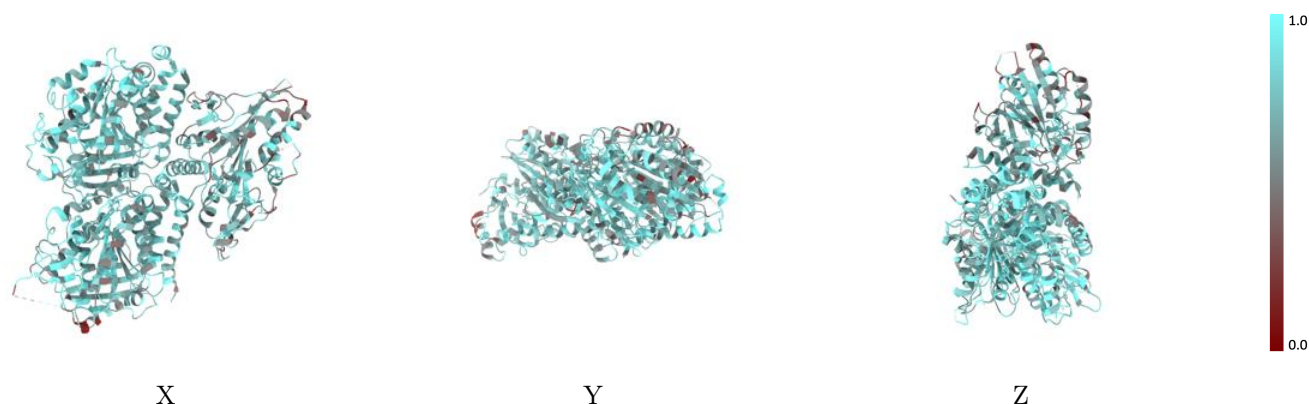
The images above show the 3D surface view of the map at the recommended contour level 0.025 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



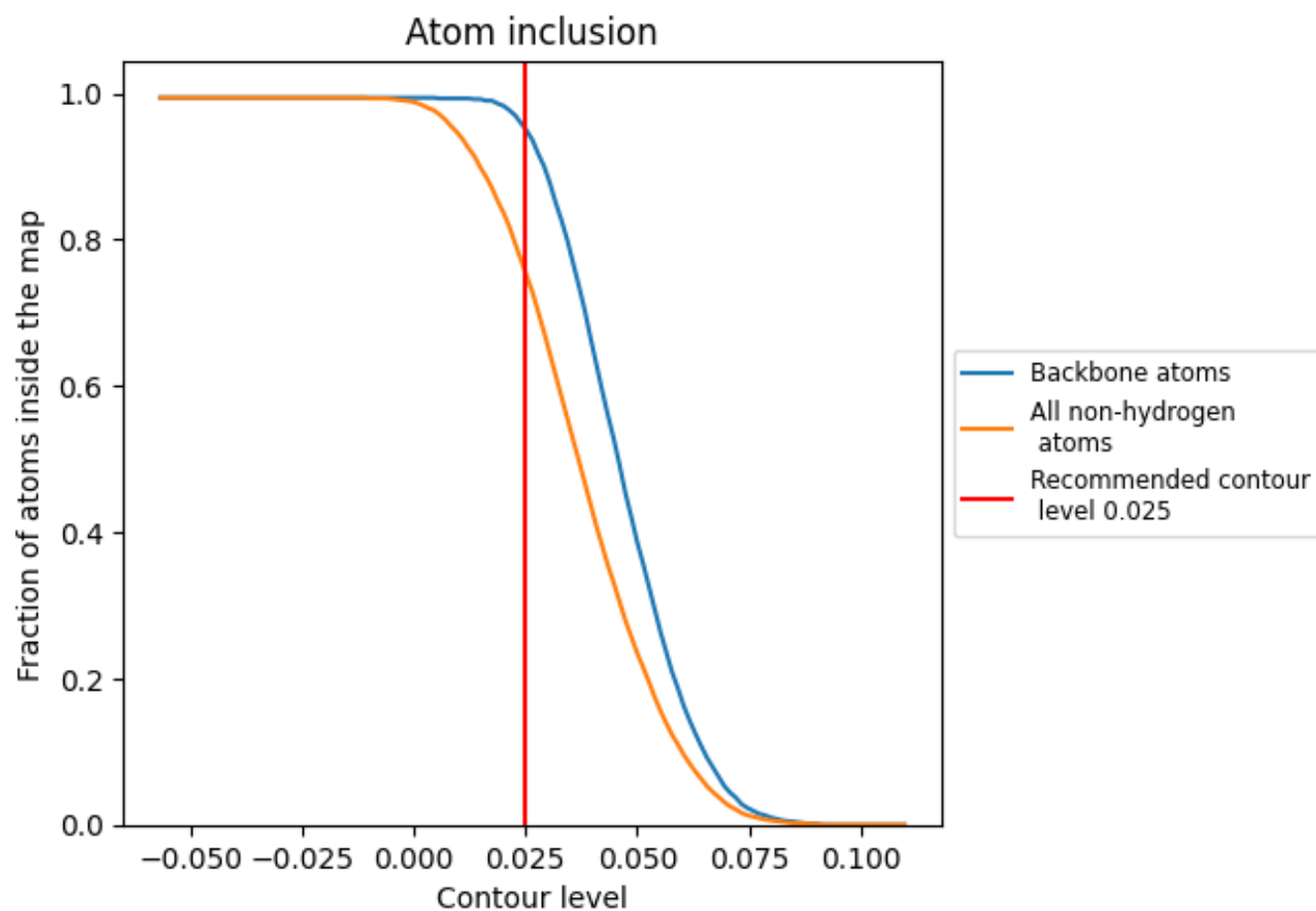
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.025).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 76% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.025) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7550	0.3240
A	0.7630	0.3250
B	0.8120	0.3410
C	0.6720	0.3000

