



## Full wwPDB EM Validation Report ⓘ

Jan 27, 2022 – 06:21 pm GMT

PDB ID : 7QFP  
EMDB ID : EMD-13946  
Title : Cryo-EM structure of Botulinum neurotoxin serotype E  
Authors : Kosenina, S.; Martinez-Carranza, M.; Davies, J.R.; Masuyer, G.; Stenmark, P.  
Deposited on : 2021-12-06  
Resolution : 3.70 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](mailto: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.

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

EMDB validation analysis : 0.0.0.dev97  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.26

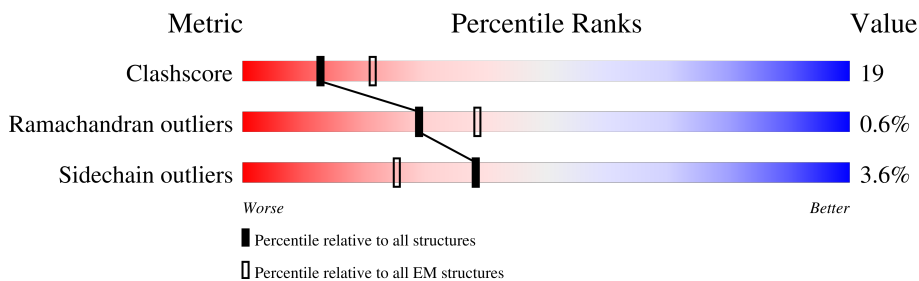
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.70 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  $\geq 3$ , 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	1269	<p>5% 69% 25% . .</p>

## 2 Entry composition i

There is only 1 type of molecule in this entry. The entry contains 9996 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 Botulinum neurotoxin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1233	9996	6363	1670	1940	23	0	0

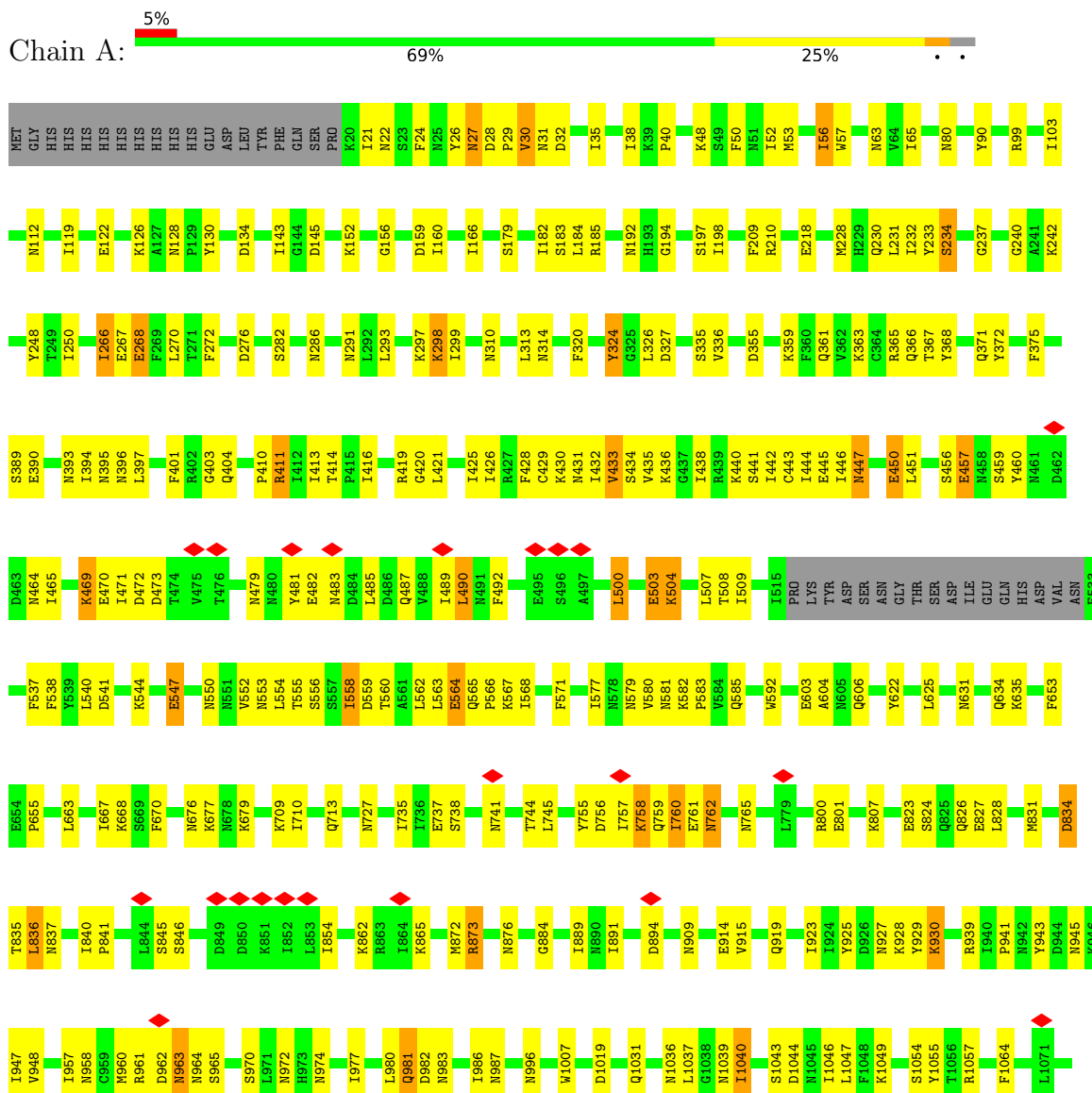
There are 21 discrepancies between the modelled and reference sequences:

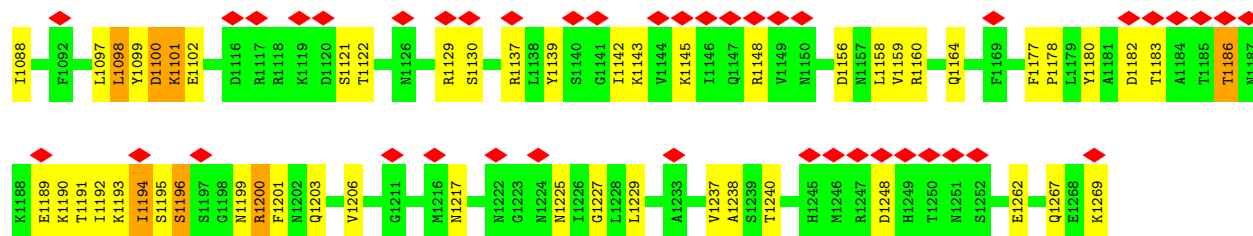
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A8Y875
A	2	GLY	-	expression tag	UNP A8Y875
A	3	HIS	-	expression tag	UNP A8Y875
A	4	HIS	-	expression tag	UNP A8Y875
A	5	HIS	-	expression tag	UNP A8Y875
A	6	HIS	-	expression tag	UNP A8Y875
A	7	HIS	-	expression tag	UNP A8Y875
A	8	HIS	-	expression tag	UNP A8Y875
A	9	HIS	-	expression tag	UNP A8Y875
A	10	HIS	-	expression tag	UNP A8Y875
A	11	HIS	-	expression tag	UNP A8Y875
A	12	GLU	-	expression tag	UNP A8Y875
A	13	ASP	-	expression tag	UNP A8Y875
A	14	LEU	-	expression tag	UNP A8Y875
A	15	TYR	-	expression tag	UNP A8Y875
A	16	PHE	-	expression tag	UNP A8Y875
A	17	GLN	-	expression tag	UNP A8Y875
A	18	SER	-	expression tag	UNP A8Y875
A	230	GLN	GLU	engineered mutation	UNP A8Y875
A	233	TYR	HIS	engineered mutation	UNP A8Y875
A	1269	LYS	-	expression tag	UNP A8Y875

### 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: Botulinum neurotoxin





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	284390	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	55	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.840	Depositor
Minimum map value	-0.397	Depositor
Average map value	-0.002	Depositor
Map value standard deviation	0.026	Depositor
Recommended contour level	0.127	Depositor
Map size (Å)	247.68001, 247.68001, 247.68001	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.86, 0.86, 0.86	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	A	0.38	0/10193	0.58	0/13808

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9996	0	9851	370	0
All	All	9996	0	9851	370	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 19.

All (370) 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:840:ILE:CD1	1:A:841:PRO:HD2	1.70	1.20
1:A:840:ILE:HD12	1:A:841:PRO:CD	1.73	1.18
1:A:758:LYS:HE3	1:A:758:LYS:H	1.20	1.03
1:A:873:ARG:HA	1:A:873:ARG:HH11	1.25	1.01
1:A:52:ILE:HB	1:A:56:ILE:HD11	1.43	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:873:ARG:HA	1:A:873:ARG:NH1	1.75	1.00
1:A:435:VAL:H	1:A:441:SER:HB2	1.29	0.96
1:A:433:VAL:HA	1:A:442:ILE:HA	1.47	0.95
1:A:930:LYS:HA	1:A:930:LYS:HE2	1.49	0.94
1:A:927:ASN:HA	1:A:1039:ASN:HA	1.51	0.93
1:A:552:VAL:HB	1:A:735:ILE:HG12	1.48	0.93
1:A:544:LYS:HD3	1:A:727:ASN:HD21	1.36	0.91
1:A:1190:LYS:HG3	1:A:1238:ALA:HB3	1.53	0.90
1:A:1098:LEU:HD23	1:A:1098:LEU:H	1.36	0.89
1:A:945:ASN:HB3	1:A:947:ILE:HD13	1.55	0.88
1:A:1100:ASP:HA	1:A:1145:LYS:HE3	1.56	0.87
1:A:504:LYS:HA	1:A:504:LYS:HZ3	1.41	0.85
1:A:504:LYS:HA	1:A:504:LYS:NZ	1.91	0.85
1:A:565:GLN:HG2	1:A:567:LYS:HE3	1.60	0.83
1:A:945:ASN:HB3	1:A:947:ILE:CD1	2.09	0.83
1:A:429:CYS:HA	1:A:445:GLU:HB2	1.61	0.83
1:A:758:LYS:H	1:A:758:LYS:CE	1.92	0.82
1:A:758:LYS:HE3	1:A:758:LYS:N	1.95	0.81
1:A:433:VAL:HG12	1:A:442:ILE:HG13	1.65	0.79
1:A:1177:PHE:HB3	1:A:1196:SER:HB2	1.65	0.78
1:A:1180:TYR:HB3	1:A:1194:ILE:HA	1.63	0.78
1:A:470:GLU:HG3	1:A:667:ILE:HG23	1.64	0.78
1:A:419:ARG:O	1:A:419:ARG:NH1	2.19	0.76
1:A:1178:PRO:HG3	1:A:1200:ARG:HE	1.51	0.75
1:A:500:LEU:HD23	1:A:500:LEU:N	1.99	0.75
1:A:1101:LYS:CE	1:A:1101:LYS:HA	2.15	0.75
1:A:554:LEU:HD11	1:A:580:VAL:HG23	1.69	0.75
1:A:267:GLU:OE2	1:A:267:GLU:N	2.17	0.75
1:A:1101:LYS:HA	1:A:1101:LYS:HE3	1.68	0.75
1:A:179:SER:HB2	1:A:233:TYR:HE2	1.51	0.73
1:A:426:ILE:HD11	1:A:451:LEU:HD21	1.70	0.73
1:A:450:GLU:CD	1:A:450:GLU:H	1.92	0.73
1:A:160:ILE:CD1	1:A:503:GLU:HA	2.18	0.72
1:A:655:PRO:HD2	1:A:710:ILE:HD11	1.71	0.72
1:A:268:GLU:HA	1:A:268:GLU:OE1	1.89	0.72
1:A:1186:THR:HG21	1:A:1189:GLU:HB2	1.72	0.71
1:A:179:SER:HB2	1:A:233:TYR:CE2	2.26	0.71
1:A:1098:LEU:HD23	1:A:1098:LEU:N	2.06	0.70
1:A:53:MET:H	1:A:56:ILE:HG12	1.56	0.70
1:A:1102:GLU:HB3	1:A:1143:LYS:HD2	1.74	0.70
1:A:433:VAL:CA	1:A:442:ILE:HA	2.23	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ARG:HD2	1:A:507:LEU:HB3	1.75	0.69
1:A:961:ARG:HD3	1:A:1043:SER:HA	1.75	0.68
1:A:872:MET:HE1	1:A:889:ILE:HG12	1.75	0.68
1:A:228:MET:HA	1:A:231:LEU:HD12	1.76	0.67
1:A:130:TYR:HA	1:A:320:PHE:HE1	1.60	0.67
1:A:555:THR:HG23	1:A:555:THR:O	1.94	0.67
1:A:807:LYS:HB2	1:A:836:LEU:HD23	1.77	0.67
1:A:421:LEU:N	1:A:421:LEU:HD12	2.10	0.66
1:A:504:LYS:H	1:A:504:LYS:HD2	1.59	0.66
1:A:555:THR:HB	1:A:568:ILE:HB	1.77	0.66
1:A:426:ILE:HB	1:A:445:GLU:OE2	1.96	0.65
1:A:1177:PHE:HB3	1:A:1196:SER:CB	2.26	0.65
1:A:426:ILE:HD12	1:A:447:ASN:HD22	1.62	0.64
1:A:230:GLN:HA	1:A:233:TYR:CE1	2.33	0.64
1:A:840:ILE:HD12	1:A:841:PRO:HD2	0.80	0.64
1:A:27:ASN:O	1:A:27:ASN:ND2	2.25	0.64
1:A:963:ASN:OD1	1:A:963:ASN:N	2.31	0.64
1:A:1200:ARG:HD2	1:A:1200:ARG:N	2.11	0.64
1:A:435:VAL:HG13	1:A:436:LYS:HD3	1.80	0.64
1:A:1190:LYS:CG	1:A:1238:ALA:HB3	2.26	0.64
1:A:371:GLN:OE1	1:A:371:GLN:N	2.25	0.64
1:A:550:ASN:HB3	1:A:738:SER:HB2	1.81	0.63
1:A:159:ASP:HB3	1:A:500:LEU:HD22	1.80	0.63
1:A:552:VAL:CB	1:A:735:ILE:HG12	2.23	0.63
1:A:436:LYS:HG2	1:A:440:LYS:HE2	1.81	0.63
1:A:440:LYS:HE3	1:A:560:THR:HG22	1.80	0.63
1:A:504:LYS:HD2	1:A:504:LYS:N	2.14	0.63
1:A:433:VAL:HA	1:A:442:ILE:CA	2.25	0.62
1:A:429:CYS:CA	1:A:445:GLU:HB2	2.29	0.62
1:A:487:GLN:O	1:A:487:GLN:NE2	2.33	0.62
1:A:421:LEU:HD12	1:A:421:LEU:H	1.65	0.61
1:A:558:ILE:HG13	1:A:560:THR:HG23	1.81	0.61
1:A:465:ILE:HG23	1:A:663:LEU:HD13	1.83	0.61
1:A:324:TYR:CD1	1:A:324:TYR:N	2.69	0.61
1:A:429:CYS:HA	1:A:445:GLU:CB	2.30	0.60
1:A:447:ASN:OD1	1:A:447:ASN:N	2.32	0.60
1:A:160:ILE:HD12	1:A:503:GLU:HA	1.83	0.60
1:A:1189:GLU:OE2	1:A:1189:GLU:N	2.35	0.60
1:A:930:LYS:HE2	1:A:930:LYS:CA	2.28	0.60
1:A:986:ILE:N	1:A:986:ILE:HD12	2.16	0.60
1:A:1206:VAL:HB	1:A:1217:ASN:HB3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:ILE:HG23	1:A:40:PRO:HD3	1.84	0.59
1:A:397:LEU:HB3	1:A:403:GLY:CA	2.32	0.59
1:A:160:ILE:HG21	1:A:504:LYS:HE2	1.84	0.59
1:A:210:ARG:NH1	1:A:218:GLU:OE1	2.36	0.59
1:A:1099:TYR:CZ	1:A:1159:VAL:HG22	2.37	0.59
1:A:430:LYS:HB3	1:A:444:ILE:H	1.68	0.59
1:A:930:LYS:NZ	1:A:930:LYS:O	2.36	0.58
1:A:919:GLN:OE1	1:A:1044:ASP:HA	2.04	0.58
1:A:420:GLY:H	1:A:421:LEU:HD12	1.68	0.58
1:A:872:MET:HG3	1:A:1064:PHE:HB3	1.85	0.58
1:A:563:LEU:C	1:A:565:GLN:N	2.56	0.58
1:A:958:ASN:HB3	1:A:1047:LEU:HB3	1.85	0.58
1:A:396:ASN:N	1:A:396:ASN:OD1	2.36	0.58
1:A:1203:GLN:HA	1:A:1203:GLN:NE2	2.19	0.57
1:A:126:LYS:HG2	1:A:492:PHE:HB2	1.85	0.57
1:A:440:LYS:HD2	1:A:560:THR:HA	1.86	0.57
1:A:485:LEU:C	1:A:485:LEU:HD23	2.24	0.57
1:A:552:VAL:HG11	1:A:735:ILE:HA	1.87	0.56
1:A:56:ILE:HG13	1:A:56:ILE:O	2.04	0.56
1:A:411:ARG:NH1	1:A:411:ARG:H	2.02	0.56
1:A:482:GLU:HG3	1:A:483:ASN:H	1.70	0.56
1:A:435:VAL:N	1:A:441:SER:HB2	2.09	0.56
1:A:26:TYR:HA	1:A:50:PHE:CZ	2.40	0.56
1:A:272:PHE:HB2	1:A:368:TYR:CZ	2.40	0.56
1:A:435:VAL:HG22	1:A:436:LYS:HD2	1.87	0.56
1:A:359:LYS:O	1:A:361:GLN:NE2	2.39	0.55
1:A:411:ARG:HH11	1:A:411:ARG:HG2	1.70	0.55
1:A:960:MET:HB3	1:A:964:ASN:HA	1.87	0.55
1:A:986:ILE:HD12	1:A:986:ILE:H	1.72	0.55
1:A:1190:LYS:HB2	1:A:1237:VAL:HG12	1.88	0.55
1:A:297:LYS:HD2	1:A:336:VAL:HG21	1.89	0.55
1:A:563:LEU:C	1:A:565:GLN:H	2.09	0.55
1:A:631:ASN:OD1	1:A:634:GLN:NE2	2.40	0.55
1:A:26:TYR:HA	1:A:50:PHE:HZ	1.71	0.55
1:A:504:LYS:NZ	1:A:504:LYS:CA	2.67	0.55
1:A:490:LEU:HD22	1:A:490:LEU:N	2.23	0.54
1:A:562:LEU:HD23	1:A:562:LEU:H	1.72	0.54
1:A:119:ILE:HG23	1:A:489:ILE:HG22	1.88	0.54
1:A:429:CYS:HA	1:A:445:GLU:CG	2.38	0.54
1:A:450:GLU:CD	1:A:450:GLU:N	2.60	0.54
1:A:53:MET:HG2	1:A:56:ILE:HD13	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:762:ASN:H	1:A:762:ASN:ND2	2.04	0.54
1:A:891:ILE:HA	1:A:915:VAL:HG12	1.90	0.54
1:A:160:ILE:HD11	1:A:503:GLU:HA	1.89	0.54
1:A:945:ASN:O	1:A:948:VAL:HG12	2.07	0.54
1:A:21:ILE:HD11	1:A:112:ASN:HD22	1.73	0.54
1:A:1040:ILE:HD12	1:A:1040:ILE:N	2.23	0.54
1:A:419:ARG:C	1:A:419:ARG:HH11	2.11	0.53
1:A:390:GLU:HG3	1:A:393:ASN:O	2.08	0.53
1:A:160:ILE:HD13	1:A:504:LYS:HD2	1.90	0.53
1:A:209:PHE:HB2	1:A:372:TYR:HB2	1.89	0.53
1:A:416:ILE:HG13	1:A:421:LEU:HD21	1.90	0.53
1:A:397:LEU:HB3	1:A:403:GLY:HA2	1.90	0.53
1:A:425:ILE:HD12	1:A:445:GLU:HG3	1.89	0.53
1:A:537:PHE:CD1	1:A:540:LEU:HD12	2.44	0.53
1:A:909:ASN:ND2	1:A:1055:TYR:O	2.42	0.53
1:A:159:ASP:CB	1:A:500:LEU:HD22	2.39	0.53
1:A:438:ILE:HB	1:A:442:ILE:HD12	1.91	0.52
1:A:250:ILE:HA	1:A:459:SER:HB3	1.90	0.52
1:A:479:ASN:ND2	1:A:481:TYR:OH	2.43	0.52
1:A:429:CYS:HA	1:A:445:GLU:HG2	1.90	0.52
1:A:577:ILE:O	1:A:581:ASN:HB2	2.10	0.52
1:A:755:TYR:HA	1:A:759:GLN:HB3	1.92	0.52
1:A:894:ASP:OD1	1:A:894:ASP:N	2.42	0.52
1:A:1019:ASP:OD2	1:A:1031:GLN:NE2	2.38	0.52
1:A:541:ASP:HA	1:A:544:LYS:HG2	1.92	0.52
1:A:919:GLN:NE2	1:A:923:ILE:O	2.42	0.52
1:A:145:ASP:HA	1:A:504:LYS:HG3	1.92	0.52
1:A:470:GLU:CG	1:A:667:ILE:HA	2.40	0.52
1:A:655:PRO:HG3	1:A:709:LYS:HE2	1.91	0.52
1:A:160:ILE:HD13	1:A:504:LYS:CD	2.39	0.52
1:A:267:GLU:HA	1:A:270:LEU:HD23	1.92	0.52
1:A:471:ILE:HG12	1:A:679:LYS:CD	2.40	0.52
1:A:469:LYS:HD2	1:A:668:LYS:HB2	1.92	0.51
1:A:862:LYS:HE3	1:A:865:LYS:HB2	1.91	0.51
1:A:930:LYS:HD3	1:A:930:LYS:C	2.31	0.51
1:A:197:SER:OG	1:A:198:ILE:N	2.44	0.51
1:A:759:GLN:O	1:A:761:GLU:N	2.43	0.51
1:A:872:MET:CE	1:A:889:ILE:HG12	2.40	0.51
1:A:1122:THR:HB	1:A:1193:LYS:NZ	2.25	0.51
1:A:431:ASN:O	1:A:443:CYS:SG	2.67	0.51
1:A:185:ARG:CD	1:A:507:LEU:HB3	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ILE:HA	1:A:459:SER:CB	2.41	0.51
1:A:552:VAL:HG23	1:A:552:VAL:O	2.10	0.51
1:A:419:ARG:O	1:A:419:ARG:CG	2.59	0.50
1:A:1180:TYR:CB	1:A:1195:SER:H	2.25	0.50
1:A:57:TRP:HB2	1:A:166:ILE:HG22	1.94	0.50
1:A:559:ASP:HA	1:A:566:PRO:HB3	1.94	0.50
1:A:737:GLU:O	1:A:741:ASN:ND2	2.44	0.50
1:A:429:CYS:SG	1:A:430:LYS:N	2.84	0.50
1:A:760:ILE:N	1:A:760:ILE:HD12	2.27	0.50
1:A:192:ASN:ND2	1:A:242:LYS:O	2.45	0.50
1:A:554:LEU:CD1	1:A:580:VAL:HG23	2.40	0.50
1:A:1101:LYS:HE3	1:A:1101:LYS:CA	2.31	0.50
1:A:130:TYR:HA	1:A:320:PHE:CE1	2.42	0.49
1:A:446:ILE:C	1:A:446:ILE:HD12	2.32	0.49
1:A:1130:SER:O	1:A:1137:ARG:NH1	2.45	0.49
1:A:395:ASN:OD1	1:A:395:ASN:N	2.45	0.49
1:A:553:ASN:HB3	1:A:564:GLU:O	2.11	0.49
1:A:1055:TYR:O	1:A:1057:ARG:NH1	2.46	0.49
1:A:1199:ASN:HB3	1:A:1201:PHE:CE2	2.46	0.49
1:A:1190:LYS:O	1:A:1237:VAL:HG11	2.13	0.49
1:A:547:GLU:H	1:A:547:GLU:CD	2.15	0.49
1:A:552:VAL:HG12	1:A:565:GLN:HE22	1.78	0.49
1:A:470:GLU:HG3	1:A:667:ILE:HA	1.94	0.49
1:A:1190:LYS:NZ	1:A:1240:THR:HG23	2.28	0.49
1:A:183:SER:OG	1:A:184:LEU:N	2.46	0.49
1:A:555:THR:HG21	1:A:577:ILE:HD13	1.95	0.49
1:A:1156:ASP:OD2	1:A:1160:ARG:NH2	2.45	0.49
1:A:429:CYS:SG	1:A:445:GLU:HB2	2.53	0.48
1:A:134:ASP:N	1:A:134:ASP:OD1	2.45	0.48
1:A:248:TYR:HB2	1:A:266:ILE:HD13	1.95	0.48
1:A:547:GLU:OE1	1:A:547:GLU:N	2.31	0.48
1:A:579:ASN:O	1:A:582:LYS:NZ	2.43	0.48
1:A:927:ASN:HA	1:A:1039:ASN:CA	2.35	0.48
1:A:160:ILE:HG21	1:A:504:LYS:CE	2.44	0.48
1:A:182:ILE:CG2	1:A:508:THR:HG23	2.43	0.48
1:A:393:ASN:HD21	1:A:404:GLN:HE22	1.62	0.48
1:A:1040:ILE:HD12	1:A:1040:ILE:H	1.77	0.48
1:A:552:VAL:HG21	1:A:735:ILE:HA	1.96	0.48
1:A:854:ILE:HD11	1:A:876:ASN:HD21	1.79	0.48
1:A:761:GLU:OE1	1:A:761:GLU:HA	2.14	0.47
1:A:119:ILE:HG23	1:A:489:ILE:CG2	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ILE:HG23	1:A:504:LYS:HZ3	1.78	0.47
1:A:152:LYS:HD2	1:A:156:GLY:HA2	1.96	0.47
1:A:428:PHE:CZ	1:A:431:ASN:HA	2.49	0.47
1:A:568:ILE:CG2	1:A:571:PHE:H	2.26	0.47
1:A:1102:GLU:H	1:A:1102:GLU:CD	2.15	0.47
1:A:962:ASP:OD2	1:A:962:ASP:N	2.45	0.47
1:A:185:ARG:HD2	1:A:507:LEU:CB	2.42	0.47
1:A:970:SER:HB3	1:A:977:ILE:HB	1.97	0.47
1:A:456:SER:OG	1:A:457:GLU:N	2.46	0.47
1:A:433:VAL:HG11	1:A:438:ILE:HG22	1.95	0.47
1:A:583:PRO:O	1:A:585:GLN:NE2	2.47	0.47
1:A:831:MET:HA	1:A:834:ASP:HB3	1.97	0.47
1:A:845:SER:OG	1:A:846:SER:N	2.47	0.47
1:A:981:GLN:HA	1:A:987:ASN:OD1	2.14	0.47
1:A:28:ASP:H	1:A:48:LYS:NZ	2.13	0.47
1:A:552:VAL:HG12	1:A:565:GLN:NE2	2.30	0.47
1:A:744:THR:OG1	1:A:745:LEU:N	2.47	0.47
1:A:957:ILE:HG23	1:A:1046:ILE:HG23	1.97	0.47
1:A:365:ARG:NH1	1:A:365:ARG:HB2	2.29	0.47
1:A:925:TYR:HA	1:A:929:TYR:OH	2.15	0.47
1:A:972:ASN:HB2	1:A:977:ILE:HD11	1.96	0.47
1:A:982:ASP:OD1	1:A:982:ASP:N	2.46	0.47
1:A:80:ASN:OD1	1:A:80:ASN:N	2.48	0.46
1:A:365:ARG:HB2	1:A:365:ARG:HH11	1.80	0.46
1:A:421:LEU:H	1:A:421:LEU:CD1	2.28	0.46
1:A:430:LYS:HE2	1:A:538:PHE:CD1	2.49	0.46
1:A:436:LYS:HB2	1:A:440:LYS:HB2	1.96	0.46
1:A:504:LYS:HA	1:A:504:LYS:HZ2	1.76	0.46
1:A:1192:ILE:HG12	1:A:1237:VAL:HG13	1.98	0.46
1:A:568:ILE:HG22	1:A:571:PHE:H	1.79	0.46
1:A:800:ARG:NH1	1:A:801:GLU:OE2	2.48	0.46
1:A:563:LEU:O	1:A:565:GLN:N	2.48	0.46
1:A:1248:ASP:OD1	1:A:1248:ASP:N	2.49	0.46
1:A:30:VAL:HG12	1:A:30:VAL:O	2.16	0.46
1:A:1190:LYS:C	1:A:1237:VAL:HG11	2.36	0.46
1:A:30:VAL:C	1:A:32:ASP:H	2.19	0.46
1:A:927:ASN:ND2	1:A:1039:ASN:HB2	2.30	0.46
1:A:30:VAL:O	1:A:32:ASP:N	2.49	0.46
1:A:945:ASN:HB3	1:A:947:ILE:HD12	1.93	0.45
1:A:122:GLU:OE1	1:A:126:LYS:NZ	2.50	0.45
1:A:507:LEU:HD23	1:A:509:ILE:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:925:TYR:HE2	1:A:928:LYS:HA	1.81	0.45
1:A:603:GLU:O	1:A:606:GLN:NE2	2.49	0.45
1:A:430:LYS:HB3	1:A:444:ILE:N	2.31	0.45
1:A:757:ILE:HD13	1:A:757:ILE:N	2.31	0.45
1:A:445:GLU:CG	1:A:445:GLU:O	2.65	0.45
1:A:421:LEU:N	1:A:421:LEU:CD1	2.79	0.45
1:A:756:ASP:OD1	1:A:756:ASP:N	2.48	0.45
1:A:537:PHE:HA	1:A:540:LEU:HB2	1.98	0.45
1:A:428:PHE:O	1:A:428:PHE:CD1	2.69	0.45
1:A:469:LYS:HD3	1:A:469:LYS:HA	1.53	0.45
1:A:291:ASN:N	1:A:291:ASN:HD22	2.15	0.45
1:A:297:LYS:NZ	1:A:336:VAL:HB	2.32	0.45
1:A:961:ARG:HD2	1:A:961:ARG:HA	1.77	0.44
1:A:559:ASP:OD2	1:A:566:PRO:HG3	2.17	0.44
1:A:571:PHE:O	1:A:571:PHE:CD2	2.70	0.44
1:A:1046:ILE:N	1:A:1046:ILE:HD12	2.33	0.44
1:A:145:ASP:OD1	1:A:145:ASP:N	2.49	0.44
1:A:974:ASN:OD1	1:A:996:ASN:ND2	2.51	0.44
1:A:1088:ILE:HD11	1:A:1158:LEU:HD13	2.00	0.44
1:A:1148:ARG:HD3	1:A:1148:ARG:HA	1.86	0.44
1:A:185:ARG:NH1	1:A:507:LEU:H	2.15	0.44
1:A:556:SER:HB2	1:A:558:ILE:HG13	1.99	0.44
1:A:826:GLN:NE2	1:A:827:GLU:OE2	2.51	0.44
1:A:1183:THR:O	1:A:1191:THR:HB	2.18	0.44
1:A:24:PHE:HB3	1:A:50:PHE:CE2	2.52	0.44
1:A:363:LYS:H	1:A:401:PHE:HZ	1.66	0.44
1:A:1054:SER:H	1:A:1057:ARG:HH22	1.65	0.44
1:A:53:MET:HB2	1:A:56:ILE:HG21	1.98	0.44
1:A:411:ARG:CZ	1:A:411:ARG:HB3	2.47	0.44
1:A:433:VAL:CG1	1:A:442:ILE:HG13	2.43	0.44
1:A:1227:GLY:HA2	1:A:1240:THR:HG22	1.98	0.44
1:A:293:LEU:O	1:A:297:LYS:HG2	2.18	0.44
1:A:276:ASP:N	1:A:276:ASP:OD1	2.51	0.43
1:A:930:LYS:CA	1:A:930:LYS:CE	2.92	0.43
1:A:90:TYR:OH	1:A:99:ARG:NH2	2.51	0.43
1:A:755:TYR:HA	1:A:759:GLN:CB	2.47	0.43
1:A:327:ASP:O	1:A:335:SER:N	2.50	0.43
1:A:552:VAL:CA	1:A:567:LYS:HZ1	2.31	0.43
1:A:556:SER:HB3	1:A:564:GLU:HG3	1.99	0.43
1:A:22:ASN:ND2	1:A:52:ILE:O	2.52	0.43
1:A:445:GLU:HG3	1:A:445:GLU:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1142:ILE:H	1:A:1142:ILE:HG13	1.54	0.43
1:A:366:GLN:OE1	1:A:366:GLN:HA	2.18	0.43
1:A:435:VAL:HG13	1:A:440:LYS:HB2	2.00	0.43
1:A:1121:SER:O	1:A:1121:SER:OG	2.32	0.43
1:A:63:ASN:HB2	1:A:65:ILE:HG22	1.99	0.43
1:A:500:LEU:N	1:A:500:LEU:CD2	2.70	0.43
1:A:824:SER:HB2	1:A:828:LEU:HD22	2.00	0.43
1:A:355:ASP:OD1	1:A:355:ASP:N	2.49	0.43
1:A:465:ILE:CG2	1:A:663:LEU:HD13	2.48	0.43
1:A:487:GLN:O	1:A:487:GLN:CD	2.57	0.43
1:A:835:THR:C	1:A:837:ASN:H	2.22	0.43
1:A:130:TYR:HE2	1:A:299:ILE:HG13	1.84	0.42
1:A:552:VAL:HA	1:A:567:LYS:HZ1	1.83	0.42
1:A:980:LEU:HD21	1:A:1037:LEU:HD11	2.00	0.42
1:A:1049:LYS:HA	1:A:1049:LYS:HD3	1.83	0.42
1:A:986:ILE:HG21	1:A:1036:ASN:O	2.19	0.42
1:A:53:MET:N	1:A:56:ILE:HG12	2.29	0.42
1:A:419:ARG:O	1:A:419:ARG:HG3	2.19	0.42
1:A:556:SER:HB3	1:A:564:GLU:CD	2.40	0.42
1:A:580:VAL:HA	1:A:592:TRP:HZ2	1.84	0.42
1:A:677:LYS:NZ	1:A:823:GLU:OE2	2.46	0.42
1:A:625:LEU:HA	1:A:635:LYS:HG2	2.01	0.42
1:A:1139:TYR:CE1	1:A:1262:GLU:HG2	2.54	0.42
1:A:326:LEU:CD2	1:A:336:VAL:HG22	2.49	0.42
1:A:389:SER:OG	1:A:390:GLU:OE1	2.35	0.42
1:A:914:GLU:HB3	1:A:1049:LYS:HB2	2.01	0.42
1:A:28:ASP:HB2	1:A:35:ILE:HD13	2.00	0.42
1:A:1097:LEU:HD23	1:A:1159:VAL:HG21	2.01	0.42
1:A:128:ASN:N	1:A:128:ASN:OD1	2.53	0.42
1:A:604:ALA:O	1:A:622:TYR:OH	2.29	0.42
1:A:653:PHE:O	1:A:713:GLN:NE2	2.53	0.42
1:A:965:SER:OG	1:A:983:ASN:HA	2.20	0.42
1:A:298:LYS:HD2	1:A:298:LYS:C	2.41	0.42
1:A:159:ASP:N	1:A:159:ASP:OD1	2.52	0.42
1:A:1164:GLN:HG2	1:A:1206:VAL:HG13	2.02	0.41
1:A:272:PHE:HB2	1:A:368:TYR:OH	2.21	0.41
1:A:103:ILE:HD13	1:A:103:ILE:HA	1.94	0.41
1:A:394:ILE:N	1:A:394:ILE:HD13	2.36	0.41
1:A:419:ARG:C	1:A:419:ARG:HD2	2.40	0.41
1:A:507:LEU:HD23	1:A:509:ILE:H	1.85	0.41
1:A:541:ASP:HA	1:A:544:LYS:CG	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1129:ARG:NH1	1:A:1137:ARG:O	2.53	0.41
1:A:1229:LEU:HA	1:A:1238:ALA:HA	2.02	0.41
1:A:1267:GLN:O	1:A:1269:LYS:NZ	2.43	0.41
1:A:52:ILE:HG22	1:A:53:MET:SD	2.61	0.41
1:A:375:PHE:HD1	1:A:375:PHE:H	1.69	0.41
1:A:504:LYS:N	1:A:504:LYS:CD	2.83	0.41
1:A:930:LYS:C	1:A:930:LYS:CD	2.88	0.41
1:A:1225:ASN:OD1	1:A:1225:ASN:N	2.48	0.41
1:A:28:ASP:H	1:A:48:LYS:HZ3	1.69	0.41
1:A:143:ILE:HG23	1:A:504:LYS:NZ	2.36	0.41
1:A:185:ARG:CD	1:A:507:LEU:HD22	2.51	0.41
1:A:282:SER:O	1:A:286:ASN:ND2	2.53	0.41
1:A:464:ASN:OD1	1:A:464:ASN:N	2.54	0.41
1:A:482:GLU:HG3	1:A:483:ASN:N	2.33	0.41
1:A:562:LEU:HD23	1:A:562:LEU:N	2.35	0.41
1:A:762:ASN:HA	1:A:765:ASN:HD22	1.85	0.41
1:A:1186:THR:CG2	1:A:1189:GLU:HB2	2.48	0.41
1:A:1192:ILE:HG12	1:A:1237:VAL:CG1	2.51	0.41
1:A:194:GLY:HA2	1:A:240:GLY:HA3	2.03	0.41
1:A:862:LYS:CE	1:A:865:LYS:HB2	2.51	0.41
1:A:228:MET:O	1:A:232:ILE:HG12	2.21	0.40
1:A:410:PRO:HD2	1:A:411:ARG:HH12	1.86	0.40
1:A:670:PHE:O	1:A:676:ASN:ND2	2.54	0.40
1:A:28:ASP:HB3	1:A:29:PRO:HD2	2.03	0.40
1:A:234:SER:O	1:A:237:GLY:N	2.55	0.40
1:A:266:ILE:H	1:A:266:ILE:HG12	1.57	0.40
1:A:272:PHE:HE2	1:A:367:THR:HA	1.86	0.40
1:A:1190:LYS:HZ2	1:A:1190:LYS:HG2	1.57	0.40
1:A:436:LYS:HE2	1:A:440:LYS:HG2	2.03	0.40
1:A:862:LYS:HG3	1:A:884:GLY:HA3	2.04	0.40
1:A:434:SER:HB3	1:A:441:SER:HB3	2.03	0.40
1:A:755:TYR:HD1	1:A:760:ILE:HD13	1.85	0.40
1:A:1180:TYR:CZ	1:A:1182:ASP:HB3	2.57	0.40
1:A:1200:ARG:HD2	1:A:1200:ARG:H	1.85	0.40
1:A:310:ASN:HB3	1:A:313:LEU:HB2	2.03	0.40
1:A:928:LYS:O	1:A:928:LYS:CG	2.70	0.40
1:A:939:ARG:HD3	1:A:1007:TRP:CE2	2.56	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	1229/1269 (97%)	1068 (87%)	154 (12%)	7 (1%)	25 62

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	473	ASP
1	A	760	ILE
1	A	472	ASP
1	A	941	PRO
1	A	564	GLU
1	A	30	VAL

### 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	1135/1169 (97%)	1094 (96%)	41 (4%)	35 63

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	56	ILE
1	A	234	SER
1	A	266	ILE

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	268	GLU
1	A	298	LYS
1	A	314	ASN
1	A	324	TYR
1	A	411	ARG
1	A	413	ILE
1	A	414	THR
1	A	432	ILE
1	A	433	VAL
1	A	447	ASN
1	A	450	GLU
1	A	457	GLU
1	A	460	TYR
1	A	469	LYS
1	A	490	LEU
1	A	500	LEU
1	A	503	GLU
1	A	504	LYS
1	A	547	GLU
1	A	558	ILE
1	A	758	LYS
1	A	762	ASN
1	A	834	ASP
1	A	836	LEU
1	A	873	ARG
1	A	930	LYS
1	A	943	TYR
1	A	963	ASN
1	A	981	GLN
1	A	1040	ILE
1	A	1098	LEU
1	A	1100	ASP
1	A	1101	LYS
1	A	1186	THR
1	A	1194	ILE
1	A	1196	SER
1	A	1200	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	192	ASN

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Mol	Chain	Res	Type
1	A	217	ASN
1	A	229	HIS
1	A	236	HIS
1	A	286	ASN
1	A	291	ASN
1	A	310	ASN
1	A	382	ASN
1	A	393	ASN
1	A	400	ASN
1	A	407	ASN
1	A	448	ASN
1	A	458	ASN
1	A	479	ASN
1	A	487	GLN
1	A	585	GLN
1	A	634	GLN
1	A	685	ASN
1	A	711	ASN
1	A	715	ASN
1	A	727	ASN
1	A	762	ASN
1	A	765	ASN
1	A	805	ASN
1	A	837	ASN
1	A	871	ASN
1	A	876	ASN
1	A	942	ASN
1	A	945	ASN
1	A	952	ASN
1	A	974	ASN
1	A	996	ASN
1	A	1025	ASN
1	A	1082	ASN
1	A	1203	GLN
1	A	1212	ASN
1	A	1224	ASN
1	A	1251	ASN

### 5.3.3 RNA

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.

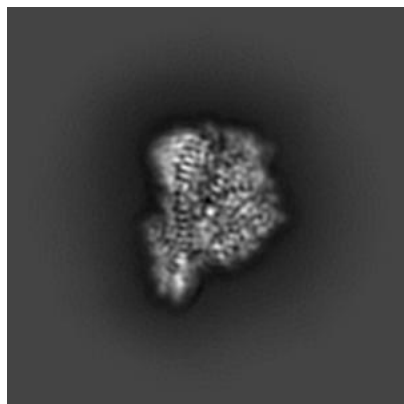
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13946. These allow visual inspection of the internal detail of the map and identification of artifacts.

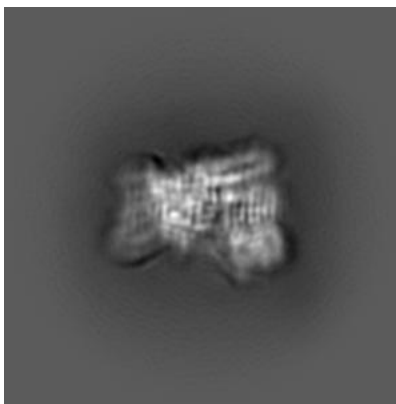
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

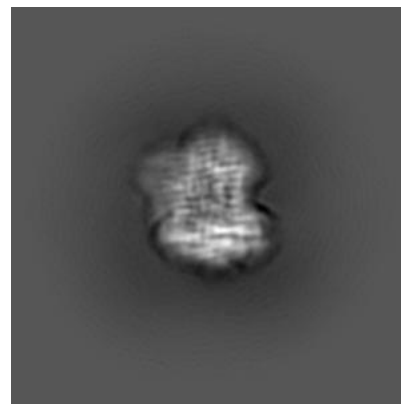
#### 6.1.1 Primary map



X

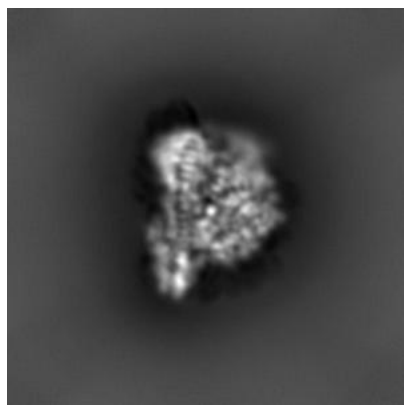


Y

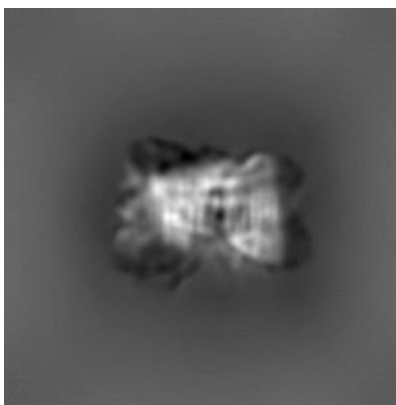


Z

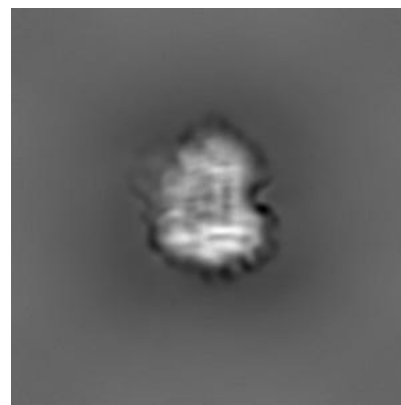
#### 6.1.2 Raw map



X



Y

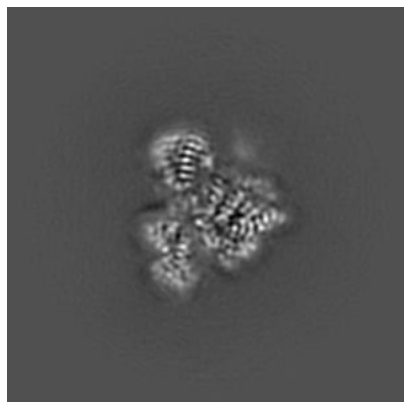


Z

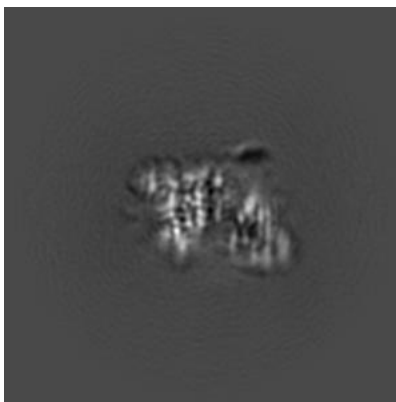
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

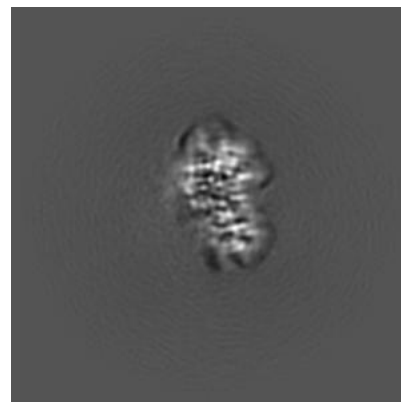
### 6.2.1 Primary map



X Index: 144

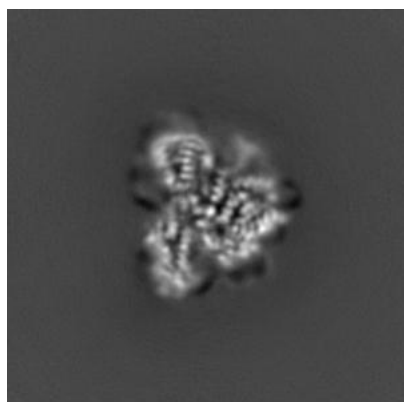


Y Index: 144



Z Index: 144

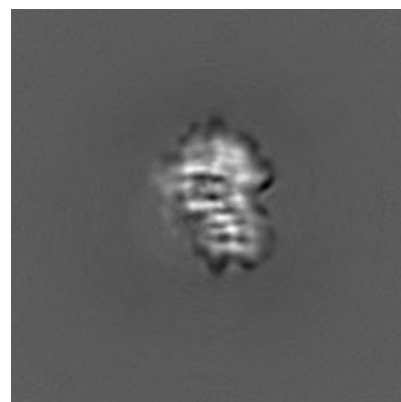
### 6.2.2 Raw map



X Index: 144



Y Index: 144

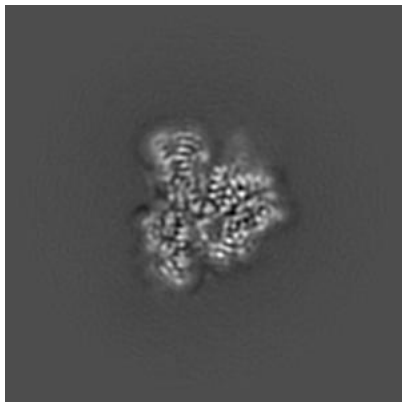


Z Index: 144

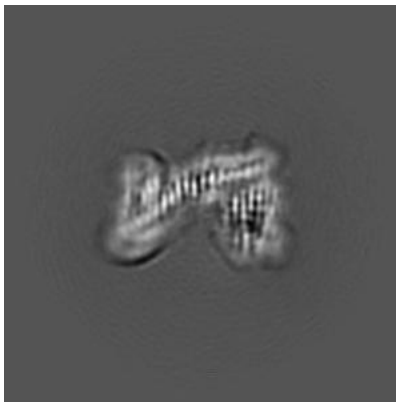
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

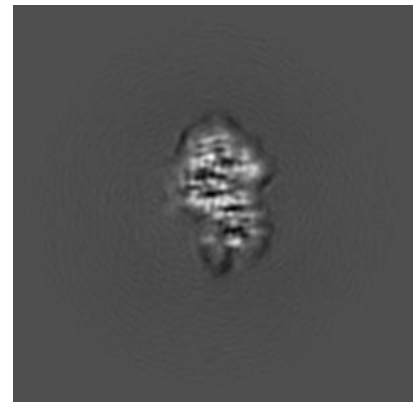
### 6.3.1 Primary map



X Index: 150

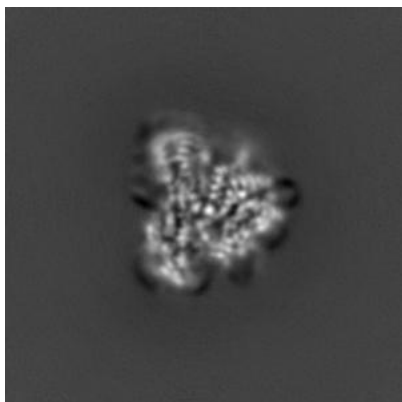


Y Index: 126

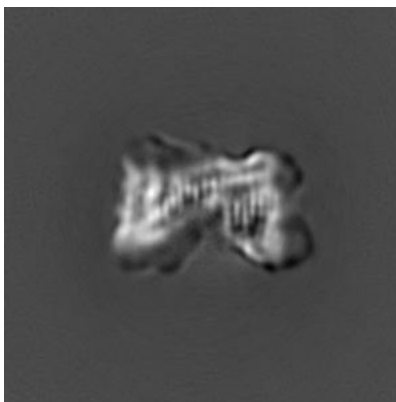


Z Index: 140

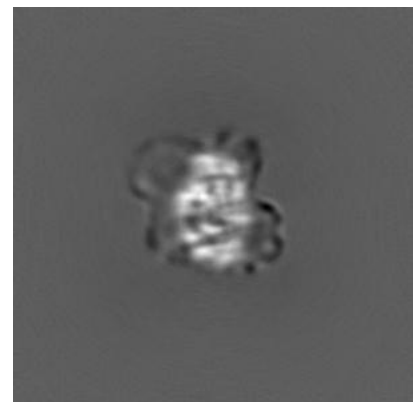
### 6.3.2 Raw map



X Index: 149



Y Index: 126

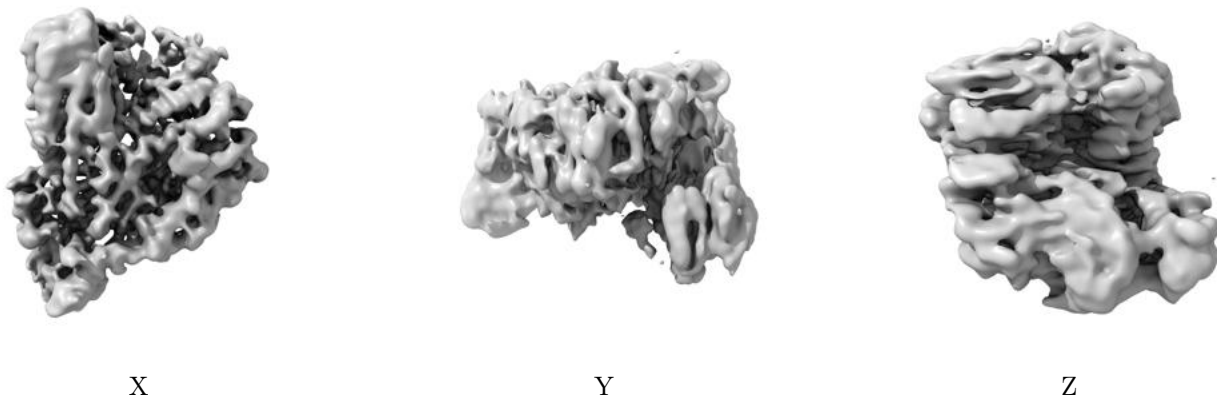


Z Index: 116

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

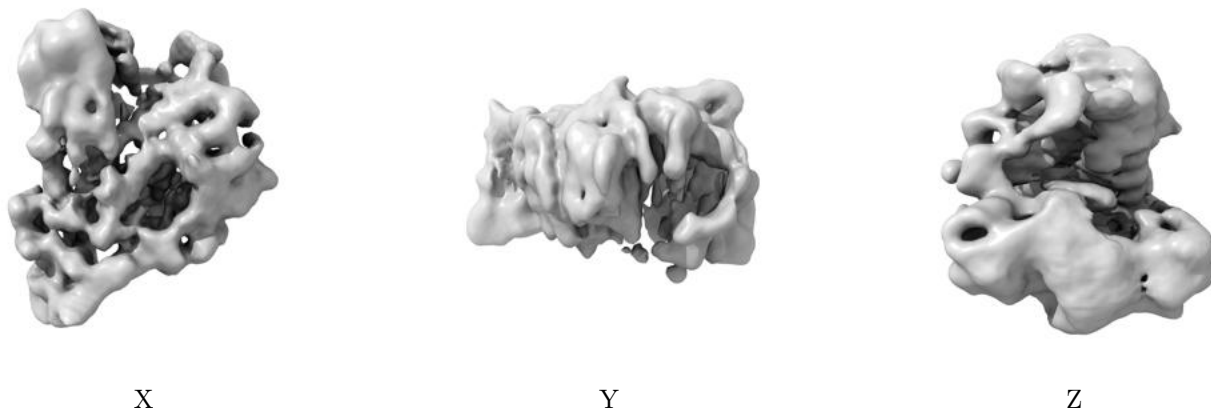
## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.127. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.4.2 Raw map



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

## 6.5 Mask visualisation [i](#)

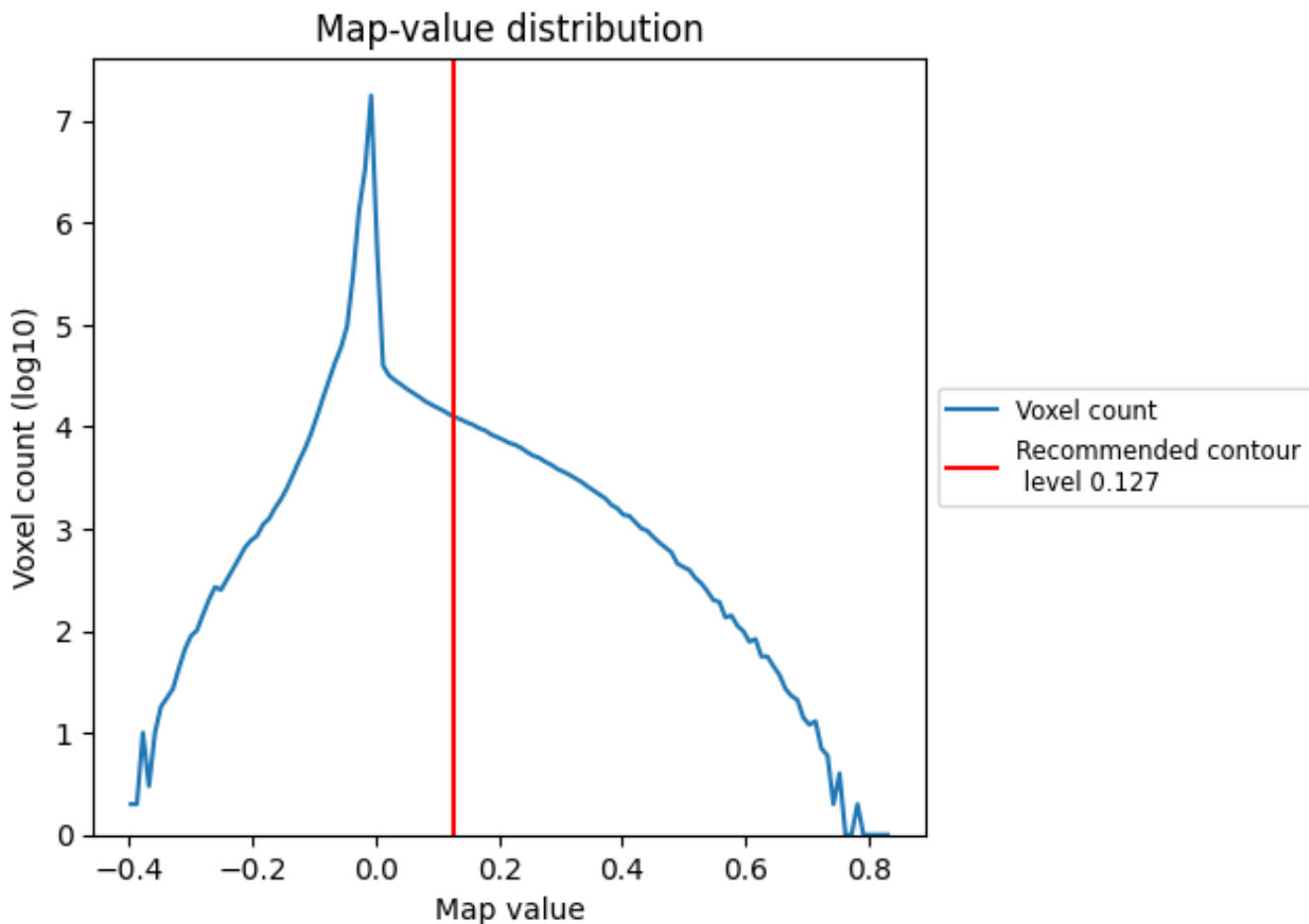
This section was not generated. No masks/segmentation were deposited.



## 7 Map analysis [i](#)

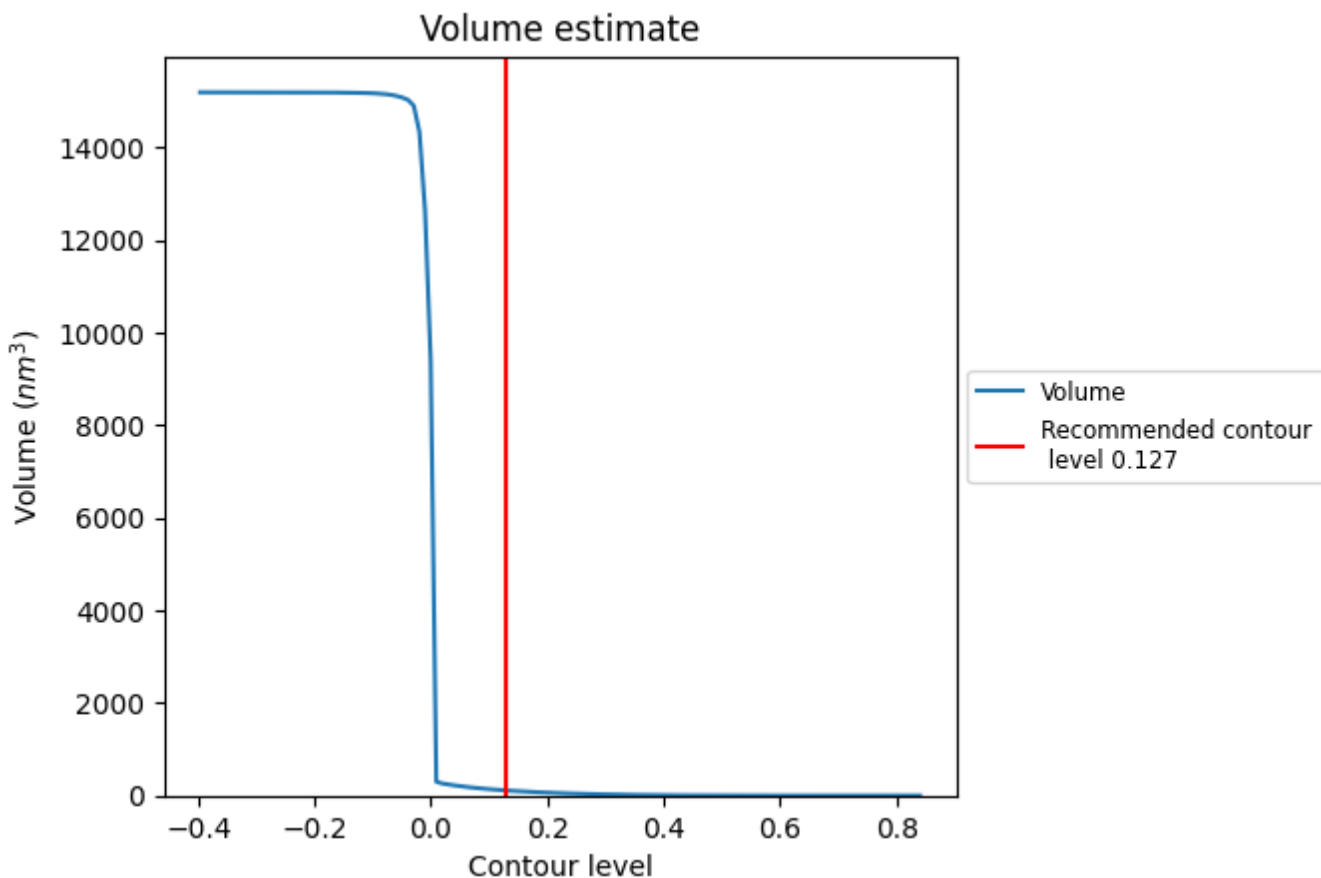
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

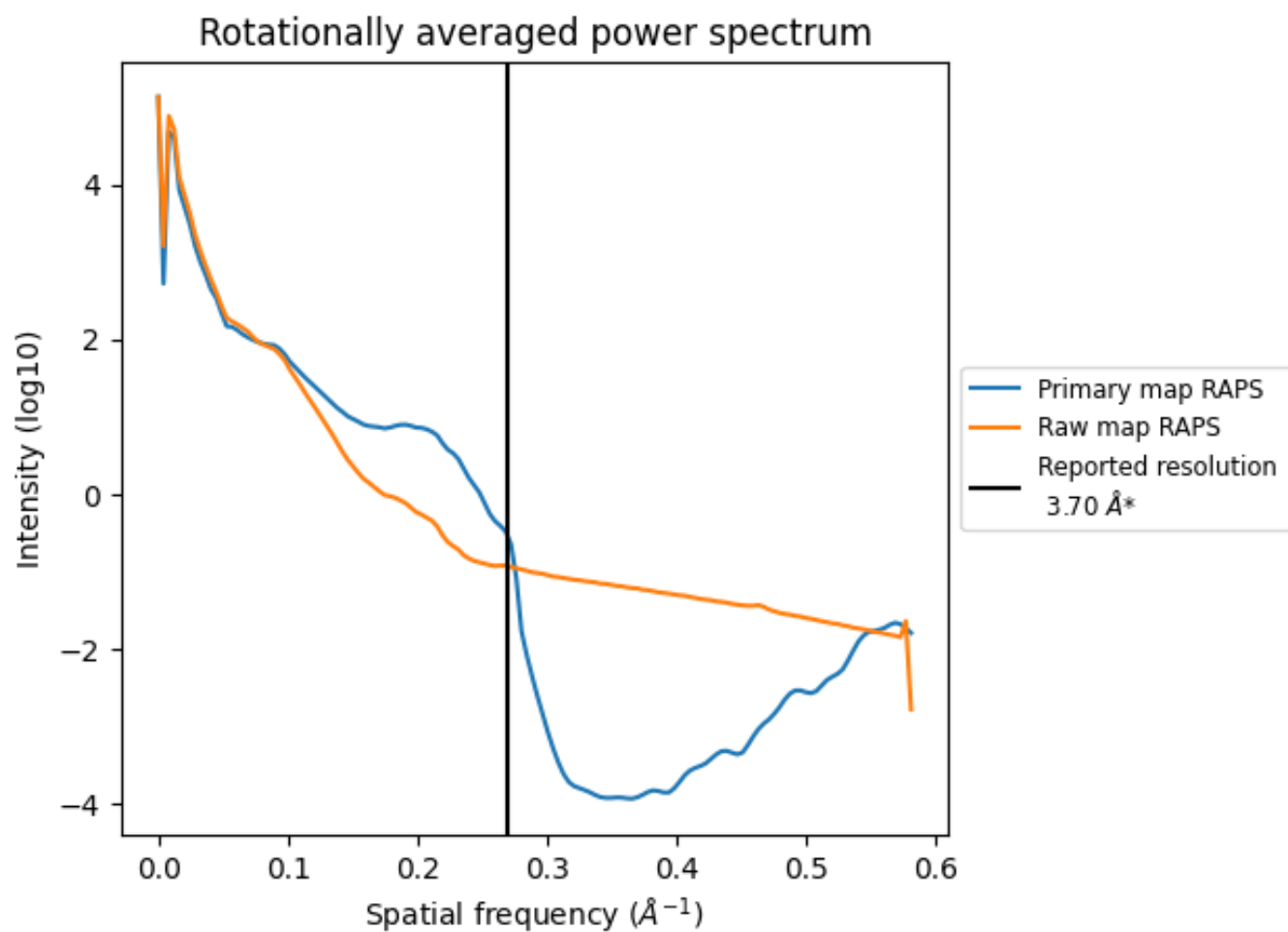
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 112 nm<sup>3</sup>; this corresponds to an approximate mass of 102 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

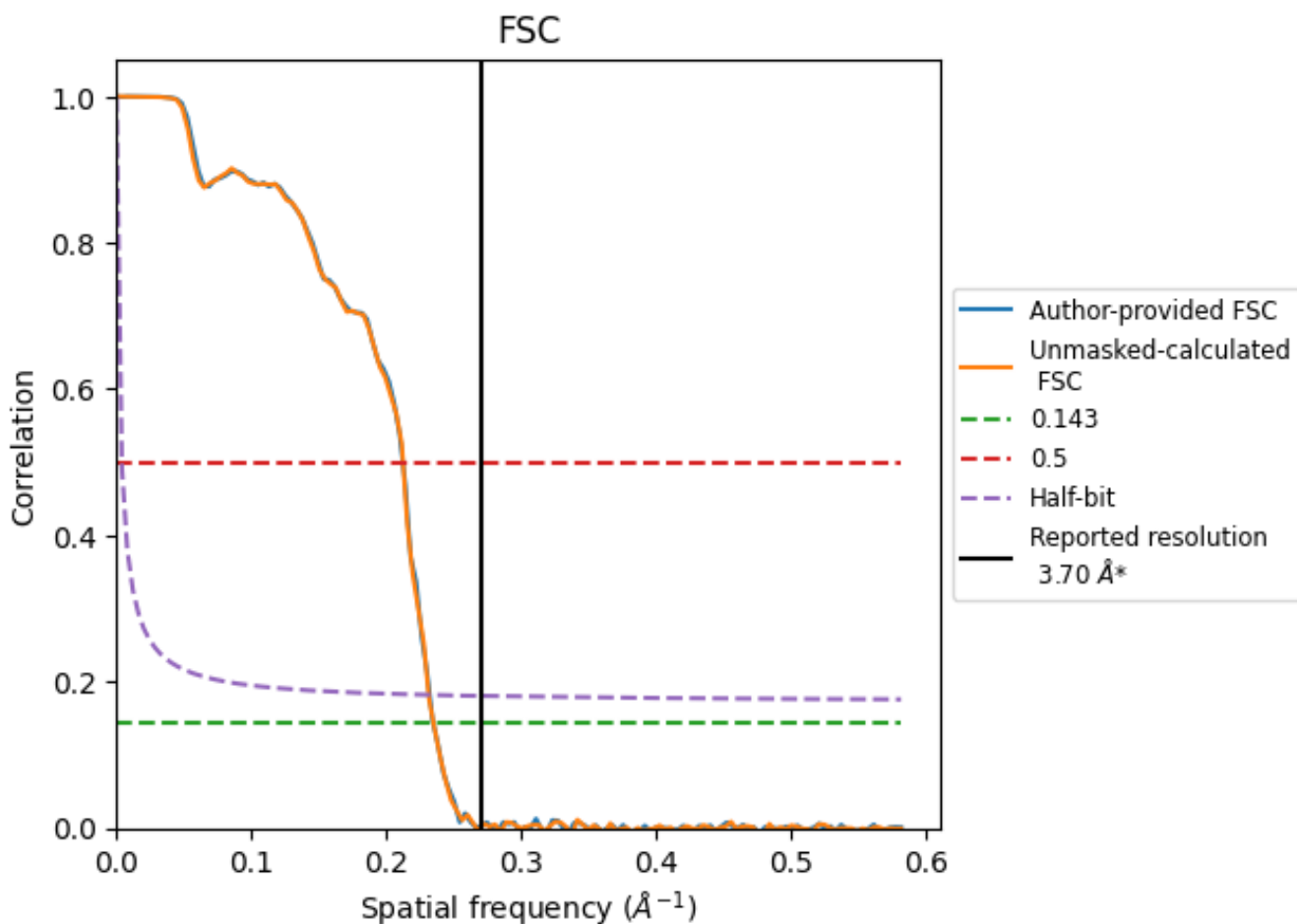


\*Reported resolution corresponds to spatial frequency of 0.270 Å<sup>-1</sup>

## 8 Fourier-Shell correlation [i](#)

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

### 8.1 FSC [i](#)



\*Reported resolution corresponds to spatial frequency of 0.270 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	4.26	4.71	4.31
Unmasked-calculated*	4.26	4.71	4.32

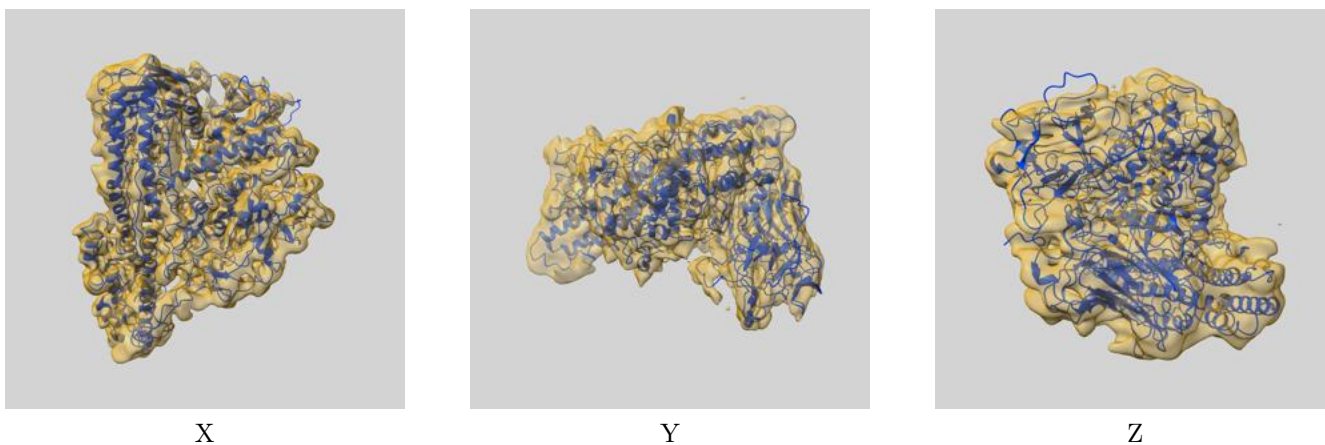
\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 4.26 differs from the reported value 3.7 by more than 10 %

The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.26 differs from the reported value 3.7 by more than 10 %

## 9 Map-model fit [i](#)

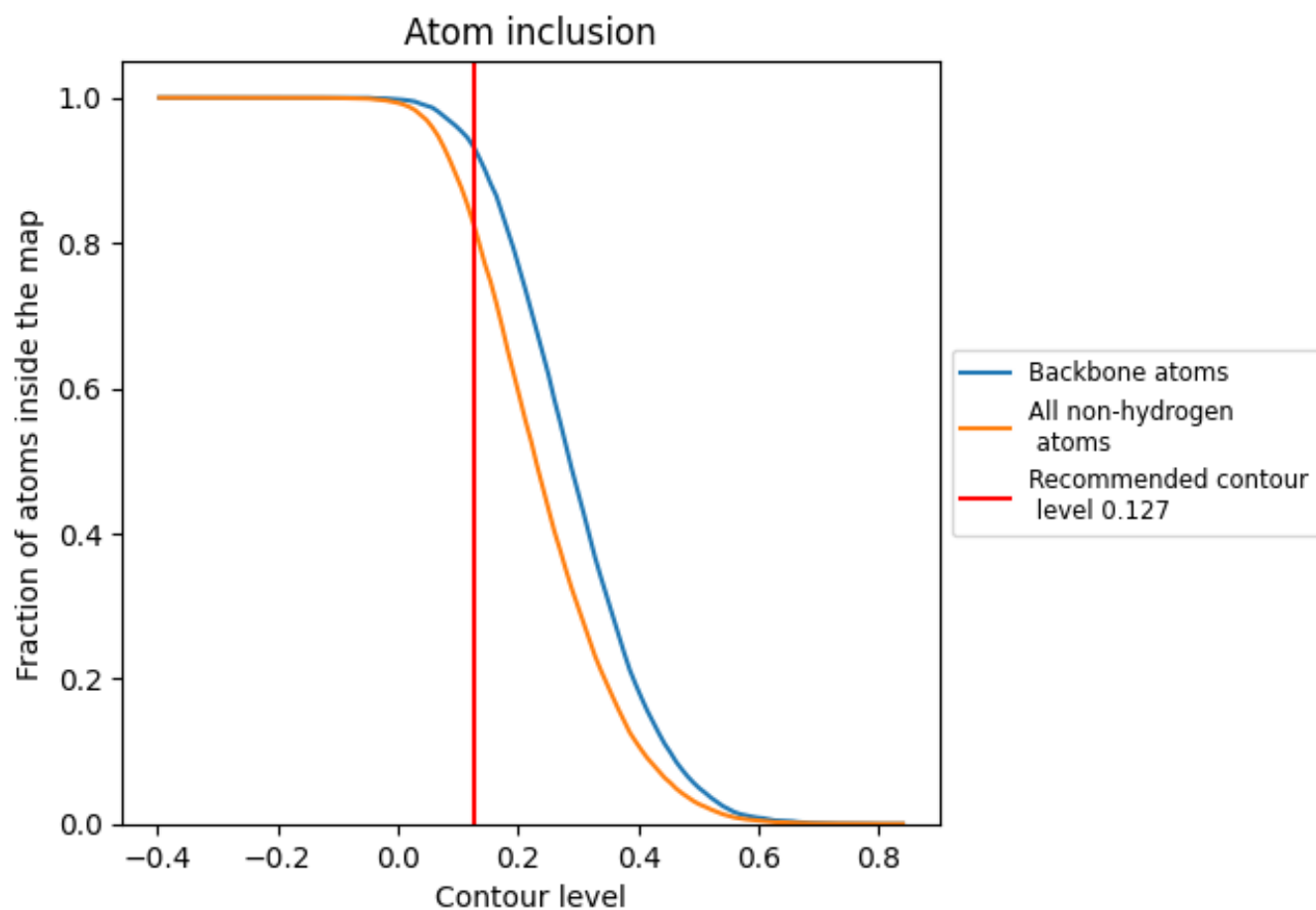
This section contains information regarding the fit between EMDB map EMD-13946 and PDB model 7QFP. Per-residue inclusion information can be found in section 3 on page 4.

### 9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.127 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 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 82% of all non-hydrogen atoms, are inside the map.