



Full wwPDB EM Validation Report ⓘ

Dec 21, 2022 – 12:49 pm GMT

PDB ID : 7QTL
EMDB ID : EMD-14144
Title : Influenza A/H7N9 polymerase elongation complex
Authors : Cusack, S.; Kouba, T.
Deposited on : 2022-01-14
Resolution : 2.48 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.dev43
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

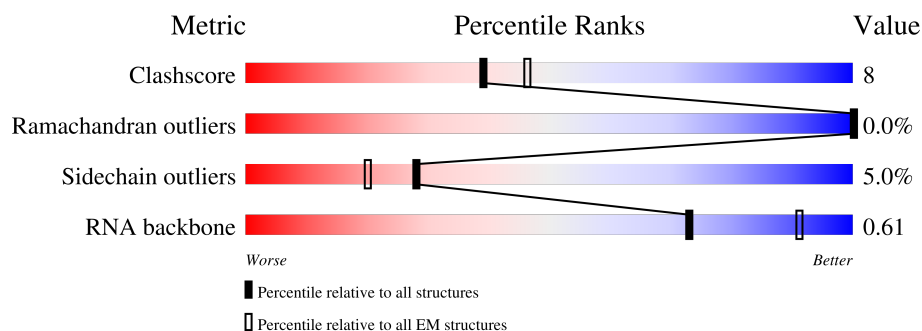
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.48 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	717	<div> <div>5%</div> <div>77%</div> <div>20%</div> <div>..</div> </div>
2	B	757	<div> <div>75%</div> <div>20%</div> <div>5%</div> </div>
3	C	759	<div> <div>73%</div> <div>20%</div> <div>..</div> </div>
4	V	14	<div> <div>7%</div> <div>50%</div> <div>43%</div> <div>7%</div> </div>
5	R	21	<div> <div>5%</div> <div>38%</div> <div>19%</div> <div>10%</div> <div>33%</div> </div>
6	M	17	<div> <div>6%</div> <div>65%</div> <div>24%</div> <div>12%</div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 18397 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 Polymerase acidic protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	709	Total	C	N	O	S	0	0
			5755	3642	979	1093	41		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP M9TI86

- Molecule 2 is a protein called RNA-directed RNA polymerase catalytic subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	717	Total	C	N	O	S	1	0
			5736	3608	1002	1081	45		

- Molecule 3 is a protein called Polymerase basic protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	726	Total	C	N	O	S	0	0
			5767	3628	1038	1061	40		

- Molecule 4 is a RNA chain called 5' vRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	V	13	Total	C	N	O	P	0	0
			284	127	57	87	13		

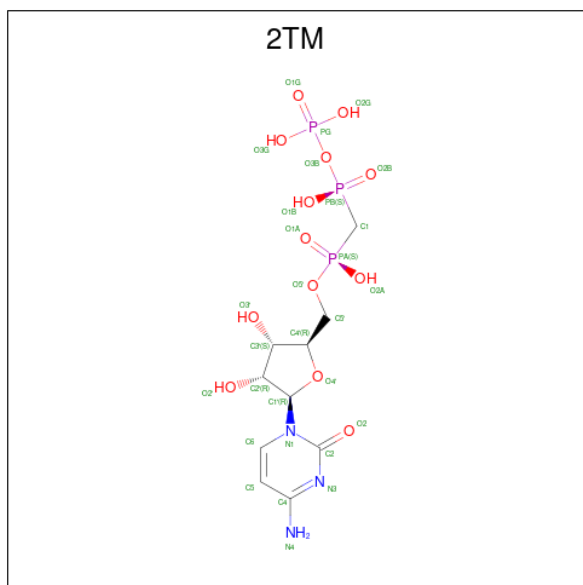
- Molecule 5 is a RNA chain called 3' vRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	R	14	Total	C	N	O	P	0	0
			288	129	41	104	14		

- Molecule 6 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	M	17	Total 365	C 164	N 69	O 115	P 17	0	0

- Molecule 7 is 5'-O-[(S)-hydroxy{[(S)-hydroxy(phosphonooxy)phosphoryl]methyl}phosphoryl]cytidine (three-letter code: 2TM) (formula: C₁₀H₁₈N₃O₁₃P₃) (labeled as "Ligand of Interest" by depositor).

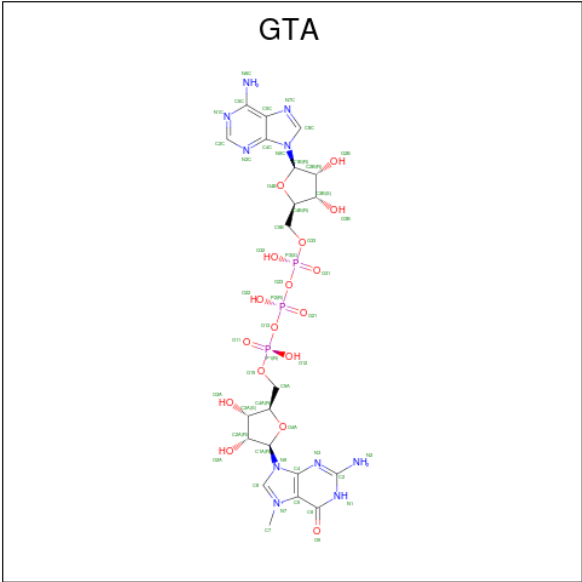


Mol	Chain	Residues	Atoms					AltConf
7	B	1	Total	C	N	O	P	0
			29	10	3	13	3	

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
8	B	2	Total	Mg	0
			2	2	

- Molecule 9 is P1-7-METHYLGUANOSINE-P3-ADENOSINE-5',5'-TRIPHOSPHATE (three-letter code: GTA) (formula: $C_{21}H_{30}N_{10}O_{17}P_3$).



Mol	Chain	Residues	Atoms					AltConf
9	M	1	Total	C	N	O	P	0
			51	21	10	17	3	

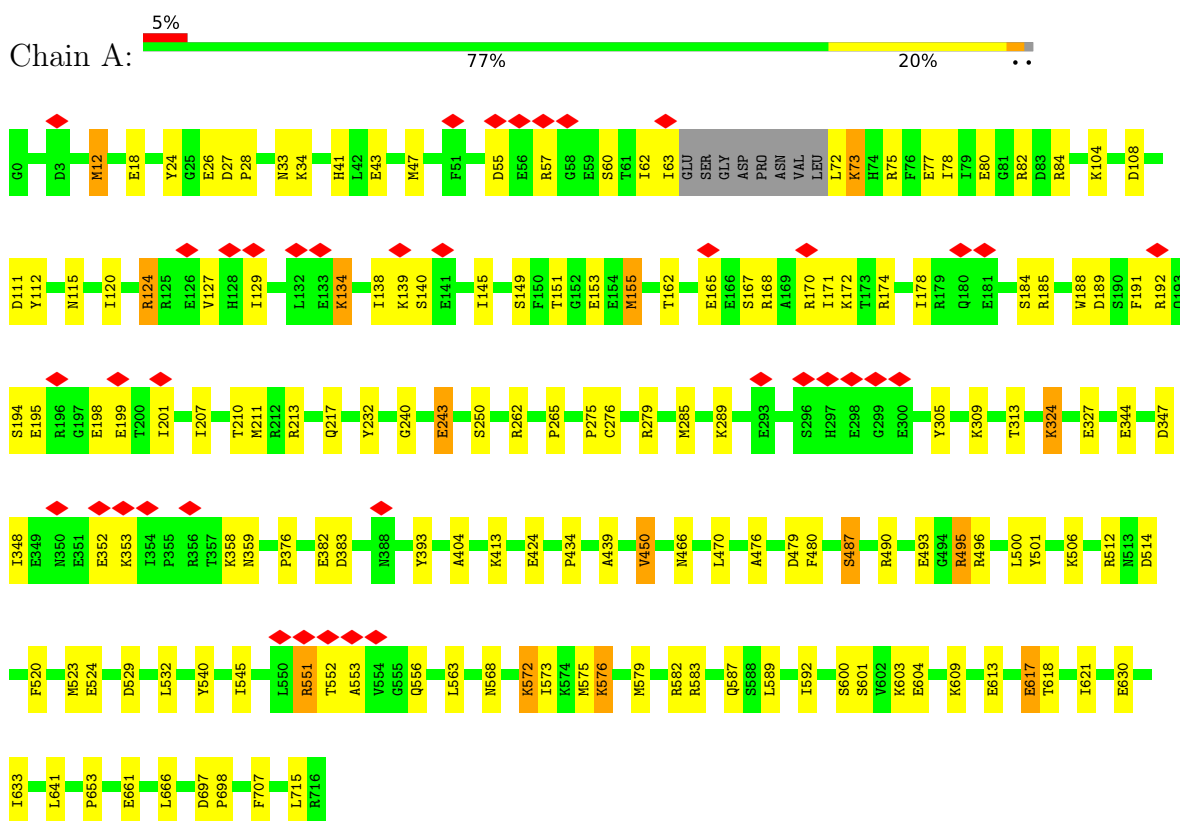
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		AltConf
10	A	28	Total	O	0
			28	28	
10	B	49	Total	O	0
			49	49	
10	C	6	Total	O	0
			6	6	
10	V	13	Total	O	0
			13	13	
10	R	12	Total	O	0
			12	12	
10	M	12	Total	O	0
			12	12	

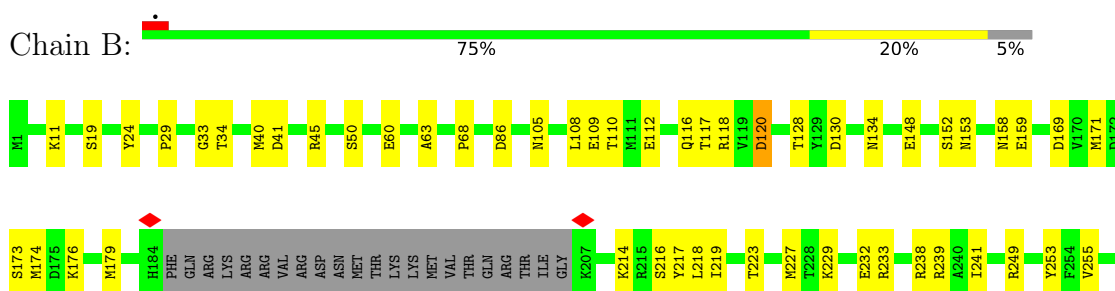
3 Residue-property plots

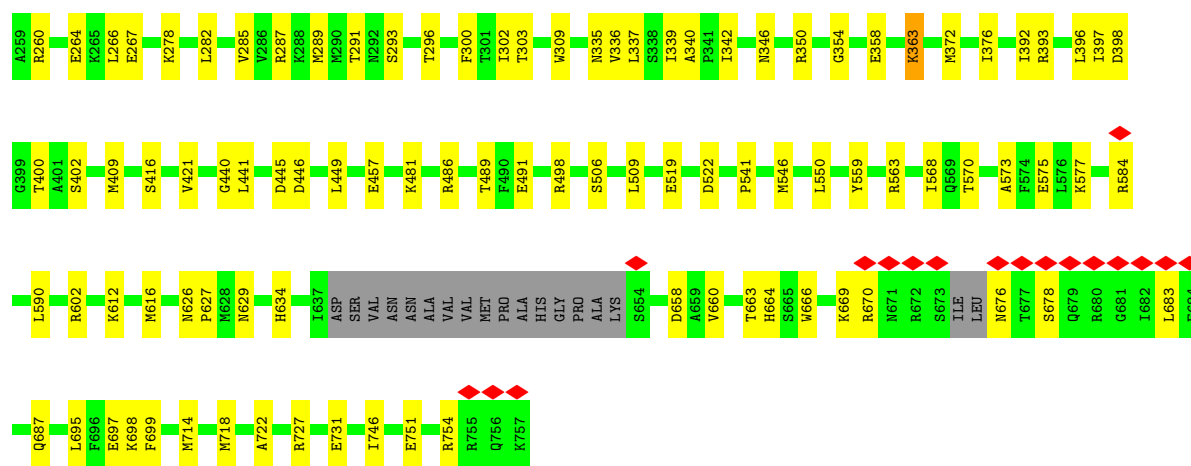
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Polymerase acidic protein

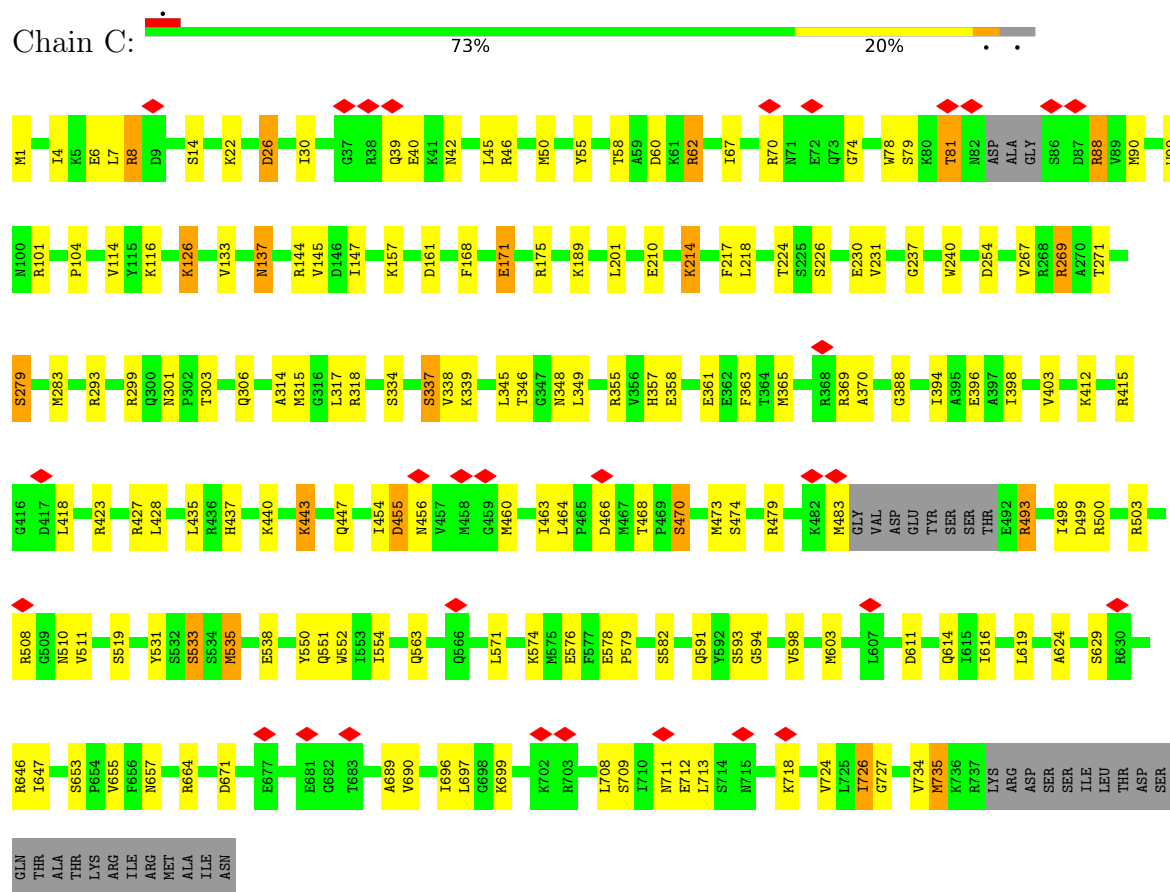


- Molecule 2: RNA-directed RNA polymerase catalytic subunit





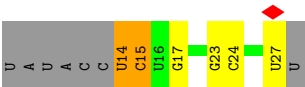
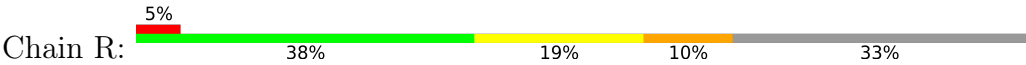
• Molecule 3: Polymerase basic protein 2



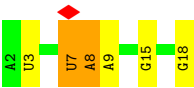
• Molecule 4: 5' vRNA



• Molecule 5: 3' vRNA



● Molecule 6: mRNA



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	116539	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 ($e^-/\text{\AA}^2$)	33.4	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	0.994	Depositor
Minimum map value	-0.289	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.019	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	357.58798, 357.58798, 357.58798	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8127, 0.8127, 0.8127	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 2TM, MG, GTA

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	0.27	0/5877	0.50	0/7920
2	B	0.27	0/5846	0.51	0/7878
3	C	0.26	0/5862	0.53	0/7902
4	V	0.68	1/319 (0.3%)	0.64	0/494
5	R	0.27	0/318	0.68	0/491
6	M	0.28	0/409	0.78	0/635
All	All	0.28	1/18631 (0.0%)	0.53	0/25320

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
2	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	V	1	A	OP3-P	-10.87	1.48	1.61

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	287	ARG	Sidechain

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	5755	0	5687	106	0
2	B	5736	0	5715	98	0
3	C	5767	0	5905	104	0
4	V	284	0	142	2	0
5	R	288	0	148	3	0
6	M	365	0	184	5	0
7	B	29	0	14	1	0
8	B	2	0	0	0	0
9	M	51	0	26	1	0
10	A	28	0	0	2	0
10	B	49	0	0	2	0
10	C	6	0	0	0	0
10	M	12	0	0	2	0
10	R	12	0	0	0	0
10	V	13	0	0	0	0
All	All	18397	0	17821	283	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 8.

All (283) 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:41:ASP:OD2	2:B:45:ARG:NH1	2.13	0.81
1:A:490:ARG:HD3	1:A:496:ARG:HE	1.46	0.81
2:B:105:ASN:HA	2:B:108:LEU:HD12	1.66	0.76
3:C:39:GLN:HG2	3:C:40:GLU:H	1.54	0.72
1:A:124:ARG:NH1	1:A:191:PHE:O	2.23	0.71
3:C:315:MET:HG3	3:C:317:LEU:HD13	1.74	0.68
3:C:147:ILE:HB	3:C:214:LYS:HD3	1.76	0.67
2:B:699:PHE:HE2	2:B:722:ALA:HB2	1.59	0.67
2:B:50:SER:HB3	2:B:68:PRO:HB3	1.76	0.67
3:C:708:LEU:HD12	3:C:712:GLU:HB3	1.78	0.66
1:A:279:ARG:NH1	1:A:697:ASP:OD1	2.29	0.66
3:C:614:GLN:NE2	3:C:689:ALA:O	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:568:ILE:HD12	2:B:568:ILE:H	1.61	0.65
2:B:153:ASN:ND2	2:B:179:MET:SD	2.71	0.64
1:A:324:LYS:NZ	1:A:540:TYR:O	2.30	0.64
1:A:524:GLU:OE1	10:A:801:HOH:O	2.14	0.64
3:C:60:ASP:HB3	3:C:62:ARG:HG2	1.80	0.64
3:C:697:LEU:HD11	3:C:735:MET:HG2	1.80	0.63
3:C:301:ASN:OD1	3:C:303:THR:OG1	2.12	0.63
1:A:376:PRO:HA	2:B:363:LYS:HE3	1.80	0.62
1:A:576:LYS:HG3	3:C:45:LEU:HB2	1.81	0.62
5:R:14:U:O2'	5:R:15:C:OP1	2.18	0.62
3:C:503:ARG:HH21	3:C:511:VAL:HG21	1.65	0.62
1:A:493:GLU:HB2	1:A:495:ARG:HH12	1.63	0.61
1:A:232:TYR:OH	2:B:86:ASP:OD1	2.18	0.61
1:A:124:ARG:HG2	1:A:192:ARG:HA	1.82	0.61
3:C:403:VAL:HG11	3:C:435:LEU:HG	1.81	0.61
3:C:428:LEU:HD22	6:M:3:U:H5''	1.82	0.61
2:B:238:ARG:NH1	10:B:1004:HOH:O	2.32	0.61
3:C:696:ILE:HG12	3:C:734:VAL:HG12	1.82	0.61
3:C:46:ARG:O	3:C:50:MET:HG3	2.02	0.60
1:A:265:PRO:HG2	1:A:715:LEU:HD11	1.84	0.59
2:B:258:LEU:HD22	2:B:337:LEU:HD13	1.83	0.59
3:C:81:THR:OG1	3:C:90:MET:SD	2.61	0.59
1:A:618:THR:HB	1:A:630:GLU:HB3	1.84	0.59
2:B:105:ASN:O	2:B:109:GLU:HG3	2.03	0.59
1:A:450:VAL:HG22	1:A:641:LEU:HD22	1.85	0.59
2:B:158:ASN:ND2	2:B:169:ASP:OD2	2.34	0.59
3:C:345:LEU:HD21	3:C:349:LEU:HD23	1.84	0.58
3:C:137:ASN:ND2	3:C:551:GLN:OE1	2.33	0.58
2:B:590:LEU:HD23	2:B:612:LYS:HD2	1.86	0.57
1:A:18:GLU:OE2	1:A:28:PRO:HG3	2.03	0.57
1:A:43:GLU:O	1:A:47:MET:HG3	2.04	0.57
2:B:446:ASP:OD2	6:M:18:G:H4'	2.04	0.56
3:C:314:ALA:HA	3:C:498:ILE:HG21	1.87	0.56
3:C:412:LYS:O	3:C:415:ARG:NH1	2.39	0.56
3:C:454:ILE:HG22	3:C:456:ASN:H	1.70	0.56
3:C:563:GLN:OE1	3:C:574:LYS:HE3	2.05	0.56
1:A:211:MET:HG3	2:B:171:MET:SD	2.46	0.56
1:A:327:GLU:OE2	10:A:802:HOH:O	2.18	0.56
3:C:454:ILE:HD11	3:C:474:SER:HB2	1.88	0.56
3:C:554:ILE:HG22	3:C:616:ILE:HD13	1.87	0.56
1:A:653:PRO:HG3	2:B:29:PRO:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:LYS:NZ	1:A:359:ASN:O	2.36	0.55
1:A:33:ASN:ND2	1:A:194:SER:OG	2.39	0.55
1:A:609:LYS:HG3	1:A:613:GLU:HG3	1.88	0.55
1:A:276:CYS:HB2	1:A:698:PRO:HG3	1.89	0.55
1:A:551:ARG:HD2	1:A:556:GLN:HB3	1.88	0.55
1:A:495:ARG:HH11	1:A:495:ARG:HG2	1.72	0.55
2:B:575:GLU:OE1	3:C:101:ARG:NE	2.31	0.55
2:B:670:ARG:NH1	3:C:46:ARG:HH21	2.03	0.55
3:C:189:LYS:NZ	3:C:210:GLU:OE1	2.35	0.55
1:A:55:ASP:HA	1:A:72:LEU:HD21	1.89	0.55
3:C:279:SER:O	3:C:283:MET:HG3	2.06	0.55
3:C:709:SER:OG	3:C:711:ASN:OD1	2.23	0.54
3:C:369:ARG:HG3	3:C:369:ARG:HH11	1.70	0.54
3:C:571:LEU:HD21	3:C:619:LEU:HD21	1.90	0.54
1:A:582:ARG:HD2	2:B:550:LEU:HD22	1.90	0.54
3:C:70:ARG:HG2	3:C:74:GLY:HA2	1.90	0.54
2:B:241:ILE:HD12	5:R:17:G:C8	2.42	0.54
1:A:592:ILE:HG23	1:A:633:ILE:HD11	1.90	0.53
1:A:62:ILE:O	1:A:63:ILE:HD13	2.09	0.53
3:C:337:SER:HA	3:C:358:GLU:O	2.09	0.53
3:C:306:GLN:NE2	9:M:101:GTA:O3A	2.41	0.53
1:A:207:ILE:HG12	2:B:339:ILE:HD11	1.91	0.53
3:C:147:ILE:HD11	3:C:218:LEU:HD11	1.90	0.53
3:C:318:ARG:HB3	3:C:493:ARG:HB3	1.90	0.53
1:A:167:SER:HA	1:A:170:ARG:HD3	1.90	0.53
1:A:487:SER:HB3	1:A:501:TYR:HE2	1.74	0.53
2:B:118:ARG:NH2	3:C:26:ASP:OD2	2.39	0.53
3:C:470:SER:OG	3:C:473:MET:O	2.25	0.52
1:A:344:GLU:O	1:A:348:ILE:HG13	2.10	0.52
1:A:529:ASP:HB3	1:A:532:LEU:HD13	1.92	0.52
2:B:489:THR:HG22	2:B:498:ARG:HA	1.92	0.52
1:A:617:GLU:OE1	2:B:11:LYS:NZ	2.38	0.52
2:B:108:LEU:O	2:B:112:GLU:HG2	2.10	0.52
3:C:346:THR:OG1	3:C:348:ASN:OD1	2.26	0.52
1:A:189:ASP:OD1	1:A:192:ARG:NH2	2.43	0.52
2:B:282:LEU:HD22	2:B:441:LEU:HD13	1.92	0.52
1:A:75:ARG:HG2	1:A:75:ARG:HH11	1.75	0.51
2:B:522:ASP:OD2	2:B:559:TYR:OH	2.15	0.51
2:B:266:LEU:HD13	2:B:421:VAL:HG11	1.92	0.51
2:B:335:ASN:O	2:B:339:ILE:HG12	2.11	0.51
1:A:24:TYR:OH	1:A:84:ARG:NH2	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:THR:HG23	1:A:153:GLU:H	1.76	0.51
3:C:116:LYS:NZ	3:C:210:GLU:OE2	2.43	0.51
1:A:520:PHE:CZ	1:A:568:ASN:HB3	2.46	0.51
1:A:666:LEU:HG	1:A:707:PHE:HE2	1.76	0.51
1:A:127:VAL:HG23	1:A:155:MET:HE2	1.93	0.50
1:A:134:LYS:HD2	1:A:145:ILE:HG23	1.93	0.50
1:A:424:GLU:OE1	1:A:490:ARG:NH1	2.45	0.50
2:B:176:LYS:O	2:B:214:LYS:HD3	2.11	0.50
3:C:603:MET:CE	3:C:611:ASP:H	2.24	0.50
2:B:63:ALA:HB1	2:B:402:SER:HB2	1.93	0.50
2:B:120:ASP:OD1	3:C:30:ILE:HG13	2.12	0.50
2:B:285:VAL:O	2:B:289:MET:HG3	2.12	0.50
3:C:370:ALA:HA	3:C:388:GLY:HA2	1.94	0.50
6:M:7:U:O2'	6:M:8:A:OP2	2.28	0.50
2:B:481:LYS:NZ	7:B:901:2TM:O2B	2.44	0.50
2:B:695:LEU:HD21	3:C:7:LEU:HA	1.93	0.50
3:C:394:ILE:O	3:C:398:ILE:HG13	2.11	0.50
2:B:669:LYS:HE2	3:C:90:MET:HB3	1.94	0.50
1:A:165:GLU:OE2	1:A:168:ARG:NH2	2.45	0.50
2:B:110:THR:HG21	2:B:258:LEU:HB2	1.94	0.50
2:B:363:LYS:O	2:B:363:LYS:HD3	2.11	0.50
2:B:112:GLU:O	2:B:116:GLN:HG2	2.12	0.49
3:C:67:ILE:HD11	3:C:99:TRP:HE3	1.76	0.49
1:A:213:ARG:HD2	2:B:60:GLU:OE2	2.12	0.49
3:C:594:GLY:O	3:C:598:VAL:HG23	2.13	0.49
1:A:26:GLU:OE2	1:A:201:ILE:HG23	2.12	0.49
1:A:285:MET:HB2	1:A:413:LYS:HD2	1.93	0.49
3:C:8:ARG:HG3	3:C:8:ARG:HH11	1.78	0.49
2:B:179:MET:HG3	2:B:217:TYR:CD1	2.48	0.49
2:B:626:ASN:ND2	3:C:104:PRO:O	2.46	0.49
1:A:41:HIS:CD2	1:A:80:GLU:HG2	2.48	0.49
1:A:552:THR:HG23	1:A:553:ALA:H	1.78	0.48
1:A:124:ARG:HG3	1:A:188:TRP:CH2	2.47	0.48
1:A:138:ILE:HG22	1:A:140:SER:H	1.78	0.48
2:B:457:GLU:OE1	2:B:457:GLU:N	2.47	0.48
1:A:587:GLN:HB2	2:B:509:LEU:HD11	1.95	0.48
1:A:174:ARG:HE	1:A:178:ILE:HG13	1.78	0.48
2:B:309:TRP:CZ2	2:B:416:SER:HB2	2.49	0.48
1:A:124:ARG:HB2	1:A:195:GLU:OE1	2.14	0.48
2:B:229:LYS:O	10:B:1001:HOH:O	2.20	0.48
2:B:629:ASN:O	2:B:634:HIS:NE2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:575:MET:O	1:A:579:MET:HG3	2.13	0.48
3:C:161:ASP:CG	3:C:175:ARG:HD3	2.34	0.47
1:A:78:ILE:O	1:A:82:ARG:NE	2.35	0.47
3:C:293:ARG:HG2	3:C:293:ARG:HH11	1.79	0.47
1:A:80:GLU:OE2	1:A:80:GLU:HA	2.14	0.47
1:A:108:ASP:OD2	1:A:120:ILE:N	2.37	0.47
1:A:393:TYR:CE1	2:B:358:GLU:HB3	2.49	0.47
1:A:545:ILE:HD13	1:A:563:LEU:HB2	1.97	0.47
1:A:583:ARG:HB3	2:B:509:LEU:HD22	1.97	0.47
2:B:117:THR:HG21	2:B:253:TYR:HE2	1.80	0.47
5:R:23:G:H2'	5:R:24:C:C6	2.50	0.47
1:A:111:ASP:O	1:A:115:ASN:N	2.47	0.47
2:B:300:PHE:CZ	2:B:486:ARG:HD2	2.50	0.47
1:A:41:HIS:HD2	1:A:80:GLU:HG2	1.78	0.46
2:B:232:GLU:OE2	2:B:239:ARG:NH2	2.35	0.46
1:A:134:LYS:HE2	1:A:134:LYS:HB2	1.52	0.46
3:C:267:VAL:O	3:C:271:THR:HG23	2.14	0.46
3:C:538:GLU:HG3	3:C:579:PRO:HG3	1.98	0.46
2:B:130:ASP:O	2:B:134:ASN:N	2.48	0.46
1:A:213:ARG:O	1:A:217:GLN:HG3	2.16	0.46
2:B:128:THR:HG23	2:B:249:ARG:NE	2.30	0.46
2:B:289:MET:HE1	2:B:440:GLY:N	2.31	0.46
2:B:698:LYS:HB2	2:B:698:LYS:HE2	1.69	0.46
3:C:563:GLN:HG2	3:C:571:LEU:HB2	1.98	0.46
3:C:724:VAL:HB	3:C:726:ILE:HG22	1.97	0.46
3:C:466:ASP:HB3	3:C:468:THR:HG23	1.98	0.46
3:C:593:SER:HB2	3:C:624:ALA:HB3	1.98	0.45
1:A:188:TRP:HE1	1:A:192:ARG:HE	1.64	0.45
3:C:46:ARG:HB3	3:C:46:ARG:CZ	2.45	0.45
1:A:149:SER:OG	1:A:151:THR:HG22	2.17	0.45
2:B:40:MET:HG2	2:B:392:ILE:HD13	1.99	0.45
3:C:46:ARG:HB3	3:C:46:ARG:NH1	2.32	0.45
1:A:199:GLU:OE1	1:A:199:GLU:N	2.46	0.45
3:C:552:TRP:CH2	3:C:576:GLU:HG3	2.52	0.45
2:B:309:TRP:HZ2	2:B:416:SER:HB2	1.82	0.45
3:C:145:VAL:HG12	3:C:224:THR:HB	1.98	0.45
3:C:293:ARG:HG2	3:C:293:ARG:NH1	2.31	0.45
1:A:138:ILE:HD12	1:A:145:ILE:HD11	1.99	0.44
2:B:506:SER:O	2:B:509:LEU:HB2	2.17	0.44
1:A:129:ILE:HD13	1:A:129:ILE:N	2.32	0.44
1:A:262:ARG:HG2	1:A:715:LEU:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:128:THR:HG23	2:B:249:ARG:HE	1.83	0.44
2:B:746:ILE:HG21	3:C:4:ILE:HG23	1.99	0.44
1:A:382:GLU:O	1:A:382:GLU:HG3	2.18	0.44
1:A:495:ARG:HG2	1:A:495:ARG:NH1	2.32	0.44
1:A:523:MET:HB3	1:A:563:LEU:HD21	1.99	0.44
3:C:463:ILE:HG13	3:C:464:LEU:N	2.33	0.44
3:C:535:MET:CE	3:C:576:GLU:HA	2.47	0.44
2:B:627:PRO:HB3	2:B:658:ASP:HB3	2.00	0.44
3:C:531:TYR:CE2	3:C:533:SER:HB2	2.52	0.44
1:A:174:ARG:HA	1:A:174:ARG:HD2	1.73	0.44
3:C:237:GLY:HA3	3:C:240:TRP:CE2	2.53	0.44
3:C:269:ARG:NH2	3:C:578:GLU:OE1	2.44	0.44
1:A:77:GLU:OE1	1:A:112:TYR:OH	2.33	0.44
3:C:690:VAL:HG11	3:C:726:ILE:HD11	2.00	0.44
1:A:476:ALA:HB1	1:A:479:ASP:HB2	2.00	0.44
1:A:572:LYS:HE2	1:A:572:LYS:HB2	1.64	0.44
2:B:296:THR:HA	3:C:727:GLY:O	2.17	0.44
1:A:439:ALA:HB2	2:B:541:PRO:HB2	1.99	0.44
1:A:603:LYS:HA	1:A:603:LYS:HD3	1.62	0.44
1:A:490:ARG:HD2	1:A:496:ARG:HA	2.00	0.43
2:B:24:TYR:CE1	2:B:233:ARG:HD2	2.54	0.43
2:B:303:THR:HG23	2:B:491:GLU:O	2.18	0.43
2:B:350:ARG:HA	2:B:400:THR:HG22	2.00	0.43
2:B:683:LEU:O	2:B:687:GLN:HG3	2.17	0.43
3:C:443:LYS:O	3:C:447:GLN:HG3	2.18	0.43
1:A:168:ARG:O	1:A:172:LYS:HG3	2.19	0.43
3:C:39:GLN:HG2	3:C:40:GLU:N	2.29	0.43
1:A:47:MET:HE3	1:A:171:ILE:HA	1.99	0.43
1:A:600:SER:O	1:A:604:GLU:N	2.50	0.43
2:B:714:MET:O	2:B:718:MET:HG3	2.19	0.43
3:C:338:VAL:O	3:C:357:HIS:HA	2.18	0.43
2:B:33:GLY:HA3	4:V:7:A:O3'	2.18	0.43
3:C:713:LEU:HD11	3:C:724:VAL:HG11	2.01	0.43
1:A:353:LYS:HB2	1:A:353:LYS:HE2	1.80	0.43
3:C:361:GLU:HG3	3:C:363:PHE:HE1	1.84	0.43
1:A:240:GLY:HA2	1:A:243:GLU:HG3	2.01	0.43
2:B:173:SER:HB2	2:B:179:MET:CE	2.49	0.43
3:C:126:LYS:HB2	3:C:126:LYS:HE2	1.40	0.43
3:C:418:LEU:HD13	3:C:437:HIS:HB2	2.01	0.43
2:B:506:SER:HA	2:B:509:LEU:HD12	2.01	0.43
2:B:663:THR:HG21	3:C:99:TRP:CD1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:THR:HG21	2:B:218[A]:LEU:HD11	2.00	0.43
1:A:621:ILE:HD12	1:A:621:ILE:HA	1.85	0.43
3:C:67:ILE:O	3:C:78:TRP:NE1	2.40	0.43
1:A:309:LYS:O	1:A:313:THR:HG23	2.19	0.43
2:B:751:GLU:O	2:B:754:ARG:HG2	2.19	0.43
3:C:4:ILE:HD12	3:C:4:ILE:HA	1.83	0.43
1:A:289:LYS:HA	1:A:500:LEU:O	2.19	0.42
1:A:470:LEU:HD22	1:A:579:MET:HE1	2.01	0.42
2:B:570:THR:HG22	2:B:573:ALA:HB2	2.00	0.42
2:B:278:LYS:O	2:B:282:LEU:HG	2.19	0.42
2:B:291:THR:O	3:C:646:ARG:NH2	2.52	0.42
2:B:302:ILE:HD12	2:B:449:LEU:HD23	2.01	0.42
3:C:168:PHE:HB3	3:C:171:GLU:HG3	2.02	0.42
1:A:34:LYS:HD2	1:A:198:GLU:HG2	2.01	0.42
2:B:568:ILE:HD12	2:B:568:ILE:N	2.32	0.42
2:B:568:ILE:H	2:B:568:ILE:CD1	2.29	0.42
2:B:669:LYS:HB2	3:C:58:THR:HG23	2.01	0.42
3:C:440:LYS:HD3	6:M:7:U:C2	2.54	0.42
2:B:260:ARG:O	2:B:264:GLU:HG3	2.19	0.42
2:B:396:LEU:C	2:B:397:ILE:HG13	2.40	0.42
3:C:647:ILE:HG12	3:C:655:VAL:HG12	2.01	0.42
1:A:589:LEU:HD13	2:B:546:MET:SD	2.60	0.42
2:B:219:ILE:O	2:B:223:THR:OG1	2.29	0.42
3:C:144:ARG:NE	10:M:201:HOH:O	2.42	0.42
1:A:275:PRO:HB3	1:A:404:ALA:HB2	2.02	0.42
2:B:660:VAL:HG21	3:C:114:VAL:HG21	2.01	0.42
1:A:551:ARG:CD	1:A:556:GLN:HB3	2.48	0.42
2:B:309:TRP:CG	2:B:445:ASP:HB3	2.54	0.42
2:B:590:LEU:HD12	2:B:664:HIS:CD2	2.55	0.42
1:A:134:LYS:HG2	1:A:145:ILE:HD13	2.01	0.41
2:B:372:MET:O	2:B:376:ILE:HB	2.20	0.41
6:M:15:G:N2	10:M:205:HOH:O	2.47	0.41
1:A:73:LYS:HE3	2:B:727:ARG:HH21	1.85	0.41
1:A:480:PHE:HA	1:A:506:LYS:O	2.21	0.41
3:C:361:GLU:HG3	3:C:363:PHE:CE1	2.55	0.41
1:A:84:ARG:HE	1:A:84:ARG:HB2	1.63	0.41
3:C:355:ARG:HD3	3:C:355:ARG:HA	1.79	0.41
1:A:434:PRO:HB2	3:C:133:VAL:HG23	2.01	0.41
3:C:8:ARG:HG3	3:C:8:ARG:NH1	2.36	0.41
2:B:397:ILE:HG22	2:B:398:ASP:OD1	2.21	0.41
1:A:466:ASN:OD1	1:A:575:MET:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:519:GLU:HG3	2:B:666:TRP:CZ2	2.55	0.41
3:C:499:ASP:OD2	3:C:503:ARG:HB2	2.21	0.41
1:A:12:MET:N	1:A:12:MET:HE3	2.36	0.41
1:A:104:LYS:HB2	1:A:104:LYS:HE3	1.94	0.41
1:A:139:LYS:HG2	1:A:139:LYS:O	2.21	0.41
1:A:514:ASP:HB3	1:A:573:ILE:HG12	2.03	0.41
1:A:551:ARG:NE	1:A:556:GLN:HB3	2.35	0.41
2:B:34:THR:HG21	2:B:354:GLY:O	2.21	0.41
2:B:255:VAL:HG21	2:B:340:ALA:HB3	2.03	0.41
3:C:22:LYS:HD2	3:C:22:LYS:HA	1.96	0.41
3:C:55:TYR:CZ	3:C:79:SER:HB3	2.56	0.41
3:C:427:ARG:HD2	3:C:427:ARG:HA	1.82	0.41
3:C:455:ASP:N	3:C:455:ASP:OD1	2.54	0.41
3:C:664:ARG:NH1	3:C:671:ASP:OD2	2.54	0.41
4:V:3:U:H2'	4:V:4:A:O4'	2.21	0.41
2:B:342:ILE:O	2:B:346:ASN:ND2	2.54	0.41
2:B:669:LYS:HB2	2:B:669:LYS:HE3	1.63	0.41
3:C:230:GLU:HG3	3:C:231:VAL:HG13	2.03	0.41
3:C:299:ARG:HA	3:C:317:LEU:HD21	2.02	0.41
2:B:612:LYS:O	2:B:616:MET:HG3	2.21	0.40
3:C:88:ARG:HA	3:C:88:ARG:HD2	1.76	0.40
3:C:550:TYR:O	3:C:554:ILE:HG23	2.22	0.40
3:C:708:LEU:HD23	3:C:708:LEU:H	1.86	0.40
3:C:708:LEU:HD11	3:C:713:LEU:HD23	2.03	0.40
3:C:254:ASP:OD1	3:C:254:ASP:N	2.53	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	705/717 (98%)	683 (97%)	22 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	710/757 (94%)	695 (98%)	14 (2%)	1 (0%)	51	71
3	C	720/759 (95%)	705 (98%)	15 (2%)	0	100	100
All	All	2135/2233 (96%)	2083 (98%)	51 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	409	MET

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	638/645 (99%)	610 (96%)	28 (4%)	28	49
2	B	633/668 (95%)	612 (97%)	21 (3%)	38	61
3	C	638/666 (96%)	591 (93%)	47 (7%)	13	25
All	All	1909/1979 (96%)	1813 (95%)	96 (5%)	28	43

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	MET
1	A	27	ASP
1	A	57	ARG
1	A	60	SER
1	A	73	LYS
1	A	124	ARG
1	A	134	LYS
1	A	155	MET
1	A	162	THR
1	A	184	SER
1	A	185	ARG
1	A	243	GLU
1	A	250	SER

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Mol	Chain	Res	Type
1	A	305	TYR
1	A	324	LYS
1	A	347	ASP
1	A	352	GLU
1	A	383	ASP
1	A	450	VAL
1	A	487	SER
1	A	495	ARG
1	A	512	ARG
1	A	551	ARG
1	A	572	LYS
1	A	576	LYS
1	A	601	SER
1	A	617	GLU
1	A	661	GLU
2	B	19	SER
2	B	120	ASP
2	B	148	GLU
2	B	152	SER
2	B	159	GLU
2	B	174	MET
2	B	216	SER
2	B	227	MET
2	B	267	GLU
2	B	293	SER
2	B	336	VAL
2	B	363	LYS
2	B	393	ARG
2	B	563	ARG
2	B	577	LYS
2	B	584	ARG
2	B	602	ARG
2	B	676	ASN
2	B	678	SER
2	B	697	GLU
2	B	731	GLU
3	C	1	MET
3	C	6	GLU
3	C	8	ARG
3	C	14	SER
3	C	26	ASP
3	C	42	ASN

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Mol	Chain	Res	Type
3	C	62	ARG
3	C	81	THR
3	C	88	ARG
3	C	126	LYS
3	C	137	ASN
3	C	157	LYS
3	C	171	GLU
3	C	201	LEU
3	C	214	LYS
3	C	217	PHE
3	C	226	SER
3	C	269	ARG
3	C	279	SER
3	C	334	SER
3	C	337	SER
3	C	339	LYS
3	C	365	MET
3	C	396	GLU
3	C	423	ARG
3	C	443	LYS
3	C	455	ASP
3	C	460	MET
3	C	470	SER
3	C	479	ARG
3	C	483	MET
3	C	493	ARG
3	C	500	ARG
3	C	508	ARG
3	C	510	ASN
3	C	519	SER
3	C	533	SER
3	C	535	MET
3	C	582	SER
3	C	591	GLN
3	C	629	SER
3	C	653	SER
3	C	657	ASN
3	C	699	LYS
3	C	718	LYS
3	C	726	ILE
3	C	735	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are

no such sidechains identified.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	V	12/14 (85%)	2 (16%)	0
5	R	14/21 (66%)	2 (14%)	1 (7%)
6	M	16/17 (94%)	2 (12%)	2 (12%)
All	All	42/52 (80%)	6 (14%)	3 (7%)

All (6) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	V	8	A
4	V	11	A
5	R	15	C
5	R	27	U
6	M	8	A
6	M	9	A

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
5	R	14	U
6	M	7	U
6	M	8	A

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are 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
9	GTA	M	101	6	46,56,56	1.67	9 (19%)	46,88,88	1.99	10 (21%)
7	2TM	B	901	8	27,30,30	0.80	2 (7%)	39,47,47	0.61	0

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
9	GTA	M	101	6	-	6/24/64/64	0/6/6/6
7	2TM	B	901	8	-	0/19/38/38	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	M	101	GTA	C2C-N3C	4.91	1.40	1.32
9	M	101	GTA	C5-C6	-4.47	1.34	1.45
9	M	101	GTA	O4B-C1B	4.08	1.46	1.41
9	M	101	GTA	C2C-N1C	3.15	1.39	1.33
9	M	101	GTA	C5C-C4C	-2.76	1.33	1.40
9	M	101	GTA	C6C-C5C	-2.68	1.33	1.43
9	M	101	GTA	C5-C4	-2.46	1.34	1.39
7	B	901	2TM	PA-O2A	-2.42	1.50	1.56
7	B	901	2TM	PB-O1B	-2.42	1.50	1.56
9	M	101	GTA	C2A-C1A	-2.12	1.50	1.53
9	M	101	GTA	C2B-C1B	-2.04	1.50	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	M	101	GTA	N3C-C2C-N1C	-6.78	118.08	128.68
9	M	101	GTA	O4A-C1A-C2A	-5.90	98.31	106.93
9	M	101	GTA	C2-N1-C6	-3.58	118.51	125.10
9	M	101	GTA	P2-O13-P1	-3.25	121.68	132.83
9	M	101	GTA	P3-O23-P2	-3.17	121.95	132.83
9	M	101	GTA	C3A-C2A-C1A	-3.16	96.23	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	M	101	GTA	C5C-C6C-N6C	-2.73	116.20	120.35
9	M	101	GTA	C2B-C3B-C4B	-2.69	97.41	102.64
9	M	101	GTA	O4B-C1B-C2B	-2.27	103.61	106.93
9	M	101	GTA	N1-C2-N3	-2.25	119.11	123.32

There are no chirality outliers.

All (6) torsion outliers are listed below:

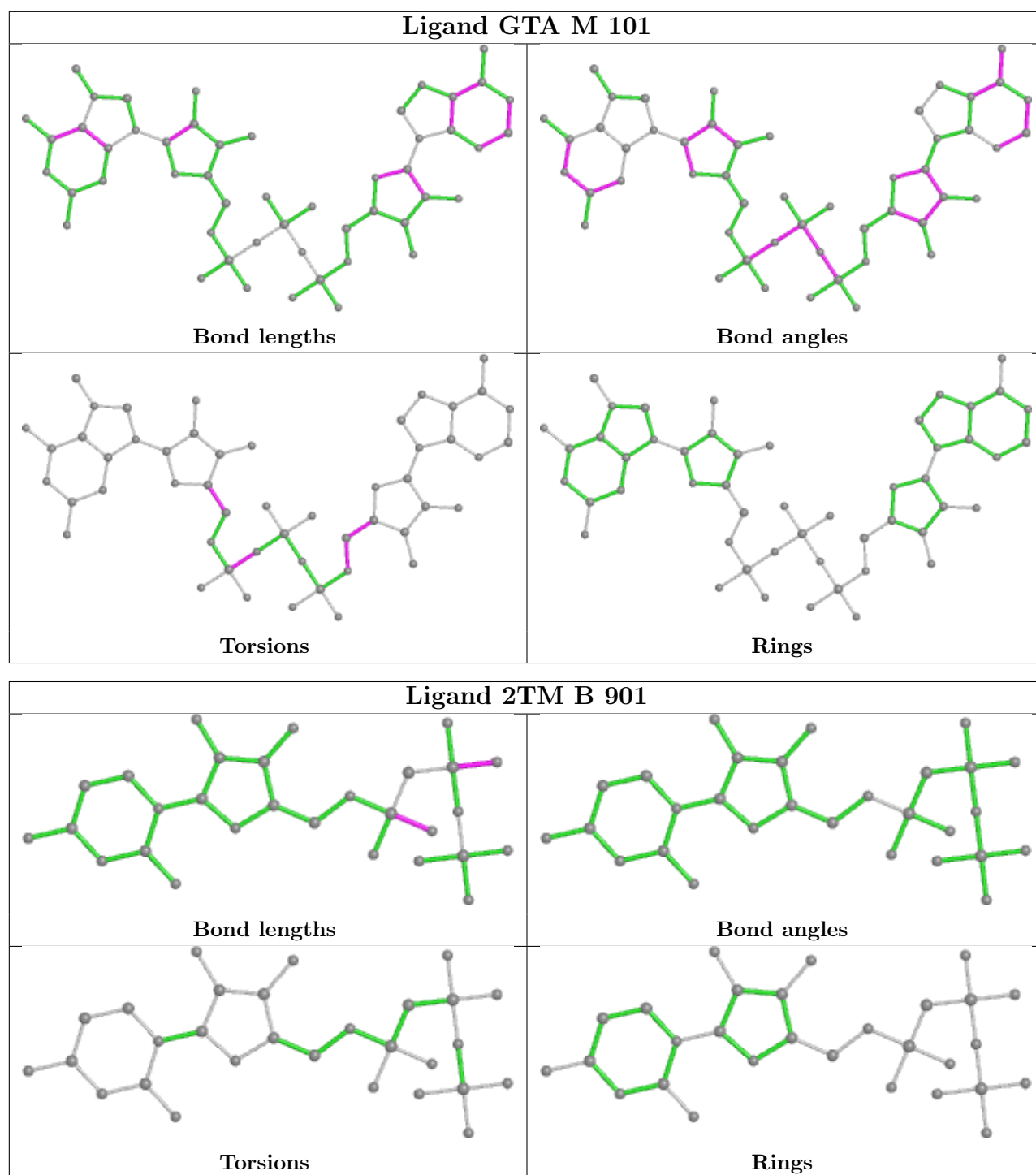
Mol	Chain	Res	Type	Atoms
9	M	101	GTA	O4B-C4B-C5B-O33
9	M	101	GTA	P2-O13-P1-O11
9	M	101	GTA	C4B-C5B-O33-P3
9	M	101	GTA	O4A-C4A-C5A-O15
9	M	101	GTA	P2-O13-P1-O12
9	M	101	GTA	C3B-C4B-C5B-O33

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	M	101	GTA	1	0
7	B	901	2TM	1	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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

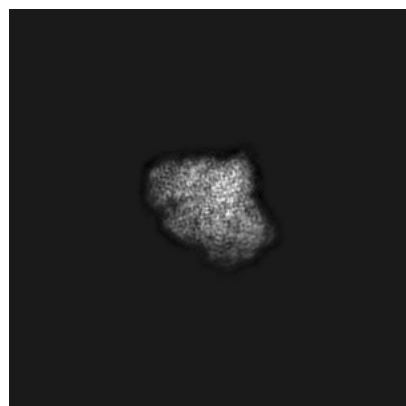
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14144. These allow visual inspection of the internal detail of the map and identification of artifacts.

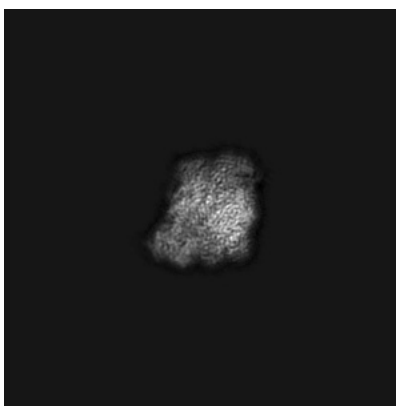
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

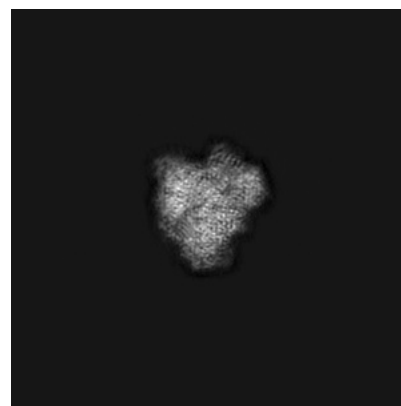
6.1.1 Primary map



X

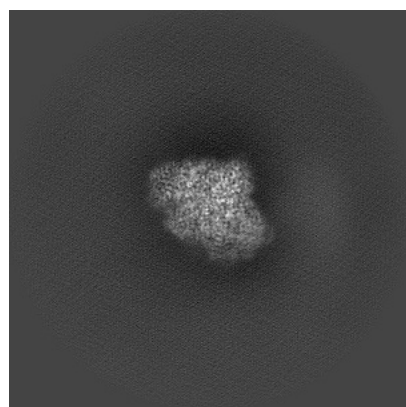


Y

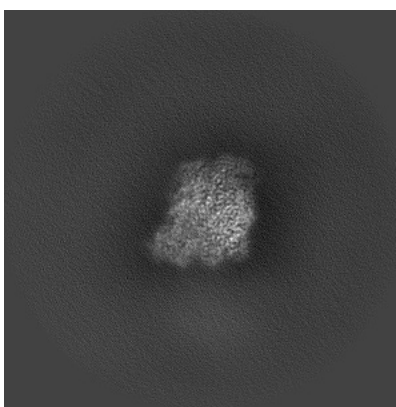


Z

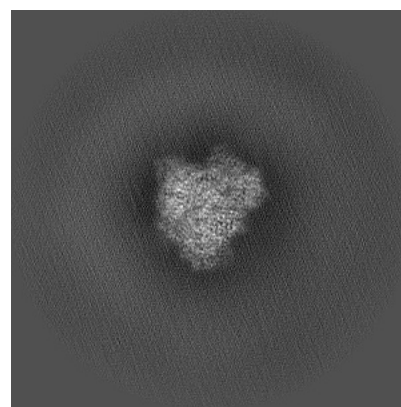
6.1.2 Raw map



X



Y

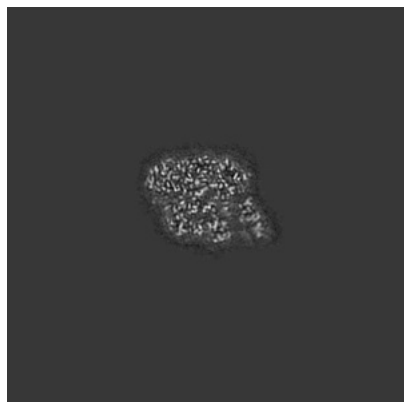


Z

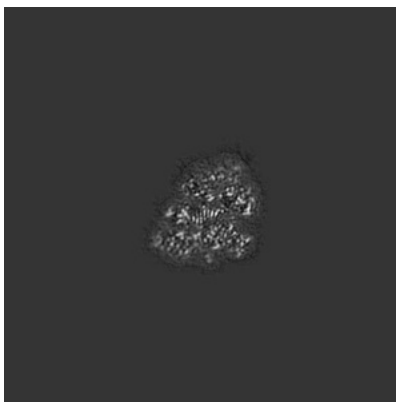
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

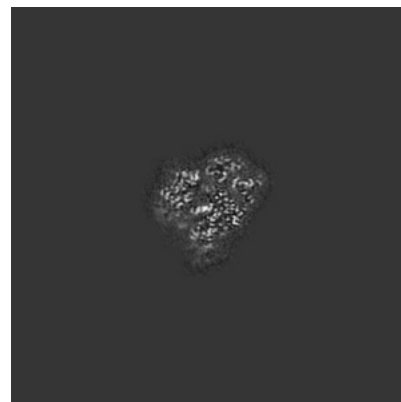
6.2.1 Primary map



X Index: 220

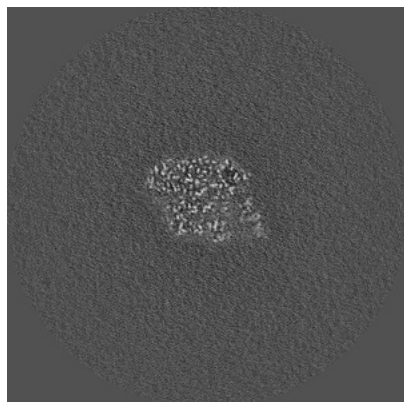


Y Index: 220

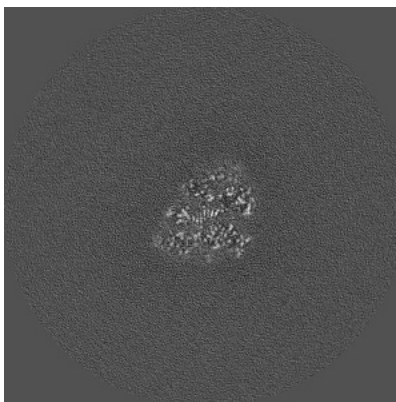


Z Index: 220

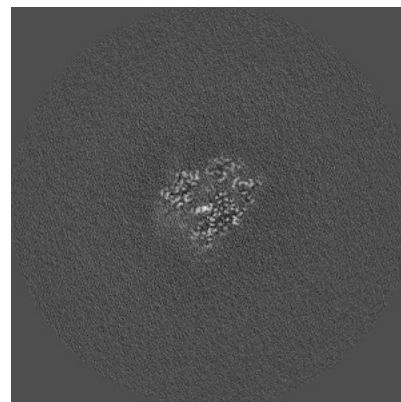
6.2.2 Raw map



X Index: 220



Y Index: 220



Z Index: 220

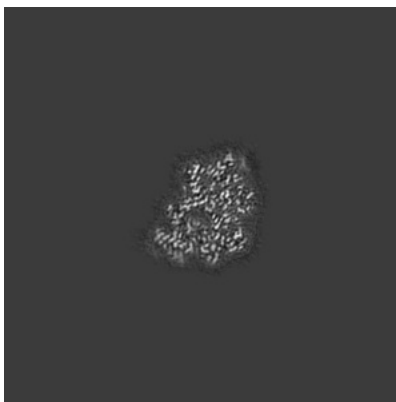
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

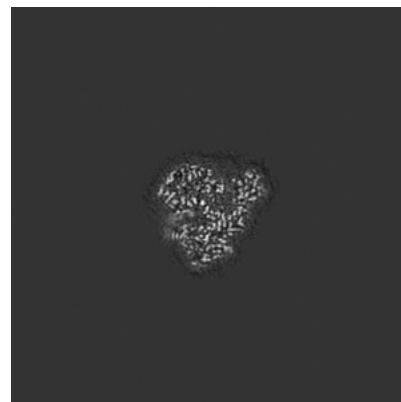
6.3.1 Primary map



X Index: 218

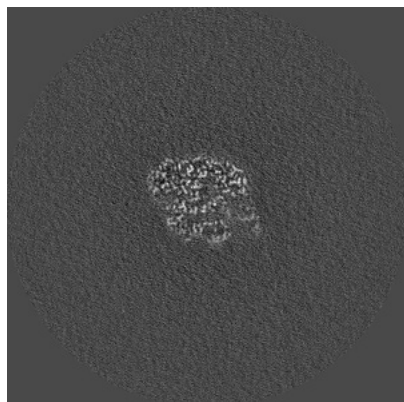


Y Index: 229

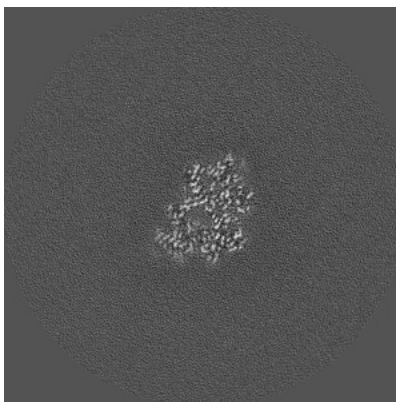


Z Index: 236

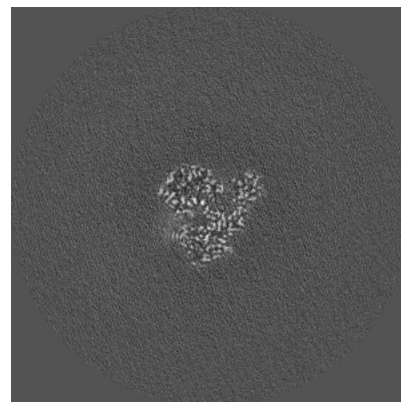
6.3.2 Raw map



X Index: 218



Y Index: 229



Z Index: 236

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

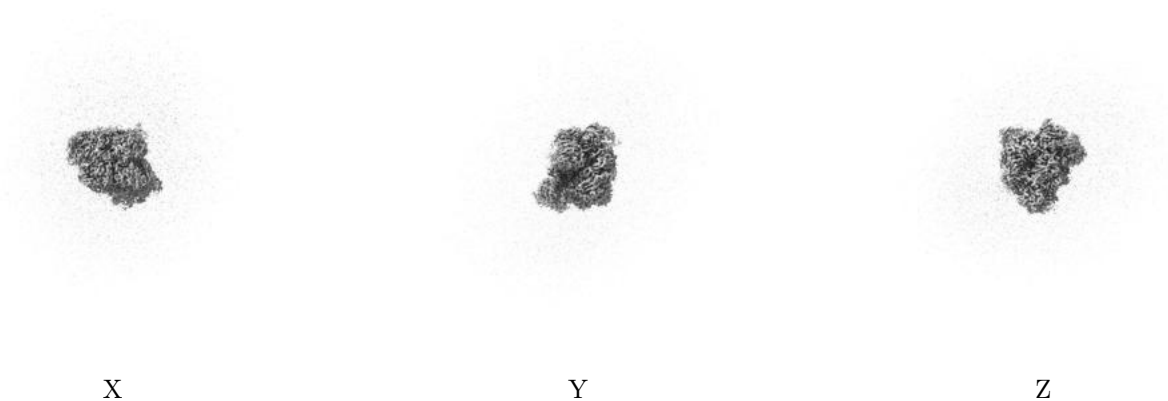
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

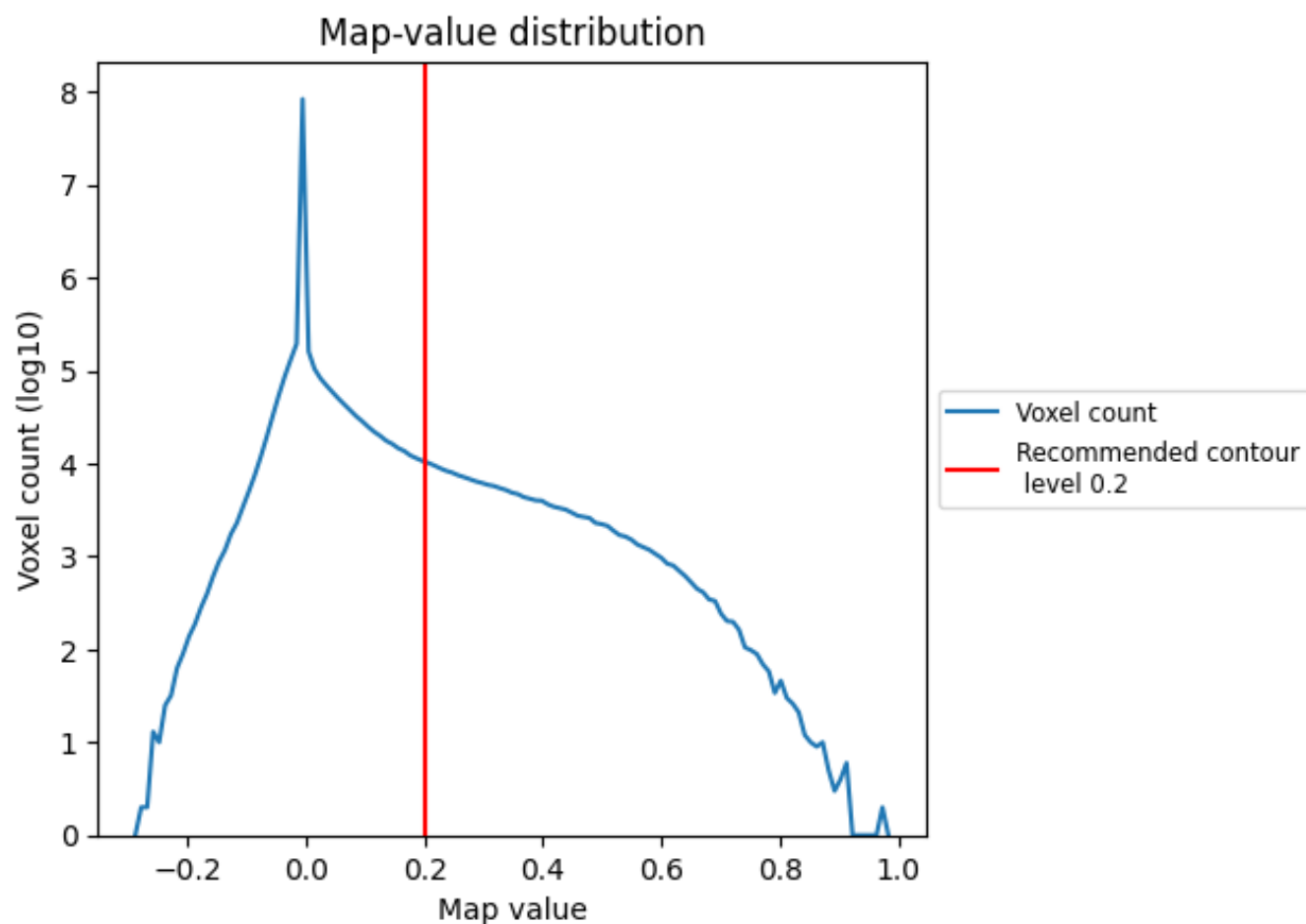
6.5 Mask visualisation [i](#)

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

7 Map analysis [i](#)

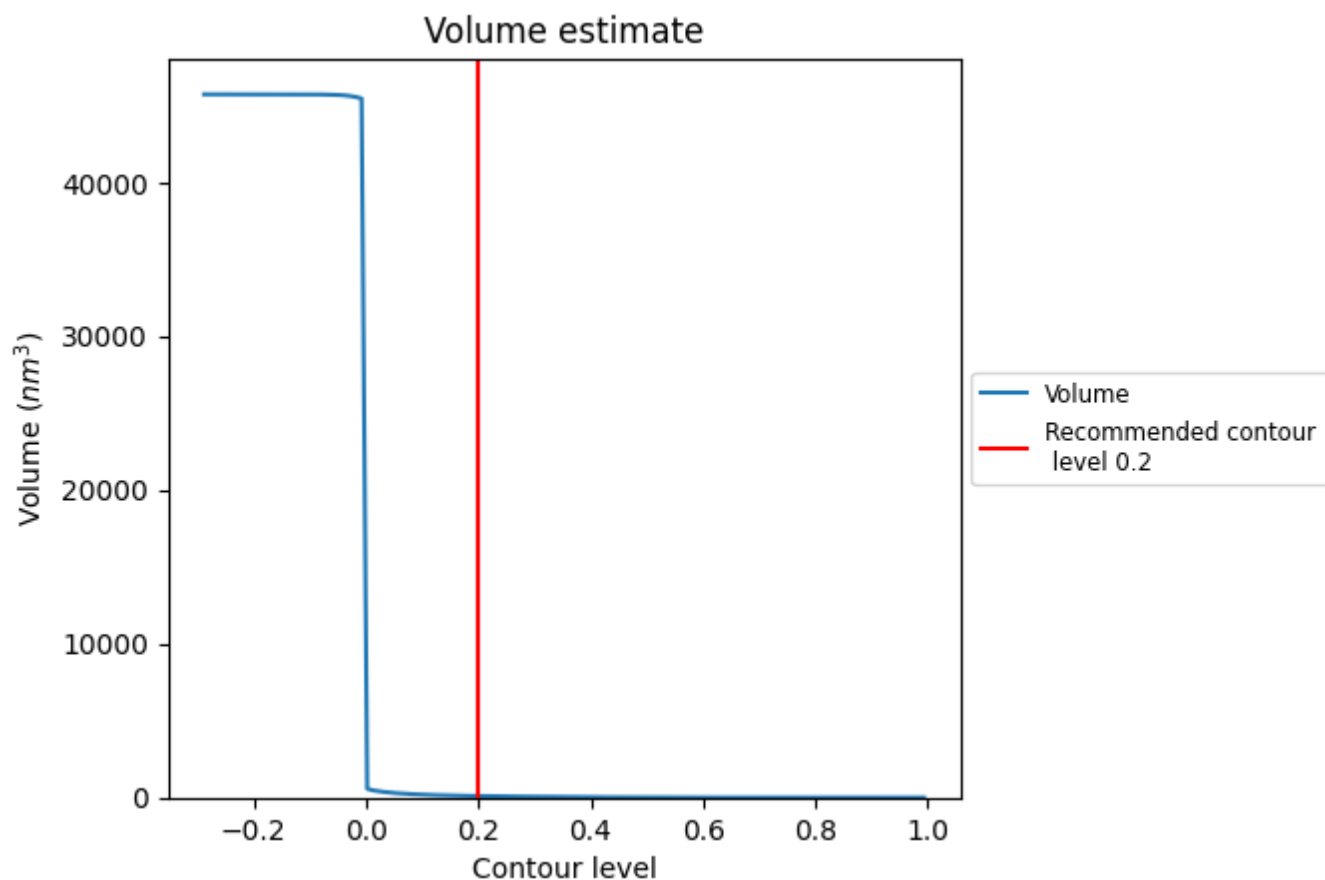
This section contains the results of statistical analysis of the map.

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.

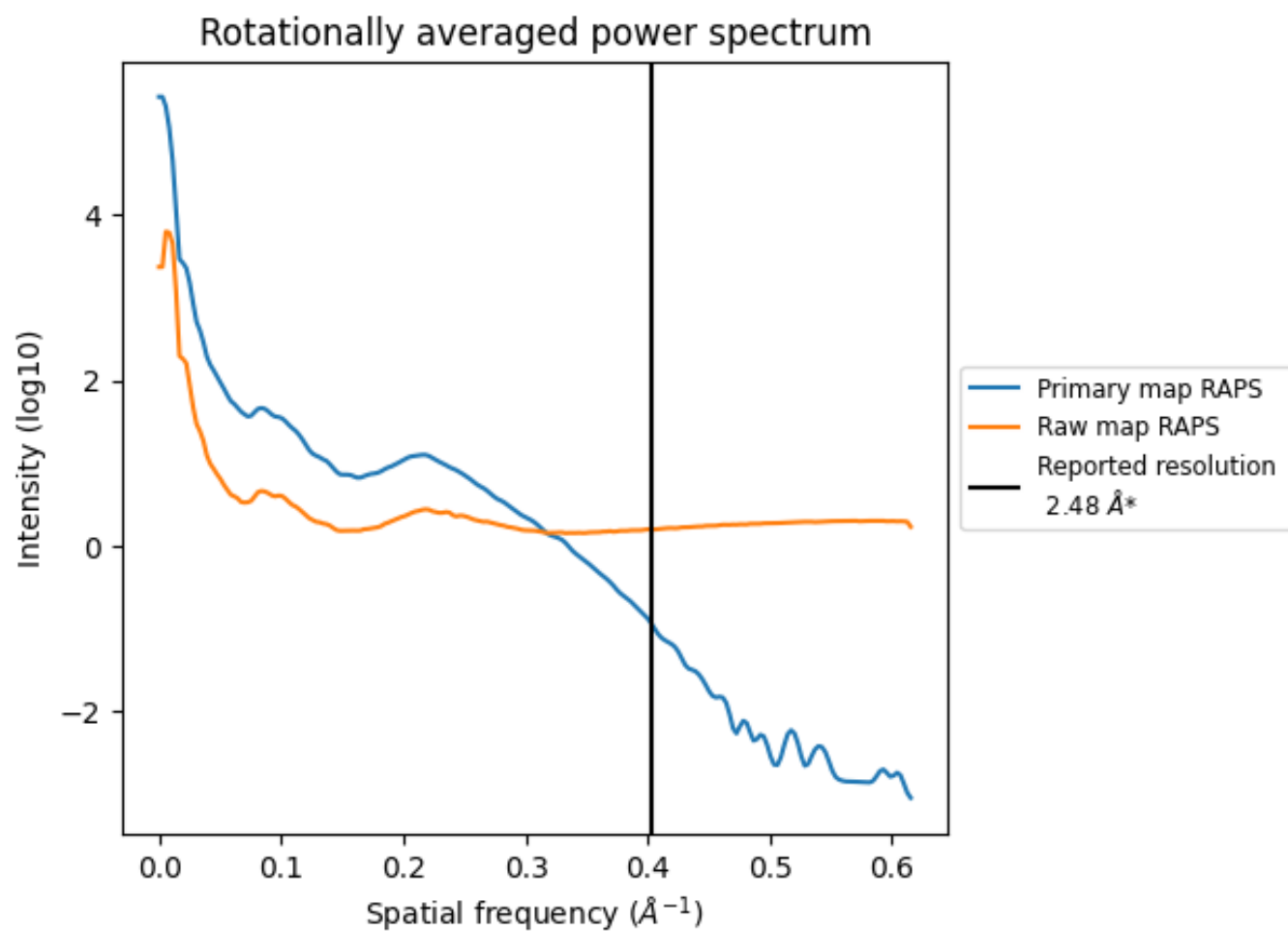
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 99 nm^3 ; this corresponds to an approximate mass of 89 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 ⓘ

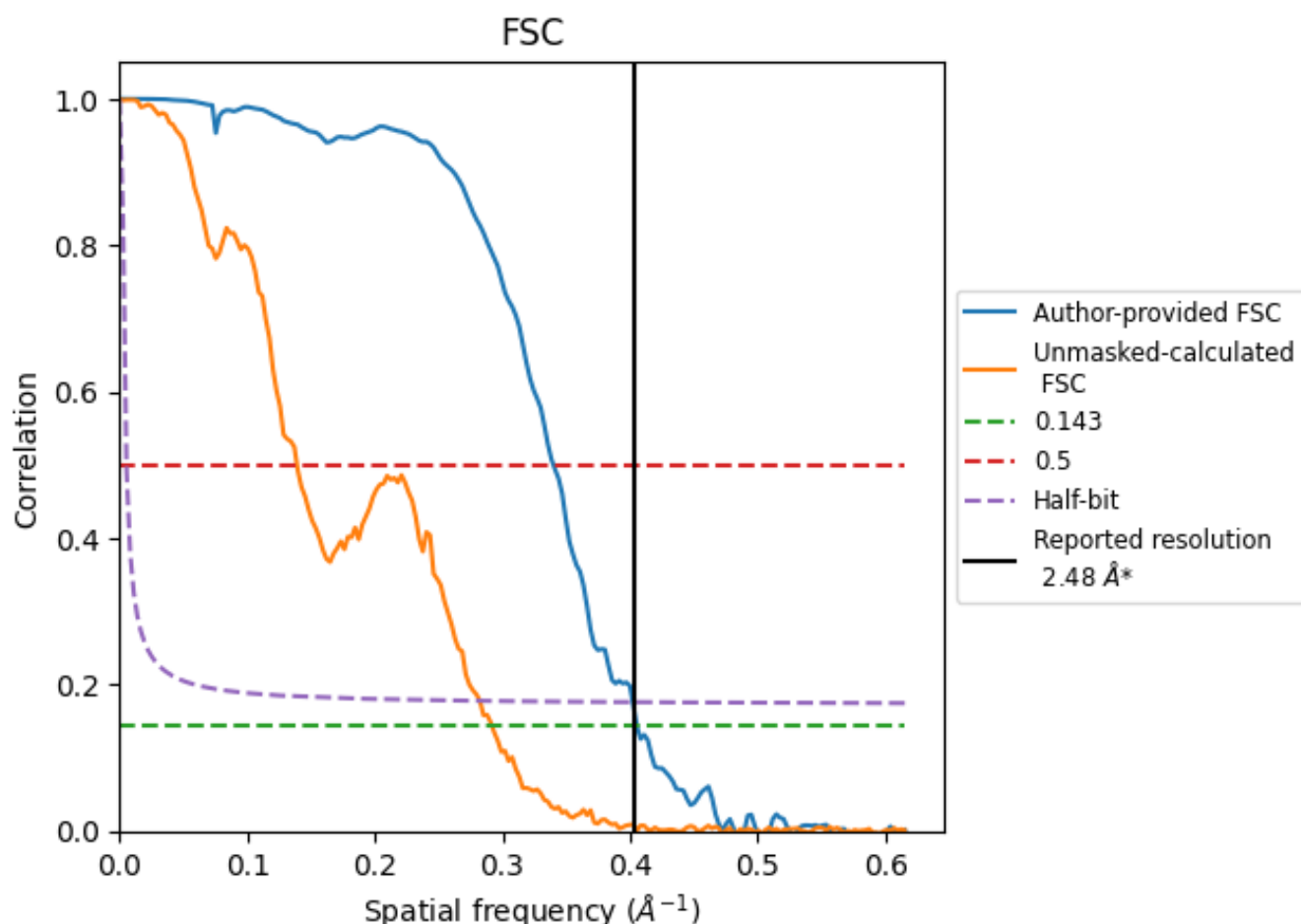


*Reported resolution corresponds to spatial frequency of 0.403 \AA^{-1}

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.403 Å⁻¹

8.2 Resolution estimates [i](#)

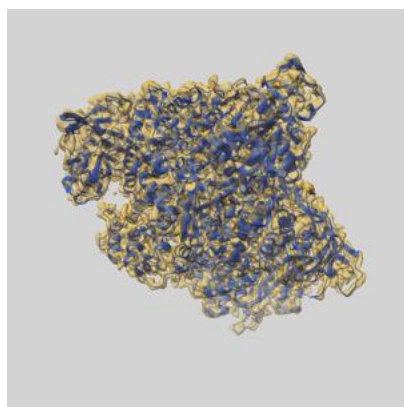
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.48	-	-
Author-provided FSC curve	2.47	2.94	2.49
Unmasked-calculated*	3.43	7.19	3.54

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.43 differs from the reported value 2.48 by more than 10 %

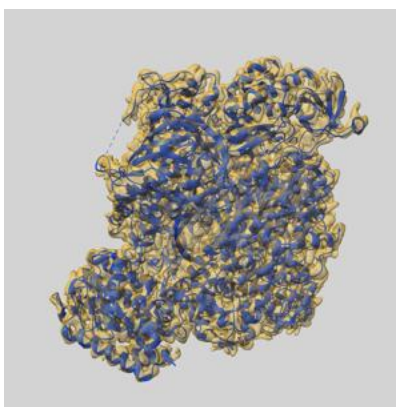
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-14144 and PDB model 7QTL. Per-residue inclusion information can be found in section [3](#) on page [6](#).

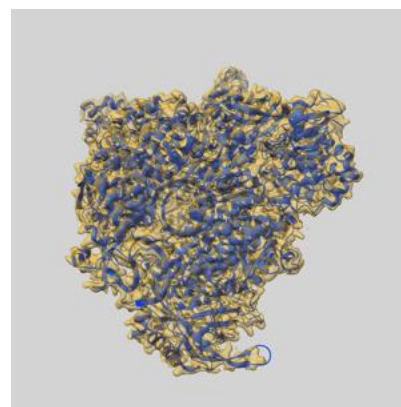
9.1 Map-model overlay [i](#)



X



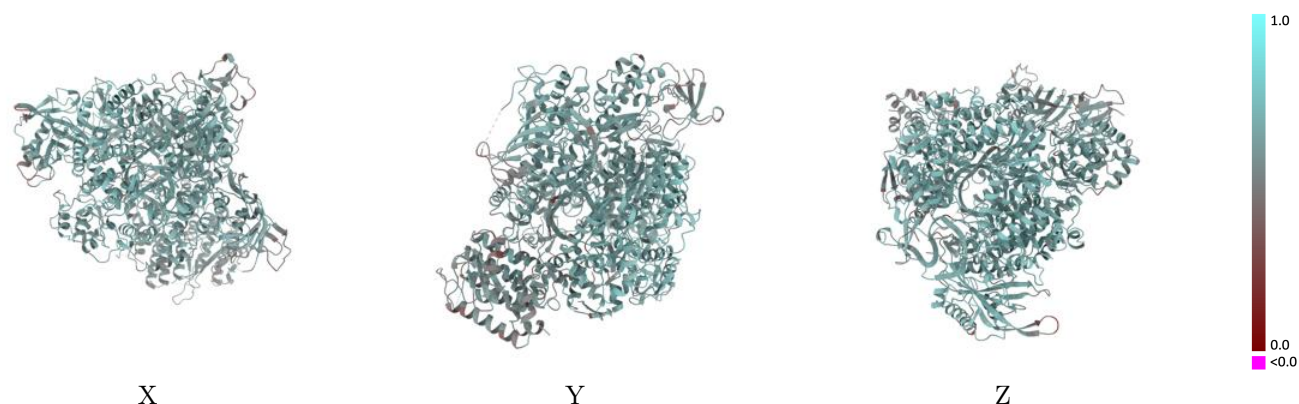
Y



Z

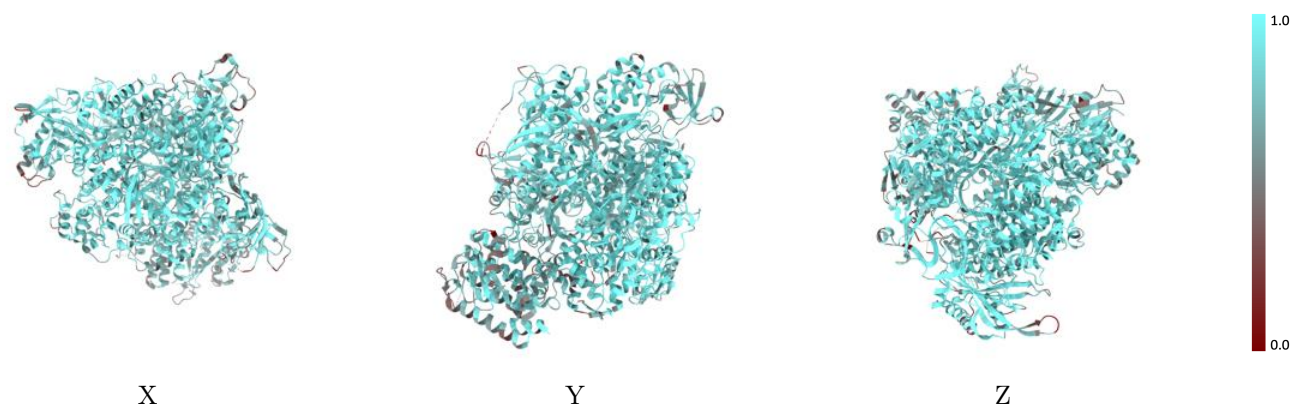
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



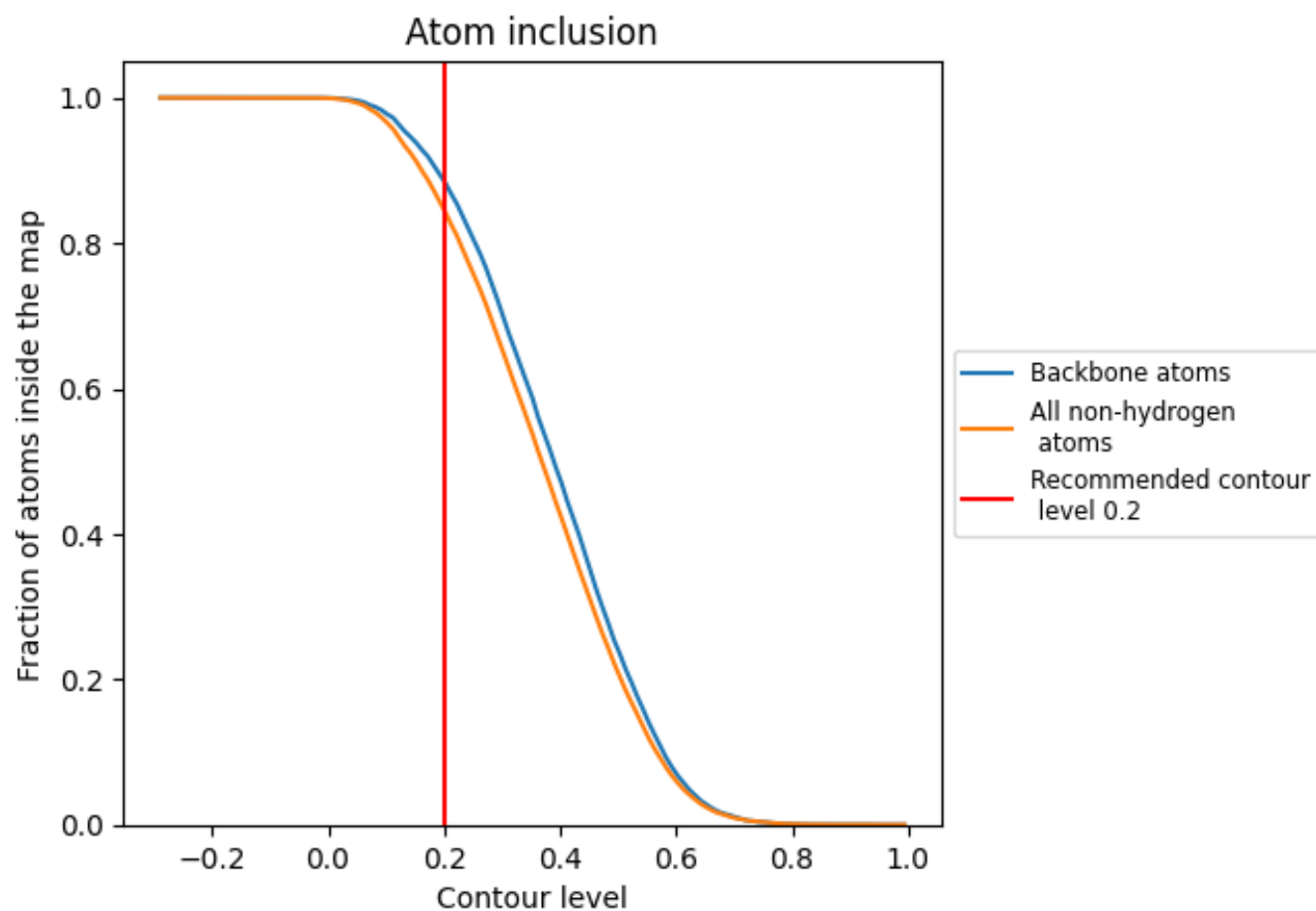
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 85% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.8459	<div><div></div></div> 0.6110
A	<div><div></div></div> 0.8298	<div><div></div></div> 0.6070
B	<div><div></div></div> 0.8916	<div><div></div></div> 0.6340
C	<div><div></div></div> 0.8213	<div><div></div></div> 0.5920
M	<div><div></div></div> 0.9111	<div><div></div></div> 0.6020
R	<div><div></div></div> 0.8889	<div><div></div></div> 0.6180
V	<div><div></div></div> 0.9120	<div><div></div></div> 0.6180

1.0

0.0

<0.0