



Full wwPDB EM Validation Report ⓘ

Nov 8, 2022 – 02:29 PM JST

PDB ID : 7W5X
EMDB ID : EMD-32323
Title : Cryo-EM structure of SoxS-dependent transcription activation complex with zwf promoter DNA
Authors : Lin, W.; Feng, Y.; Shi, J.
Deposited on : 2021-11-30
Resolution : 3.40 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:

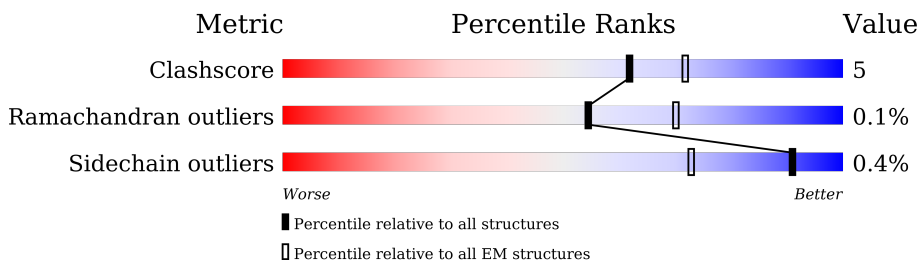
EMDB validation analysis : 0.0.1.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

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 3.40 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	1	75	
2	2	75	
3	C	1342	
4	D	1407	
5	E	91	
6	F	613	
7	K	107	
8	A	329	

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Mol	Chain	Length	Quality of chain
8	B	329	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a red segment on the left labeled '30%', a green segment labeled '79%', a yellow segment labeled '15%', and a grey segment on the far right labeled '6%'. The segments are stacked horizontally, with the red segment starting from the left and the grey segment ending on the right.</p>

2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 34265 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 DNA chain called zwf promoter DNA forward strand.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	1	75	1548	736	293	445	74	0	0

- Molecule 2 is a DNA chain called zwf promoter DNA reverse strand.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	2	75	1537	732	279	451	75	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	1340	10560	6627	1840	2050	43	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	516	VAL	ASP	variant	UNP P0A8V2

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	1334	10369	6514	1849	1956	50	0	0

- Molecule 5 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	79	627	382	118	126	1	0	0

- Molecule 6 is a protein called RNA polymerase sigma factor RpoD.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	F	495	3979	2487	707	763	22	0	0

- Molecule 7 is a protein called Regulatory protein SoxS.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	K	107	912	579	170	160	3	0	0

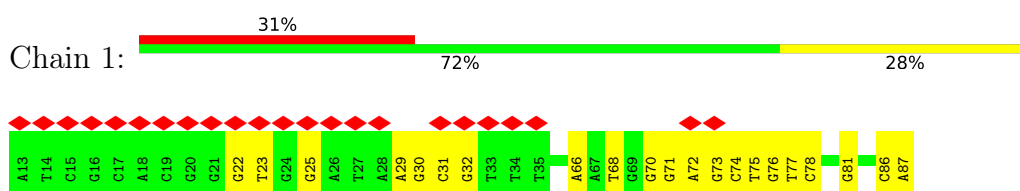
- Molecule 8 is a protein called DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	A	301	2334	1459	410	457	8	0	0
8	B	309	2399	1500	420	471	8	0	0

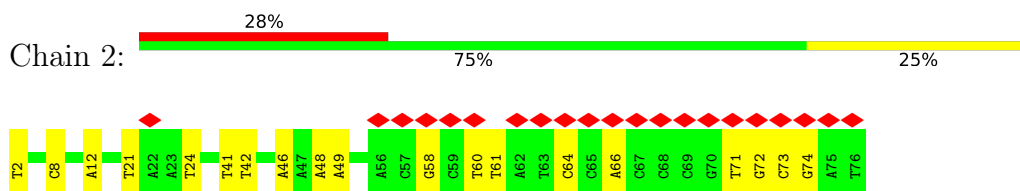
3 Residue-property plots [i](#)

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.

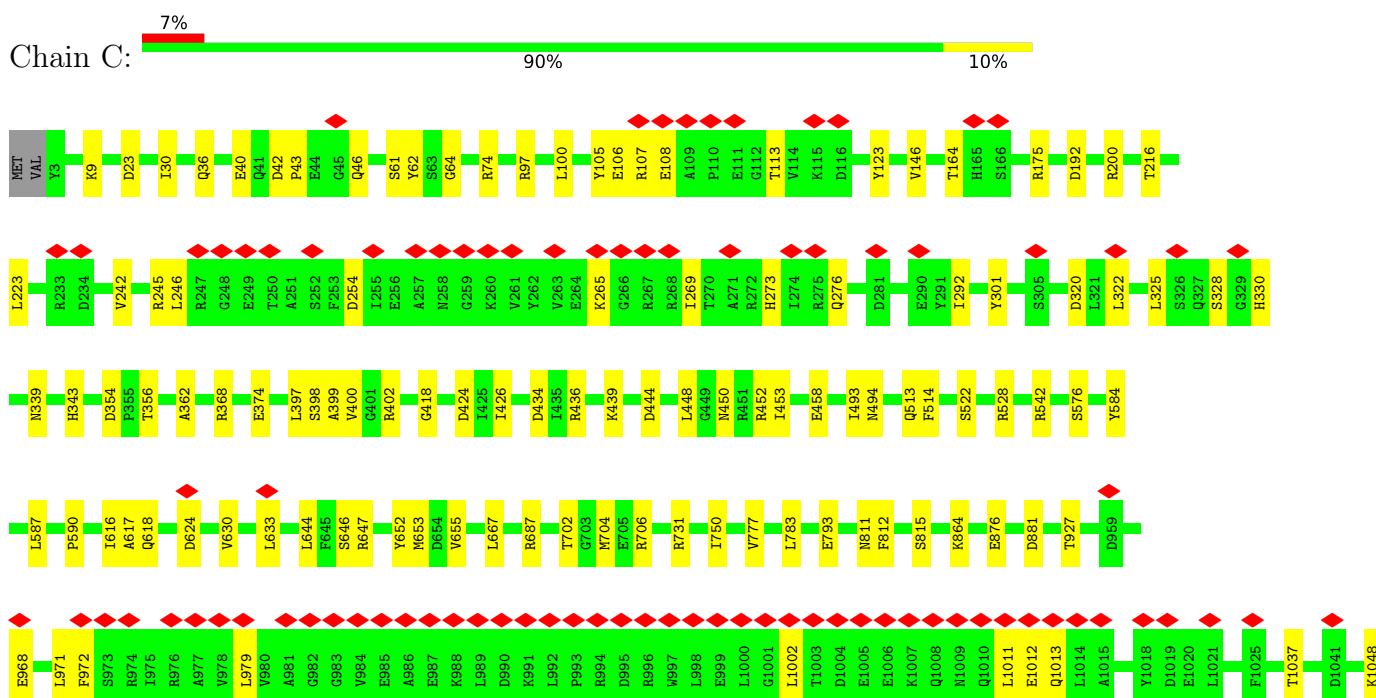
- Molecule 1: zwf promoter DNA forward strand

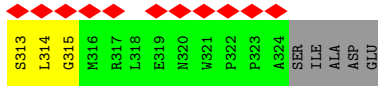


- Molecule 2: zwf promoter DNA reverse strand

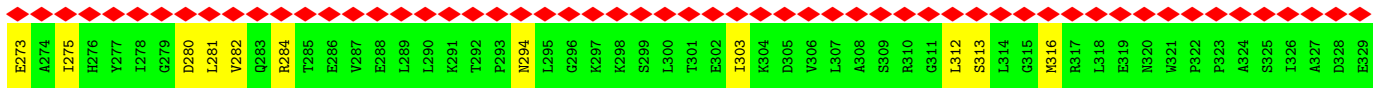
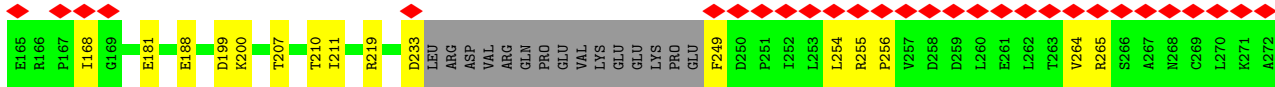
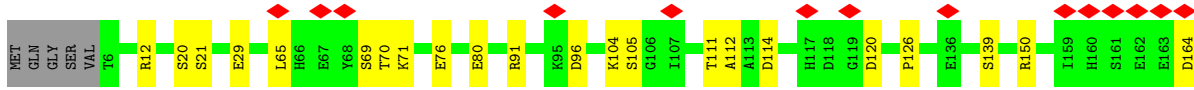
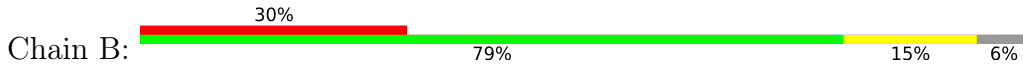


- Molecule 3: DNA-directed RNA polymerase subunit beta





• Molecule 8: DNA-directed RNA polymerase subunit alpha



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	116760	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	52	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.173	Depositor
Minimum map value	-0.093	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.02	Depositor
Map size (\AA)	264.0, 264.0, 264.0	wwPDB
Map dimensions	240, 240, 240	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.1, 1.1, 1.1	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	1	0.61	0/1740	0.91	0/2687
2	2	0.60	0/1723	0.95	0/2657
3	C	0.33	0/10728	0.51	0/14474
4	D	0.32	0/10521	0.53	0/14196
5	E	0.27	0/629	0.44	0/847
6	F	0.28	0/4029	0.51	0/5418
7	K	0.25	0/932	0.49	0/1256
8	A	0.30	0/2363	0.53	0/3205
8	B	0.27	0/2430	0.52	0/3293
All	All	0.35	0/35095	0.58	0/48033

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	1	1548	0	845	22	0
2	2	1537	0	846	15	0
3	C	10560	0	10573	81	0
4	D	10369	0	10590	107	0
5	E	627	0	634	4	0
6	F	3979	0	4019	37	0
7	K	912	0	917	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	2334	0	2375	20	0
8	B	2399	0	2443	29	0
All	All	34265	0	33242	298	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 5.

All (298) 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)
1:1:78:DC:H5'	3:C:542:ARG:O	1.72	0.90
1:1:70:DG:H2''	1:1:71:DG:C4'	2.15	0.77
1:1:70:DG:H2''	1:1:71:DG:O4'	1.87	0.73
4:D:1104:LYS:HB2	4:D:1124:ILE:HG23	1.72	0.71
1:1:78:DC:OP2	1:1:78:DC:H2'	1.91	0.70
8:A:7:GLU:HB2	8:B:150:ARG:HH12	1.62	0.64
1:1:66:DA:H8	6:F:429:THR:HG22	1.61	0.64
1:1:81:DG:H1	2:2:8:DC:H42	1.47	0.63
4:D:968:ASN:HD21	4:D:972:LYS:HB2	1.62	0.63
4:D:1028:ILE:HG22	4:D:1120:THR:HB	1.83	0.61
3:C:146:VAL:HG21	3:C:513:GLN:HE21	1.66	0.61
4:D:803:VAL:HG21	4:D:1309:ILE:HG22	1.81	0.61
4:D:952:VAL:HG12	4:D:1013:GLY:HA2	1.82	0.60
8:B:255:ARG:HH11	8:B:256:PRO:HD2	1.66	0.60
7:K:7:ILE:HG13	7:K:45:VAL:HG21	1.83	0.59
4:D:1108:GLN:HE21	4:D:1123:ARG:HH21	1.49	0.59
3:C:667:LEU:HA	3:C:702:THR:HG21	1.84	0.59
7:K:52:ASP:OD2	7:K:55:ARG:NH2	2.35	0.59
8:B:275:ILE:HG23	8:B:280:ASP:HB3	1.83	0.59
3:C:9:LYS:HG2	3:C:1171:ARG:HE	1.67	0.58
8:A:131:CYS:SG	8:A:132:HIS:N	2.76	0.58
8:A:252:ILE:HA	8:A:255:ARG:HE	1.67	0.58
8:B:313:SER:HB2	8:B:316:MET:HB2	1.85	0.58
3:C:1117:LEU:HD13	3:C:1195:ILE:HG12	1.85	0.58
7:K:71:ARG:HB3	8:A:287:VAL:HG22	1.86	0.57
3:C:292:ILE:HG21	3:C:322:LEU:HD21	1.86	0.57
4:D:799:ARG:O	4:D:803:VAL:HG23	2.05	0.57
4:D:1239:ASP:OD2	4:D:1242:ARG:NH2	2.36	0.57
6:F:583:THR:HG23	6:F:587:ILE:HG12	1.85	0.57
4:D:475:GLU:OE2	5:E:28:ARG:NH2	2.37	0.57
3:C:968:GLU:HA	3:C:971:LEU:HB2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1246:ARG:NH1	3:C:1265:PHE:O	2.37	0.57
3:C:452:ARG:NH2	3:C:458:GLU:OE2	2.36	0.57
6:F:278:ASP:HA	6:F:281:ARG:HB2	1.86	0.57
8:B:188:GLU:HB2	8:B:200:LYS:HG2	1.87	0.57
3:C:453:ILE:HD12	3:C:587:LEU:HD21	1.87	0.56
4:D:824:PRO:HB2	4:D:826:ILE:HG23	1.87	0.56
7:K:16:GLU:HB2	8:B:264:VAL:HB	1.87	0.56
1:1:72:DA:OP1	1:1:73:DG:N2	2.39	0.56
3:C:36:GLN:NE2	3:C:40:GLU:OE1	2.38	0.56
3:C:444:ASP:O	3:C:450:ASN:ND2	2.38	0.56
6:F:334:SER:O	6:F:338:HIS:NE2	2.38	0.56
1:1:23:DT:OP1	7:K:94:ARG:NH2	2.38	0.55
1:1:29:DA:H2'	1:1:30:DG:H8	1.71	0.55
1:1:68:DT:H3	6:F:383:ASN:HD22	1.53	0.55
6:F:586:ARG:HA	6:F:589:GLN:HB2	1.88	0.55
3:C:400:VAL:HG13	3:C:584:TYR:HB3	1.89	0.55
3:C:731:ARG:HD2	3:C:750:ILE:HD11	1.88	0.54
4:D:495:ASN:HD22	4:D:1247:LYS:HB3	1.72	0.54
4:D:674:THR:HG23	4:D:677:GLU:H	1.72	0.54
4:D:876:SER:HA	4:D:990:ARG:HH21	1.71	0.54
4:D:875:ASN:O	4:D:990:ARG:NH2	2.41	0.54
4:D:968:ASN:HD22	4:D:974:VAL:HG13	1.73	0.54
2:2:61:DT:O4	7:K:86:GLN:NE2	2.40	0.54
4:D:799:ARG:NH1	4:D:1146:GLU:OE2	2.41	0.54
3:C:633:LEU:HB3	3:C:644:LEU:HB3	1.90	0.54
3:C:1253:LEU:HA	6:F:525:ASP:HB2	1.89	0.54
4:D:69:GLU:HG2	4:D:76:LYS:HG2	1.88	0.54
4:D:955:LYS:HA	4:D:1012:ALA:HA	1.89	0.54
8:B:104:LYS:NZ	8:B:105:SER:O	2.40	0.53
1:1:71:DG:N2	3:C:374:GLU:OE1	2.40	0.53
3:C:617:ALA:HB3	3:C:653:MET:HG3	1.89	0.53
4:D:1172:LYS:HE3	4:D:1191:PRO:HG3	1.89	0.53
3:C:811:ASN:HA	3:C:815:SER:HB2	1.90	0.53
3:C:1090:ASN:O	8:A:182:ARG:NH1	2.41	0.53
5:E:44:ASP:OD2	5:E:52:ARG:NH2	2.41	0.53
4:D:983:LYS:NZ	4:D:994:SER:OG	2.42	0.53
3:C:74:ARG:HH22	3:C:97:ARG:HD2	1.72	0.53
4:D:1024:THR:HA	4:D:1125:PRO:HA	1.90	0.53
7:K:46:THR:O	7:K:47:HIS:ND1	2.41	0.53
6:F:407:GLU:HG2	6:F:442:SER:HB3	1.91	0.53
3:C:1192:GLU:OE1	4:D:764:ARG:NH1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:317:THR:HA	4:D:323:PRO:HA	1.92	0.52
3:C:702:THR:HG23	3:C:704:MET:H	1.74	0.52
4:D:885:VAL:HG11	4:D:1255:VAL:HG23	1.91	0.52
4:D:1261:LEU:HD21	4:D:1306:LEU:HD12	1.90	0.52
4:D:1215:GLU:HG3	4:D:1220:ILE:HD11	1.90	0.52
3:C:444:ASP:N	3:C:444:ASP:OD1	2.41	0.52
3:C:864:LYS:NZ	3:C:881:ASP:OD2	2.43	0.52
3:C:1142:ARG:NH1	3:C:1161:LEU:O	2.43	0.52
3:C:106:GLU:HG3	3:C:108:GLU:H	1.75	0.52
4:D:1173:ARG:HD3	4:D:1192:LYS:HD3	1.91	0.52
8:A:57:THR:HG22	8:A:58:GLU:HG3	1.90	0.52
4:D:850:LYS:NZ	4:D:855:ASP:O	2.43	0.52
3:C:23:ASP:OD1	3:C:23:ASP:N	2.42	0.52
4:D:851:PRO:HG3	4:D:876:SER:HB2	1.92	0.52
8:B:282:VAL:HG11	8:B:312:LEU:HD13	1.91	0.52
4:D:1211:SER:OG	4:D:1212:ASP:N	2.43	0.51
8:B:249:PHE:HE2	8:B:254:LEU:HB2	1.75	0.51
1:I:77:DT:O4	3:C:175:ARG:NE	2.39	0.51
3:C:1058:ARG:NH1	3:C:1240:ASP:OD2	2.44	0.51
3:C:368:ARG:NH2	6:F:90:GLU:O	2.44	0.51
7:K:20:GLN:O	7:K:58:ARG:NH2	2.43	0.51
3:C:30:ILE:HD11	3:C:528:ARG:HA	1.93	0.51
4:D:309:ASN:HB2	4:D:326:SER:HB3	1.91	0.51
4:D:1206:ARG:NH2	4:D:1223:LEU:O	2.44	0.51
7:K:10:LEU:HA	7:K:13:TRP:HB2	1.93	0.51
8:A:230:ALA:HA	8:A:234:LEU:HB2	1.93	0.51
8:B:29:GLU:HB3	8:B:200:LYS:HB2	1.93	0.51
8:B:164:ASP:N	8:B:164:ASP:OD1	2.44	0.51
3:C:61:SER:OG	3:C:62:TYR:N	2.44	0.51
3:C:105:TYR:HD1	3:C:113:THR:HA	1.75	0.51
3:C:1274:GLU:HG2	4:D:424:ASN:HD21	1.75	0.50
3:C:254:ASP:HB2	3:C:265:LYS:HE2	1.93	0.50
4:D:606:ASN:OD1	4:D:610:ARG:NH1	2.44	0.50
2:2:60:DT:OP2	7:K:85:GLN:NE2	2.45	0.50
4:D:1034:PHE:HA	4:D:1114:GLN:HA	1.93	0.50
4:D:202:ARG:NH2	4:D:225:GLU:OE2	2.45	0.50
7:K:74:PHE:HA	7:K:85:GLN:HB3	1.94	0.50
8:B:199:ASP:OD1	8:B:199:ASP:N	2.44	0.50
3:C:397:LEU:HB2	3:C:418:GLY:HA2	1.93	0.49
6:F:122:ARG:NH2	6:F:378:GLU:OE1	2.41	0.49
4:D:417:ARG:HG2	4:D:418:GLU:HG2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:285:THR:HG22	8:A:315:GLY:HA2	1.94	0.49
6:F:479:THR:HG22	6:F:481:GLU:H	1.78	0.49
4:D:818:GLU:HA	4:D:845:ALA:HB1	1.94	0.49
8:B:20:SER:OG	8:B:21:SER:N	2.46	0.49
3:C:528:ARG:NH2	3:C:576:SER:O	2.38	0.49
4:D:363:LEU:HD23	4:D:618:VAL:HG13	1.93	0.49
4:D:514:THR:HG21	4:D:596:LEU:HD12	1.95	0.49
2:2:21:DT:O4	4:D:320:ASN:ND2	2.46	0.48
3:C:246:LEU:HD22	3:C:269:ILE:HD13	1.95	0.48
4:D:839:VAL:HG12	4:D:864:LEU:HD22	1.95	0.48
6:F:154:GLU:OE1	6:F:157:ARG:NH2	2.44	0.48
4:D:948:SER:OG	4:D:949:SER:N	2.46	0.48
3:C:192:ASP:OD2	3:C:436:ARG:NH2	2.46	0.48
3:C:876:GLU:OE1	3:C:927:THR:OG1	2.30	0.48
4:D:1175:LEU:HG	4:D:1190:ILE:HG21	1.95	0.48
4:D:110:PRO:HB3	4:D:240:THR:HG22	1.96	0.48
4:D:1221:LEU:HD22	4:D:1306:LEU:HB2	1.96	0.48
8:B:96:ASP:OD1	8:B:96:ASP:N	2.46	0.48
4:D:214:ARG:NH1	4:D:218:THR:OG1	2.47	0.48
4:D:826:ILE:HG22	4:D:831:VAL:HG22	1.95	0.48
2:2:48:DA:H2''	2:2:49:DA:H5''	1.96	0.47
4:D:586:GLY:HA3	4:D:612:LEU:HD11	1.95	0.47
6:F:142:THR:HA	6:F:145:LEU:HB2	1.97	0.47
6:F:216:LEU:HA	6:F:219:GLU:HB3	1.96	0.47
4:D:970:SER:HB2	4:D:972:LYS:HE2	1.96	0.47
4:D:1028:ILE:O	4:D:1099:TYR:OH	2.31	0.47
8:A:255:ARG:HB2	8:A:278:ILE:HD12	1.97	0.47
3:C:522:SER:OG	3:C:687:ARG:O	2.32	0.47
3:C:646:SER:OG	3:C:647:ARG:N	2.47	0.47
4:D:141:PHE:HA	4:D:180:MET:HG2	1.97	0.47
4:D:1046:ILE:HD13	4:D:1059:LEU:HD13	1.97	0.47
8:B:69:SER:OG	8:B:70:THR:N	2.48	0.47
4:D:679:TYR:OH	4:D:754:ILE:O	2.33	0.47
1:1:70:DG:O6	6:F:99:ARG:HG2	2.15	0.47
6:F:606:VAL:O	6:F:609:SER:OG	2.31	0.47
6:F:561:MET:HB3	6:F:571:TYR:HD2	1.80	0.46
3:C:402:ARG:NH1	3:C:424:ASP:OD2	2.42	0.46
6:F:267:ASP:O	6:F:271:ASN:ND2	2.48	0.46
7:K:7:ILE:HA	7:K:10:LEU:HB2	1.98	0.46
8:B:265:ARG:NH2	8:B:294:ASN:O	2.48	0.46
3:C:320:ASP:N	3:C:320:ASP:OD1	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:1347:LEU:HD22	4:D:1357:ILE:HG23	1.97	0.46
8:A:269:CYS:SG	8:A:292:THR:OG1	2.67	0.46
8:B:120:ASP:OD1	8:B:120:ASP:N	2.43	0.46
4:D:1028:ILE:HA	4:D:1120:THR:HA	1.97	0.46
4:D:1034:PHE:O	4:D:1081:VAL:N	2.41	0.46
6:F:586:ARG:O	6:F:590:ILE:N	2.44	0.46
7:K:77:ALA:HB1	7:K:82:TYR:HB2	1.98	0.46
3:C:1121:ALA:HB2	3:C:1182:ILE:HD12	1.98	0.46
1:1:86:DC:H2''	1:1:87:DA:H5'	1.98	0.46
3:C:1223:ARG:NH2	4:D:721:SER:OG	2.49	0.46
1:1:70:DG:H2''	1:1:71:DG:H4'	1.95	0.46
3:C:223:LEU:HD13	3:C:426:ILE:HG21	1.98	0.46
6:F:268:TYR:HA	6:F:271:ASN:HD22	1.81	0.45
3:C:434:ASP:HB3	3:C:439:LYS:HG3	1.98	0.45
3:C:1012:GLU:OE1	3:C:1013:GLN:NE2	2.49	0.45
4:D:803:VAL:HG21	4:D:1309:ILE:CG2	2.45	0.45
4:D:1154:ALA:N	4:D:1214:PRO:O	2.46	0.45
1:1:75:DT:H2''	1:1:76:DG:H5''	1.98	0.45
3:C:242:VAL:HB	3:C:245:ARG:HG3	1.97	0.45
3:C:812:PHE:O	4:D:504:GLN:NE2	2.49	0.45
4:D:968:ASN:HD21	4:D:972:LYS:HE3	1.82	0.45
5:E:25:ARG:NH2	5:E:68:GLU:OE1	2.42	0.45
2:2:24:DT:OP1	3:C:494:ASN:ND2	2.49	0.45
3:C:624:ASP:HB2	3:C:630:VAL:HG12	1.97	0.45
7:K:46:THR:HG23	7:K:48:GLN:HG2	1.98	0.45
7:K:65:GLU:O	7:K:69:THR:OG1	2.26	0.45
3:C:301:TYR:HB3	3:C:325:LEU:HD11	1.99	0.45
3:C:400:VAL:HG11	3:C:452:ARG:HD2	1.99	0.45
3:C:64:GLY:O	3:C:107:ARG:NH1	2.50	0.45
4:D:806:ASP:HB3	4:D:1346:GLY:HA2	1.98	0.45
4:D:1061:VAL:HB	4:D:1105:ALA:HB3	1.99	0.45
1:1:25:DG:H1	2:2:64:DC:H42	1.63	0.45
2:2:58:DG:H5'	8:A:290:LEU:HD22	1.99	0.45
3:C:328:SER:HB2	3:C:330:HIS:HD2	1.82	0.45
3:C:1341:ASP:HA	4:D:17:PHE:HA	1.99	0.45
8:B:281:LEU:HD21	8:B:303:ILE:HG23	1.98	0.45
3:C:398:SER:O	3:C:400:VAL:N	2.50	0.45
4:D:1000:GLY:HA3	4:D:1026:PRO:HG2	1.99	0.45
7:K:58:ARG:HG2	7:K:80:LEU:HD22	1.99	0.45
3:C:354:ASP:OD1	3:C:356:THR:OG1	2.32	0.44
3:C:200:ARG:HD3	3:C:200:ARG:HA	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:1021:ASP:HB3	4:D:1024:THR:HB	1.99	0.44
6:F:280:VAL:O	6:F:284:GLU:N	2.45	0.44
8:A:161:SER:OG	8:A:162:GLU:N	2.48	0.44
4:D:126:LEU:HD23	4:D:223:LEU:HD22	1.98	0.44
4:D:654:ILE:O	4:D:658:GLU:N	2.47	0.44
7:K:64:VAL:O	7:K:68:THR:OG1	2.33	0.44
1:1:29:DA:H2'	1:1:30:DG:C8	2.51	0.44
4:D:884:SER:OG	4:D:885:VAL:N	2.51	0.44
3:C:100:LEU:HD12	3:C:493:ILE:HD11	1.98	0.44
4:D:147:ILE:N	4:D:177:ASP:O	2.48	0.44
2:2:73:DC:H2'	2:2:74:DG:H8	1.83	0.44
4:D:256:ASP:OD1	4:D:256:ASP:N	2.51	0.44
3:C:1271:GLY:N	3:C:1274:GLU:OE1	2.50	0.44
6:F:263:PRO:HA	6:F:266:PHE:HB3	2.00	0.44
6:F:583:THR:O	6:F:584:ARG:NH1	2.46	0.44
2:2:41:DT:H2''	2:2:42:DT:H5''	1.99	0.43
4:D:635:SER:OG	4:D:636:GLY:N	2.51	0.43
4:D:1079:LYS:HE2	4:D:1081:VAL:HG12	1.99	0.43
4:D:1191:PRO:HB2	4:D:1194:ARG:HD3	1.99	0.43
3:C:42:ASP:OD2	3:C:42:ASP:N	2.52	0.43
4:D:91:GLU:OE1	4:D:101:ARG:NH2	2.40	0.43
4:D:438:GLU:HA	4:D:439:PRO:HD3	1.91	0.43
4:D:1038:THR:HG21	4:D:1079:LYS:HB3	2.01	0.43
2:2:46:DA:H5'	6:F:588:ARG:HD2	2.00	0.43
3:C:706:ARG:NH2	3:C:793:GLU:OE2	2.52	0.43
4:D:1109:LEU:HD22	4:D:1113:VAL:HG21	2.00	0.43
6:F:273:MET:HB3	6:F:362:ASN:HD21	1.83	0.43
6:F:233:ASP:HA	6:F:236:LYS:HB2	2.01	0.43
6:F:426:LYS:O	6:F:429:THR:OG1	2.34	0.43
1:1:22:DG:O6	2:2:66:DA:N6	2.52	0.43
4:D:1357:ILE:HD13	4:D:1357:ILE:HA	1.86	0.43
8:A:263:THR:O	8:A:266:SER:OG	2.32	0.43
8:A:286:GLU:HB2	8:A:314:LEU:HD13	2.01	0.43
3:C:448:LEU:HD23	3:C:448:LEU:HA	1.88	0.42
3:C:1048:LYS:HB3	3:C:1048:LYS:HE2	1.84	0.42
4:D:972:LYS:HB2	4:D:972:LYS:HE3	1.89	0.42
6:F:402:LEU:HD23	6:F:405:ILE:HD12	2.02	0.42
4:D:147:ILE:HB	4:D:177:ASP:HB3	2.01	0.42
4:D:1199:PHE:H	4:D:1202:GLU:HG3	1.84	0.42
8:A:120:ASP:OD1	8:A:120:ASP:N	2.50	0.42
3:C:43:PRO:O	3:C:46:GLN:NE2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:618:GLN:HE21	4:D:770:LEU:HD22	1.84	0.42
4:D:294:ASN:HD22	6:F:406:GLN:HE21	1.68	0.42
4:D:816:THR:HG22	4:D:883:ARG:HH21	1.84	0.42
6:F:165:PHE:HB3	6:F:166:VAL:H	1.70	0.42
2:2:2:DT:H6	2:2:2:DT:H2'	1.68	0.42
3:C:1002:LEU:HD23	3:C:1002:LEU:HA	1.88	0.42
4:D:1307:LEU:HB2	4:D:1312:ALA:HB2	2.00	0.42
3:C:123:TYR:HE2	6:F:471:LEU:HD22	1.84	0.42
4:D:847:ASP:N	4:D:847:ASP:OD1	2.52	0.42
2:2:71:DT:H2'	2:2:72:DG:C8	2.55	0.42
3:C:616:ILE:HG12	3:C:652:TYR:HB2	2.02	0.42
3:C:1293:VAL:HG11	3:C:1315:MET:HG3	2.01	0.42
8:B:211:ILE:HD11	8:B:219:ARG:HD2	2.00	0.42
8:B:273:GLU:O	8:B:284:ARG:NH2	2.53	0.42
3:C:979:LEU:HD21	3:C:1011:LEU:HD21	2.02	0.42
8:B:71:LYS:NZ	8:B:139:SER:O	2.52	0.42
8:B:111:THR:OG1	8:B:126:PRO:O	2.34	0.42
8:B:112:ALA:HB3	8:B:126:PRO:HA	2.01	0.42
3:C:273:HIS:HA	3:C:276:GLN:HB2	2.01	0.42
6:F:574:GLU:O	6:F:578:LYS:N	2.53	0.42
8:B:65:LEU:HD12	8:B:168:ILE:HD11	2.02	0.42
1:1:68:DT:H1'	6:F:385:ARG:HB2	2.02	0.41
6:F:151:VAL:HB	6:F:157:ARG:HG2	2.02	0.41
6:F:559:LEU:HD23	6:F:559:LEU:HA	1.91	0.41
7:K:29:LYS:HD2	7:K:29:LYS:HA	1.91	0.41
4:D:153:ASN:HD22	4:D:172:PHE:HZ	1.67	0.41
4:D:473:THR:OG1	4:D:474:LEU:N	2.53	0.41
4:D:1076:PRO:HG2	4:D:1101:LEU:HB2	2.01	0.41
8:B:91:ARG:HB2	8:B:210:THR:HG22	2.01	0.41
3:C:339:ASN:HB3	3:C:343:HIS:H	1.85	0.41
4:D:1107:VAL:HA	4:D:1122:ALA:HA	2.03	0.41
8:A:307:LEU:HD13	8:A:313:SER:HA	2.03	0.41
3:C:777:VAL:HG11	3:C:783:LEU:HD21	2.02	0.41
4:D:1164:SER:N	4:D:1176:VAL:O	2.43	0.41
8:B:76:GLU:HB3	8:B:80:GLU:HB2	2.01	0.41
4:D:1149:ARG:HD2	4:D:1150:PRO:HD2	2.03	0.41
8:A:218:ARG:NH1	8:B:233:ASP:O	2.53	0.41
1:1:74:DC:OP1	4:D:321:LYS:NZ	2.54	0.41
4:D:222:LYS:HA	4:D:222:LYS:HD2	1.87	0.41
4:D:795:TYR:OH	4:D:1326:GLN:NE2	2.45	0.41
4:D:858:VAL:HG13	4:D:860:ARG:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:K:5:LYS:HB2	7:K:5:LYS:HE3	1.82	0.41
1:1:31:DC:H2'	1:1:32:DG:C8	2.55	0.41
2:2:12:DA:OP1	4:D:334:LYS:NZ	2.42	0.41
4:D:301:GLU:HA	4:D:312:ARG:HH21	1.86	0.41
4:D:566:LYS:HE3	4:D:566:LYS:HB2	1.89	0.41
4:D:1036:ARG:HG3	4:D:1081:VAL:HG21	2.02	0.41
6:F:584:ARG:HA	6:F:584:ARG:HD3	1.73	0.41
8:A:100:LEU:HD13	8:A:115:ILE:HG21	2.02	0.41
8:B:181:GLU:H	8:B:207:THR:HA	1.85	0.41
3:C:590:PRO:HB2	3:C:655:VAL:HG21	2.03	0.41
3:C:1150:ASP:N	3:C:1150:ASP:OD1	2.54	0.41
3:C:356:THR:HG21	3:C:362:ALA:HA	2.02	0.40
4:D:986:ASP:OD1	4:D:986:ASP:N	2.42	0.40
8:A:45:ARG:O	8:A:49:SER:OG	2.37	0.40
8:A:142:MET:HE3	8:A:142:MET:HB3	1.94	0.40
4:D:768:ASN:H	4:D:771:GLN:HB2	1.86	0.40
4:D:1048:ARG:HA	4:D:1048:ARG:HD2	1.91	0.40
4:D:479:GLU:HG3	5:E:20:VAL:HG11	2.02	0.40
4:D:1115:ILE:HB	4:D:1119:ASP:HB3	2.03	0.40
8:B:114:ASP:N	8:B:114:ASP:OD1	2.54	0.40
6:F:315:TRP:HZ2	6:F:341:LEU:HD11	1.85	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	1338/1342 (100%)	1269 (95%)	68 (5%)	1 (0%)	51	82
4	D	1322/1407 (94%)	1238 (94%)	84 (6%)	0	100	100
5	E	77/91 (85%)	77 (100%)	0	0	100	100
6	F	487/613 (79%)	450 (92%)	36 (7%)	1 (0%)	47	78

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	K	105/107 (98%)	99 (94%)	6 (6%)	0	100	100
8	A	295/329 (90%)	284 (96%)	11 (4%)	0	100	100
8	B	305/329 (93%)	284 (93%)	21 (7%)	0	100	100
All	All	3929/4218 (93%)	3701 (94%)	226 (6%)	2 (0%)	54	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	399	ALA
6	F	505	ILE

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	1152/1157 (100%)	1146 (100%)	6 (0%)	88	94
4	D	1115/1168 (96%)	1111 (100%)	4 (0%)	91	95
5	E	67/75 (89%)	67 (100%)	0	100	100
6	F	431/540 (80%)	430 (100%)	1 (0%)	93	98
7	K	99/99 (100%)	97 (98%)	2 (2%)	55	77
8	A	260/286 (91%)	260 (100%)	0	100	100
8	B	267/286 (93%)	266 (100%)	1 (0%)	91	95
All	All	3391/3611 (94%)	3377 (100%)	14 (0%)	91	95

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	164	THR
3	C	216	THR
3	C	514	PHE
3	C	972	PHE
3	C	1037	THR

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Mol	Chain	Res	Type
3	C	1149	TYR
4	D	709	ARG
4	D	744	ARG
4	D	1123	ARG
4	D	1224	ARG
6	F	509	THR
7	K	43	ARG
7	K	49	THR
8	B	12	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	120	GLN
3	C	276	GLN
3	C	330	HIS
3	C	447	HIS
3	C	513	GLN
3	C	622	ASN
3	C	952	GLN
3	C	965	GLN
3	C	1136	GLN
3	C	1257	GLN
4	D	294	ASN
4	D	424	ASN
4	D	477	GLN
4	D	560	ASN
4	D	1084	GLN
4	D	1108	GLN
4	D	1227	HIS
4	D	1235	ASN
4	D	1238	GLN
4	D	1326	GLN
6	F	271	ASN
6	F	362	ASN
6	F	383	ASN
6	F	461	ASN
7	K	20	GLN
7	K	86	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.

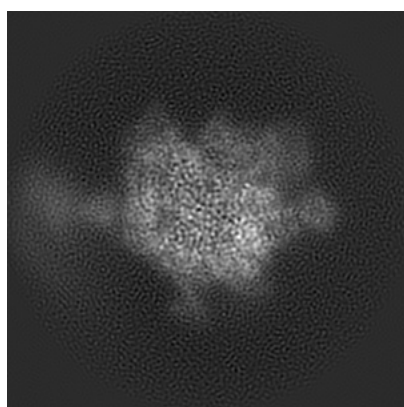
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32323. These allow visual inspection of the internal detail of the map and identification of artifacts.

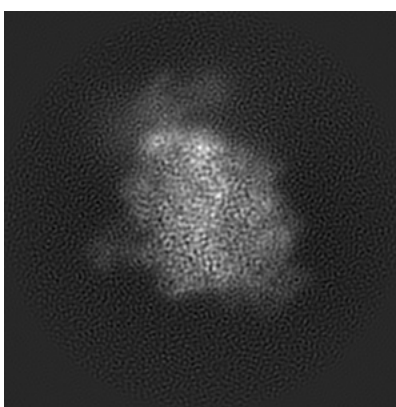
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

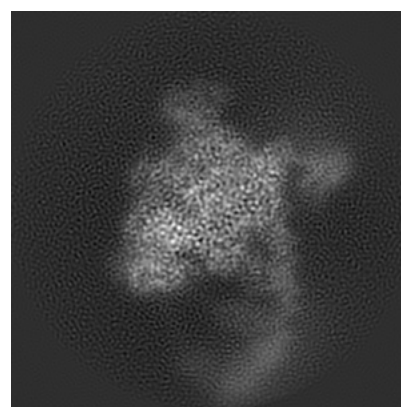
6.1.1 Primary map



X



Y

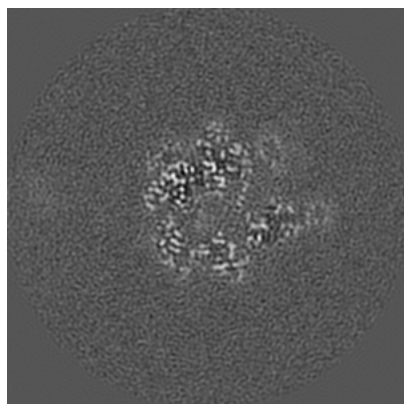


Z

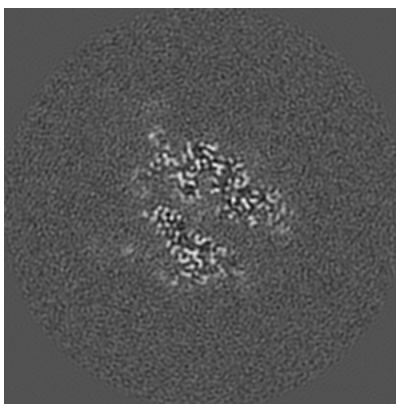
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

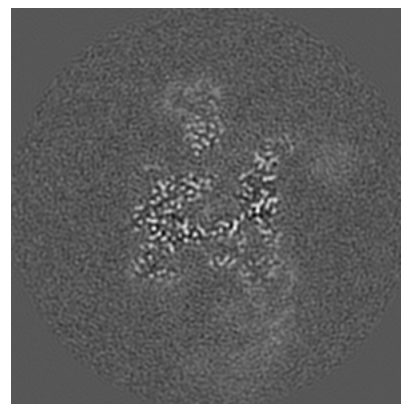
6.2.1 Primary map



X Index: 120



Y Index: 120

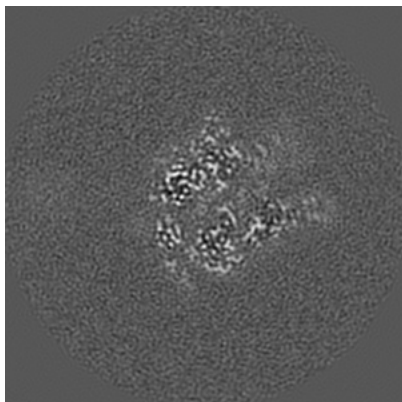


Z Index: 120

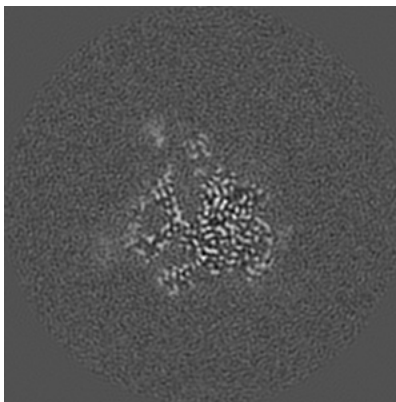
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

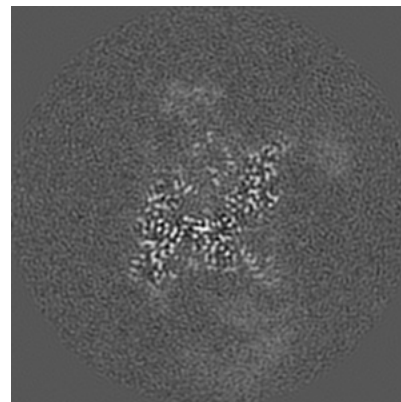
6.3.1 Primary map



X Index: 116



Y Index: 104

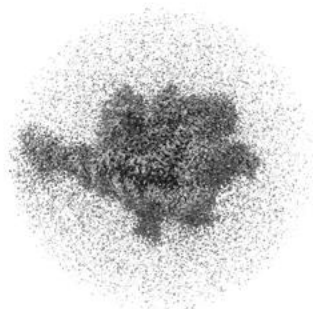


Z Index: 125

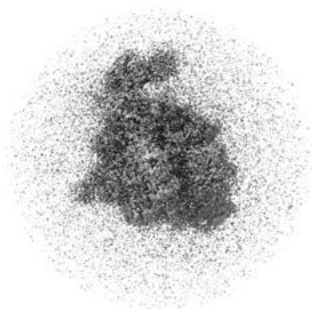
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

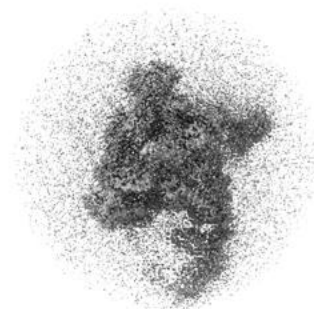
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.02. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

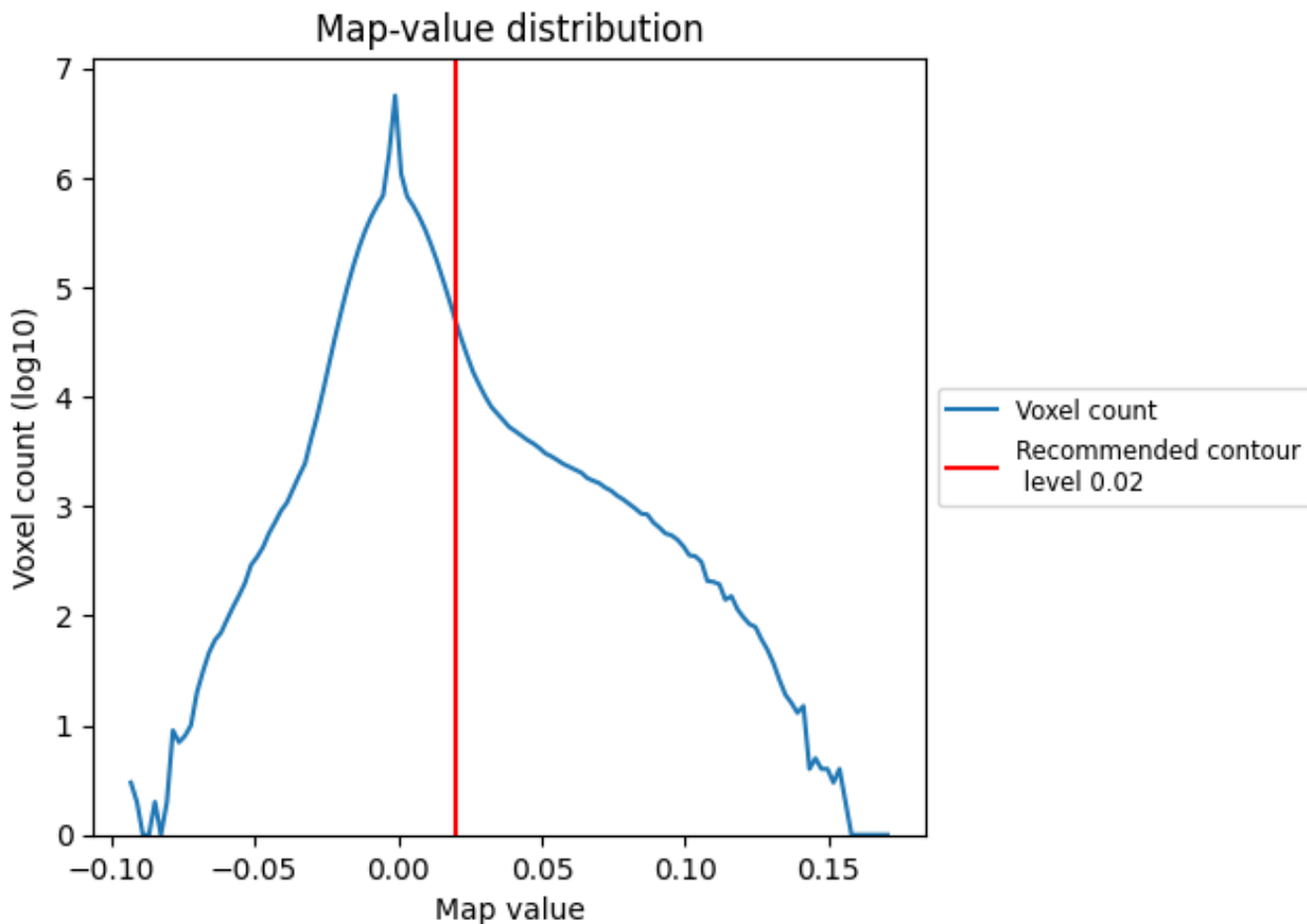
6.5 Mask visualisation

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

7 Map analysis [i](#)

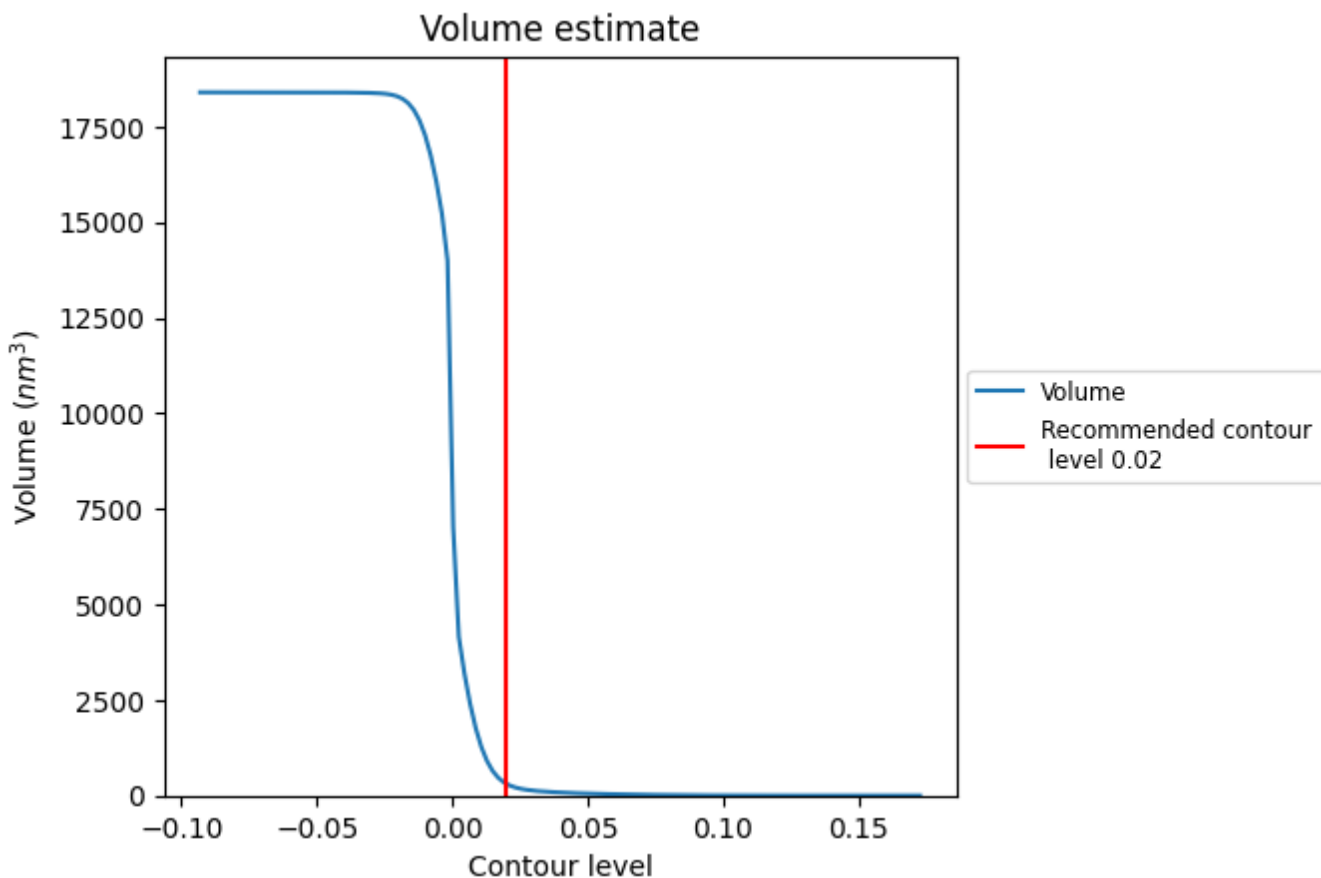
This section contains the results of statistical analysis of the map.

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.

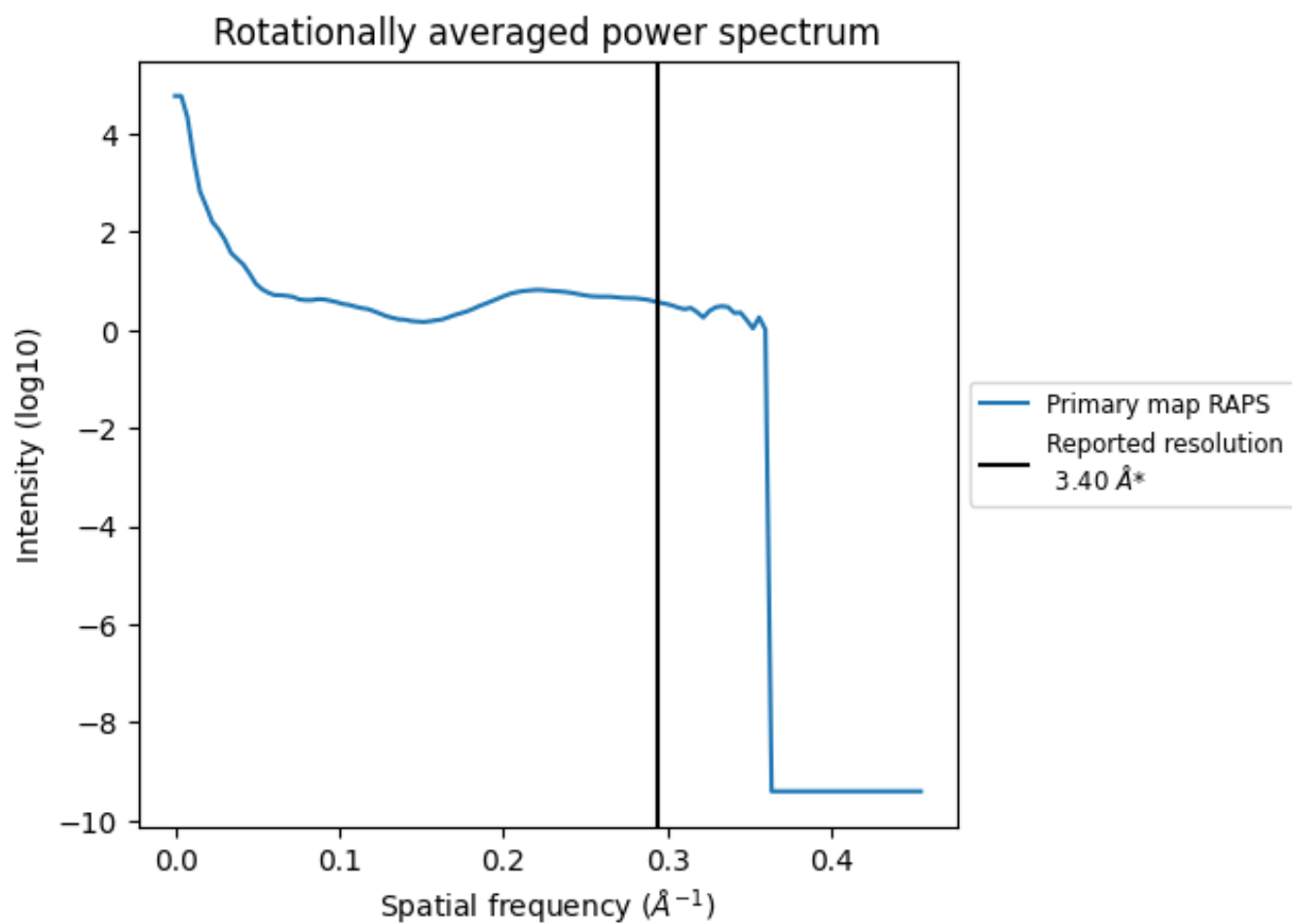
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 308 nm^3 ; this corresponds to an approximate mass of 279 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 [\(i\)](#)



*Reported resolution corresponds to spatial frequency of 0.294\AA^{-1}

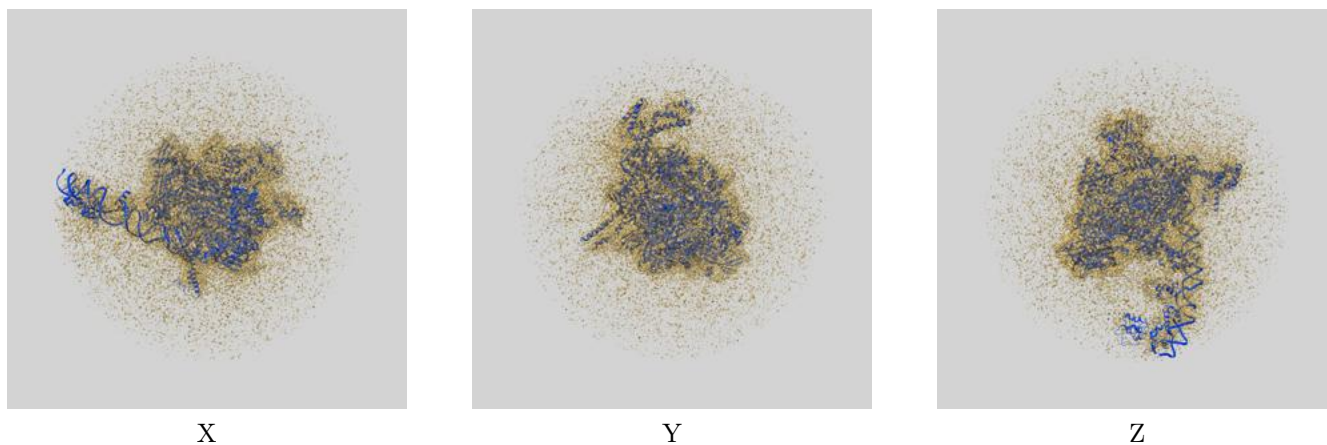
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

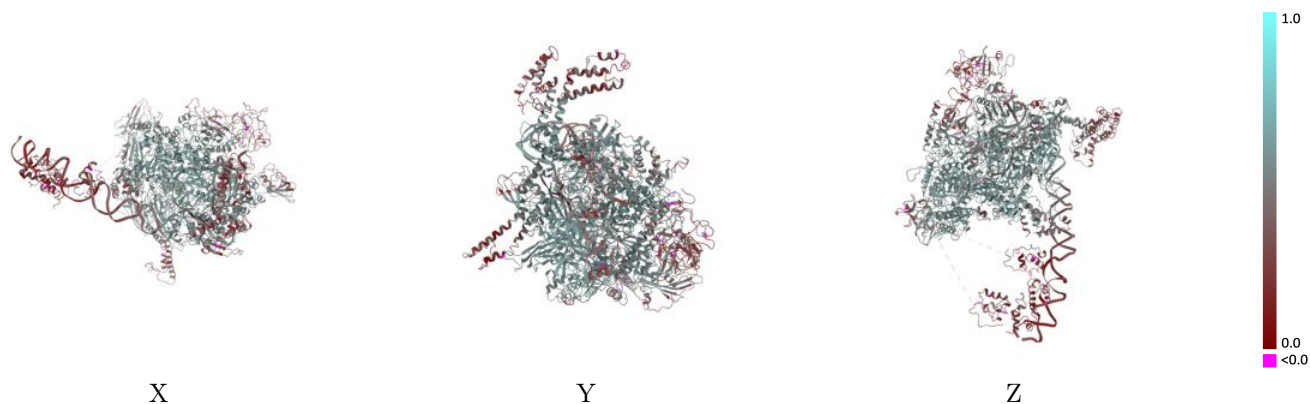
This section contains information regarding the fit between EMDB map EMD-32323 and PDB model 7W5X. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



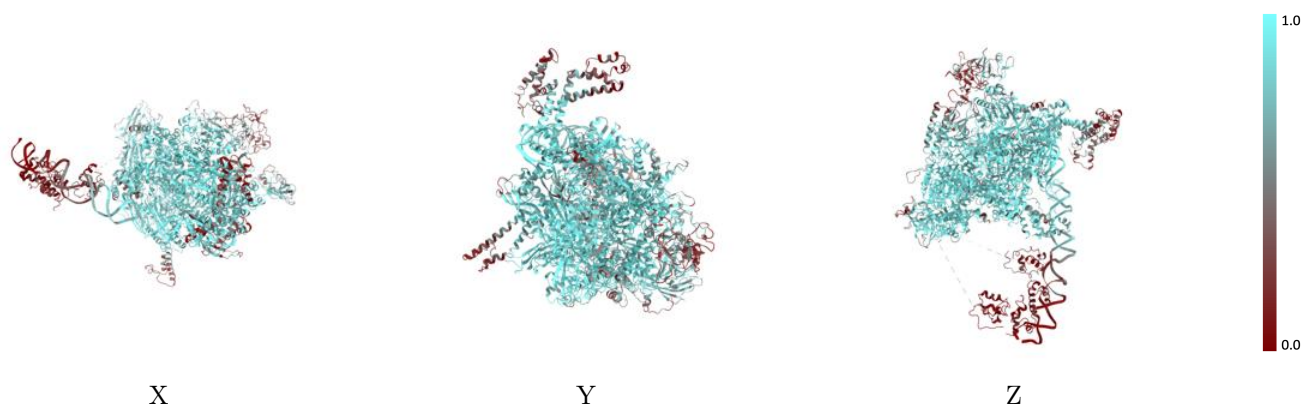
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



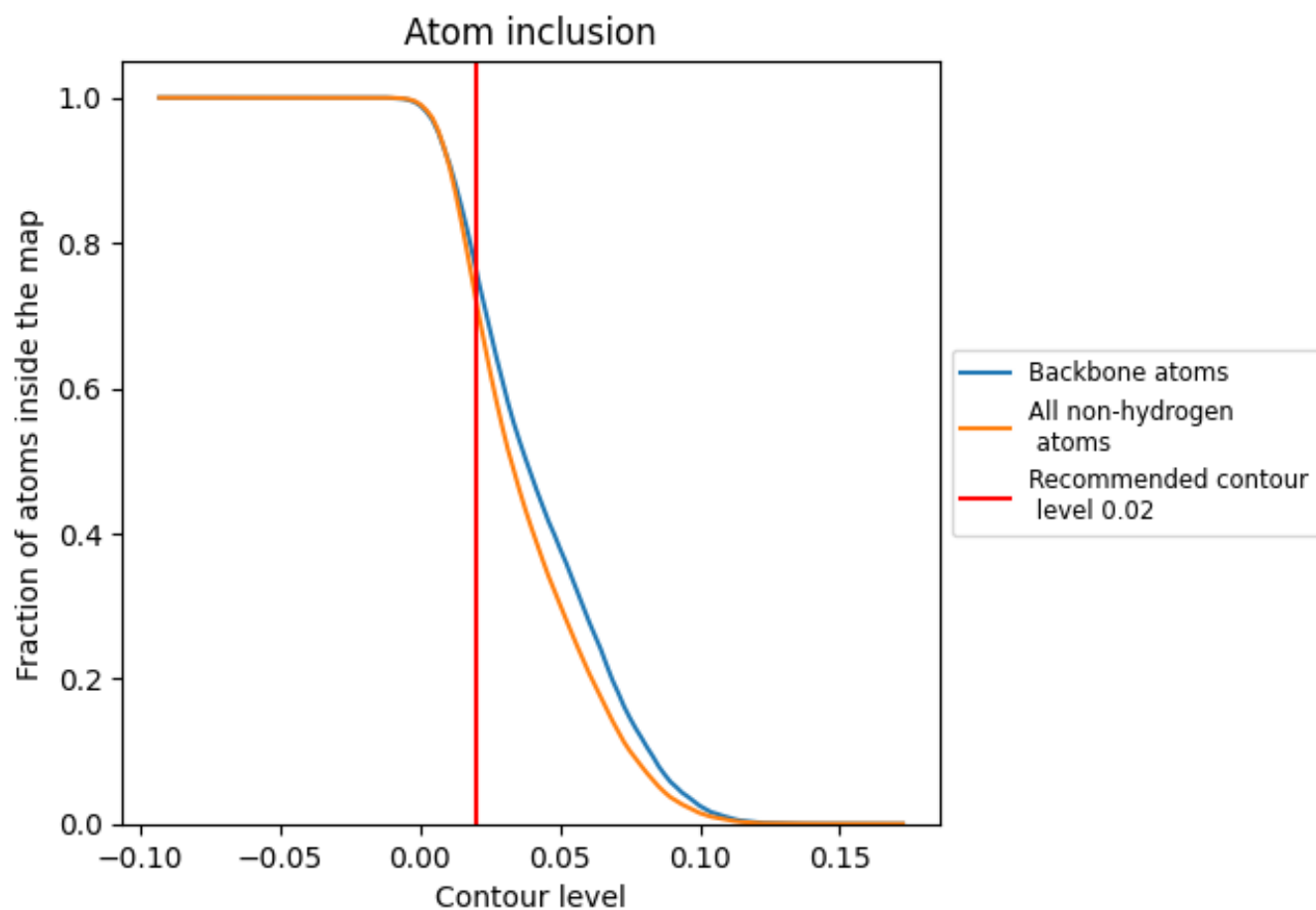
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.02).





















9.4 Atom inclusion [i](#)



At the recommended contour level, 77% of all backbone atoms, 72% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7213	 0.4700
1	 0.6072	 0.3550
2	 0.5856	 0.3590
A	 0.6955	 0.4690
B	 0.5666	 0.4150
C	 0.8353	 0.5290
D	 0.7727	 0.4970
E	 0.7921	 0.5250
F	 0.6209	 0.4130
K	 0.1013	 0.2300

