



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 03:12 AM UTC

PDB ID : 7EGH / pdb_00007egh
EMDB ID : EMD-31117
Title : TFIID lobe C subcomplex
Authors : Chen, X.; Wu, Z.; Li, J.; Zhao, D.; Xu, Y.
Deposited on : 2021-03-24
Resolution : 3.04 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

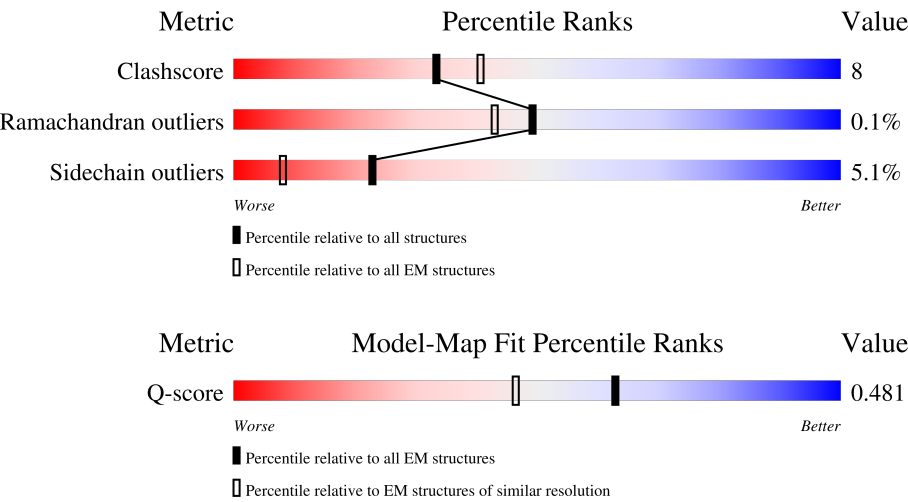
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.04 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	13952 (2.54 - 3.54)

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	1872	
2	B	1199	
3	F	677	
3	f	677	

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Mol	Chain	Length	Quality of chain
4	G	349	36% 59%
5	H	310	18% 5% 77%
6	X	72	57% 6% 38%
7	Y	94	41% 6% 52%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 20543 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 Transcription initiation factor TFIID subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	630	Total	C	N	O	S	0	0
			5199	3303	923	944	29		

- Molecule 2 is a protein called Transcription initiation factor TFIID subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	963	Total	C	N	O	S	0	0
			7796	5011	1315	1412	58		

- Molecule 3 is a protein called Transcription initiation factor TFIID subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	F	267	Total	C	N	O	S	0	0
			2006	1265	362	368	11		
3	f	260	Total	C	N	O	S	0	0
			1946	1227	348	360	11		

- Molecule 4 is a protein called Transcription initiation factor TFIID subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	144	Total	C	N	O	S	0	0
			1171	742	215	210	4		

- Molecule 5 is a protein called Transcription initiation factor TFIID subunit 8.

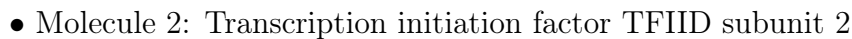
Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	72	Total	C	N	O	S	0	0
			580	370	100	109	1		

- Molecule 6 is a DNA chain called DNA (45-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	X	45	Total	C	N	O	P	0	0
			918	434	172	267	45		

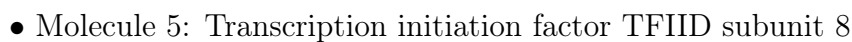
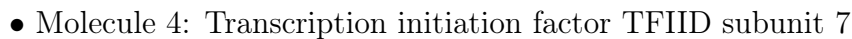
- Molecule 7 is a DNA chain called DNA (45-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	Y	45	Total	C	N	O	P	0	0
			927	438	171	273	45		



Device Type	Percentage
Smartphones	68%
Tablets	11%
Other mobile devices	20%





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	121285	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 ($e^-/\text{\AA}^2$)	50	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	5.396	Depositor
Minimum map value	-3.106	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.046	Depositor
Recommended contour level	0.222	Depositor
Map size (Å)	540.16, 540.16, 540.16	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.055, 1.055, 1.055	Depositor

5 Model quality

5.1 Standard geometry

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	0.64	0/5320	0.97	5/7165 (0.1%)
2	B	0.45	1/7993 (0.0%)	0.68	7/10836 (0.1%)
3	F	0.72	1/2040 (0.0%)	1.14	7/2775 (0.3%)
3	f	0.55	0/1979	0.90	4/2694 (0.1%)
4	G	0.67	0/1190	0.94	1/1601 (0.1%)
5	H	0.24	0/592	0.53	0/803
6	X	0.30	0/1029	0.45	0/1584
7	Y	0.29	0/1039	0.43	0/1603
All	All	0.54	2/21182 (0.0%)	0.83	24/29061 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	61	ASN	C-O	-6.02	1.18	1.24
3	F	395	ARG	C-O	5.61	1.31	1.24

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	491	HIS	N-CA-C	-9.47	100.96	111.28
3	F	271	VAL	N-CA-C	-7.16	106.03	112.90
1	A	715	PRO	N-CA-C	-7.11	100.08	111.03
2	B	495	GLN	N-CA-C	-7.04	104.69	113.28
1	A	396	LEU	N-CA-C	-6.72	104.83	113.16
3	f	326	HIS	N-CA-C	-6.67	104.24	112.38
3	f	319	LEU	N-CA-C	-6.54	103.98	112.68
1	A	467	ILE	N-CA-C	-6.42	104.00	111.00
1	A	1205	PHE	CA-CB-CG	-6.25	107.55	113.80
3	f	255	MET	N-CA-C	-5.96	104.86	114.09
3	F	316	GLN	CB-CA-C	5.93	120.27	110.78
2	B	181	GLY	CA-C-O	-5.76	118.17	122.37
2	B	583	GLU	CB-CA-C	-5.72	101.29	110.79
3	F	310	THR	CB-CA-C	5.71	120.27	110.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	330	ARG	N-CA-C	-5.62	105.15	111.28
2	B	264	GLU	N-CA-C	-5.56	106.16	113.17
3	F	324	ASP	N-CA-C	-5.56	105.32	111.71
3	f	329	LEU	N-CA-C	-5.35	106.45	112.87
3	F	272	VAL	N-CA-C	-5.35	105.16	110.62
4	G	183	ILE	N-CA-C	-5.33	102.64	109.30
2	B	491	HIS	CB-CA-C	5.23	119.48	110.79
1	A	1067	GLN	CB-CA-C	5.14	119.32	110.79
2	B	541	ARG	CB-CA-C	-5.03	101.68	110.09
3	F	438	LEU	N-CA-C	5.03	118.49	111.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5199	0	5158	134	0
2	B	7796	0	7751	89	0
3	F	2006	0	1954	71	0
3	f	1946	0	1881	48	0
4	G	1171	0	1222	15	0
5	H	580	0	591	15	0
6	X	918	0	503	3	0
7	Y	927	0	506	5	0
All	All	20543	0	19566	332	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 (332) 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:A:1200:THR:O	1:A:1204:GLU:HG2	1.52	1.08
1:A:339:ALA:HB3	1:A:355:VAL:HG21	1.42	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:666:ARG:HH22	2:B:509:ASN:HA	1.25	1.00
1:A:725:CYS:O	1:A:725:CYS:SG	2.22	0.97
1:A:467:ILE:H	3:F:221:GLN:NE2	1.65	0.95
4:G:153:LYS:HD3	4:G:153:LYS:H	1.36	0.88
1:A:475:LEU:HD11	3:f:351:ILE:HD11	1.57	0.87
1:A:511:LEU:H	1:A:511:LEU:HD22	1.37	0.87
1:A:1023:TRP:HB2	6:X:44:DT:H5'	1.59	0.84
3:f:447:ALA:HA	3:f:456:GLY:HA3	1.61	0.83
1:A:1065:GLU:HA	1:A:1065:GLU:OE1	1.78	0.83
1:A:1217:ARG:H	1:A:1217:ARG:HD3	1.44	0.81
1:A:821:ARG:HA	1:A:821:ARG:CZ	2.10	0.81
3:F:324:ASP:HA	3:F:327:TRP:HB3	1.63	0.80
1:A:1220:MET:HE3	1:A:1220:MET:O	1.82	0.79
3:F:274:ASN:ND2	3:F:316:GLN:HA	1.99	0.78
3:F:273:GLN:H	3:F:273:GLN:NE2	1.81	0.78
3:F:429:LYS:O	3:F:433:PRO:HD2	1.85	0.75
3:F:359:PHE:HB3	3:F:380:LEU:HG	1.68	0.73
1:A:467:ILE:H	3:F:221:GLN:HE21	1.33	0.73
1:A:488:ALA:HB1	3:f:271:VAL:HG23	1.71	0.72
3:F:418:ILE:HD12	3:F:418:ILE:H	1.54	0.72
1:A:639:LEU:HD11	1:A:774:ARG:HG2	1.70	0.72
1:A:666:ARG:NH2	2:B:509:ASN:HA	2.03	0.72
4:G:129:LYS:O	4:G:129:LYS:NZ	2.23	0.71
1:A:1217:ARG:H	1:A:1217:ARG:CD	2.04	0.70
2:B:241:PRO:HG2	2:B:496:MET:HE1	1.72	0.70
1:A:353:LEU:HD21	3:f:272:VAL:HG21	1.72	0.70
6:X:38:DC:H2''	6:X:39:DC:H5'	1.73	0.70
2:B:490:SER:HA	2:B:493:TRP:HB2	1.75	0.69
1:A:568:SER:HB2	2:B:389:ASN:HB2	1.75	0.69
3:f:447:ALA:CA	3:f:456:GLY:HA3	2.23	0.68
1:A:396:LEU:HD21	3:f:394:PRO:HB2	1.77	0.67
3:f:239:ARG:HD3	3:f:284:ARG:HH11	1.60	0.66
3:F:274:ASN:HB2	3:F:317:LEU:H	1.58	0.66
1:A:339:ALA:CB	1:A:355:VAL:HG21	2.22	0.65
1:A:1168:TYR:HB2	4:G:180:ARG:HB3	1.78	0.65
1:A:821:ARG:HA	1:A:821:ARG:NH1	2.11	0.65
1:A:869:ARG:HH21	2:B:582:GLU:HG2	1.62	0.65
3:f:320:ARG:NH1	3:f:320:ARG:HB3	2.12	0.65
1:A:970:ASN:OD1	1:A:970:ASN:N	2.21	0.64
2:B:361:ARG:HD2	2:B:367:GLU:HB2	1.80	0.63
2:B:583:GLU:CD	2:B:583:GLU:H	2.06	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1217:ARG:HD3	1:A:1217:ARG:N	2.13	0.62
1:A:342:ARG:O	1:A:342:ARG:HG2	1.99	0.61
3:f:219:GLU:OE1	3:f:219:GLU:N	2.26	0.61
2:B:203:LYS:HD3	2:B:237:MET:HE3	1.83	0.60
3:f:330:ARG:HH22	3:f:371:THR:HB	1.66	0.60
5:H:194:MET:HE1	3:f:226:GLU:HG3	1.83	0.60
1:A:475:LEU:HD13	3:f:349:ASN:OD1	2.02	0.60
1:A:1193:ALA:HB3	4:G:166:VAL:HG21	1.83	0.60
5:H:192:ARG:NH1	5:H:209:SER:O	2.34	0.60
1:A:470:ILE:O	1:A:470:ILE:HG22	2.02	0.59
1:A:1242:LYS:C	1:A:1242:LYS:HD2	2.28	0.59
2:B:27:LYS:HE3	2:B:490:SER:HB3	1.84	0.59
3:F:312:ILE:HG13	3:F:334:ALA:CB	2.33	0.58
1:A:511:LEU:HD22	1:A:511:LEU:N	2.11	0.58
2:B:769:LYS:HG3	2:B:769:LYS:O	2.03	0.58
3:F:207:ARG:HD2	3:F:207:ARG:N	2.18	0.58
3:F:335:ARG:NE	3:F:382:GLU:OE1	2.33	0.58
5:H:171:ASP:HB3	5:H:174:VAL:HG12	1.86	0.58
3:f:349:ASN:HB3	3:f:351:ILE:HG13	1.84	0.58
1:A:929:THR:HG22	1:A:954:ALA:HA	1.86	0.58
3:f:320:ARG:CB	3:f:320:ARG:HH11	2.17	0.58
1:A:334:THR:HA	1:A:489:GLN:O	2.04	0.57
1:A:1020:LEU:HD12	1:A:1020:LEU:O	2.05	0.57
1:A:1063:LYS:C	1:A:1063:LYS:HD3	2.29	0.57
3:F:335:ARG:HH21	3:F:378:ALA:HB1	1.69	0.57
2:B:837:LEU:HD22	5:H:213:LEU:HD22	1.86	0.57
1:A:340:GLU:N	1:A:497:PRO:HG2	2.18	0.57
1:A:402:PHE:HA	1:A:407:GLN:NE2	2.20	0.57
1:A:1222:LYS:HD2	1:A:1222:LYS:O	2.04	0.57
2:B:560:TYR:CD2	2:B:581:ILE:HD11	2.39	0.56
3:F:418:ILE:HD12	3:F:418:ILE:N	2.20	0.56
3:F:315:ARG:CZ	3:F:369:PRO:HB3	2.35	0.56
3:f:223:TYR:HD2	3:f:255:MET:HE1	1.71	0.56
2:B:27:LYS:CE	2:B:490:SER:HB3	2.35	0.56
1:A:401:ASN:HD22	1:A:401:ASN:H	1.53	0.56
2:B:259:ASP:HB3	2:B:262:MET:O	2.05	0.56
3:f:280:ILE:O	3:f:284:ARG:HG3	2.06	0.56
2:B:61:ASN:O	2:B:130:VAL:HG23	2.06	0.55
2:B:544:LEU:HD23	2:B:590:ILE:HD11	1.89	0.55
2:B:248:SER:OG	2:B:322:MET:SD	2.65	0.55
3:f:352:GLN:O	3:f:356:THR:OG1	2.25	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:406:VAL:HG11	3:F:420:ALA:HB2	1.88	0.55
3:F:265:GLU:HA	3:F:265:GLU:OE2	2.06	0.55
1:A:463:PRO:HD3	3:F:222:LEU:HB2	1.89	0.55
1:A:1052:ARG:HD2	1:A:1052:ARG:O	2.07	0.54
3:F:312:ILE:HG13	3:F:334:ALA:HB3	1.88	0.54
3:F:403:ILE:O	3:F:406:VAL:HG22	2.08	0.54
3:f:253:TYR:O	3:f:254:GLN:HB3	2.06	0.54
1:A:395:ASP:N	1:A:395:ASP:OD1	2.41	0.54
3:f:390:THR:HG23	3:f:391:LEU:HD22	1.89	0.54
1:A:355:VAL:HG23	1:A:355:VAL:O	2.07	0.54
1:A:489:GLN:H	1:A:489:GLN:NE2	2.05	0.54
2:B:66:ASN:HD21	2:B:119:PRO:HB3	1.73	0.53
2:B:571:LEU:HD21	2:B:619:MET:HE1	1.89	0.53
2:B:820:ASP:OD1	2:B:820:ASP:N	2.37	0.53
3:f:318:CYS:SG	3:f:326:HIS:HB3	2.49	0.53
2:B:628:ILE:O	2:B:644:GLN:NE2	2.41	0.53
2:B:489:GLN:OE1	2:B:489:GLN:N	2.34	0.53
3:F:433:PRO:HA	3:F:436:ALA:HB3	1.89	0.53
1:A:339:ALA:HA	1:A:497:PRO:HG3	1.91	0.53
1:A:675:ASP:OD1	4:G:78:LYS:NZ	2.37	0.53
2:B:629:ARG:NH1	2:B:658:ASP:OD2	2.37	0.53
3:f:447:ALA:O	3:f:453:GLY:HA2	2.09	0.52
4:G:181:TRP:O	4:G:181:TRP:CE3	2.63	0.52
1:A:851:ASP:OD1	1:A:851:ASP:N	2.38	0.52
2:B:345:CYS:HA	2:B:348:GLN:HE21	1.73	0.52
2:B:937:THR:HG23	2:B:939:ASN:H	1.74	0.52
3:F:335:ARG:HD2	3:F:382:GLU:HB2	1.90	0.52
4:G:149:ARG:O	4:G:149:ARG:HG3	2.09	0.52
1:A:843:ARG:NH1	7:Y:18:DC:OP2	2.43	0.52
3:F:273:GLN:H	3:F:273:GLN:CD	2.17	0.52
2:B:66:ASN:OD1	2:B:187:ARG:NH1	2.43	0.52
2:B:529:VAL:HG11	2:B:560:TYR:HB2	1.91	0.52
2:B:808:PRO:HA	2:B:864:ASN:HB3	1.90	0.52
2:B:90:THR:O	2:B:968:ARG:NH2	2.43	0.51
3:f:320:ARG:NH1	3:f:320:ARG:CB	2.73	0.51
1:A:828:GLU:HA	1:A:831:LYS:HB2	1.92	0.51
1:A:510:ILE:HB	1:A:511:LEU:HD13	1.93	0.51
1:A:1200:THR:C	1:A:1204:GLU:HG2	2.30	0.51
2:B:202:TRP:HB2	2:B:238:LEU:HB3	1.92	0.51
2:B:780:LYS:O	5:H:183:ARG:NH1	2.44	0.51
2:B:983:ARG:HG3	2:B:983:ARG:HH11	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1222:LYS:HD2	1:A:1222:LYS:C	2.35	0.51
1:A:338:VAL:HG12	1:A:497:PRO:HD3	1.91	0.51
2:B:656:GLU:HG2	2:B:661:ALA:HB3	1.92	0.51
3:f:283:MET:HG3	3:f:329:LEU:HD11	1.91	0.51
1:A:804:ARG:NH1	6:X:64:DC:OP1	2.44	0.50
2:B:274:LEU:HG	2:B:278:LYS:HE2	1.93	0.50
2:B:983:ARG:HG3	2:B:983:ARG:NH1	2.26	0.50
1:A:1163:ARG:HB3	4:G:183:ILE:HG22	1.92	0.50
3:F:273:GLN:CD	3:F:273:GLN:N	2.70	0.50
2:B:739:ASN:ND2	2:B:782:ASN:OD1	2.44	0.50
2:B:818:THR:OG1	2:B:819:LEU:N	2.45	0.50
1:A:994:ASP:N	1:A:994:ASP:OD1	2.45	0.50
3:F:243:LEU:HD21	3:F:284:ARG:HB3	1.92	0.50
3:F:257:PRO:O	3:F:261:THR:OG1	2.26	0.50
4:G:94:MET:HE3	4:G:96:VAL:HG22	1.94	0.50
1:A:468:PHE:CE1	3:F:216:LEU:HD12	2.47	0.50
1:A:821:ARG:HA	1:A:821:ARG:NE	2.25	0.49
2:B:152:LEU:HD23	2:B:155:PRO:HB3	1.94	0.49
3:F:316:GLN:NE2	3:F:319:LEU:O	2.45	0.49
1:A:638:PRO:HB3	4:G:39:LEU:HD23	1.95	0.49
4:G:181:TRP:CD2	4:G:181:TRP:C	2.87	0.49
1:A:467:ILE:N	3:F:221:GLN:NE2	2.48	0.49
2:B:570:GLU:HA	2:B:625:LEU:HA	1.95	0.49
1:A:337:ARG:HH11	1:A:338:VAL:H	1.61	0.49
1:A:927:VAL:O	1:A:933:ASN:ND2	2.44	0.48
1:A:1022:ARG:HG3	7:Y:47:DG:H4'	1.95	0.48
1:A:639:LEU:O	1:A:643:ILE:HG13	2.13	0.48
2:B:598:ARG:HH11	2:B:619:MET:HE3	1.77	0.48
1:A:338:VAL:HG13	1:A:360:SER:HB2	1.96	0.48
2:B:560:TYR:HD2	2:B:581:ILE:HD11	1.78	0.48
1:A:485:ILE:HB	3:f:310:THR:HG23	1.94	0.48
2:B:882:HIS:HE1	5:H:216:ALA:HB1	1.77	0.48
3:F:312:ILE:HG22	3:F:313:VAL:HG13	1.94	0.48
5:H:197:THR:HB	3:f:238:LYS:HG2	1.95	0.48
2:B:217:ASN:ND2	2:B:320:ALA:O	2.35	0.48
3:F:215:GLU:O	3:F:254:GLN:NE2	2.39	0.48
1:A:475:LEU:O	3:f:354:ARG:NH2	2.47	0.48
2:B:329:LEU:HB3	2:B:342:THR:HG23	1.95	0.48
2:B:953:LEU:HD12	2:B:979:PHE:HE2	1.79	0.48
3:F:415:ILE:HG12	3:F:415:ILE:O	2.14	0.48
4:G:48:HIS:HD2	4:G:54:GLY:HA2	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:636:VAL:HG23	2:B:638:ARG:HB3	1.96	0.48
2:B:781:TYR:O	5:H:176:ARG:NH2	2.47	0.48
3:f:257:PRO:O	3:f:261:THR:OG1	2.31	0.48
1:A:569:ASN:HB3	1:A:572:TYR:HD2	1.78	0.48
1:A:1014:GLU:O	1:A:1018:LYS:HB2	2.13	0.48
3:f:254:GLN:O	3:f:258:ARG:NH2	2.47	0.48
3:F:409:GLY:O	3:F:417:ARG:NH2	2.46	0.48
1:A:575:PRO:HD3	2:B:608:ASN:HD22	1.79	0.47
1:A:995:LEU:O	1:A:998:LEU:HB2	2.14	0.47
1:A:501:THR:O	1:A:502:LEU:C	2.58	0.47
3:F:317:LEU:N	3:F:317:LEU:CD2	2.77	0.47
1:A:824:ARG:HB3	1:A:863:VAL:HG12	1.95	0.47
2:B:184:ASN:N	2:B:184:ASN:ND2	2.63	0.47
3:f:326:HIS:O	3:f:330:ARG:HG2	2.15	0.47
2:B:135:TRP:HA	2:B:138:VAL:HG12	1.95	0.47
3:F:317:LEU:N	3:F:317:LEU:HD23	2.30	0.47
4:G:41:ASP:N	4:G:41:ASP:OD1	2.48	0.47
2:B:564:LEU:HB2	2:B:581:ILE:HD13	1.97	0.46
3:F:418:ILE:H	3:F:418:ILE:CD1	2.26	0.46
2:B:559:LYS:HB2	2:B:559:LYS:HZ3	1.80	0.46
2:B:71:ARG:HD3	2:B:73:TYR:CE1	2.50	0.46
3:F:411:VAL:O	3:F:411:VAL:HG13	2.15	0.46
1:A:857:MET:HE3	1:A:858:ASP:H	1.81	0.46
2:B:559:LYS:HB2	2:B:559:LYS:NZ	2.31	0.46
5:H:186:VAL:HG21	3:f:220:GLN:HB3	1.97	0.46
3:f:328:ALA:C	3:f:330:ARG:N	2.72	0.46
1:A:397:LEU:HD21	3:f:361:LYS:HB3	1.97	0.46
1:A:1063:LYS:HD3	1:A:1063:LYS:O	2.16	0.46
3:f:330:ARG:NH1	3:f:371:THR:O	2.47	0.46
1:A:1165:LEU:HB2	1:A:1191:ILE:HG12	1.97	0.45
3:f:326:HIS:ND1	3:f:326:HIS:N	2.63	0.45
1:A:404:MET:O	1:A:408:LEU:HB2	2.16	0.45
1:A:1219:GLU:CD	1:A:1219:GLU:C	2.84	0.45
1:A:475:LEU:HD23	1:A:480:TRP:HE1	1.82	0.45
2:B:568:VAL:HG22	2:B:628:ILE:HG23	1.98	0.45
3:F:424:GLN:O	3:F:428:LEU:HG	2.15	0.45
3:F:426:LEU:C	3:F:426:LEU:HD23	2.42	0.45
1:A:1224:ARG:O	1:A:1227:ILE:HG22	2.16	0.45
3:f:317:LEU:HD12	3:f:329:LEU:HD23	1.98	0.45
1:A:1053:PHE:HB2	1:A:1057:GLU:HB2	1.99	0.45
3:f:327:TRP:O	3:f:330:ARG:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:194:MET:HB2	5:H:194:MET:HE2	1.85	0.44
1:A:690:LEU:HD22	1:A:747:LEU:HD22	1.99	0.44
1:A:470:ILE:O	3:f:299:LYS:NZ	2.49	0.44
3:F:350:ASN:OD1	3:F:350:ASN:N	2.51	0.44
1:A:352:MET:HE3	1:A:352:MET:HB3	1.81	0.44
1:A:463:PRO:O	3:F:218:VAL:HG23	2.18	0.44
1:A:1182:CYS:O	1:A:1182:CYS:SG	2.75	0.44
3:F:361:LYS:HA	3:F:364:VAL:HG22	1.99	0.44
1:A:899:LYS:HE3	1:A:899:LYS:HB2	1.81	0.44
2:B:492:MET:HE3	2:B:492:MET:HB2	1.85	0.44
2:B:839:MET:HE1	2:B:843:LEU:HD13	2.00	0.44
3:F:233:GLY:HA2	3:F:239:ARG:HE	1.82	0.44
1:A:1239:GLU:HA	1:A:1239:GLU:OE1	2.18	0.44
3:F:328:ALA:O	3:F:332:PHE:N	2.51	0.44
3:F:400:GLY:O	3:F:404:ARG:HG2	2.18	0.44
1:A:464:TRP:CD1	1:A:464:TRP:C	2.96	0.44
3:F:223:TYR:HH	3:F:245:SER:HG	1.62	0.44
3:f:317:LEU:HG	3:f:330:ARG:HE	1.83	0.44
1:A:401:ASN:HD22	1:A:401:ASN:N	2.14	0.44
2:B:414:LEU:HB2	2:B:523:VAL:HG13	1.99	0.43
2:B:887:ARG:NH2	2:B:917:ASP:OD2	2.49	0.43
1:A:1022:ARG:NH2	7:Y:46:DT:O2	2.45	0.43
3:F:370:TRP:CZ3	3:F:420:ALA:HA	2.53	0.43
1:A:592:SER:OG	1:A:695:GLY:O	2.31	0.43
1:A:639:LEU:O	1:A:639:LEU:CD1	2.67	0.43
1:A:730:PHE:CD1	1:A:730:PHE:N	2.82	0.43
1:A:996:ARG:HA	1:A:1026:ILE:HD13	1.99	0.43
3:f:223:TYR:CD2	3:f:255:MET:HE1	2.52	0.43
1:A:400:GLU:HG3	1:A:401:ASN:N	2.32	0.43
2:B:328:ASN:HA	5:H:227:LEU:HD21	2.00	0.43
3:F:402:ARG:O	3:F:406:VAL:HG13	2.19	0.43
1:A:854:ARG:NH2	7:Y:17:DT:O2	2.49	0.43
1:A:1052:ARG:O	1:A:1052:ARG:CG	2.67	0.43
2:B:766:LEU:HA	2:B:807:THR:HG21	2.00	0.43
3:F:283:MET:HE1	3:F:308:VAL:HG12	2.00	0.43
1:A:925:ASP:N	1:A:925:ASP:OD1	2.52	0.43
2:B:562:GLY:O	2:B:581:ILE:HG12	2.19	0.43
2:B:831:GLU:OE1	2:B:835:ARG:NH2	2.52	0.43
3:F:312:ILE:HG13	3:F:334:ALA:HB1	2.00	0.43
3:f:320:ARG:HH11	3:f:320:ARG:HB2	1.83	0.43
2:B:653:LEU:HG	2:B:684:ILE:HG13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:f:447:ALA:CB	3:f:456:GLY:HA3	2.49	0.43
1:A:727:THR:HG23	1:A:727:THR:O	2.19	0.42
1:A:825:ILE:HD12	1:A:830:ILE:HD11	2.01	0.42
2:B:882:HIS:CE1	5:H:216:ALA:HB1	2.54	0.42
4:G:181:TRP:O	4:G:181:TRP:CD2	2.72	0.42
3:f:432:ALA:HB1	3:f:462:GLN:CB	2.50	0.42
3:F:312:ILE:HG22	3:F:313:VAL:CG1	2.50	0.42
3:F:324:ASP:CG	3:F:327:TRP:HD1	2.28	0.42
7:Y:14:DA:H2''	7:Y:15:DC:H5''	2.01	0.42
3:F:401:GLU:CD	3:F:401:GLU:C	2.87	0.42
3:f:428:LEU:HD13	3:f:458:LEU:CB	2.49	0.42
1:A:847:LYS:HE2	1:A:847:LYS:HB3	1.77	0.42
1:A:1218:GLU:C	1:A:1218:GLU:CD	2.88	0.42
2:B:216:SER:OG	2:B:217:ASN:N	2.52	0.42
2:B:594:SER:OG	2:B:595:LYS:N	2.52	0.42
1:A:604:PRO:O	1:A:889:TYR:OH	2.34	0.42
1:A:1217:ARG:HG2	1:A:1217:ARG:HH11	1.85	0.42
3:F:397:GLN:H	3:F:397:GLN:HG3	1.60	0.42
3:F:410:PRO:O	3:F:411:VAL:C	2.63	0.42
1:A:666:ARG:NH2	2:B:508:SER:O	2.52	0.42
2:B:298:ARG:HH11	2:B:298:ARG:HD3	1.69	0.42
2:B:957:MET:O	2:B:968:ARG:NH1	2.53	0.42
1:A:620:LEU:HD11	1:A:766:ARG:HD3	2.02	0.42
1:A:645:LYS:HD2	1:A:645:LYS:HA	1.76	0.42
3:F:396:LEU:HA	3:F:399:GLU:OE1	2.20	0.42
1:A:468:PHE:HE1	3:F:221:GLN:HG3	1.85	0.41
2:B:203:LYS:HZ3	2:B:235:HIS:HB3	1.85	0.41
1:A:464:TRP:CD1	1:A:464:TRP:O	2.74	0.41
1:A:502:LEU:HD22	1:A:502:LEU:HA	1.85	0.41
1:A:639:LEU:O	1:A:639:LEU:HD13	2.20	0.41
1:A:819:LYS:HZ3	1:A:819:LYS:H	1.67	0.41
1:A:1007:LEU:HD22	1:A:1012:VAL:HG21	2.02	0.41
2:B:639:LYS:NZ	2:B:641:GLU:OE2	2.42	0.41
2:B:749:LEU:HA	2:B:752:THR:HG22	2.03	0.41
1:A:353:LEU:HD22	3:f:273:GLN:HE22	1.85	0.41
1:A:472:ASN:ND2	3:f:299:LYS:O	2.53	0.41
2:B:176:HIS:HE1	2:B:310:ASP:H	1.68	0.41
2:B:544:LEU:HD22	2:B:625:LEU:HD22	2.02	0.41
1:A:1241:GLU:OE1	1:A:1241:GLU:HA	2.21	0.41
2:B:55:PRO:HB3	2:B:60:LEU:HD22	2.02	0.41
3:F:219:GLU:CD	3:F:219:GLU:C	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:GLU:H	1:A:497:PRO:HG2	1.83	0.41
1:A:468:PHE:CE1	3:F:221:GLN:HG3	2.54	0.41
1:A:1018:LYS:C	1:A:1020:LEU:N	2.78	0.41
2:B:881:GLY:HA3	5:H:218:PRO:HB3	2.02	0.41
3:f:402:ARG:HH22	3:f:403:ILE:HD12	1.86	0.41
1:A:338:VAL:HG22	1:A:360:SER:HB2	2.03	0.41
2:B:102:ASN:OD1	2:B:102:ASN:N	2.54	0.41
3:F:277:ALA:O	3:F:278:LEU:C	2.64	0.41
2:B:100:GLN:O	2:B:101:ARG:NH1	2.46	0.41
2:B:464:ILE:HG12	2:B:513:LYS:HE2	2.02	0.41
2:B:984:PRO:HG2	2:B:987:LEU:HD22	2.02	0.41
1:A:511:LEU:N	1:A:511:LEU:HD13	2.35	0.41
1:A:725:CYS:SG	1:A:734:LEU:HD12	2.61	0.41
2:B:295:LEU:HD13	2:B:477:LEU:HD11	2.03	0.41
3:F:388:ILE:HA	3:F:392:ILE:HG12	2.03	0.41
3:F:426:LEU:O	3:F:429:LYS:HG2	2.21	0.41
1:A:1165:LEU:HD22	1:A:1165:LEU:HA	1.87	0.41
1:A:1216:HIS:HB3	1:A:1219:GLU:HB3	2.03	0.41
1:A:1221:ARG:HG2	1:A:1224:ARG:HH11	1.86	0.41
2:B:62:ARG:H	2:B:62:ARG:HG2	1.58	0.41
2:B:218:GLY:HA2	2:B:238:LEU:HD13	2.03	0.41
3:F:329:LEU:HA	3:F:332:PHE:HB3	2.02	0.41
3:F:374:TYR:CE1	3:F:422:HIS:HB3	2.56	0.41
1:A:337:ARG:HD3	1:A:338:VAL:H	1.85	0.41
5:H:166:ARG:HD2	5:H:166:ARG:HA	1.87	0.41
3:f:216:LEU:HD21	3:f:254:GLN:O	2.21	0.41
1:A:1061:ARG:HD3	1:A:1061:ARG:HA	1.63	0.40
3:F:231:CYS:HB3	3:F:285:MET:HE2	2.02	0.40
3:F:427:LEU:HD12	3:F:427:LEU:HA	1.72	0.40
1:A:567:LEU:HD22	2:B:745:GLN:HG2	2.04	0.40
3:F:233:GLY:CA	3:F:239:ARG:HE	2.34	0.40
2:B:77:ILE:HG12	2:B:146:ILE:HG23	2.03	0.40
4:G:129:LYS:HD2	4:G:129:LYS:HA	1.76	0.40
3:f:446:ASP:O	3:f:450:ALA:N	2.54	0.40
1:A:463:PRO:HD2	3:F:218:VAL:HG22	2.03	0.40
1:A:1019:LYS:HB2	1:A:1019:LYS:NZ	2.37	0.40
2:B:180:CYS:HB2	2:B:313:TYR:HB2	2.04	0.40
2:B:319:TYR:O	2:B:353:GLN:NE2	2.49	0.40
3:F:370:TRP:CH2	3:F:402:ARG:HB3	2.56	0.40
3:F:433:PRO:C	3:F:435:LEU:N	2.78	0.40
1:A:1068:ARG:HD2	1:A:1068:ARG:C	2.47	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:396:LEU:HA	3:F:399:GLU:CD	2.46	0.40
5:H:178:LYS:HA	5:H:178:LYS:HD3	1.82	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	610/1872 (33%)	583 (96%)	25 (4%)	2 (0%)	36	67
2	B	959/1199 (80%)	911 (95%)	48 (5%)	0	100	100
3	F	265/677 (39%)	251 (95%)	13 (5%)	1 (0%)	30	61
3	f	258/677 (38%)	252 (98%)	6 (2%)	0	100	100
4	G	138/349 (40%)	135 (98%)	3 (2%)	0	100	100
5	H	70/310 (23%)	67 (96%)	3 (4%)	0	100	100
All	All	2300/5084 (45%)	2199 (96%)	98 (4%)	3 (0%)	49	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	502	LEU
3	F	411	VAL
1	A	998	LEU

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	564/1665 (34%)	511 (91%)	53 (9%)	8	29
2	B	876/1083 (81%)	862 (98%)	14 (2%)	55	76
3	F	202/574 (35%)	184 (91%)	18 (9%)	9	32
3	f	195/574 (34%)	185 (95%)	10 (5%)	21	52
4	G	132/322 (41%)	124 (94%)	8 (6%)	17	46
5	H	64/270 (24%)	64 (100%)	0	100	100
All	All	2033/4488 (45%)	1930 (95%)	103 (5%)	23	52

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	337	ARG
1	A	338	VAL
1	A	348	LEU
1	A	353	LEU
1	A	395	ASP
1	A	397	LEU
1	A	400	GLU
1	A	401	ASN
1	A	403	LEU
1	A	404	MET
1	A	408	LEU
1	A	415	ILE
1	A	416	TRP
1	A	419	GLU
1	A	470	ILE
1	A	475	LEU
1	A	481	GLU
1	A	483	ASN
1	A	491	MET
1	A	499	VAL
1	A	500	LEU
1	A	501	THR
1	A	502	LEU
1	A	505	ASN
1	A	511	LEU
1	A	639	LEU
1	A	661	GLU
1	A	667	THR

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Mol	Chain	Res	Type
1	A	711	ASP
1	A	727	THR
1	A	730	PHE
1	A	819	LYS
1	A	821	ARG
1	A	828	GLU
1	A	925	ASP
1	A	943	LYS
1	A	970	ASN
1	A	994	ASP
1	A	1020	LEU
1	A	1023	TRP
1	A	1025	VAL
1	A	1029	VAL
1	A	1052	ARG
1	A	1058	HIS
1	A	1062	TYR
1	A	1068	ARG
1	A	1073	GLN
1	A	1165	LEU
1	A	1203	GLU
1	A	1217	ARG
1	A	1219	GLU
1	A	1227	ILE
1	A	1237	ASN
2	B	21	GLU
2	B	71	ARG
2	B	140	GLU
2	B	184	ASN
2	B	262	MET
2	B	266	THR
2	B	293	GLU
2	B	559	LYS
2	B	603	LYS
2	B	640	VAL
2	B	746	SER
2	B	771	VAL
2	B	818	THR
2	B	987	LEU
3	F	219	GLU
3	F	261	THR
3	F	269	VAL

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Mol	Chain	Res	Type
3	F	271	VAL
3	F	280	ILE
3	F	295	LEU
3	F	301	VAL
3	F	312	ILE
3	F	317	LEU
3	F	326	HIS
3	F	340	ILE
3	F	348	THR
3	F	354	ARG
3	F	368	THR
3	F	397	GLN
3	F	406	VAL
3	F	415	ILE
3	F	427	LEU
4	G	41	ASP
4	G	129	LYS
4	G	131	ILE
4	G	143	VAL
4	G	153	LYS
4	G	181	TRP
4	G	182	GLU
4	G	183	ILE
3	f	215	GLU
3	f	253	TYR
3	f	261	THR
3	f	272	VAL
3	f	322	ASP
3	f	323	VAL
3	f	326	HIS
3	f	356	THR
3	f	366	GLU
3	f	421	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	401	ASN
1	A	472	ASN
1	A	489	GLN
1	A	590	GLN
1	A	630	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	687	ASN
1	A	693	GLN
1	A	702	ASN
1	A	755	HIS
1	A	860	ASN
1	A	896	GLN
1	A	1073	GLN
1	A	1237	ASN
1	A	1238	GLN
2	B	30	HIS
2	B	36	ASN
2	B	123	ASN
2	B	137	HIS
2	B	176	HIS
2	B	183	GLN
2	B	235	HIS
2	B	348	GLN
2	B	439	HIS
2	B	450	GLN
2	B	509	ASN
2	B	521	GLN
2	B	577	HIS
2	B	644	GLN
2	B	652	GLN
2	B	750	GLN
2	B	813	ASN
2	B	864	ASN
2	B	882	HIS
2	B	908	GLN
2	B	912	ASN
2	B	963	HIS
3	F	221	GLN
3	F	273	GLN
3	F	275	ASN
3	F	326	HIS
4	G	10	HIS
4	G	48	HIS
4	G	142	ASN
3	f	273	GLN
3	f	325	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](#)

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.

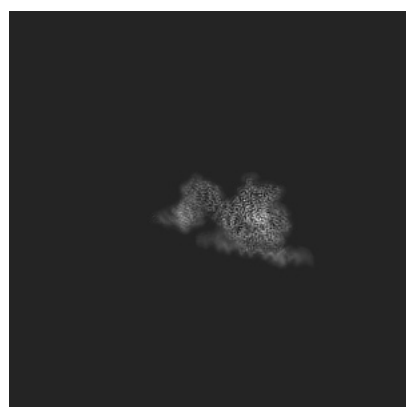
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31117. These allow visual inspection of the internal detail of the map and identification of artifacts.

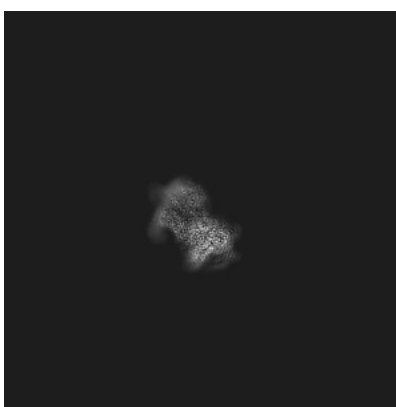
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

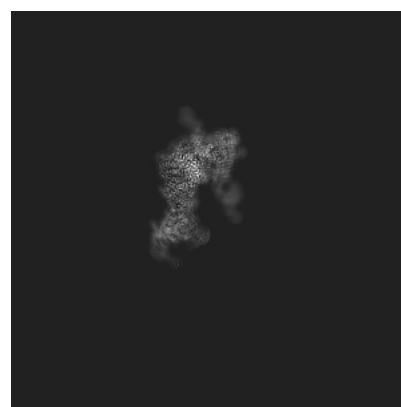
6.1.1 Primary map



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: 256



Y Index: 256



Z Index: 256

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: 231



Y Index: 321

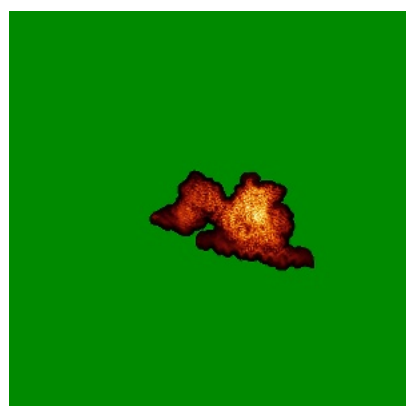


Z Index: 248

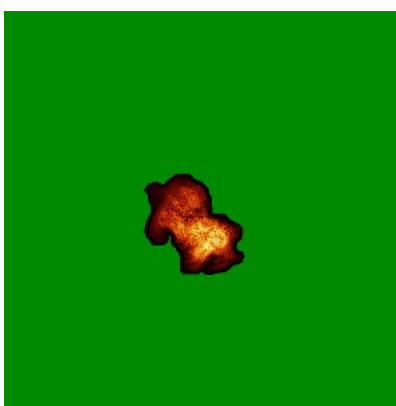
The images above show the largest variance slices of the map in three orthogonal directions.

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

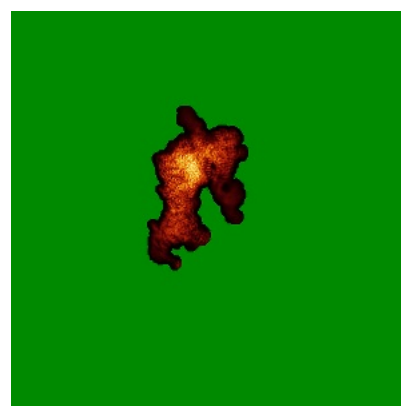
6.4.1 Primary map



X



Y

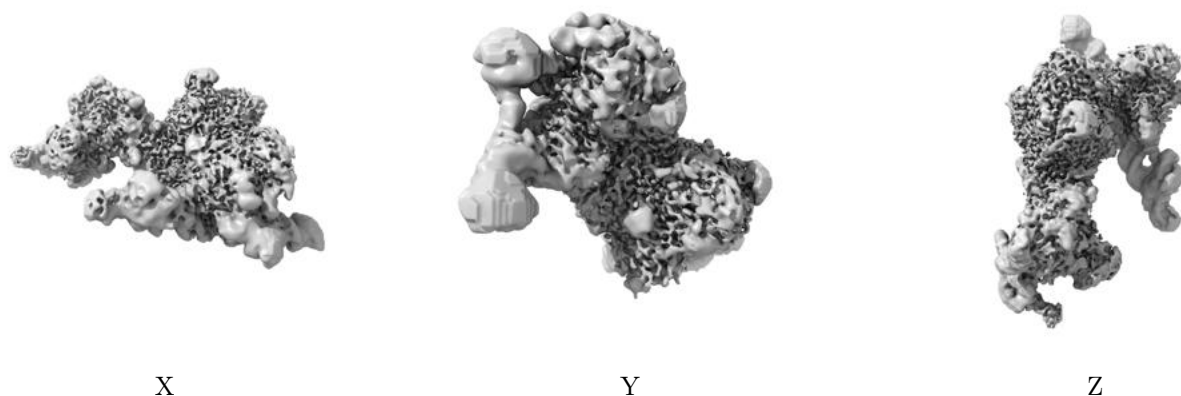


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.222. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

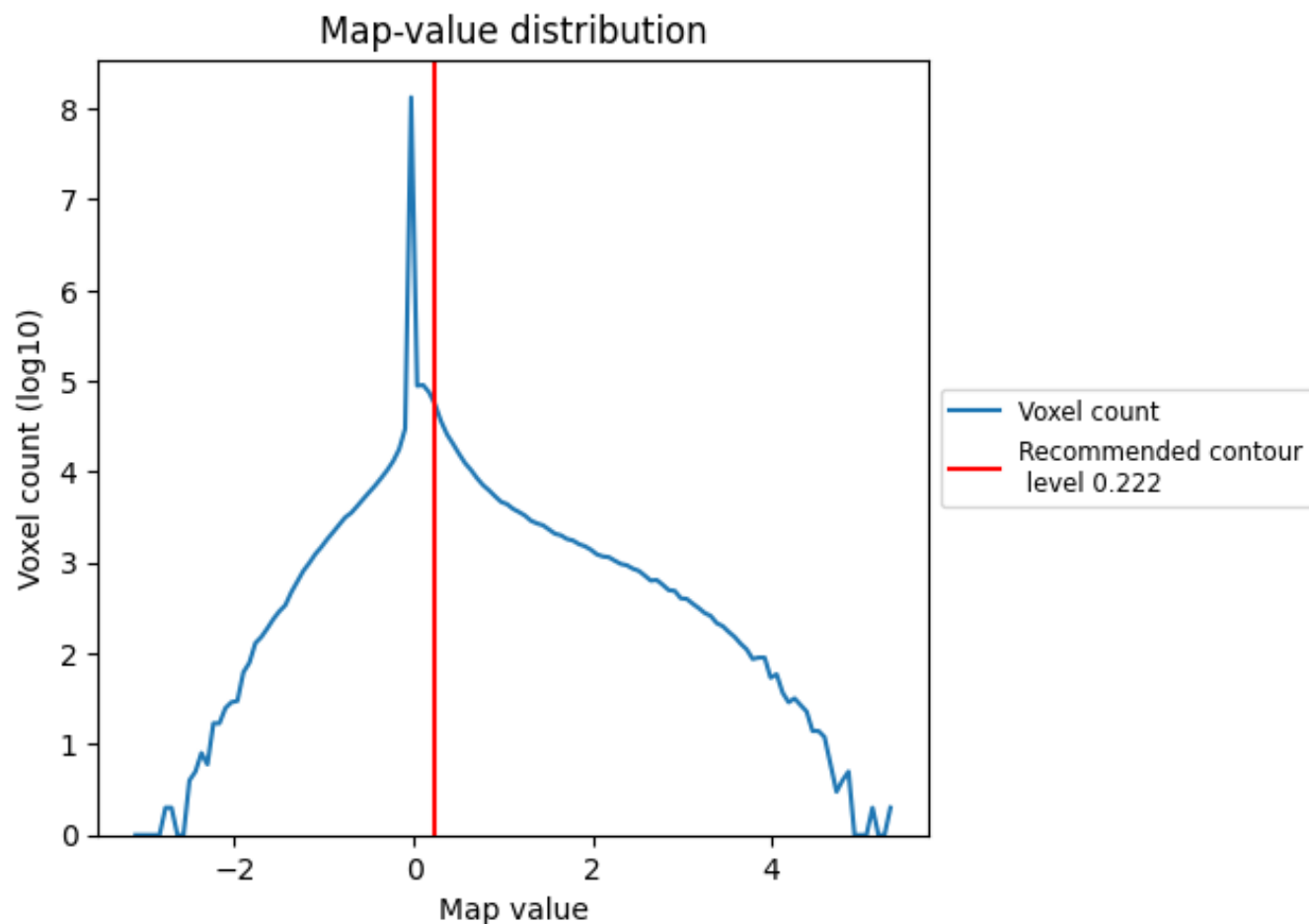
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

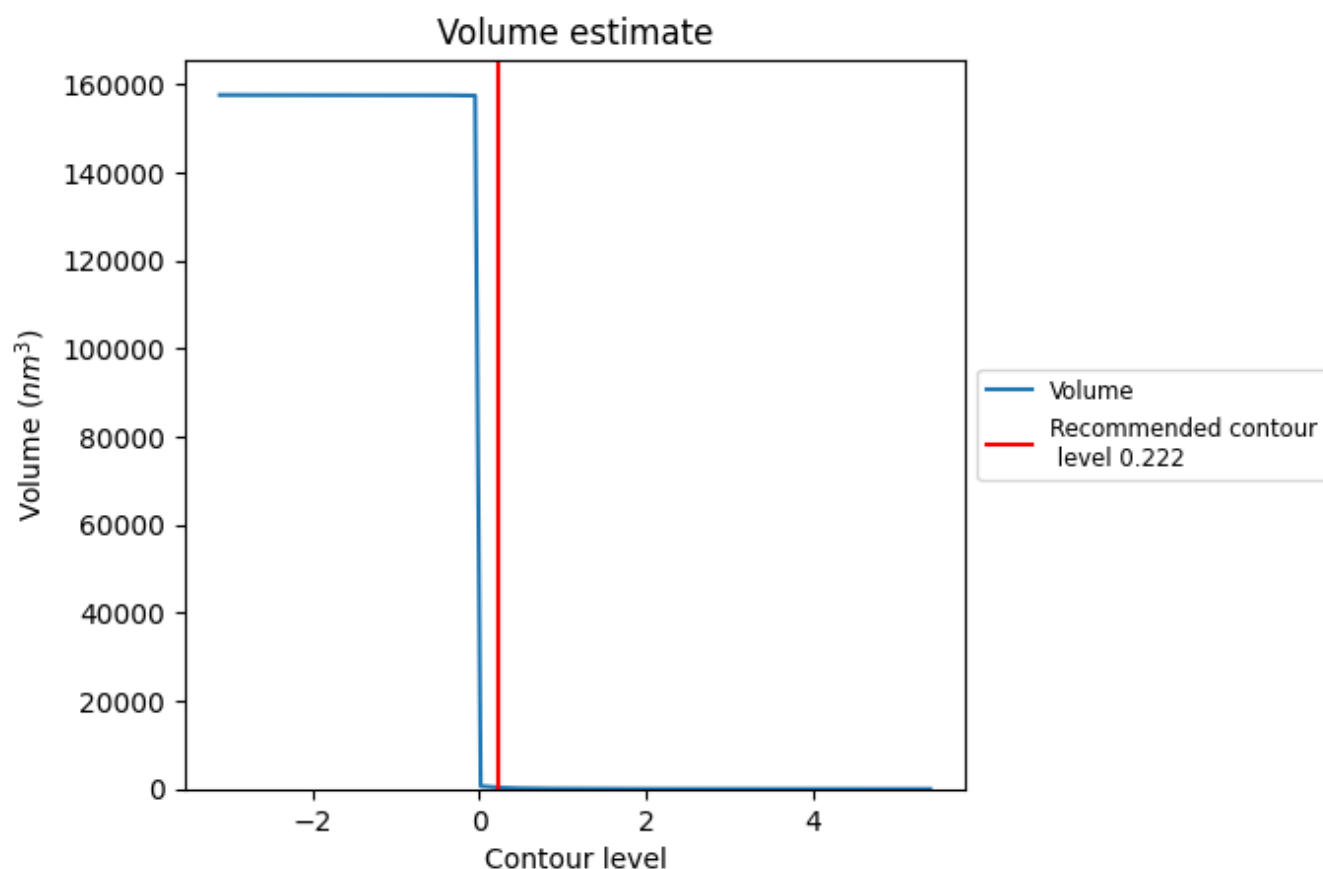
This section contains the results of statistical analysis of the map.

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.

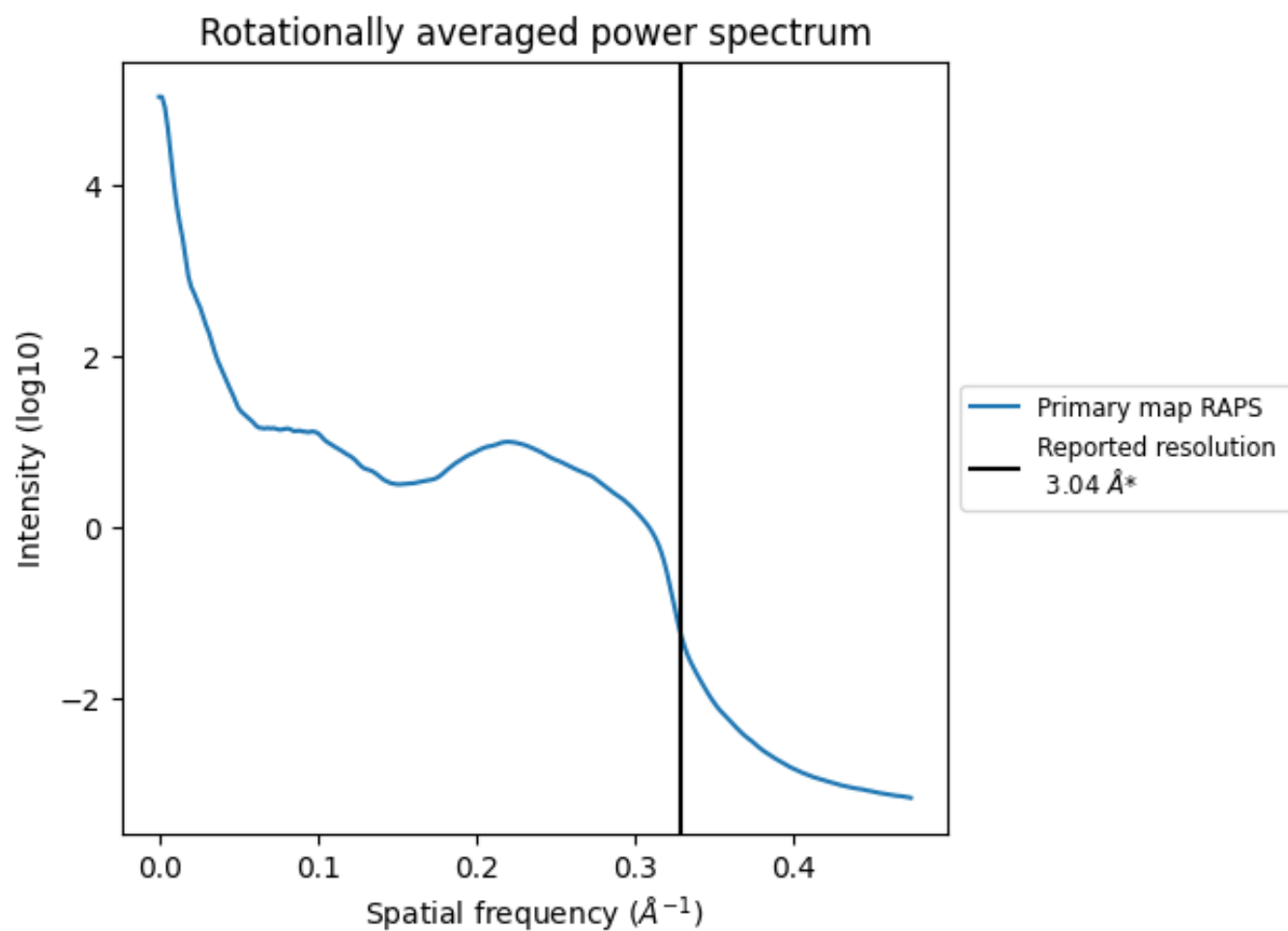
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 331 nm^3 ; this corresponds to an approximate mass of 299 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 0.329 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

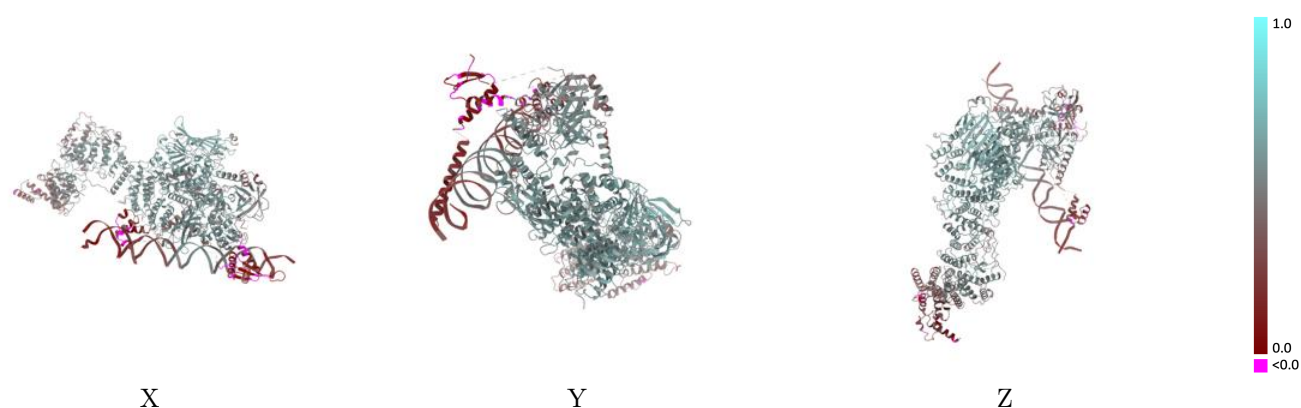
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-31117 and PDB model 7EGH. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

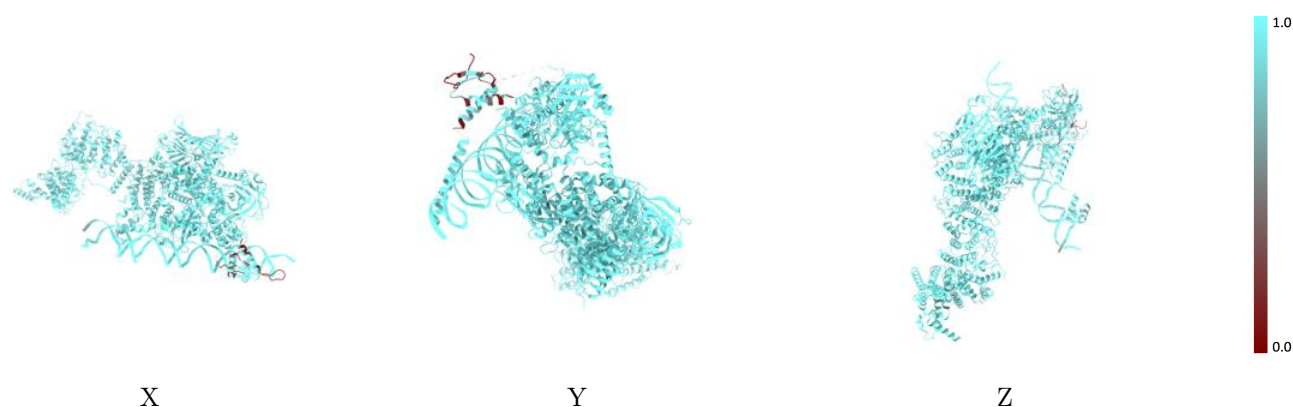
This section was not generated.

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



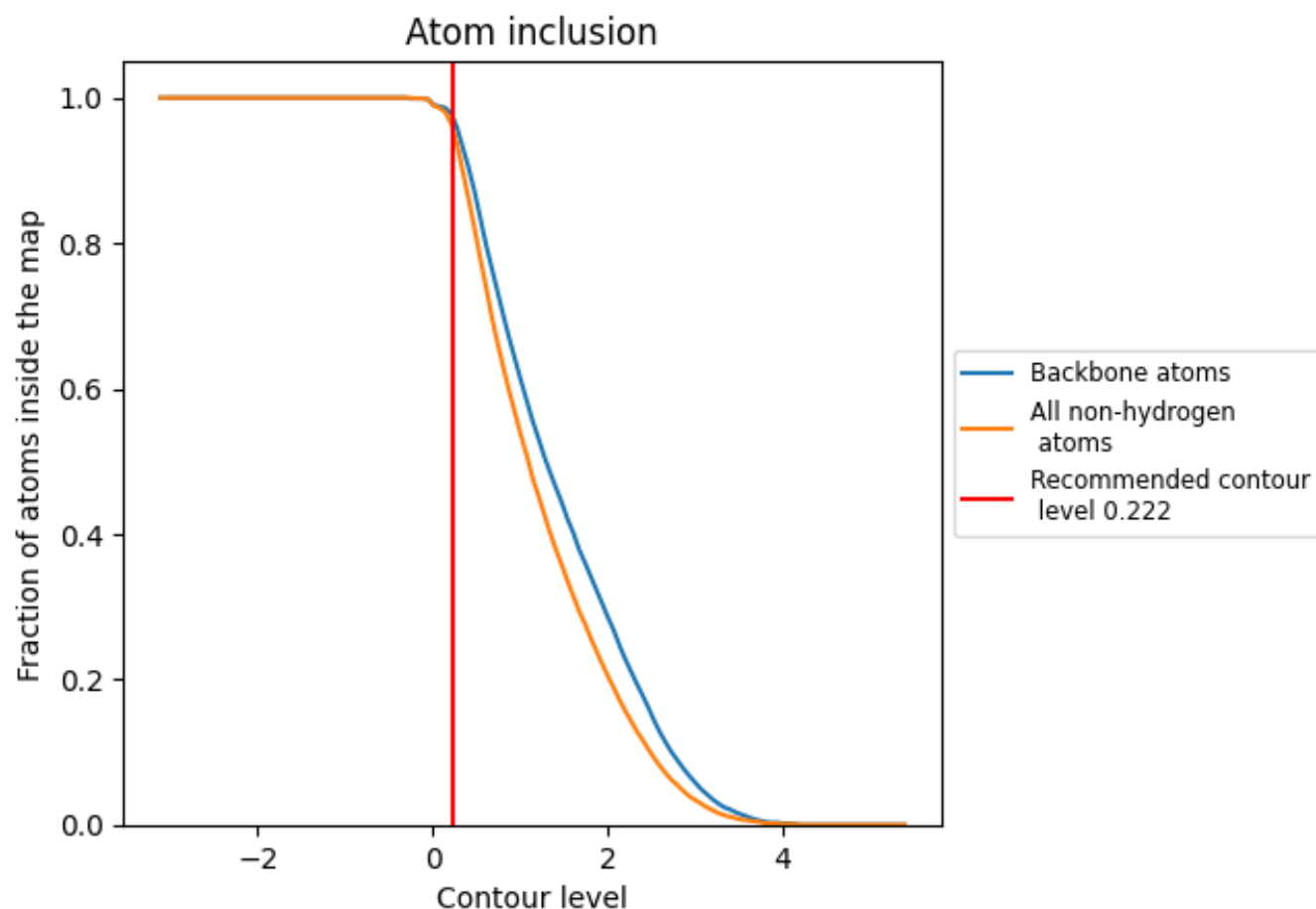
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.222).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 96% 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 (0.222) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9630	0.4810
A	0.9340	0.4360
B	0.9850	0.5790
F	0.9530	0.3690
G	0.8900	0.4040
H	0.9770	0.5400
X	0.9890	0.3420
Y	0.9800	0.3520
f	0.9790	0.4790

