



Full wwPDB EM Validation Report ⓘ

Aug 5, 2026 – 07:59 AM EDT

PDB ID : 7KIF / pdb_00007kif
EMDB ID : EMD-22886
Title : Mycobacterium tuberculosis WT RNAP transcription open promoter complex
with WhiB7 transcription factor
Authors : Lilic, M.; Darst, S.A.; Campbell, E.A.
Deposited on : 2020-10-23
Resolution : 2.94 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:

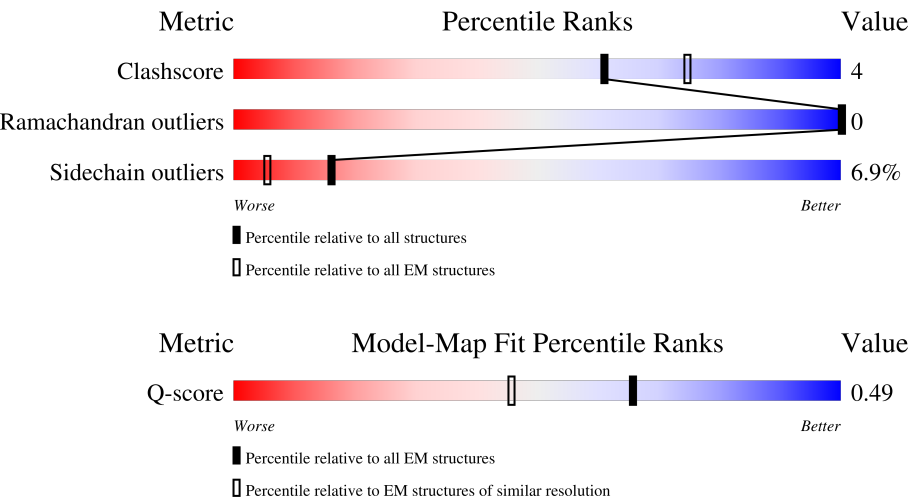
EMDB validation analysis : 0.0.1.dev133
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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) : 2.50

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.94 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.



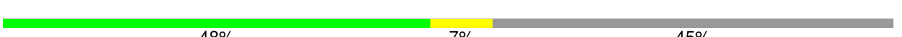
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13068 (2.44 - 3.44)

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	347	52%11%•35%
1	B	347	56%12%•32%
2	C	1172	78%15%•5%
3	D	1318	84%11%••

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Mol	Chain	Length	Quality of chain
4	E	110	
5	F	528	
6	J	111	
7	M	162	
8	O	100	
9	P	100	
10	Z	92	

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
13	SF4	Z	101	-	-	X	-

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 30322 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	225	Total	C	N	O	S	0	0
			1716	1080	296	338	2		
1	B	237	Total	C	N	O	S	0	0
			1765	1115	301	346	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	1111	Total	C	N	O	S	0	0
			8606	5392	1511	1664	39		

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	1270	Total	C	N	O	S	0	0
			9914	6208	1802	1862	42		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	GLY	-	expression tag	UNP A0A045J9E2
D	0	ALA	-	expression tag	UNP A0A045J9E2

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	E	83	Total	C	N	O	0	0
			649	414	108	127		

- Molecule 5 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	322	Total	C	N	O	S	0	0
			2548	1588	461	490	9		

- Molecule 6 is a protein called RNA polymerase-binding protein RbpA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	J	109	Total	C	N	O	S	0	0
			880	543	166	168	3		

- Molecule 7 is a protein called RNA polymerase-binding transcription factor CarD.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	M	159	Total	C	N	O	S	0	0
			1241	777	224	239	1		

- Molecule 8 is a DNA chain called DNA (63-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	O	63	Total	C	N	O	P	0	0
			1282	611	229	380	62		

- Molecule 9 is a DNA chain called DNA (55-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
9	P	55	Total	C	N	O	P	0	0
			1129	534	216	324	55		

- Molecule 10 is a protein called Probable transcriptional regulator WhiB7.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	Z	76	Total	C	N	O	S	0	0
			581	369	110	98	4		

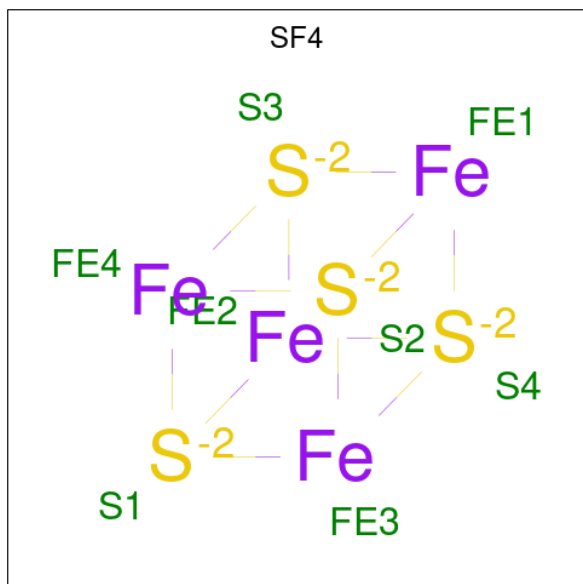
- Molecule 11 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
11	D	2	Total	Zn	0
			2	2	

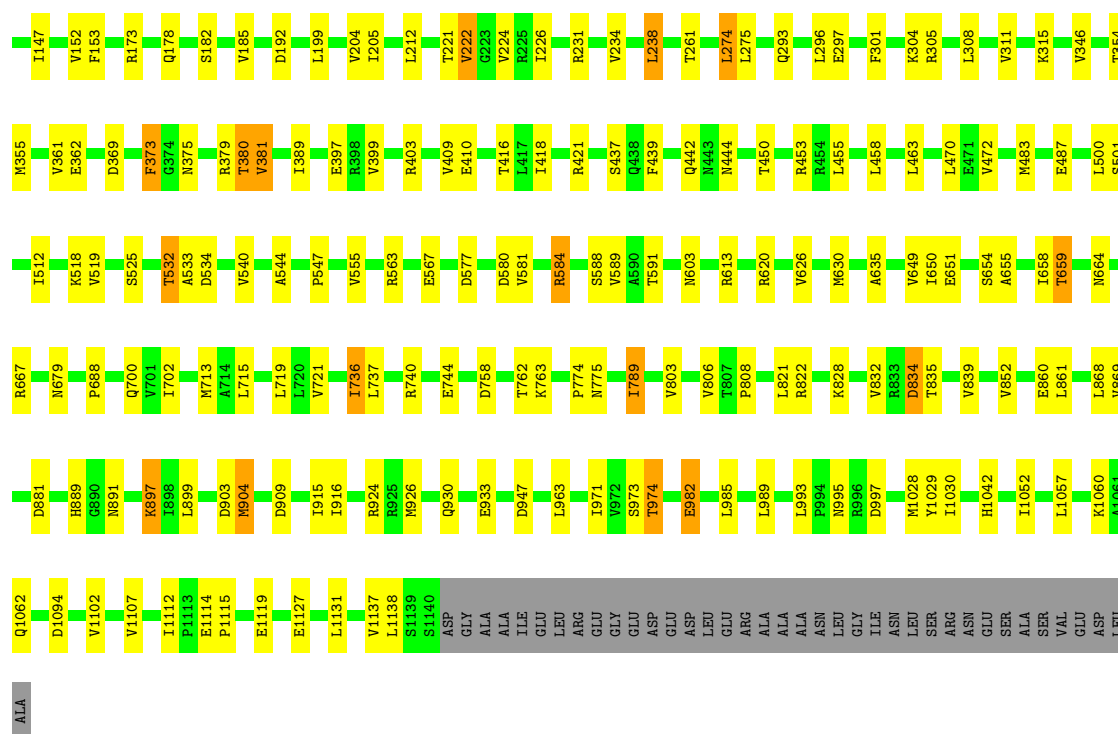
- Molecule 12 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
12	D	1	Total	Mg	0
			1	1	

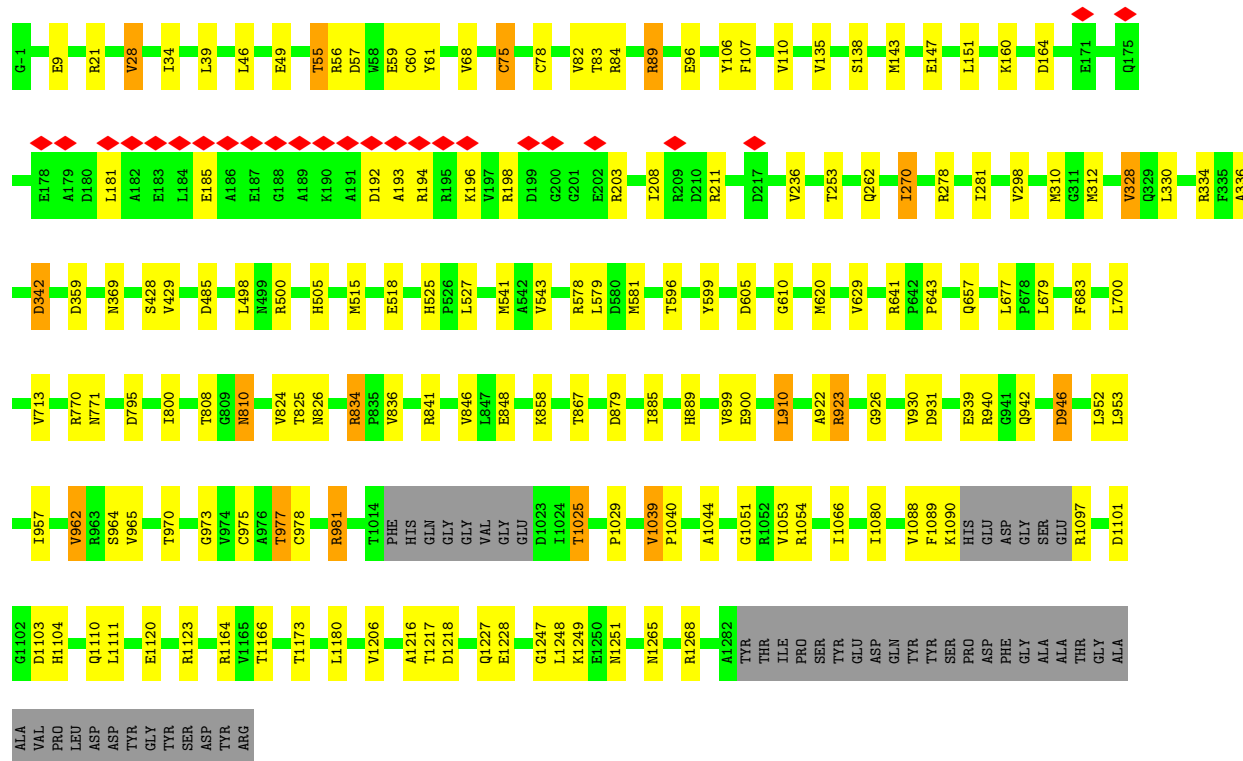
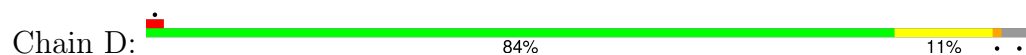
- Molecule 13 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
13	Z	1	Total	Fe	S	0
			8	4	4	

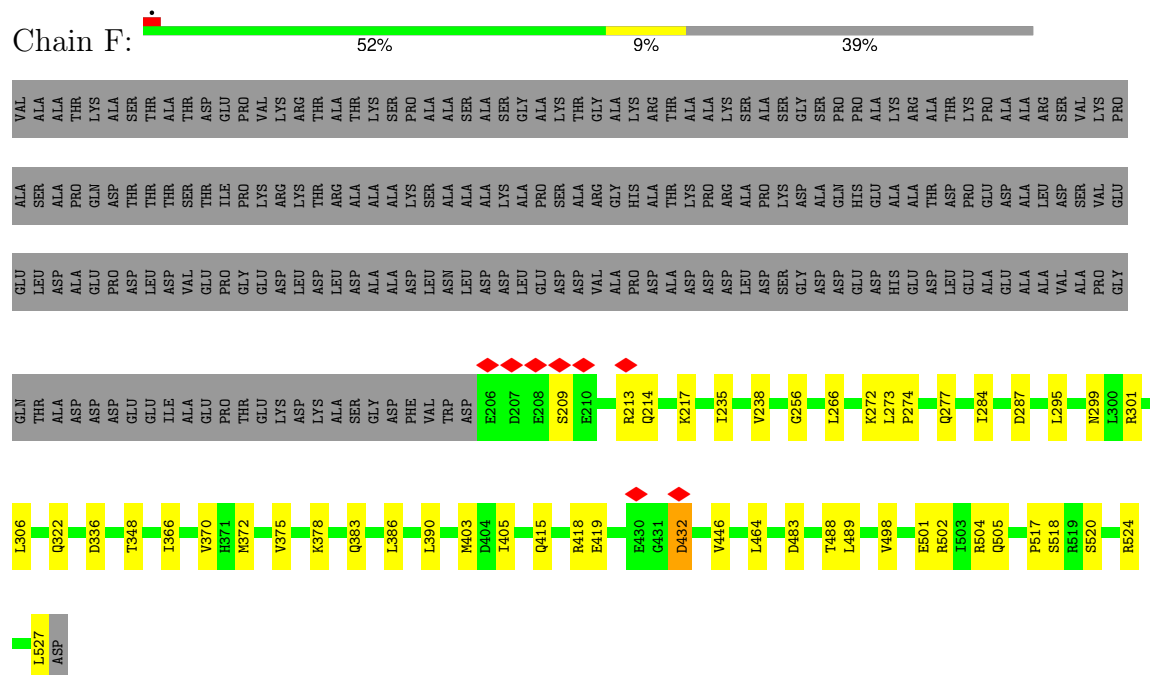


• Molecule 3: DNA-directed RNA polymerase subunit beta'

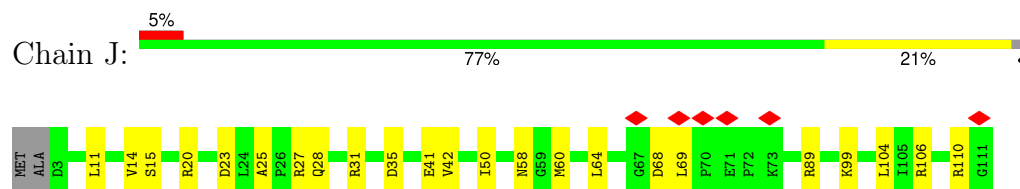


• Molecule 4: DNA-directed RNA polymerase subunit omega

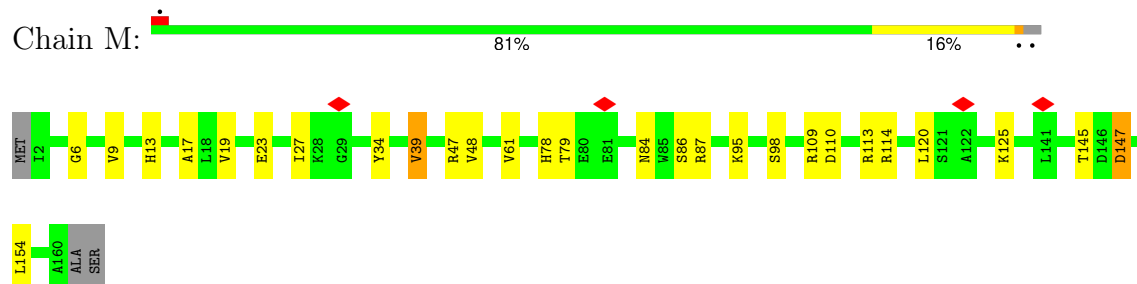
- Molecule 5: RNA polymerase sigma factor SigA



- Molecule 6: RNA polymerase-binding protein RbpA

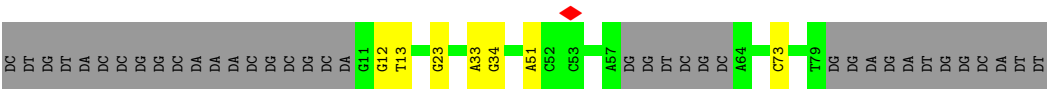


- Molecule 7: RNA polymerase-binding transcription factor CarD



- Molecule 8: DNA (63-MER)





● Molecule 9: DNA (55-MER)



● Molecule 10: Probable transcriptional regulator WhiB7



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	87743	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 ($e^-/\text{\AA}^2$)	66	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	44.350	Depositor
Minimum map value	-19.151	Depositor
Average map value	0.001	Depositor
Map value standard deviation	1.008	Depositor
Recommended contour level	3.6	Depositor
Map size ($\text{\AA}$)	330.0, 330.0, 330.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, MG, ZN

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	0.37	0/1742	0.44	0/2370
1	B	0.33	0/1792	0.53	0/2442
2	C	0.40	0/8765	0.51	0/11885
3	D	0.39	0/10078	0.54	4/13624 (0.0%)
4	E	0.37	0/662	0.50	0/901
5	F	0.26	0/2579	0.45	0/3477
6	J	0.25	0/896	0.48	0/1210
7	M	0.21	0/1257	0.57	2/1700 (0.1%)
8	O	0.29	0/1434	0.54	0/2208
9	P	0.31	0/1267	0.53	0/1950
10	Z	0.28	0/598	0.55	0/815
All	All	0.36	0/31070	0.52	6/42582 (0.0%)

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
3	D	0	1
5	F	0	1
All	All	0	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	M	147	ASP	CA-C-N	6.25	133.47	121.54
7	M	147	ASP	C-N-CA	6.25	133.47	121.54
3	D	578	ARG	CA-C-N	6.15	133.28	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	578	ARG	C-N-CA	6.15	133.28	121.54
3	D	834	ARG	CA-C-N	5.75	140.81	127.00
3	D	834	ARG	C-N-CA	5.75	140.81	127.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	579	LEU	Mainchain
5	F	273	LEU	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	1716	0	1756	18	0
1	B	1765	0	1794	18	0
2	C	8606	0	8544	92	0
3	D	9914	0	9986	77	0
4	E	649	0	645	7	0
5	F	2548	0	2569	28	0
6	J	880	0	852	13	0
7	M	1241	0	1259	15	0
8	O	1282	0	712	6	0
9	P	1129	0	616	6	0
10	Z	581	0	574	7	0
11	D	2	0	0	0	0
12	D	1	0	0	0	0
13	Z	8	0	0	3	0
All	All	30322	0	29307	262	0

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

All (262) 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)
3:D:60:CYS:SG	3:D:61:TYR:N	2.65	0.70
3:D:147:GLU:O	3:D:151:LEU:HB2	1.94	0.68
3:D:181:LEU:O	3:D:185:GLU:HB2	1.93	0.68
6:J:28:GLN:HE22	6:J:50:ILE:HD11	1.59	0.67
2:C:75:ALA:HA	2:C:79:ASP:HB3	1.78	0.64
2:C:891:ASN:HD21	2:C:1028:MET:HE1	1.65	0.62
5:F:274:PRO:HB2	5:F:277:GLN:HB2	1.83	0.61
5:F:432:ASP:OD1	5:F:432:ASP:N	2.35	0.60
2:C:563:ARG:HB2	2:C:567:GLU:HG3	1.83	0.60
2:C:1057:LEU:HD23	2:C:1062:GLN:HG2	1.82	0.60
2:C:1094:ASP:HB3	2:C:1119:GLU:H	1.67	0.59
3:D:500:ARG:HB2	3:D:541:MET:HE2	1.84	0.59
2:C:453:ARG:NH2	2:C:501:SER:O	2.35	0.59
3:D:181:LEU:HA	3:D:198:ARG:HH21	1.67	0.59
6:J:20:ARG:NH1	6:J:23:ASP:O	2.35	0.59
2:C:635:ALA:HB2	2:C:713:MET:HG2	1.85	0.58
5:F:209:SER:O	5:F:213:ARG:HB2	2.04	0.58
6:J:31:ARG:HG2	6:J:41:GLU:HG2	1.85	0.57
3:D:952:LEU:HD22	3:D:957:ILE:HD11	1.85	0.57
3:D:1164:ARG:NH2	3:D:1216:ALA:O	2.38	0.57
2:C:532:THR:OG1	2:C:533:ALA:N	2.37	0.57
1:A:53:SER:OG	1:A:161:ARG:NH2	2.38	0.57
2:C:173:ARG:NH1	2:C:437:SER:O	2.38	0.57
3:D:56:ARG:HB2	3:D:59:GLU:HB3	1.87	0.57
3:D:525:HIS:HE1	3:D:527:LEU:HD12	1.70	0.56
5:F:504:ARG:NH2	9:P:127:DA:OP2	2.39	0.56
3:D:39:LEU:HB3	6:J:11:LEU:HD13	1.87	0.56
2:C:891:ASN:ND2	2:C:930:GLN:OE1	2.36	0.56
9:P:126:DC:H2''	9:P:127:DA:H2'	1.87	0.56
4:E:89:GLU:OE2	4:E:97:ARG:NH1	2.39	0.56
2:C:1119:GLU:OE2	3:D:89:ARG:NH1	2.39	0.56
3:D:75:CYS:HB3	3:D:78:CYS:SG	2.46	0.55
2:C:995:ASN:N	2:C:995:ASN:OD1	2.39	0.55
3:D:1247:GLY:O	3:D:1251:ASN:ND2	2.38	0.55
1:A:223:ARG:HH21	1:B:213:LYS:HB2	1.69	0.55
2:C:899:LEU:HD12	2:C:904:MET:HE3	1.88	0.55
3:D:770:ARG:NH1	3:D:771:ASN:OD1	2.39	0.55
5:F:295:LEU:O	5:F:299:ASN:ND2	2.39	0.55
3:D:138:SER:HG	3:D:253:THR:HG1	1.55	0.55
5:F:518:SER:HA	10:Z:21:HIS:HB3	1.89	0.55
1:A:139:VAL:HG21	1:A:161:ARG:HH22	1.71	0.55
2:C:380:THR:OG1	2:C:381:VAL:N	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:GLU:OE2	2:C:620:ARG:NH1	2.40	0.55
3:D:1265:ASN:OD1	3:D:1268:ARG:NH2	2.40	0.55
1:B:24:GLU:HB3	1:B:191:LYS:HG3	1.88	0.55
7:M:109:ARG:NH1	7:M:110:ASP:OD1	2.40	0.55
3:D:931:ASP:OD1	3:D:931:ASP:N	2.39	0.54
1:A:142:ARG:NH2	1:B:230:GLU:OE1	2.40	0.54
3:D:1025:THR:HG21	3:D:1029:PRO:HB2	1.90	0.54
2:C:222:VAL:HG22	2:C:261:THR:HG21	1.90	0.54
3:D:143:MET:SD	3:D:143:MET:N	2.81	0.54
3:D:1054:ARG:NH2	3:D:1101:ASP:OD1	2.41	0.54
1:A:206:ASP:OD1	1:B:226:ASN:ND2	2.41	0.54
3:D:1089:PHE:O	3:D:1097:ARG:N	2.41	0.54
5:F:502:ARG:HD3	8:O:23:DG:H2'	1.89	0.53
2:C:221:THR:OG1	2:C:231:ARG:NH2	2.42	0.53
1:B:97:LEU:HB3	1:B:110:ILE:HG22	1.90	0.53
2:C:588:SER:OG	2:C:589:VAL:N	2.40	0.53
3:D:194:ARG:O	3:D:198:ARG:NH1	2.42	0.53
2:C:455:LEU:HD21	2:C:500:LEU:HG	1.90	0.53
5:F:235:ILE:O	5:F:301:ARG:NH1	2.42	0.53
2:C:654:SER:OG	2:C:655:ALA:N	2.42	0.52
2:C:721:VAL:HG23	2:C:915:ILE:HG13	1.91	0.52
2:C:891:ASN:OD1	2:C:891:ASN:N	2.43	0.52
5:F:277:GLN:OE1	6:J:106:ARG:NH1	2.42	0.52
3:D:334:ARG:NH1	5:F:419:GLU:O	2.43	0.52
3:D:879:ASP:OD2	3:D:1249:LYS:NZ	2.35	0.52
3:D:160:LYS:NZ	3:D:164:ASP:OD1	2.43	0.52
3:D:1090:LYS:HD3	3:D:1097:ARG:HE	1.74	0.51
2:C:808:PRO:HA	2:C:832:VAL:HG12	1.91	0.51
3:D:926:GLY:O	3:D:940:ARG:NH2	2.42	0.51
2:C:737:LEU:HD23	2:C:915:ILE:HG22	1.92	0.51
3:D:83:THR:OG1	3:D:84:ARG:N	2.43	0.51
3:D:1180:LEU:HD22	3:D:1206:VAL:HG11	1.93	0.51
5:F:483:ASP:N	5:F:483:ASP:OD1	2.44	0.51
2:C:444:ASN:N	2:C:444:ASN:OD1	2.43	0.51
7:M:87:ARG:NH2	8:O:51:DA:OP2	2.43	0.51
2:C:719:LEU:HD12	2:C:1030:ILE:HD11	1.92	0.51
2:C:487:GLU:OE2	2:C:613:ARG:NH1	2.42	0.51
2:C:1042:HIS:HB3	2:C:1060:LYS:HG3	1.92	0.51
1:A:94:THR:HG22	1:A:139:VAL:HG22	1.93	0.51
3:D:922:ALA:HB1	3:D:981:ARG:HG2	1.93	0.51
5:F:524:ARG:O	5:F:524:ARG:NH2	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:M:78:HIS:O	7:M:109:ARG:NH1	2.45	0.50
3:D:278:ARG:HA	3:D:281:ILE:HG12	1.92	0.50
10:Z:31:ASP:OD1	10:Z:31:ASP:N	2.43	0.50
2:C:304:LYS:HD3	2:C:305:ARG:HG2	1.92	0.50
4:E:48:SER:OG	4:E:49:SER:N	2.45	0.50
2:C:46:GLU:OE2	2:C:584:ARG:NH2	2.45	0.50
2:C:822:ARG:NH1	2:C:828:LYS:O	2.45	0.50
2:C:577:ASP:OD1	2:C:577:ASP:N	2.45	0.50
2:C:915:ILE:HD13	2:C:1030:ILE:HD13	1.94	0.50
3:D:641:ARG:HA	3:D:657:GLN:HG3	1.94	0.50
2:C:740:ARG:NH1	2:C:744:GLU:OE1	2.45	0.49
7:M:147:ASP:OD1	7:M:147:ASP:N	2.45	0.49
1:B:25:PRO:HB2	1:B:189:PHE:HD2	1.77	0.49
2:C:659:THR:O	2:C:659:THR:OG1	2.30	0.49
2:C:120:ASP:OD1	2:C:120:ASP:N	2.44	0.49
2:C:947:ASP:OD1	2:C:947:ASP:N	2.43	0.49
3:D:946:ASP:OD1	3:D:946:ASP:N	2.32	0.49
3:D:143:MET:O	3:D:147:GLU:HB2	2.13	0.49
10:Z:52:CYS:HB3	13:Z:101:SF4:S4	2.53	0.49
2:C:182:SER:OG	2:C:315:LYS:NZ	2.46	0.49
2:C:658:ILE:HD11	2:C:688:PRO:HB3	1.95	0.49
5:F:214:GLN:HA	5:F:217:LYS:HZ3	1.77	0.49
2:C:38:ARG:HG2	2:C:973:SER:HB3	1.95	0.48
2:C:758:ASP:HB3	2:C:868:LEU:HD12	1.94	0.48
4:E:60:ARG:NH2	4:E:98:GLU:OE2	2.46	0.48
3:D:49:GLU:OE2	3:D:55:THR:OG1	2.31	0.48
5:F:266:LEU:HD13	5:F:272:LYS:HE3	1.96	0.48
7:M:23:GLU:N	7:M:34:TYR:O	2.46	0.48
2:C:982:GLU:OE1	3:D:841:ARG:NH2	2.46	0.48
2:C:1138:LEU:HB2	3:D:9:GLU:HB2	1.95	0.48
4:E:47:VAL:HG21	4:E:53:LEU:HB2	1.95	0.48
2:C:442:GLN:NE2	2:C:679:ASN:OD1	2.37	0.48
3:D:975:CYS:SG	3:D:978:CYS:HB2	2.54	0.48
1:A:155:SER:OG	1:A:156:GLY:N	2.45	0.48
2:C:178:GLN:OE1	2:C:379:ARG:NH2	2.47	0.48
2:C:410:GLU:O	7:M:13:HIS:NE2	2.47	0.48
3:D:1217:THR:OG1	3:D:1218:ASP:N	2.46	0.48
1:A:104:GLU:OE2	1:A:124:HIS:ND1	2.44	0.47
1:B:124:HIS:NE2	1:B:126:ALA:O	2.47	0.47
1:B:99:LYS:NZ	1:B:104:GLU:O	2.43	0.47
7:M:79:THR:OG1	7:M:113:ARG:NH1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:515:MET:HE2	3:D:515:MET:HB3	1.76	0.47
3:D:342:ASP:N	3:D:342:ASP:OD1	2.47	0.47
3:D:1120:GLU:OE2	3:D:1123:ARG:NH2	2.39	0.47
1:A:57:ASP:OD1	1:A:57:ASP:N	2.38	0.47
2:C:881:ASP:OD1	2:C:881:ASP:N	2.43	0.47
2:C:835:THR:O	2:C:835:THR:OG1	2.32	0.47
1:A:71:GLU:HB3	1:A:75:GLU:HB3	1.97	0.47
2:C:518:LYS:O	2:C:525:SER:OG	2.30	0.47
3:D:939:GLU:OE2	3:D:942:GLN:NE2	2.48	0.47
6:J:25:ALA:O	6:J:27:ARG:NH2	2.47	0.47
2:C:974:THR:O	2:C:974:THR:OG1	2.33	0.46
7:M:114:ARG:HB3	7:M:120:LEU:HD11	1.97	0.46
1:A:130:ASP:OD1	1:A:130:ASP:N	2.48	0.46
2:C:762:THR:OG1	2:C:763:LYS:N	2.42	0.46
5:F:238:VAL:O	5:F:301:ARG:NH2	2.46	0.46
1:B:22:VAL:HG12	1:B:193:ILE:HG23	1.98	0.46
1:B:167:ILE:HG23	3:D:620:MET:HE1	1.98	0.46
1:B:198:THR:OG1	1:B:199:LYS:O	2.27	0.46
3:D:1227:GLN:HG2	3:D:1228:GLU:HG3	1.97	0.46
3:D:605:ASP:OD1	3:D:605:ASP:N	2.49	0.46
5:F:256:GLY:HA2	5:F:284:ILE:HG22	1.96	0.46
5:F:378:LYS:HE3	5:F:403:MET:HE3	1.98	0.45
2:C:274:LEU:HD11	2:C:293:GLN:HA	1.98	0.45
5:F:336:ASP:OD1	6:J:89:ARG:NH2	2.48	0.45
5:F:502:ARG:NE	8:O:23:DG:OP2	2.38	0.45
5:F:415:GLN:OE1	5:F:418:ARG:NH1	2.44	0.45
1:A:64:THR:OG1	1:A:65:THR:N	2.49	0.45
2:C:821:LEU:HD23	5:F:527:LEU:HB2	1.98	0.45
2:C:889:HIS:NE2	2:C:933:GLU:OE1	2.42	0.45
1:B:165:ASP:OD1	1:B:165:ASP:N	2.49	0.45
2:C:98:ASP:OD1	2:C:98:ASP:N	2.47	0.45
2:C:1102:VAL:HG23	2:C:1112:ILE:HG23	1.97	0.45
10:Z:52:CYS:CB	13:Z:101:SF4:S4	3.05	0.45
2:C:373:PHE:HE2	2:C:512:ILE:HD12	1.81	0.45
3:D:57:ASP:HB2	6:J:15:SER:H	1.81	0.45
1:A:93:VAL:HG21	1:A:116:VAL:HG11	1.99	0.45
2:C:775:ASN:N	2:C:775:ASN:OD1	2.50	0.45
3:D:106:TYR:HB3	3:D:312:MET:HE2	1.98	0.45
3:D:328:VAL:HG13	3:D:336:ALA:HB3	1.99	0.45
2:C:713:MET:HE3	2:C:713:MET:HB3	1.77	0.45
10:Z:67:GLY:N	13:Z:101:SF4:S2	2.86	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:LYS:HD3	1:B:105:VAL:HG22	1.99	0.44
2:C:354:THR:OG1	2:C:362:GLU:OE1	2.31	0.44
3:D:21:ARG:NH2	3:D:96:GLU:OE2	2.46	0.44
1:B:129:ASN:OD1	1:B:129:ASN:N	2.51	0.44
3:D:910:LEU:HD13	3:D:910:LEU:HA	1.87	0.44
5:F:306:LEU:HD11	5:F:348:THR:HG23	1.99	0.44
2:C:924:ARG:HD3	3:D:808:THR:HB	1.99	0.44
3:D:1044:ALA:HB3	3:D:1110:GLN:HE21	1.81	0.44
9:P:80:DT:H2''	9:P:81:DG:H5''	1.98	0.44
2:C:224:VAL:HB	2:C:234:VAL:HG22	2.00	0.44
2:C:664:ASN:OD1	2:C:664:ASN:N	2.51	0.44
1:B:80:LEU:HD22	1:B:138:LEU:HD21	2.00	0.44
2:C:403:ARG:NH1	2:C:416:THR:O	2.45	0.44
3:D:643:PRO:HD3	3:D:683:PHE:HB3	1.99	0.44
6:J:58:ASN:HB2	6:J:60:MET:HG2	1.99	0.44
2:C:547:PRO:HB2	2:C:555:VAL:HB	2.00	0.43
1:A:90:ASP:OD1	1:A:90:ASP:N	2.36	0.43
1:B:177:LYS:HE3	1:B:177:LYS:HB2	1.87	0.43
2:C:926:MET:HE3	2:C:926:MET:HB3	1.81	0.43
2:C:985:LEU:HD23	2:C:985:LEU:HA	1.85	0.43
3:D:193:ALA:HA	3:D:196:LYS:HG2	2.00	0.43
3:D:208:ILE:HG12	3:D:211:ARG:HH22	1.82	0.43
7:M:145:THR:O	7:M:145:THR:OG1	2.35	0.43
9:P:81:DG:H2'	9:P:82:DG:C8	2.54	0.43
1:A:123:MET:HE3	1:A:123:MET:HB2	1.89	0.43
2:C:651:GLU:OE1	2:C:667:ARG:NH1	2.51	0.43
3:D:836:VAL:HG13	3:D:848:GLU:HB3	2.01	0.43
6:J:68:ASP:OD1	6:J:68:ASP:N	2.50	0.43
7:M:95:LYS:O	7:M:98:SER:OG	2.36	0.43
2:C:238:LEU:HD13	2:C:238:LEU:HA	1.91	0.43
3:D:369:ASN:OD1	5:F:322:GLN:NE2	2.51	0.43
1:A:72:ASP:OD1	1:A:72:ASP:N	2.46	0.42
2:C:421:ARG:NH2	5:F:383:GLN:OE1	2.52	0.42
2:C:736:ILE:HG13	2:C:916:ILE:HB	2.01	0.42
5:F:403:MET:HE2	5:F:403:MET:HB3	1.78	0.42
2:C:897:LYS:HE2	2:C:897:LYS:HB2	1.94	0.42
2:C:1052:ILE:O	3:D:89:ARG:NH2	2.53	0.42
3:D:1039:VAL:HA	3:D:1040:PRO:HD3	1.91	0.42
5:F:517:PRO:HA	5:F:520:SER:HB3	2.00	0.42
3:D:834:ARG:HD2	3:D:834:ARG:HA	1.87	0.42
7:M:125:LYS:HE2	7:M:125:LYS:HB3	1.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:62:GLU:HG3	2:C:70:TRP:HB2	2.01	0.42
1:B:59:VAL:HG21	1:B:66:VAL:HG22	2.02	0.42
2:C:774:PRO:HD2	2:C:834:ASP:HB2	2.02	0.42
3:D:970:THR:OG1	3:D:973:GLY:O	2.37	0.42
3:D:1051:GLY:O	3:D:1104:HIS:ND1	2.52	0.42
8:O:73:DC:N4	9:P:81:DG:O6	2.53	0.42
3:D:923:ARG:HD3	3:D:962:VAL:HG11	2.01	0.42
6:J:110:ARG:HD3	6:J:110:ARG:HA	1.91	0.42
2:C:544:ALA:HB2	2:C:580:ASP:HB2	2.01	0.41
3:D:270:ILE:H	3:D:270:ILE:HG12	1.60	0.41
3:D:810:ASN:OD1	3:D:810:ASN:N	2.53	0.41
2:C:369:ASP:O	2:C:375:ASN:ND2	2.44	0.41
2:C:903:ASP:OD1	2:C:903:ASP:N	2.48	0.41
3:D:262:GLN:HG3	3:D:310:MET:HE1	2.01	0.41
3:D:826:ASN:O	3:D:858:LYS:NZ	2.43	0.41
4:E:60:ARG:HH22	4:E:80:GLY:HA3	1.86	0.41
2:C:86:LEU:HD21	2:C:389:ILE:HD13	2.01	0.41
6:J:35:ASP:OD1	6:J:35:ASP:N	2.40	0.41
7:M:84:ASN:OD1	7:M:86:SER:OG	2.32	0.41
8:O:12:DG:H2"	8:O:13:DT:H72	2.02	0.41
2:C:963:LEU:HD23	2:C:963:LEU:HA	1.90	0.41
3:D:599:TYR:HA	3:D:610:GLY:HA3	2.02	0.41
5:F:501:GLU:HA	5:F:504:ARG:HG2	2.03	0.41
2:C:79:ASP:OD1	2:C:79:ASP:N	2.54	0.41
6:J:23:ASP:OD2	6:J:23:ASP:N	2.52	0.41
5:F:372:MET:HA	5:F:375:VAL:HG22	2.03	0.41
2:C:132:PRO:HB3	2:C:153:PHE:HE1	1.86	0.41
2:C:789:ILE:HG23	2:C:803:VAL:HG22	2.03	0.41
3:D:203:ARG:HA	3:D:203:ARG:HD2	1.90	0.41
3:D:310:MET:HB3	3:D:310:MET:HE3	1.85	0.41
3:D:498:LEU:HG	3:D:543:VAL:HG22	2.03	0.41
3:D:889:HIS:O	3:D:977:THR:OG1	2.38	0.41
8:O:33:DA:H2"	8:O:34:DG:H8	1.85	0.41
1:A:95:MET:HB3	1:A:113:PRO:HD3	2.03	0.41
2:C:34:GLY:HA3	2:C:700:GLN:HG3	2.03	0.41
2:C:152:VAL:HG21	2:C:418:ILE:HD12	2.04	0.41
2:C:50:VAL:HG21	2:C:630:MET:HE1	2.02	0.40
2:C:715:LEU:N	2:C:1029:TYR:OH	2.54	0.40
4:E:36:THR:O	4:E:36:THR:OG1	2.35	0.40
4:E:68:TYR:HD1	4:E:68:TYR:HA	1.79	0.40
2:C:192:ASP:HB2	2:C:199:LEU:HD11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:964:SER:OG	3:D:965:VAL:N	2.53	0.40
1:B:183:VAL:HG12	1:B:188:ASP:HA	2.04	0.40
3:D:28:VAL:HG21	3:D:46:LEU:HD23	2.03	0.40
2:C:297:GLU:HA	2:C:301:PHE:HD2	1.86	0.40
2:C:789:ILE:H	2:C:789:ILE:HG13	1.41	0.40
2:C:1114:GLU:HA	2:C:1115:PRO:HD3	1.93	0.40
3:D:1103:ASP:OD1	3:D:1104:HIS:N	2.54	0.40
7:M:6:GLY:H	7:M:19:VAL:HG23	1.87	0.40
7:M:23:GLU:OE1	7:M:47:ARG:NH1	2.48	0.40
10:Z:38:VAL:O	10:Z:42:LEU:HB2	2.22	0.40
10:Z:50:ARG:HD3	10:Z:50:ARG:HA	1.90	0.40
2:C:113:ASP:HB2	2:C:132:PRO:HD2	2.04	0.40
3:D:192:ASP:O	3:D:196:LYS:N	2.55	0.40
7:M:17:ALA:HB2	7:M:39:VAL:HB	2.02	0.40
9:P:122:DT:H2'	9:P:123:DG:C8	2.57	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	223/347 (64%)	204 (92%)	19 (8%)	0	100	100
1	B	235/347 (68%)	211 (90%)	24 (10%)	0	100	100
2	C	1109/1172 (95%)	1023 (92%)	86 (8%)	0	100	100
3	D	1264/1318 (96%)	1192 (94%)	72 (6%)	0	100	100
4	E	81/110 (74%)	77 (95%)	4 (5%)	0	100	100
5	F	320/528 (61%)	304 (95%)	16 (5%)	0	100	100
6	J	107/111 (96%)	97 (91%)	10 (9%)	0	100	100
7	M	157/162 (97%)	146 (93%)	11 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	Z	74/92 (80%)	64 (86%)	10 (14%)	0	100	100
All	All	3570/4187 (85%)	3318 (93%)	252 (7%)	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	194/297 (65%)	171 (88%)	23 (12%)	5	14
1	B	195/297 (66%)	177 (91%)	18 (9%)	8	22
2	C	937/993 (94%)	868 (93%)	69 (7%)	13	30
3	D	1048/1095 (96%)	991 (95%)	57 (5%)	20	42
4	E	69/90 (77%)	63 (91%)	6 (9%)	9	24
5	F	268/427 (63%)	255 (95%)	13 (5%)	22	47
6	J	92/97 (95%)	86 (94%)	6 (6%)	15	35
7	M	129/131 (98%)	123 (95%)	6 (5%)	23	48
10	Z	60/76 (79%)	53 (88%)	7 (12%)	5	15
All	All	2992/3503 (85%)	2787 (93%)	205 (7%)	16	33

All (205) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	22	VAL
1	A	30	PHE
1	A	43	LEU
1	A	51	VAL
1	A	56	ILE
1	A	59	VAL
1	A	65	THR
1	A	84	VAL

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Mol	Chain	Res	Type
1	A	111	VAL
1	A	116	VAL
1	A	123	MET
1	A	127	THR
1	A	130	ASP
1	A	136	VAL
1	A	147	VAL
1	A	150	VAL
1	A	167	ILE
1	A	203	SER
1	A	215	LEU
1	A	217	GLU
1	A	218	LEU
1	A	221	LEU
1	B	33	THR
1	B	34	LEU
1	B	43	LEU
1	B	56	ILE
1	B	74	THR
1	B	84	VAL
1	B	111	VAL
1	B	125	ILE
1	B	131	LYS
1	B	147	VAL
1	B	150	VAL
1	B	162	ILE
1	B	165	ASP
1	B	175	THR
1	B	178	VAL
1	B	198	THR
1	B	216	VAL
1	B	227	VAL
2	C	56	VAL
2	C	64	LEU
2	C	83	VAL
2	C	90	LEU
2	C	124	ASP
2	C	147	ILE
2	C	185	VAL
2	C	204	VAL
2	C	205	ILE
2	C	212	LEU

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Mol	Chain	Res	Type
2	C	222	VAL
2	C	226	ILE
2	C	238	LEU
2	C	274	LEU
2	C	275	LEU
2	C	296	LEU
2	C	308	LEU
2	C	311	VAL
2	C	346	VAL
2	C	355	MET
2	C	361	VAL
2	C	373	PHE
2	C	380	THR
2	C	381	VAL
2	C	397	GLU
2	C	399	VAL
2	C	409	VAL
2	C	439	PHE
2	C	450	THR
2	C	458	LEU
2	C	463	LEU
2	C	470	LEU
2	C	472	VAL
2	C	483	MET
2	C	519	VAL
2	C	532	THR
2	C	534	ASP
2	C	540	VAL
2	C	581	VAL
2	C	584	ARG
2	C	591	THR
2	C	603	ASN
2	C	626	VAL
2	C	649	VAL
2	C	650	ILE
2	C	659	THR
2	C	702	ILE
2	C	736	ILE
2	C	789	ILE
2	C	806	VAL
2	C	834	ASP
2	C	839	VAL

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Mol	Chain	Res	Type
2	C	852	VAL
2	C	860	GLU
2	C	861	LEU
2	C	869	VAL
2	C	897	LYS
2	C	904	MET
2	C	909	ASP
2	C	971	ILE
2	C	974	THR
2	C	982	GLU
2	C	989	LEU
2	C	993	LEU
2	C	997	ASP
2	C	1107	VAL
2	C	1127	GLU
2	C	1131	LEU
2	C	1137	VAL
3	D	28	VAL
3	D	34	ILE
3	D	55	THR
3	D	68	VAL
3	D	75	CYS
3	D	82	VAL
3	D	89	ARG
3	D	107	PHE
3	D	110	VAL
3	D	135	VAL
3	D	236	VAL
3	D	270	ILE
3	D	298	VAL
3	D	328	VAL
3	D	330	LEU
3	D	342	ASP
3	D	359	ASP
3	D	428	SER
3	D	429	VAL
3	D	485	ASP
3	D	505	HIS
3	D	518	GLU
3	D	581	MET
3	D	596	THR
3	D	629	VAL

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Mol	Chain	Res	Type
3	D	677	LEU
3	D	679	LEU
3	D	700	LEU
3	D	713	VAL
3	D	795	ASP
3	D	800	ILE
3	D	810	ASN
3	D	824	VAL
3	D	825	THR
3	D	846	VAL
3	D	867	THR
3	D	885	ILE
3	D	899	VAL
3	D	900	GLU
3	D	910	LEU
3	D	923	ARG
3	D	930	VAL
3	D	946	ASP
3	D	953	LEU
3	D	962	VAL
3	D	977	THR
3	D	981	ARG
3	D	1025	THR
3	D	1039	VAL
3	D	1053	VAL
3	D	1066	ILE
3	D	1080	ILE
3	D	1088	VAL
3	D	1111	LEU
3	D	1166	THR
3	D	1173	THR
3	D	1248	LEU
4	E	31	THR
4	E	33	LEU
4	E	35	ILE
4	E	49	SER
4	E	50	LYS
4	E	79	VAL
5	F	287	ASP
5	F	366	ILE
5	F	370	VAL
5	F	386	LEU

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Mol	Chain	Res	Type
5	F	390	LEU
5	F	405	ILE
5	F	432	ASP
5	F	446	VAL
5	F	464	LEU
5	F	488	THR
5	F	489	LEU
5	F	498	VAL
5	F	505	GLN
6	J	14	VAL
6	J	42	VAL
6	J	64	LEU
6	J	69	LEU
6	J	99	LYS
6	J	104	LEU
7	M	9	VAL
7	M	27	ILE
7	M	39	VAL
7	M	48	VAL
7	M	61	VAL
7	M	154	LEU
10	Z	17	VAL
10	Z	22	VAL
10	Z	24	ASP
10	Z	32	THR
10	Z	36	LEU
10	Z	63	TRP
10	Z	77	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
3	D	369	ASN
3	D	415	GLN
3	D	544	HIS
3	D	1125	GLN
3	D	1160	GLN
5	F	234	GLN
5	F	242	ASN
5	F	322	GLN
5	F	325	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
6	J	28	GLN
7	M	31	GLN
7	M	52	ASN

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, 3 are monoatomic - leaving 1 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	$\# Z > 2$
13	SF4	Z	101	10	0,12,12	-	-	-		

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
13	SF4	Z	101	10	-	-	0/6/5/5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

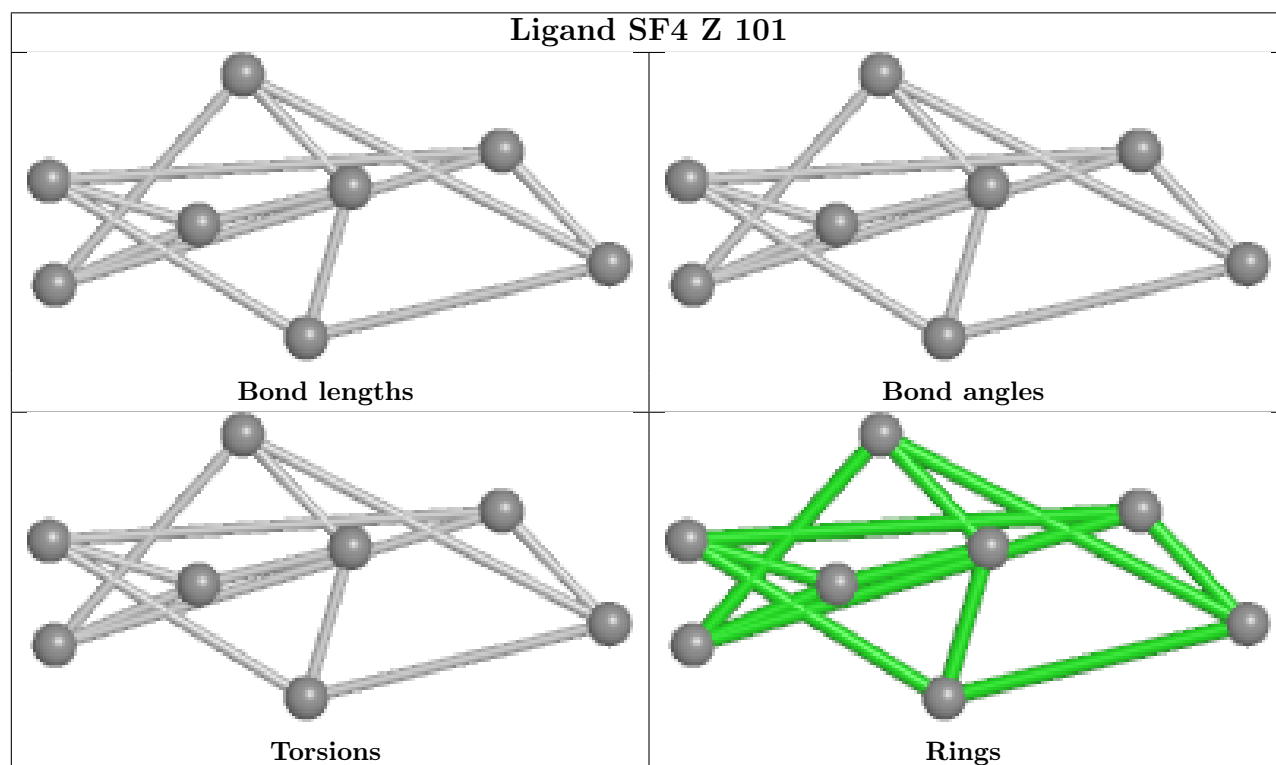
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	Z	101	SF4	3	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.

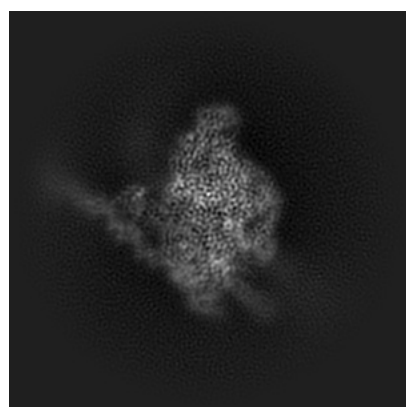
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22886. These allow visual inspection of the internal detail of the map and identification of artifacts.

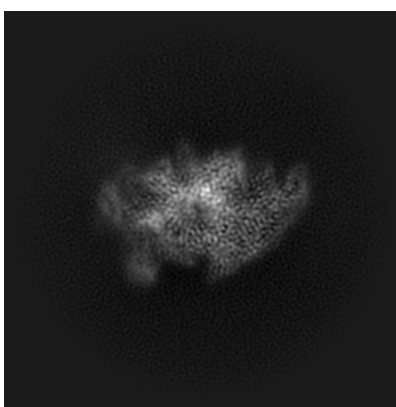
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

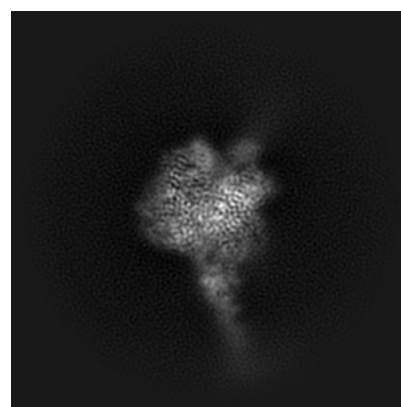
6.1.1 Primary map



X



Y

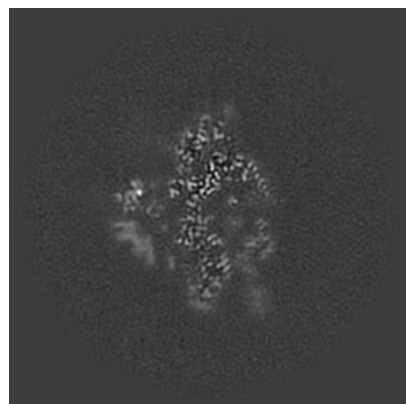


Z

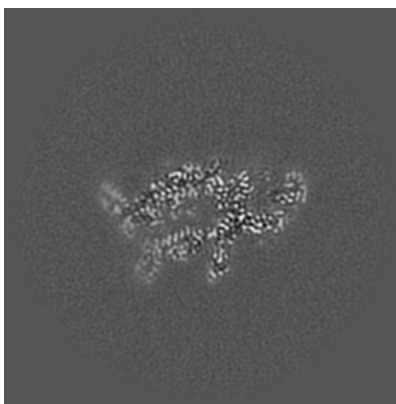
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

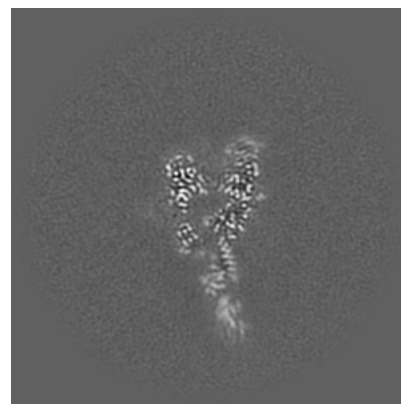
6.2.1 Primary map



X Index: 150



Y Index: 150

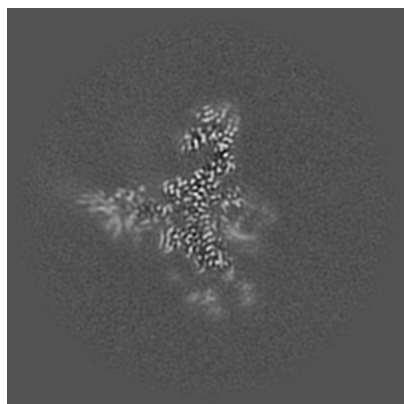


Z Index: 150

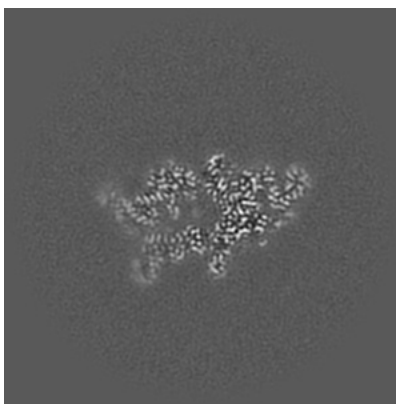
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

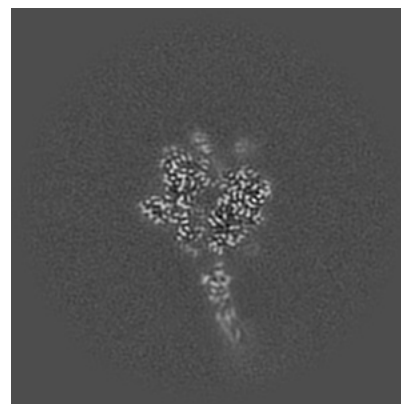
6.3.1 Primary map



X Index: 160



Y Index: 154

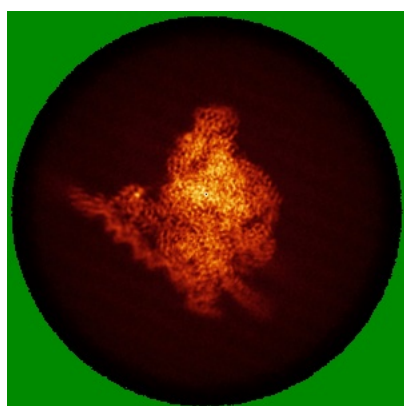


Z Index: 158

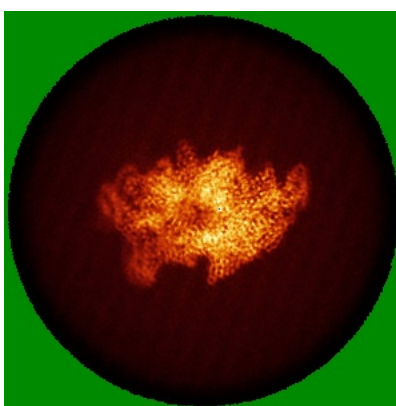
The images above show the largest variance slices of the map in three orthogonal directions.

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

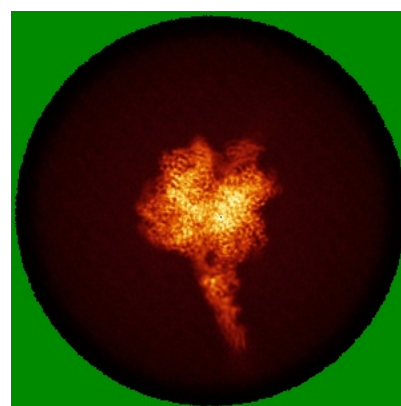
6.4.1 Primary map



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 3.6. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

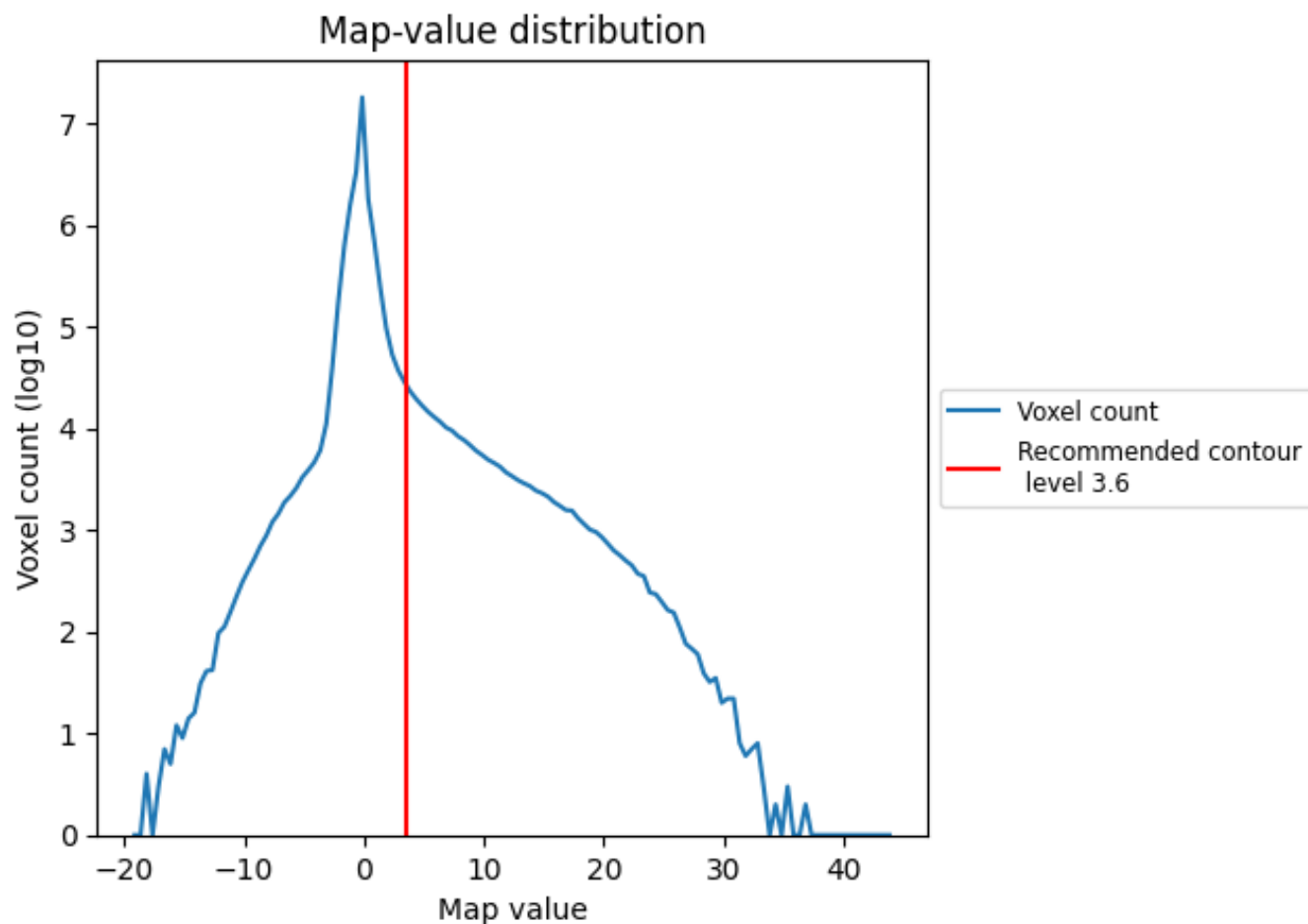
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

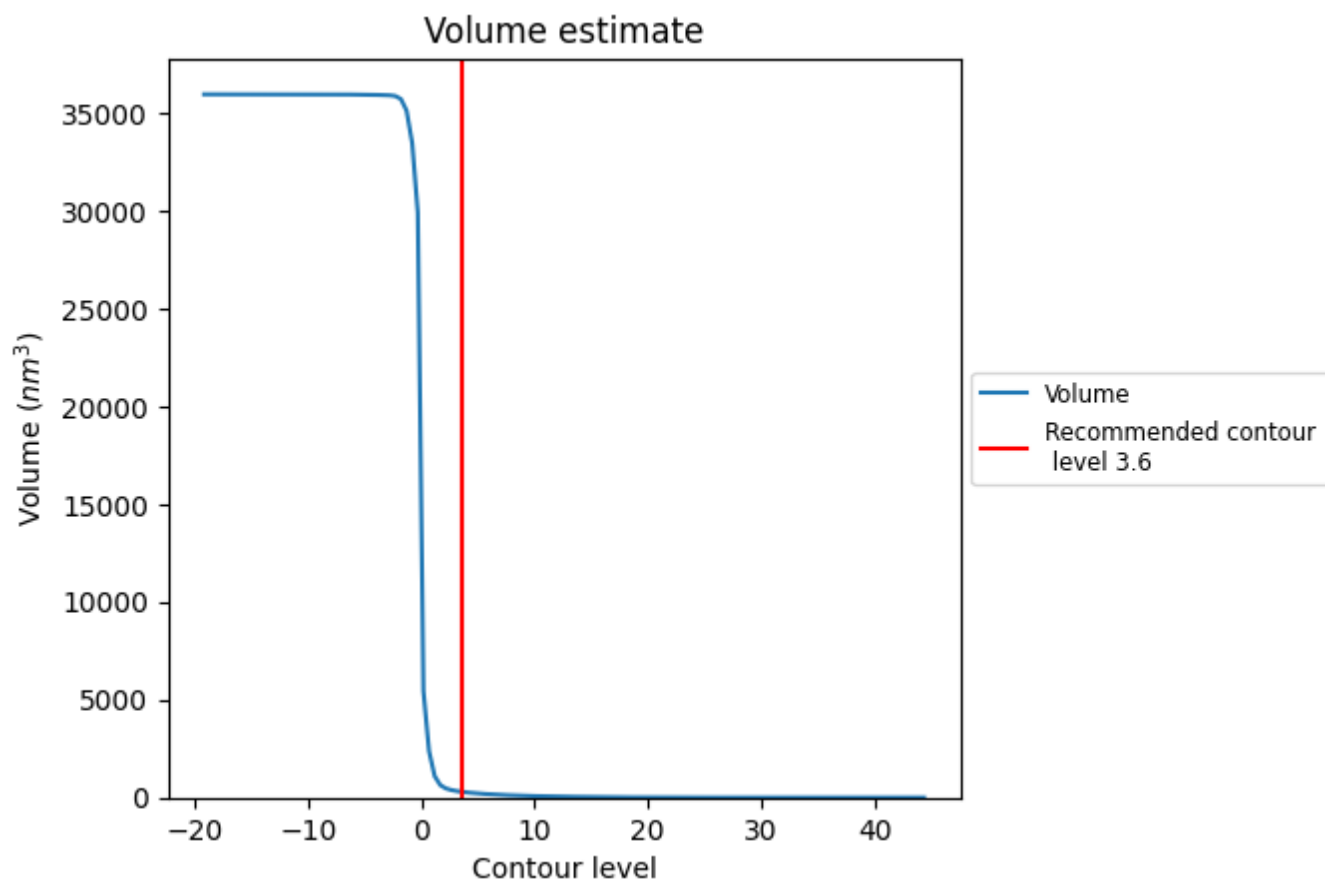
This section contains the results of statistical analysis of the map.

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.

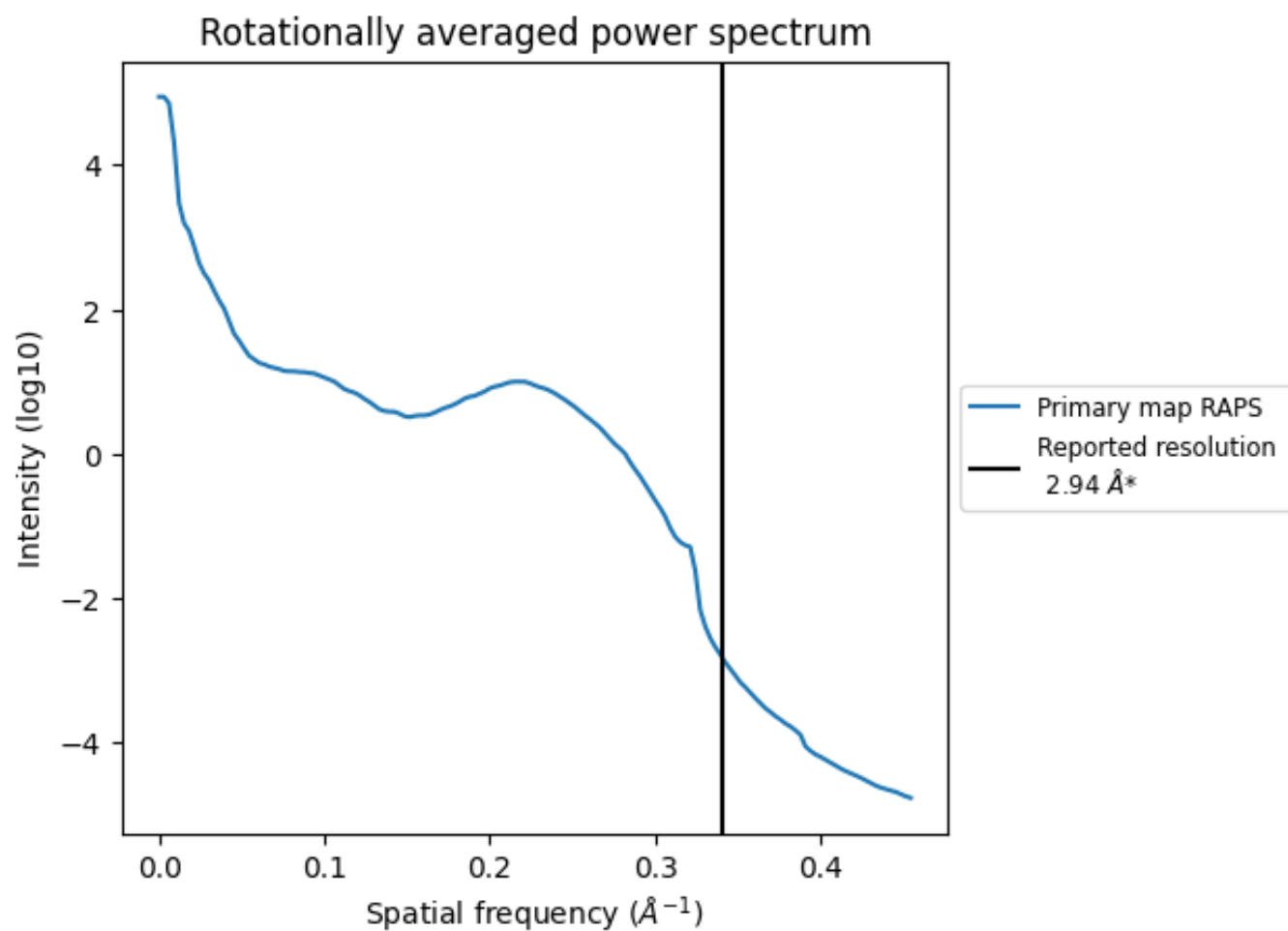
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 292 nm^3 ; this corresponds to an approximate mass of 264 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 0.340 Å⁻¹

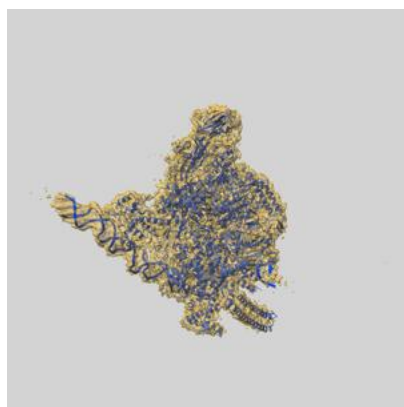
8 Fourier-Shell correlation ⓘ

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

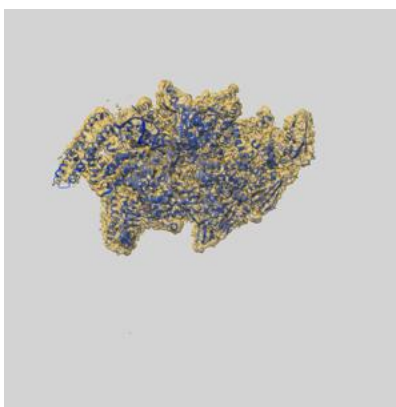
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-22886 and PDB model 7KIF. Per-residue inclusion information can be found in section [3](#) on page [7](#).

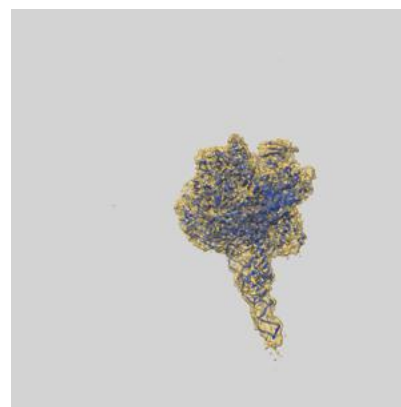
9.1 Map-model overlay [i](#)



X



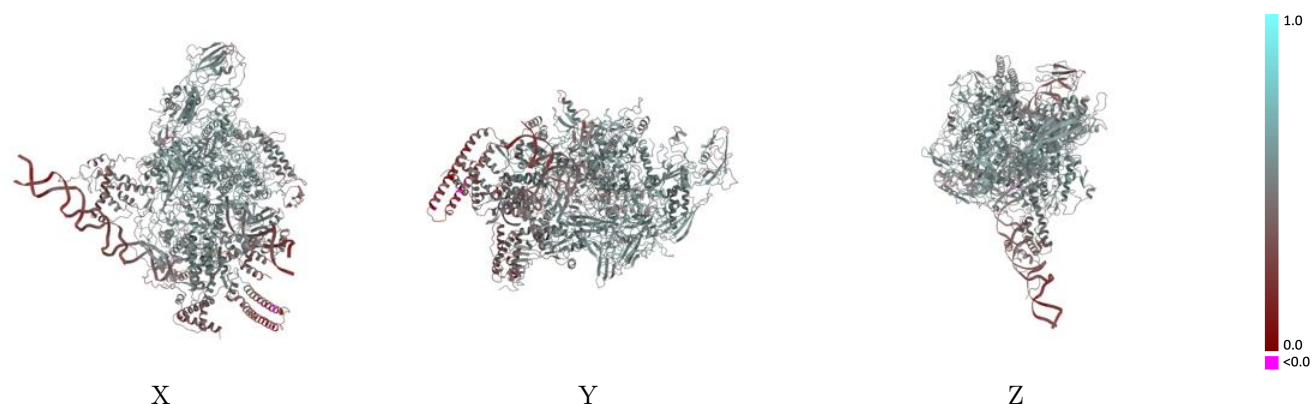
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 3.6 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](#)



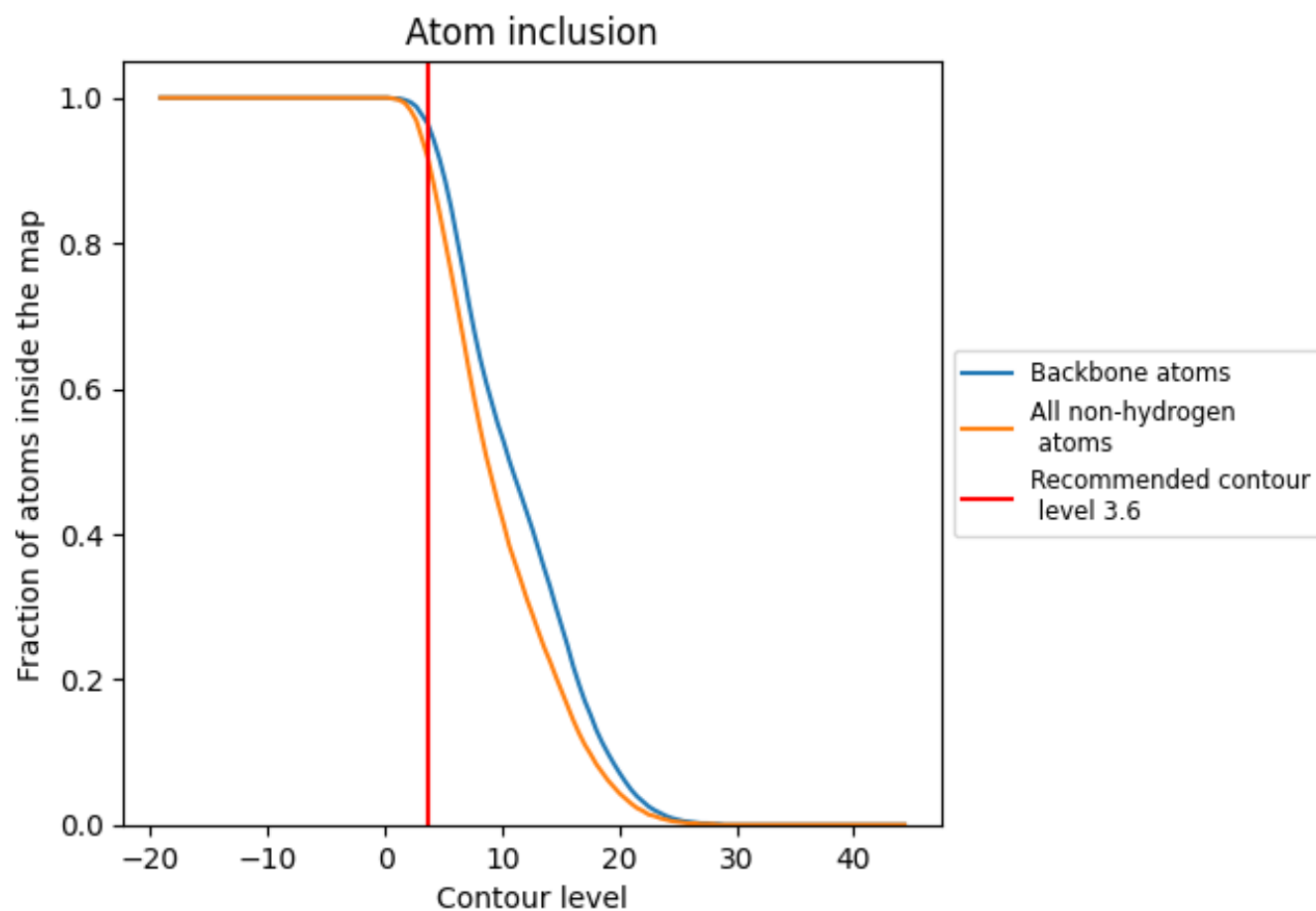
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (3.6).

9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 92% 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 (3.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9210	 0.4900
A	 0.9510	 0.5400
B	 0.9310	 0.5090
C	 0.9390	 0.5300
D	 0.9220	 0.5140
E	 0.9320	 0.5270
F	 0.8860	 0.4560
J	 0.8580	 0.4450
M	 0.8310	 0.3790
O	 0.9120	 0.3370
P	 0.9290	 0.3080
Z	 0.9440	 0.4060

