



## Full wwPDB EM Validation Report ⓘ

Nov 19, 2022 – 08:30 AM EST

PDB ID : 7LGU  
EMDB ID : EMD-23329  
Title : Structure of human prestin in the presence of NaCl  
Authors : Ge, J.; Gouaux, E.  
Deposited on : 2021-01-21  
Resolution : 2.30 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.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev43  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
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

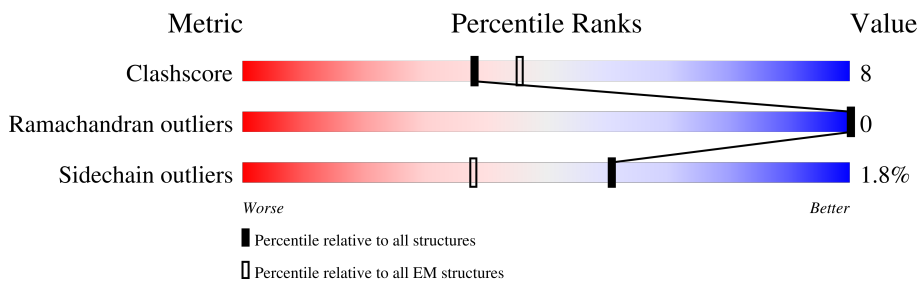
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.30 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	750	
1	B	750	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CLR	A	802	-	-	X	-
5	D10	A	839	-	-	X	-
5	D10	B	839	-	-	X	-
8	HP6	A	841	-	-	X	-

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<b>Mol</b>	<b>Type</b>	<b>Chain</b>	<b>Res</b>	<b>Chirality</b>	<b>Geometry</b>	<b>Clashes</b>	<b>Electron density</b>
8	HP6	B	841	-	-	X	-

## 2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 11597 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 Prestin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	680	5226	3431	845	926	24	0	0
1	B	680	5226	3431	845	926	24	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	745	LEU	-	expression tag	UNP P58743
A	746	GLU	-	expression tag	UNP P58743
A	747	VAL	-	expression tag	UNP P58743
A	748	LEU	-	expression tag	UNP P58743
A	749	PHE	-	expression tag	UNP P58743
A	750	GLN	-	expression tag	UNP P58743
B	745	LEU	-	expression tag	UNP P58743
B	746	GLU	-	expression tag	UNP P58743
B	747	VAL	-	expression tag	UNP P58743
B	748	LEU	-	expression tag	UNP P58743
B	749	PHE	-	expression tag	UNP P58743
B	750	GLN	-	expression tag	UNP P58743

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

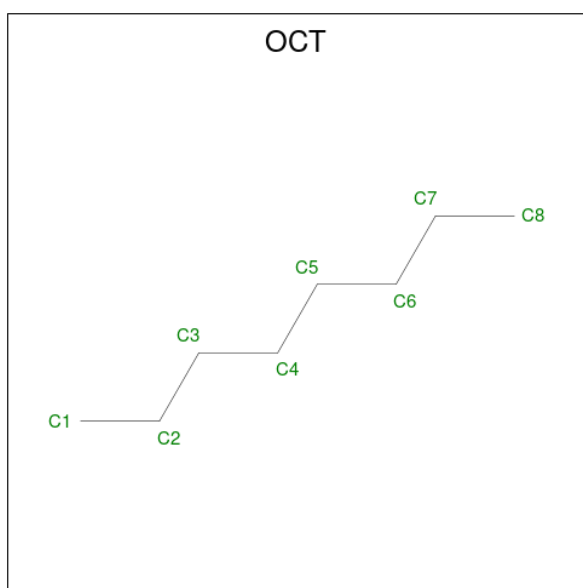
Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total	Cl	0
			1	1	
2	B	1	Total	Cl	0
			1	1	

- Molecule 3 is CHOLESTEROL (three-letter code: CLR) (formula: C<sub>27</sub>H<sub>46</sub>O).



Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	C O	0
			28	27 1	
3	B	1	Total	C O	0
			28	27 1	

- Molecule 4 is N-OCTANE (three-letter code: OCT) (formula:  $C_8H_{18}$ ).



Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	C	0
			200	200	
4	A	1	Total	C	0
			200	200	

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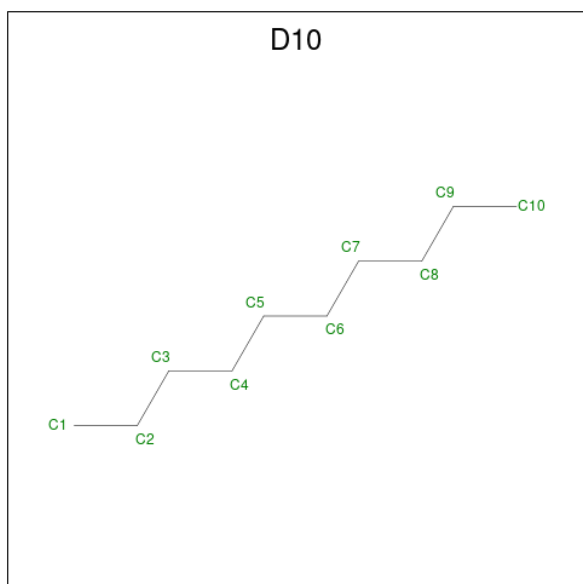




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Mol	Chain	Residues	Atoms	AltConf
4	B	1	Total C 200 200	0
4	B	1	Total C 200 200	0
4	B	1	Total C 200 200	0
4	B	1	Total C 200 200	0
4	B	1	Total C 200 200	0
4	B	1	Total C 200 200	0

- Molecule 5 is DECANE (three-letter code: D10) (formula:  $C_{10}H_{22}$ ).



Mol	Chain	Residues	Atoms	AltConf
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0

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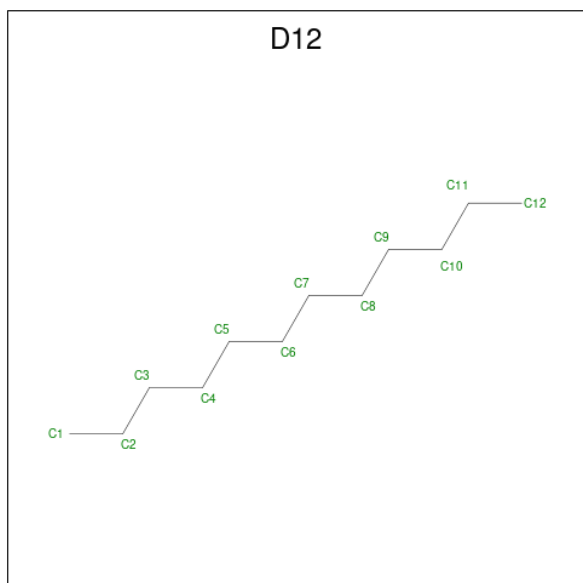
Mol	Chain	Residues	Atoms	AltConf
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	A	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0

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Mol	Chain	Residues	Atoms	AltConf
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0
5	B	1	Total C 180 180	0

- Molecule 6 is DODECANE (three-letter code: D12) (formula:  $C_{12}H_{26}$ ).



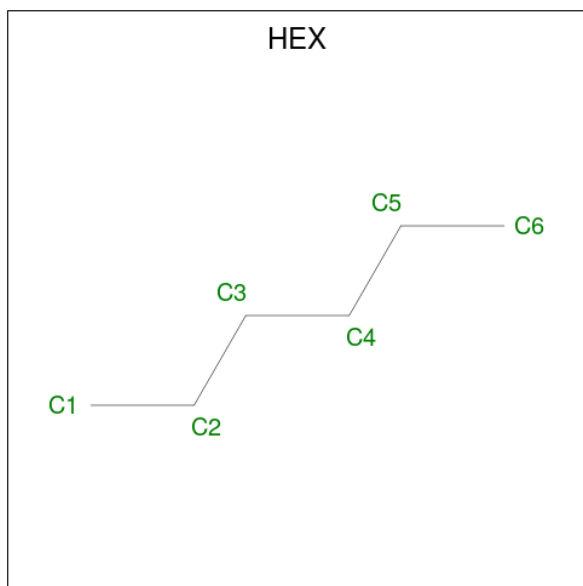
Mol	Chain	Residues	Atoms	AltConf
6	A	1	Total C 60 60	0

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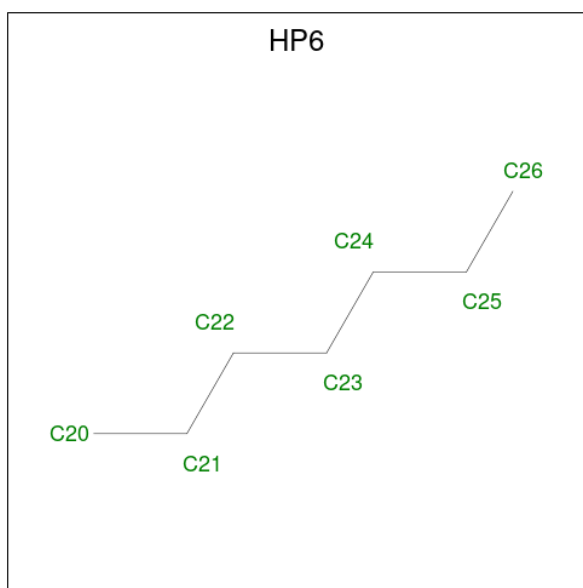
Mol	Chain	Residues	Atoms	AltConf
6	A	1	Total C 60 60	0
6	A	1	Total C 60 60	0
6	A	1	Total C 60 60	0
6	A	1	Total C 60 60	0
6	B	1	Total C 60 60	0
6	B	1	Total C 60 60	0
6	B	1	Total C 60 60	0
6	B	1	Total C 60 60	0
6	B	1	Total C 60 60	0

- Molecule 7 is HEXANE (three-letter code: HEX) (formula: C<sub>6</sub>H<sub>14</sub>).



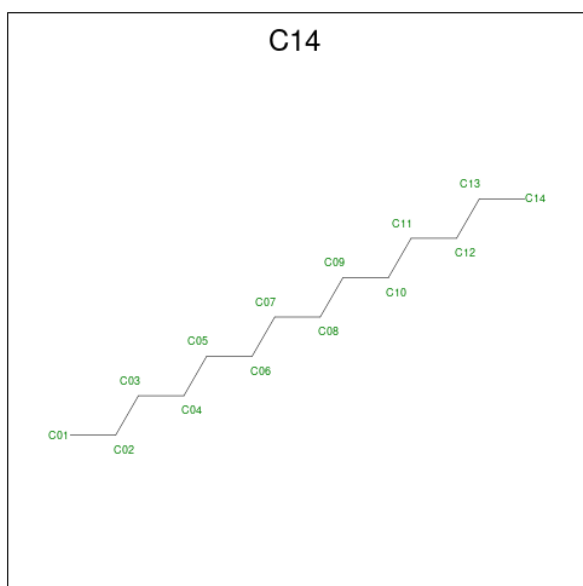
Mol	Chain	Residues	Atoms	AltConf
7	A	1	Total C 6 6	0
7	B	1	Total C 6 6	0

- Molecule 8 is HEPTANE (three-letter code: HP6) (formula:  $C_7H_{16}$ ).



Mol	Chain	Residues	Atoms	AltConf
8	A	1	Total C 14 14	0
8	A	1	Total C 14 14	0
8	B	1	Total C 14 14	0
8	B	1	Total C 14 14	0

- Molecule 9 is TETRADECANE (three-letter code: C14) (formula:  $C_{14}H_{30}$ ).



Mol	Chain	Residues	Atoms	AltConf
9	A	1	Total C 28 28	0
9	A	1	Total C 28 28	0
9	B	1	Total C 28 28	0
9	B	1	Total C 28 28	0

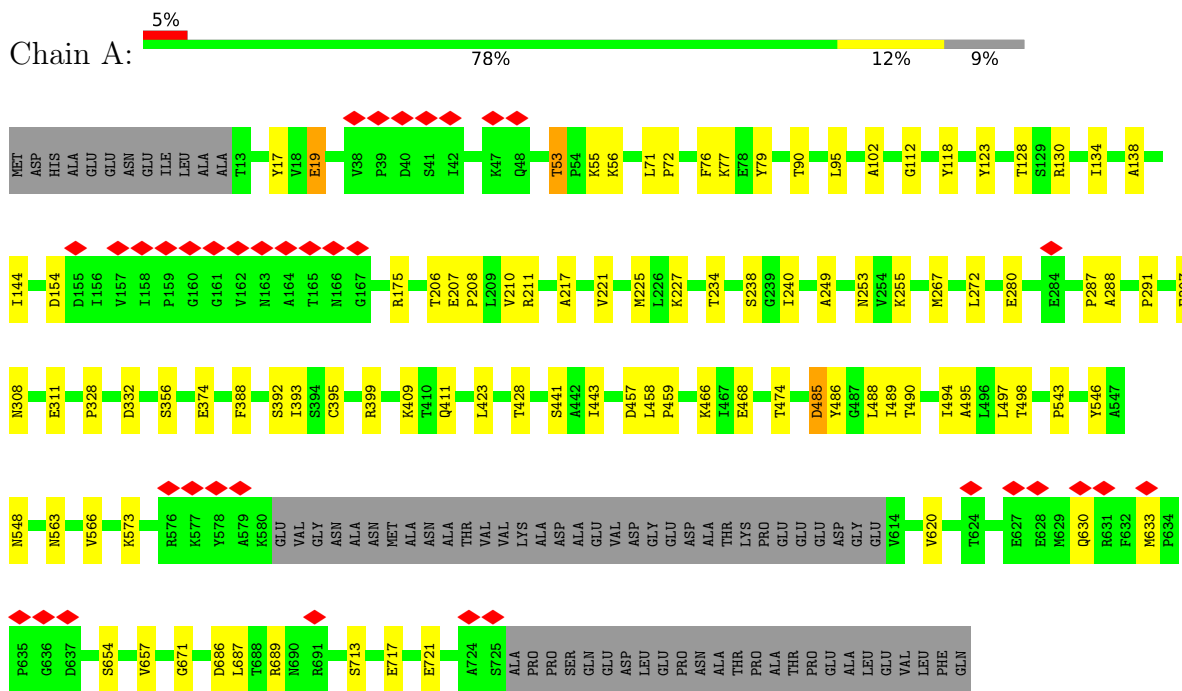
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	AltConf
10	A	55	Total O 55 55	0
10	B	56	Total O 56 56	0

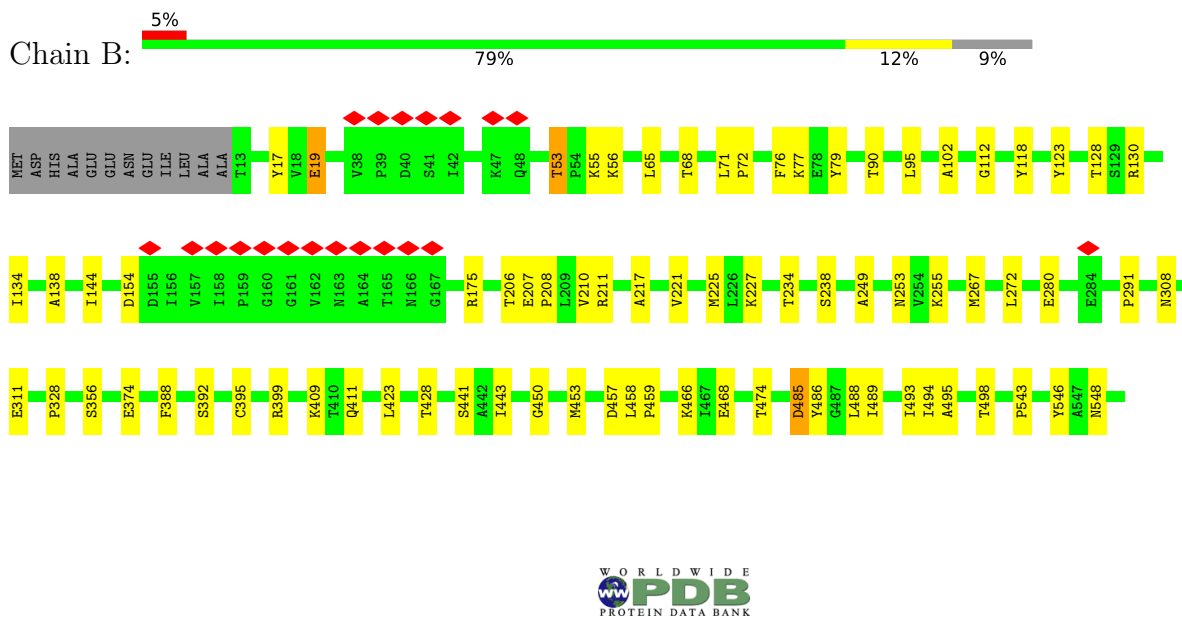
### 3 Residue-property plots

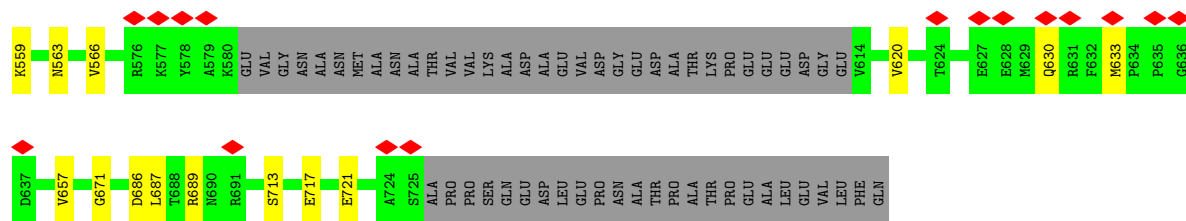
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Prestin



- Molecule 1: Prestin





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	234000	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 K3 (6k x 4k)	Depositor
Maximum map value	1.059	Depositor
Minimum map value	-0.624	Depositor
Average map value	0.006	Depositor
Map value standard deviation	0.048	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	166.5792, 166.5792, 166.5792	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.6507, 0.6507, 0.6507	Depositor



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CLR, D12, D10, HP6, CL, HEX, C14, OCT

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	0.26	0/5344	0.39	0/7266
1	B	0.26	0/5344	0.39	0/7266
All	All	0.26	0/10688	0.39	0/14532

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

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	5226	0	5384	88	0
1	B	5226	0	5384	64	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	28	0	46	34	0
3	B	28	0	46	9	0
4	A	200	0	450	12	0
4	B	200	0	450	9	0
5	A	180	0	396	16	0
5	B	180	0	396	31	0
6	A	60	0	130	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	60	0	130	3	0
7	A	6	0	14	0	0
7	B	6	0	14	0	0
8	A	14	0	32	11	0
8	B	14	0	32	24	0
9	A	28	0	60	3	0
9	B	28	0	60	3	0
10	A	55	0	0	1	0
10	B	56	0	0	1	0
All	All	11597	0	13024	197	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 (197) 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)
5:B:839:D10:H101	8:B:841:HP6:C23	1.49	1.40
1:A:494:ILE:CG1	3:A:802:CLR:H261	1.63	1.29
5:B:839:D10:C10	8:B:841:HP6:H232	1.67	1.24
5:B:839:D10:H101	8:B:841:HP6:C22	1.75	1.17
1:A:267:MET:HE2	8:A:841:HP6:H202	1.21	1.13
1:A:267:MET:CE	8:A:841:HP6:H202	1.83	1.07
5:B:839:D10:C10	8:B:841:HP6:C22	2.32	1.06
5:B:839:D10:C10	8:B:841:HP6:H212	1.85	1.06
1:A:494:ILE:HG12	3:A:802:CLR:H261	1.08	1.04
5:B:839:D10:H102	8:B:841:HP6:H212	1.40	1.03
3:A:802:CLR:H272	3:A:802:CLR:H221	1.39	1.03
5:B:839:D10:H92	8:B:841:HP6:H221	1.42	1.00
5:B:839:D10:C10	8:B:841:HP6:C21	2.45	0.94
5:B:839:D10:H101	8:B:841:HP6:H232	0.95	0.94
1:A:494:ILE:HG12	3:A:802:CLR:C26	1.98	0.92
1:A:490:THR:HG23	3:A:802:CLR:H272	1.52	0.92
1:A:486:TYR:CE1	3:A:802:CLR:H21	2.06	0.89
5:B:839:D10:C10	8:B:841:HP6:C23	2.32	0.89
5:B:839:D10:H101	8:B:841:HP6:C21	2.01	0.88
1:A:490:THR:HG23	3:A:802:CLR:C27	2.06	0.85
1:A:494:ILE:HG13	3:A:802:CLR:H261	1.54	0.85
5:B:839:D10:C9	8:B:841:HP6:H221	2.06	0.85
1:A:267:MET:HE2	8:A:841:HP6:C20	2.06	0.85
5:A:839:D10:H103	8:A:841:HP6:H232	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:486:TYR:CZ	3:A:802:CLR:H21	2.15	0.81
5:B:839:D10:H102	8:B:841:HP6:C21	2.11	0.81
1:A:494:ILE:CG1	3:A:802:CLR:C26	2.57	0.78
1:B:486:TYR:CZ	3:B:802:CLR:H12	2.20	0.77
5:B:839:D10:C10	8:B:841:HP6:H221	2.15	0.73
5:B:839:D10:C9	8:B:841:HP6:C23	2.67	0.72
1:B:486:TYR:CE2	3:B:802:CLR:H12	2.25	0.71
1:B:489:ILE:HB	5:B:813:D10:H62	1.74	0.70
1:A:489:ILE:HB	5:A:813:D10:H62	1.74	0.69
5:A:839:D10:H103	8:A:841:HP6:H212	1.74	0.69
3:A:802:CLR:H272	3:A:802:CLR:C22	2.21	0.68
5:A:839:D10:H103	8:A:841:HP6:C23	2.24	0.66
5:B:839:D10:C9	8:B:841:HP6:C22	2.68	0.66
1:A:486:TYR:CE1	3:A:802:CLR:C2	2.79	0.65
3:A:802:CLR:H151	5:A:808:D10:H62	1.78	0.65
1:A:686:ASP:OD1	1:A:689:ARG:NH1	2.30	0.65
1:A:227:LYS:HE2	1:A:234:THR:HB	1.79	0.65
5:B:839:D10:C9	8:B:841:HP6:H232	2.25	0.65
1:B:227:LYS:HE2	1:B:234:THR:HB	1.79	0.64
1:A:486:TYR:CE2	3:A:802:CLR:H11	2.33	0.64
1:B:686:ASP:OD1	1:B:689:ARG:NH1	2.29	0.63
1:B:53:THR:HG23	1:B:56:LYS:HB2	1.82	0.62
1:A:53:THR:HG23	1:A:56:LYS:HB2	1.82	0.61
1:A:76:PHE:HB3	4:A:848:OCT:H22	1.82	0.61
1:A:486:TYR:CD1	3:A:802:CLR:H12	2.34	0.61
1:B:154:ASP:OD1	1:B:175:ARG:NH2	2.33	0.61
1:B:76:PHE:HB3	4:B:848:OCT:H22	1.82	0.60
1:A:494:ILE:HG13	3:A:802:CLR:H271	1.83	0.60
1:B:486:TYR:OH	3:B:802:CLR:H21	2.02	0.60
1:A:490:THR:CG2	3:A:802:CLR:C27	2.80	0.60
3:A:802:CLR:H221	3:A:802:CLR:C27	2.22	0.60
5:A:839:D10:C10	8:A:841:HP6:H232	2.30	0.60
1:A:77:LYS:HE3	4:A:848:OCT:H31	1.83	0.59
1:B:77:LYS:HE3	4:B:848:OCT:H31	1.83	0.59
1:A:486:TYR:CZ	3:A:802:CLR:C1	2.85	0.59
1:A:486:TYR:CZ	3:A:802:CLR:C2	2.84	0.59
1:A:154:ASP:OD1	1:A:175:ARG:NH2	2.33	0.59
1:B:308:ASN:ND2	1:B:311:GLU:OE1	2.35	0.59
1:B:495:ALA:O	1:B:498:THR:OG1	2.20	0.59
1:A:138:ALA:HB3	1:A:221:VAL:HG22	1.85	0.58
1:B:138:ALA:HB3	1:B:221:VAL:HG22	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:839:D10:H103	8:A:841:HP6:C21	2.35	0.57
1:A:399:ARG:NH2	10:A:904:HOH:O	2.37	0.56
1:A:466:LYS:HB2	9:A:831:C14:H052	1.88	0.55
1:A:308:ASN:ND2	1:A:311:GLU:OE1	2.35	0.55
6:A:824:D12:H71	4:A:825:OCT:H82	1.88	0.55
1:A:144:ILE:HD11	1:A:392:SER:HB3	1.89	0.55
1:B:249:ALA:O	1:B:253:ASN:ND2	2.40	0.55
6:B:824:D12:H71	4:B:825:OCT:H82	1.89	0.55
1:A:249:ALA:O	1:A:253:ASN:ND2	2.40	0.55
1:B:466:LYS:HB2	9:B:831:C14:H052	1.88	0.54
1:A:494:ILE:HG13	3:A:802:CLR:C26	2.31	0.54
1:B:102:ALA:HB2	1:B:112:GLY:HA3	1.90	0.54
1:B:144:ILE:HD11	1:B:392:SER:HB3	1.89	0.54
4:A:851:OCT:H62	4:A:853:OCT:H62	1.89	0.54
1:A:102:ALA:HB2	1:A:112:GLY:HA3	1.90	0.53
4:B:851:OCT:H62	4:B:853:OCT:H62	1.89	0.53
1:B:543:PRO:O	1:B:548:ASN:ND2	2.41	0.53
1:A:543:PRO:O	1:A:548:ASN:ND2	2.41	0.53
6:A:816:D12:H52	5:A:817:D10:H81	1.91	0.53
5:B:839:D10:H101	8:B:841:HP6:H212	1.65	0.53
1:B:399:ARG:NH2	10:B:904:HOH:O	2.37	0.53
1:A:497:LEU:HD12	3:A:802:CLR:H263	1.90	0.53
1:A:495:ALA:O	1:A:498:THR:OG1	2.20	0.52
1:A:485:ASP:OD1	1:A:485:ASP:N	2.35	0.52
1:A:486:TYR:CE1	3:A:802:CLR:C1	2.91	0.52
1:B:291:PRO:HD3	4:B:853:OCT:H51	1.91	0.52
6:B:816:D12:H52	5:B:817:D10:H81	1.91	0.52
1:A:291:PRO:HD3	4:A:853:OCT:H51	1.91	0.52
1:A:19:GLU:HG2	1:B:17:TYR:HD1	1.75	0.52
1:B:267:MET:HE2	8:B:841:HP6:C20	2.40	0.51
1:B:630:GLN:HA	1:B:633:MET:HB2	1.92	0.51
3:A:802:CLR:C27	3:A:802:CLR:C22	2.85	0.51
1:A:630:GLN:HA	1:A:633:MET:HB2	1.92	0.51
1:B:428:THR:HG21	4:B:820:OCT:H13	1.93	0.51
1:A:211:ARG:NE	1:A:457:ASP:OD2	2.27	0.51
1:A:428:THR:HG21	4:A:820:OCT:H13	1.93	0.51
4:B:850:OCT:H12	4:B:851:OCT:H13	1.94	0.50
4:A:850:OCT:H12	4:A:851:OCT:H13	1.94	0.50
5:A:814:D10:H102	3:B:802:CLR:C26	2.42	0.50
5:B:839:D10:H91	8:B:841:HP6:H241	1.94	0.50
1:A:208:PRO:HG2	1:A:468:GLU:HG2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:ARG:NE	1:B:457:ASP:OD2	2.27	0.50
1:A:128:THR:OG1	1:A:374:GLU:OE2	2.30	0.50
1:B:208:PRO:HG2	1:B:468:GLU:HG2	1.93	0.49
1:A:17:TYR:HD1	1:B:19:GLU:HG2	1.75	0.49
1:A:486:TYR:CZ	3:A:802:CLR:H11	2.48	0.49
1:B:466:LYS:HD2	9:B:830:C14:H031	1.93	0.49
5:A:839:D10:C10	8:A:841:HP6:C23	2.91	0.49
1:B:128:THR:OG1	1:B:374:GLU:OE2	2.30	0.49
1:A:466:LYS:HD2	9:A:830:C14:H031	1.94	0.48
1:B:79:TYR:CD1	1:B:409:LYS:HG2	2.48	0.48
1:B:65:LEU:O	1:B:68:THR:OG1	2.29	0.48
1:B:474:THR:HG21	1:B:494:ILE:HG21	1.96	0.48
1:A:474:THR:HG21	1:A:494:ILE:HG21	1.96	0.48
1:A:332:ASP:OD1	1:A:332:ASP:N	2.36	0.47
5:A:839:D10:C10	8:A:841:HP6:H212	2.41	0.47
1:A:79:TYR:CD1	1:A:409:LYS:HG2	2.48	0.47
1:A:95:LEU:HD22	4:A:820:OCT:H72	1.96	0.47
3:A:802:CLR:H72	5:A:808:D10:H42	1.96	0.47
1:A:240:ILE:HD11	3:A:802:CLR:H3	1.95	0.47
1:A:486:TYR:CE1	3:A:802:CLR:H12	2.50	0.47
1:B:95:LEU:HD22	4:B:820:OCT:H72	1.96	0.47
1:A:488:LEU:HD23	5:A:813:D10:H81	1.97	0.47
5:B:839:D10:C9	8:B:841:HP6:H241	2.45	0.47
1:B:493:ILE:CG2	3:B:802:CLR:H273	2.45	0.47
1:A:717:GLU:O	1:A:721:GLU:HG3	2.15	0.46
1:B:488:LEU:HD23	5:B:813:D10:H81	1.97	0.46
1:B:717:GLU:O	1:B:721:GLU:HG3	2.15	0.46
1:B:486:TYR:CZ	3:B:802:CLR:C1	2.95	0.46
1:B:657:VAL:HG22	1:B:687:LEU:HD23	1.98	0.45
1:B:210:VAL:HG13	1:B:356:SER:HB3	1.98	0.45
6:B:812:D12:H72	6:B:812:D12:H41	1.80	0.45
1:A:494:ILE:CG1	3:A:802:CLR:H271	2.46	0.45
1:B:225:MET:O	1:B:441:SER:OG	2.31	0.44
1:B:328:PRO:HG3	1:B:388:PHE:CE1	2.52	0.44
4:A:806:OCT:H52	5:A:817:D10:H41	1.99	0.44
1:B:563:ASN:HB3	1:B:566:VAL:HB	1.99	0.44
1:A:328:PRO:HG3	1:A:388:PHE:CE1	2.52	0.44
1:B:272:LEU:HD11	1:B:443:ILE:HG12	1.99	0.44
1:A:563:ASN:HB3	1:A:566:VAL:HB	1.99	0.44
1:A:573:LYS:HE2	1:A:573:LYS:HB3	1.78	0.44
4:B:806:OCT:H52	5:B:817:D10:H41	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:VAL:HG13	1:A:356:SER:HB3	1.98	0.44
1:A:307:PHE:CG	8:A:841:HP6:H263	2.52	0.44
1:A:620:VAL:HG13	1:A:671:GLY:HA2	1.99	0.44
1:B:71:LEU:HB3	1:B:72:PRO:HD3	2.00	0.44
1:B:620:VAL:HG13	1:B:671:GLY:HA2	1.99	0.43
1:B:90:THR:HG23	1:B:399:ARG:HG2	2.00	0.43
1:A:486:TYR:CG	3:A:802:CLR:H12	2.52	0.43
1:A:657:VAL:HG22	1:A:687:LEU:HD23	1.98	0.43
1:A:272:LEU:HD11	1:A:443:ILE:HG12	1.99	0.43
5:B:839:D10:H92	8:B:841:HP6:C22	2.24	0.43
1:A:71:LEU:HB3	1:A:72:PRO:HD3	2.00	0.43
1:A:123:TYR:CG	1:A:134:ILE:HG13	2.54	0.43
1:A:486:TYR:CE2	3:A:802:CLR:C1	3.02	0.43
1:B:423:LEU:HB2	5:B:821:D10:H92	2.00	0.43
1:A:423:LEU:HB2	5:A:821:D10:H92	2.00	0.43
1:A:90:THR:HG23	1:A:399:ARG:HG2	2.00	0.42
4:A:845:OCT:H71	4:A:845:OCT:H42	1.83	0.42
1:B:466:LYS:HB2	9:B:831:C14:H021	2.01	0.42
1:A:130:ARG:HA	1:A:411:GLN:HG2	2.01	0.42
3:A:802:CLR:H152	5:B:814:D10:H41	2.02	0.42
1:A:458:LEU:HB2	1:A:459:PRO:HD3	2.02	0.42
1:B:130:ARG:HA	1:B:411:GLN:HG2	2.01	0.42
1:A:466:LYS:HB2	9:A:831:C14:H021	2.01	0.42
1:A:494:ILE:CD1	3:A:802:CLR:H271	2.50	0.42
1:A:206:THR:HB	1:B:546:TYR:CD2	2.55	0.42
1:B:458:LEU:HB2	1:B:459:PRO:HD3	2.02	0.42
1:B:493:ILE:HG21	3:B:802:CLR:H273	2.01	0.42
1:B:559:LYS:HE2	1:B:559:LYS:HB3	1.88	0.42
1:A:217:ALA:O	1:A:221:VAL:HG23	2.20	0.41
1:A:546:TYR:CG	1:B:206:THR:HB	2.55	0.41
1:B:123:TYR:CG	1:B:134:ILE:HG13	2.54	0.41
1:B:118:TYR:HE1	5:B:823:D10:H81	1.85	0.41
1:A:206:THR:HB	1:B:546:TYR:CG	2.56	0.41
1:A:546:TYR:CD2	1:B:206:THR:HB	2.55	0.41
3:B:802:CLR:C15	5:B:808:D10:H62	2.50	0.41
1:A:490:THR:CG2	3:A:802:CLR:H271	2.48	0.41
1:B:217:ALA:O	1:B:221:VAL:HG23	2.20	0.41
1:A:118:TYR:HE1	5:A:823:D10:H81	1.85	0.41
6:A:812:D12:H41	6:A:812:D12:H72	1.80	0.41
5:B:839:D10:C9	8:B:841:HP6:C24	2.99	0.41
1:A:225:MET:O	1:A:441:SER:OG	2.31	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:MET:CE	8:B:841:HP6:H211	2.51	0.40
1:B:486:TYR:CE1	3:B:802:CLR:H3	2.57	0.40
1:A:393:ILE:H	1:A:393:ILE:HG13	1.77	0.40
1:A:287:PRO:HB2	4:A:851:OCT:H12	2.03	0.40
1:A:288:ALA:HB3	4:A:853:OCT:H42	2.02	0.40
1:A:654:SER:OG	1:B:468:GLU:OE2	2.26	0.40
5:B:836:D10:H52	5:B:836:D10:H81	1.94	0.40
1:B:450:GLY:O	1:B:453:MET:HB2	2.21	0.40
1:B:485:ASP:OD1	1:B:485:ASP:N	2.35	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	676/750 (90%)	666 (98%)	10 (2%)	0	100	100
1	B	676/750 (90%)	665 (98%)	11 (2%)	0	100	100
All	All	1352/1500 (90%)	1331 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	567/631 (90%)	557 (98%)	10 (2%)	59	75
1	B	567/631 (90%)	557 (98%)	10 (2%)	59	75
All	All	1134/1262 (90%)	1114 (98%)	20 (2%)	61	75

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	53	THR
1	A	55	LYS
1	A	207	GLU
1	A	238	SER
1	A	255	LYS
1	A	280	GLU
1	A	395	CYS
1	A	485	ASP
1	A	713	SER
1	B	19	GLU
1	B	53	THR
1	B	55	LYS
1	B	207	GLU
1	B	238	SER
1	B	255	LYS
1	B	280	GLU
1	B	395	CYS
1	B	485	ASP
1	B	713	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	454	GLN
1	B	454	GLN

### 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](#)

Of 110 ligands modelled in this entry, 2 are monoatomic - leaving 108 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	OCT	B	809	-	7,7,7	0.29	0	6,6,6	0.74	0
4	OCT	A	828	-	7,7,7	0.29	0	6,6,6	0.73	0
4	OCT	A	848	-	7,7,7	0.31	0	6,6,6	0.72	0
4	OCT	A	844	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	A	832	-	9,9,9	0.28	0	8,8,8	0.81	0
3	CLR	A	802	-	31,31,31	0.28	0	48,48,48	0.39	0
4	OCT	A	853	-	7,7,7	0.30	0	6,6,6	0.71	0
4	OCT	B	828	-	7,7,7	0.29	0	6,6,6	0.73	0
4	OCT	B	818	-	7,7,7	0.29	0	6,6,6	0.74	0
6	D12	A	811	-	11,11,11	0.30	0	10,10,10	0.82	0
4	OCT	B	806	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	845	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	A	821	-	9,9,9	0.30	0	8,8,8	0.80	0
8	HP6	B	827	-	6,6,6	0.29	0	5,5,5	0.68	0
4	OCT	B	850	-	7,7,7	0.30	0	6,6,6	0.73	0
7	HEX	B	826	-	5,5,5	0.31	0	4,4,4	0.57	0
5	D10	A	819	-	9,9,9	0.28	0	8,8,8	0.82	0
5	D10	A	836	-	9,9,9	0.29	0	8,8,8	0.78	0
5	D10	A	823	-	9,9,9	0.29	0	8,8,8	0.81	0
5	D10	B	823	-	9,9,9	0.29	0	8,8,8	0.82	0
4	OCT	A	847	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	B	837	-	9,9,9	0.29	0	8,8,8	0.81	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	OCT	B	805	-	7,7,7	0.29	0	6,6,6	0.72	0
4	OCT	A	846	-	7,7,7	0.30	0	6,6,6	0.72	0
4	OCT	A	852	-	7,7,7	0.30	0	6,6,6	0.71	0
5	D10	B	836	-	9,9,9	0.30	0	8,8,8	0.79	0
4	OCT	B	838	-	7,7,7	0.29	0	6,6,6	0.72	0
5	D10	A	808	-	9,9,9	0.30	0	8,8,8	0.80	0
6	D12	B	816	-	11,11,11	0.28	0	10,10,10	0.85	0
4	OCT	A	806	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	849	-	7,7,7	0.30	0	6,6,6	0.74	0
6	D12	A	812	-	11,11,11	0.29	0	10,10,10	0.84	0
9	C14	B	830	-	13,13,13	0.28	0	12,12,12	0.86	0
9	C14	B	831	-	13,13,13	0.27	0	12,12,12	0.86	0
4	OCT	A	804	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	A	803	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	A	833	-	9,9,9	0.29	0	8,8,8	0.80	0
6	D12	B	811	-	11,11,11	0.29	0	10,10,10	0.82	0
9	C14	A	830	-	13,13,13	0.28	0	12,12,12	0.87	0
4	OCT	A	825	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	B	853	-	7,7,7	0.30	0	6,6,6	0.71	0
6	D12	B	812	-	11,11,11	0.29	0	10,10,10	0.84	0
5	D10	B	817	-	9,9,9	0.29	0	8,8,8	0.80	0
4	OCT	B	844	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	855	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	847	-	7,7,7	0.30	0	6,6,6	0.74	0
9	C14	A	831	-	13,13,13	0.27	0	12,12,12	0.87	0
4	OCT	A	820	-	7,7,7	0.29	0	6,6,6	0.73	0
4	OCT	A	810	-	7,7,7	0.28	0	6,6,6	0.77	0
4	OCT	B	807	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	A	855	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	B	808	-	9,9,9	0.29	0	8,8,8	0.80	0
4	OCT	A	818	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	B	835	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	B	839	-	9,9,9	0.10	0	8,8,8	0.06	0
5	D10	A	835	-	9,9,9	0.29	0	8,8,8	0.79	0
4	OCT	A	807	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	A	854	-	7,7,7	0.29	0	6,6,6	0.73	0
3	CLR	B	802	-	31,31,31	0.29	0	48,48,48	0.38	0
5	D10	A	822	-	9,9,9	0.29	0	8,8,8	0.81	0
7	HEX	A	826	-	5,5,5	0.31	0	4,4,4	0.57	0
5	D10	B	819	-	9,9,9	0.28	0	8,8,8	0.82	0
4	OCT	B	848	-	7,7,7	0.30	0	6,6,6	0.71	0
4	OCT	B	820	-	7,7,7	0.29	0	6,6,6	0.74	0
4	OCT	B	851	-	7,7,7	0.29	0	6,6,6	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	OCT	B	840	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	B	829	-	9,9,9	0.29	0	8,8,8	0.79	0
6	D12	B	842	-	11,11,11	0.29	0	10,10,10	0.84	0
4	OCT	A	850	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	846	-	7,7,7	0.30	0	6,6,6	0.72	0
5	D10	B	843	-	9,9,9	0.30	0	8,8,8	0.78	0
4	OCT	A	840	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	A	839	-	9,9,9	0.10	0	8,8,8	0.06	0
5	D10	B	814	-	9,9,9	0.29	0	8,8,8	0.81	0
5	D10	A	829	-	9,9,9	0.29	0	8,8,8	0.79	0
5	D10	B	834	-	9,9,9	0.29	0	8,8,8	0.82	0
5	D10	B	813	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	A	817	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	A	834	-	9,9,9	0.29	0	8,8,8	0.82	0
4	OCT	B	854	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	A	805	-	7,7,7	0.29	0	6,6,6	0.72	0
5	D10	A	814	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	A	809	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	B	822	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	B	810	-	7,7,7	0.28	0	6,6,6	0.77	0
6	D12	A	816	-	11,11,11	0.28	0	10,10,10	0.85	0
8	HP6	A	841	-	6,6,6	0.11	0	5,5,5	0.10	0
4	OCT	B	852	-	7,7,7	0.31	0	6,6,6	0.71	0
5	D10	A	837	-	9,9,9	0.29	0	8,8,8	0.81	0
6	D12	A	842	-	11,11,11	0.29	0	10,10,10	0.83	0
4	OCT	B	825	-	7,7,7	0.30	0	6,6,6	0.74	0
6	D12	B	824	-	11,11,11	0.29	0	10,10,10	0.84	0
5	D10	B	815	-	9,9,9	0.29	0	8,8,8	0.79	0
5	D10	B	821	-	9,9,9	0.29	0	8,8,8	0.80	0
4	OCT	A	838	-	7,7,7	0.30	0	6,6,6	0.72	0
8	HP6	B	841	-	6,6,6	0.11	0	5,5,5	0.09	0
4	OCT	A	851	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	A	849	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	804	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	A	845	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	B	832	-	9,9,9	0.28	0	8,8,8	0.82	0
6	D12	A	824	-	11,11,11	0.29	0	10,10,10	0.84	0
8	HP6	A	827	-	6,6,6	0.30	0	5,5,5	0.68	0
4	OCT	B	803	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	A	815	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	A	843	-	9,9,9	0.30	0	8,8,8	0.78	0
5	D10	B	833	-	9,9,9	0.29	0	8,8,8	0.81	0
5	D10	A	813	-	9,9,9	0.29	0	8,8,8	0.81	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	OCT	B	809	-	-	0/5/5/5	-
4	OCT	A	828	-	-	0/5/5/5	-
4	OCT	A	848	-	-	0/5/5/5	-
4	OCT	A	844	-	-	0/5/5/5	-
5	D10	A	832	-	-	2/7/7/7	-
3	CLR	A	802	-	-	0/10/68/68	0/4/4/4
4	OCT	A	853	-	-	0/5/5/5	-
4	OCT	B	828	-	-	0/5/5/5	-
4	OCT	B	818	-	-	0/5/5/5	-
6	D12	A	811	-	-	3/9/9/9	-
4	OCT	B	806	-	-	0/5/5/5	-
4	OCT	B	845	-	-	0/5/5/5	-
5	D10	A	821	-	-	0/7/7/7	-
8	HP6	B	827	-	-	0/4/4/4	-
4	OCT	B	850	-	-	0/5/5/5	-
7	HEX	B	826	-	-	0/3/3/3	-
5	D10	A	819	-	-	0/7/7/7	-
5	D10	A	836	-	-	0/7/7/7	-
5	D10	A	823	-	-	0/7/7/7	-
5	D10	B	823	-	-	0/7/7/7	-
4	OCT	A	847	-	-	0/5/5/5	-
5	D10	B	837	-	-	0/7/7/7	-
4	OCT	B	805	-	-	1/5/5/5	-
4	OCT	A	846	-	-	0/5/5/5	-
4	OCT	A	852	-	-	0/5/5/5	-
5	D10	B	836	-	-	0/7/7/7	-
4	OCT	B	838	-	-	0/5/5/5	-
5	D10	A	808	-	-	0/7/7/7	-
6	D12	B	816	-	-	0/9/9/9	-
4	OCT	A	806	-	-	0/5/5/5	-
4	OCT	B	849	-	-	0/5/5/5	-
6	D12	A	812	-	-	2/9/9/9	-
9	C14	B	830	-	-	0/11/11/11	-
9	C14	B	831	-	-	0/11/11/11	-
4	OCT	A	804	-	-	0/5/5/5	-
4	OCT	A	803	-	-	0/5/5/5	-
5	D10	A	833	-	-	0/7/7/7	-
6	D12	B	811	-	-	3/9/9/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	C14	A	830	-	-	0/11/11/11	-
4	OCT	A	825	-	-	0/5/5/5	-
4	OCT	B	853	-	-	0/5/5/5	-
6	D12	B	812	-	-	2/9/9/9	-
5	D10	B	817	-	-	0/7/7/7	-
4	OCT	B	844	-	-	0/5/5/5	-
4	OCT	B	855	-	-	0/5/5/5	-
4	OCT	B	847	-	-	0/5/5/5	-
9	C14	A	831	-	-	0/11/11/11	-
4	OCT	A	820	-	-	0/5/5/5	-
4	OCT	A	810	-	-	0/5/5/5	-
4	OCT	B	807	-	-	0/5/5/5	-
4	OCT	A	855	-	-	0/5/5/5	-
5	D10	B	808	-	-	0/7/7/7	-
4	OCT	A	818	-	-	0/5/5/5	-
5	D10	B	835	-	-	0/7/7/7	-
5	D10	B	839	-	-	0/7/7/7	-
5	D10	A	835	-	-	0/7/7/7	-
4	OCT	A	807	-	-	0/5/5/5	-
4	OCT	A	854	-	-	0/5/5/5	-
3	CLR	B	802	-	-	0/10/68/68	0/4/4/4
5	D10	A	822	-	-	0/7/7/7	-
7	HEX	A	826	-	-	0/3/3/3	-
5	D10	B	819	-	-	0/7/7/7	-
4	OCT	B	848	-	-	0/5/5/5	-
4	OCT	B	820	-	-	0/5/5/5	-
4	OCT	B	851	-	-	0/5/5/5	-
4	OCT	B	840	-	-	0/5/5/5	-
5	D10	B	829	-	-	0/7/7/7	-
6	D12	B	842	-	-	1/9/9/9	-
4	OCT	A	850	-	-	0/5/5/5	-
4	OCT	B	846	-	-	0/5/5/5	-
5	D10	B	843	-	-	1/7/7/7	-
4	OCT	A	840	-	-	0/5/5/5	-
5	D10	A	839	-	-	0/7/7/7	-
5	D10	B	814	-	-	0/7/7/7	-
5	D10	A	829	-	-	0/7/7/7	-
5	D10	B	834	-	-	0/7/7/7	-
5	D10	B	813	-	-	0/7/7/7	-
5	D10	A	817	-	-	0/7/7/7	-
5	D10	A	834	-	-	0/7/7/7	-
4	OCT	B	854	-	-	0/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	OCT	A	805	-	-	1/5/5/5	-
5	D10	A	814	-	-	0/7/7/7	-
4	OCT	A	809	-	-	0/5/5/5	-
5	D10	B	822	-	-	0/7/7/7	-
4	OCT	B	810	-	-	0/5/5/5	-
6	D12	A	816	-	-	0/9/9/9	-
8	HP6	A	841	-	-	0/4/4/4	-
4	OCT	B	852	-	-	0/5/5/5	-
5	D10	A	837	-	-	0/7/7/7	-
6	D12	A	842	-	-	1/9/9/9	-
4	OCT	B	825	-	-	0/5/5/5	-
6	D12	B	824	-	-	0/9/9/9	-
5	D10	B	815	-	-	0/7/7/7	-
5	D10	B	821	-	-	0/7/7/7	-
4	OCT	A	838	-	-	0/5/5/5	-
8	HP6	B	841	-	-	0/4/4/4	-
4	OCT	A	851	-	-	0/5/5/5	-
4	OCT	A	849	-	-	0/5/5/5	-
4	OCT	B	804	-	-	0/5/5/5	-
4	OCT	A	845	-	-	0/5/5/5	-
5	D10	B	832	-	-	2/7/7/7	-
6	D12	A	824	-	-	0/9/9/9	-
8	HP6	A	827	-	-	0/4/4/4	-
4	OCT	B	803	-	-	0/5/5/5	-
5	D10	A	815	-	-	0/7/7/7	-
5	D10	A	843	-	-	1/7/7/7	-
5	D10	B	833	-	-	0/7/7/7	-
5	D10	A	813	-	-	0/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	811	D12	C3-C4-C5-C6
6	B	811	D12	C3-C4-C5-C6
6	A	812	D12	C7-C8-C9-C10
6	B	812	D12	C7-C8-C9-C10
6	A	811	D12	C4-C5-C6-C7
6	B	811	D12	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
6	A	811	D12	C2-C3-C4-C5
6	A	842	D12	C11-C10-C9-C8
6	B	842	D12	C11-C10-C9-C8
6	B	811	D12	C2-C3-C4-C5
6	A	812	D12	C3-C4-C5-C6
6	B	812	D12	C3-C4-C5-C6
4	B	805	OCT	C4-C5-C6-C7
4	A	805	OCT	C4-C5-C6-C7
5	A	832	D10	C4-C5-C6-C7
5	B	832	D10	C4-C5-C6-C7
5	A	832	D10	C5-C6-C7-C8
5	B	832	D10	C5-C6-C7-C8
5	A	843	D10	C3-C4-C5-C6
5	B	843	D10	C3-C4-C5-C6

There are no ring outliers.

44 monomers are involved in 118 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	848	OCT	2	0
3	A	802	CLR	34	0
4	A	853	OCT	3	0
4	B	806	OCT	1	0
5	A	821	D10	1	0
4	B	850	OCT	1	0
5	A	823	D10	1	0
5	B	823	D10	1	0
5	B	836	D10	1	0
5	A	808	D10	2	0
6	B	816	D12	1	0
4	A	806	OCT	1	0
6	A	812	D12	1	0
9	B	830	C14	1	0
9	B	831	C14	2	0
9	A	830	C14	1	0
4	A	825	OCT	1	0
4	B	853	OCT	2	0
6	B	812	D12	1	0
5	B	817	D10	2	0
9	A	831	C14	2	0
4	A	820	OCT	2	0
5	B	808	D10	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	839	D10	22	0
3	B	802	CLR	9	0
4	B	848	OCT	2	0
4	B	820	OCT	2	0
4	B	851	OCT	2	0
4	A	850	OCT	1	0
5	A	839	D10	7	0
5	B	814	D10	1	0
5	B	813	D10	2	0
5	A	817	D10	2	0
5	A	814	D10	1	0
6	A	816	D12	1	0
8	A	841	HP6	11	0
4	B	825	OCT	1	0
6	B	824	D12	1	0
5	B	821	D10	1	0
8	B	841	HP6	24	0
4	A	851	OCT	3	0
4	A	845	OCT	1	0
6	A	824	D12	1	0
5	A	813	D10	2	0

## 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.



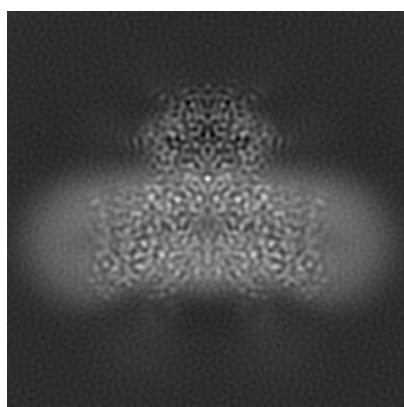
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23329. These allow visual inspection of the internal detail of the map and identification of artifacts.

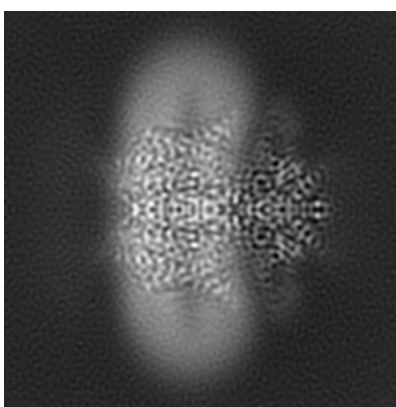
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

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

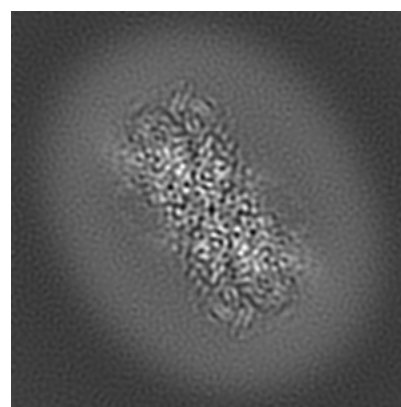
#### 6.1.1 Primary map



X



Y

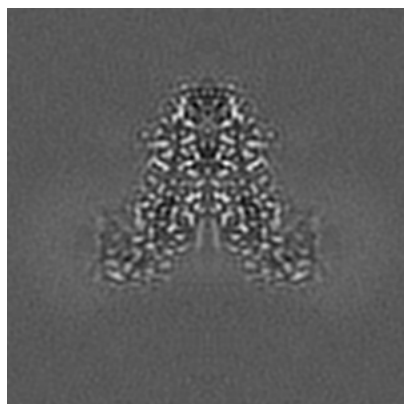


Z

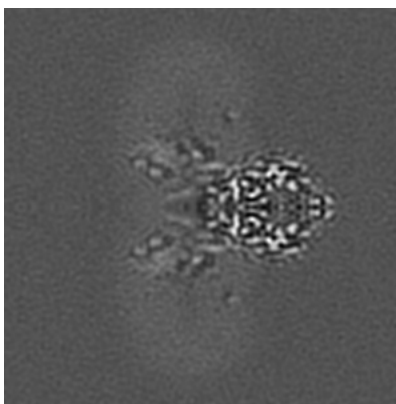
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

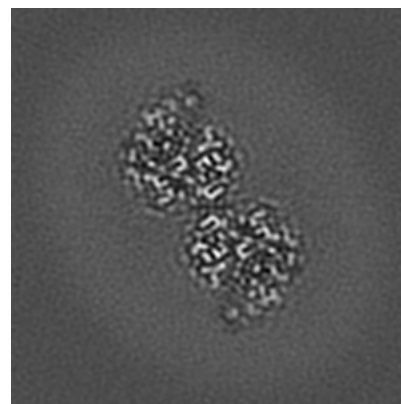
#### 6.2.1 Primary map



X Index: 128



Y Index: 128

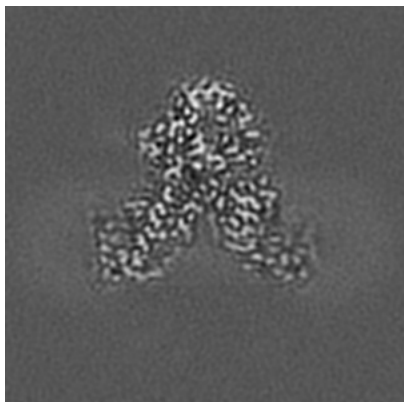


Z Index: 128

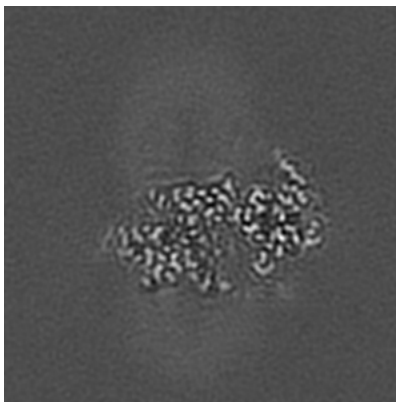
The images above show central slices of the map in three orthogonal directions.

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

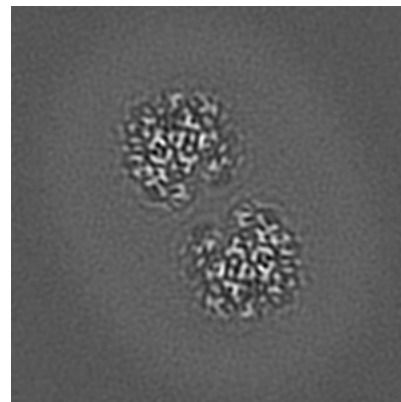
### 6.3.1 Primary map



X Index: 131



Y Index: 151

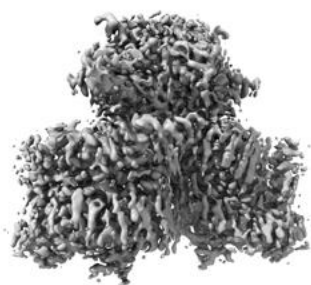


Z Index: 99

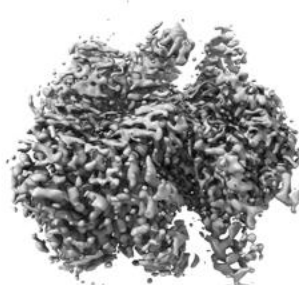
The images above show the largest variance slices of the map in three orthogonal directions.

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

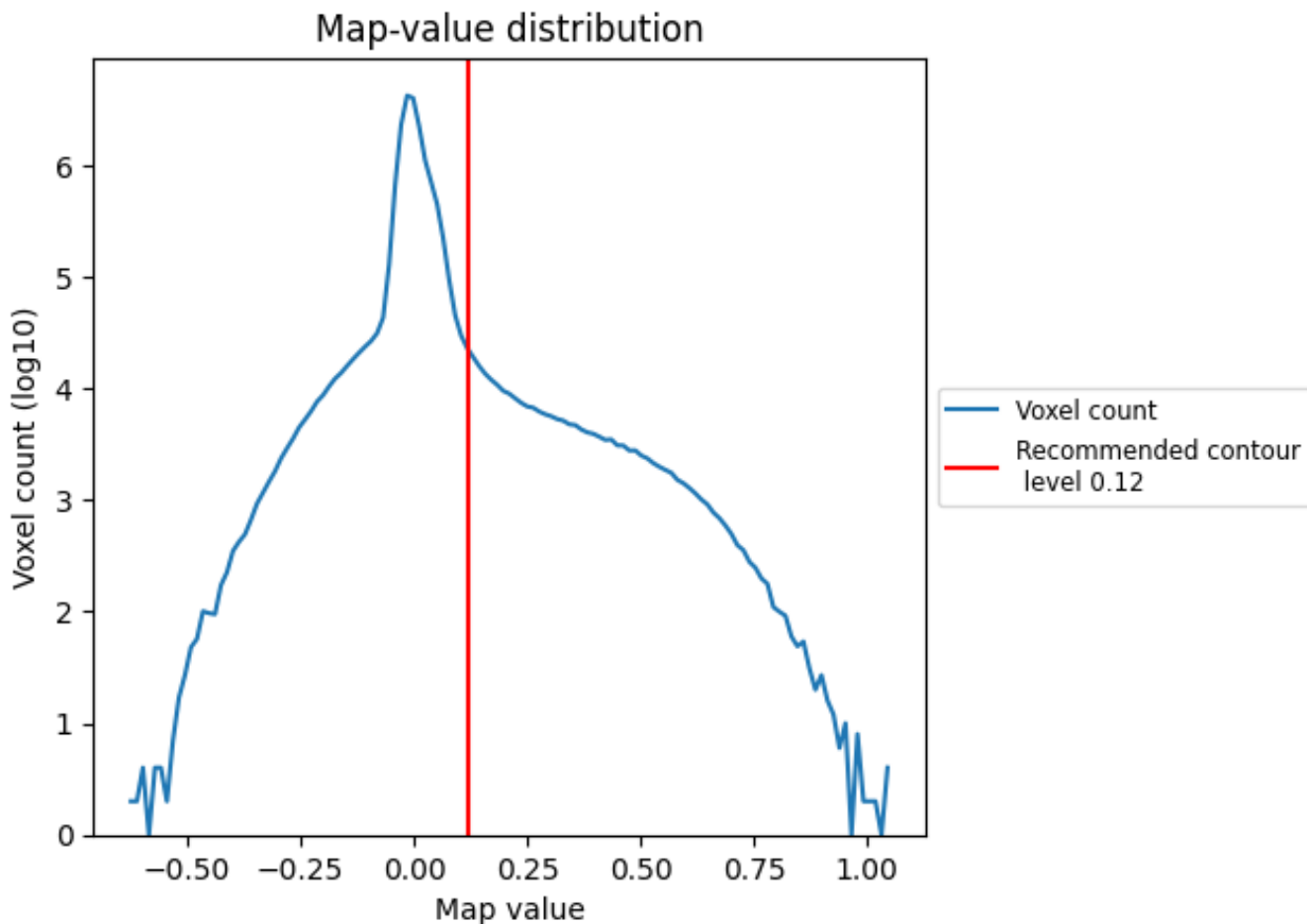
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

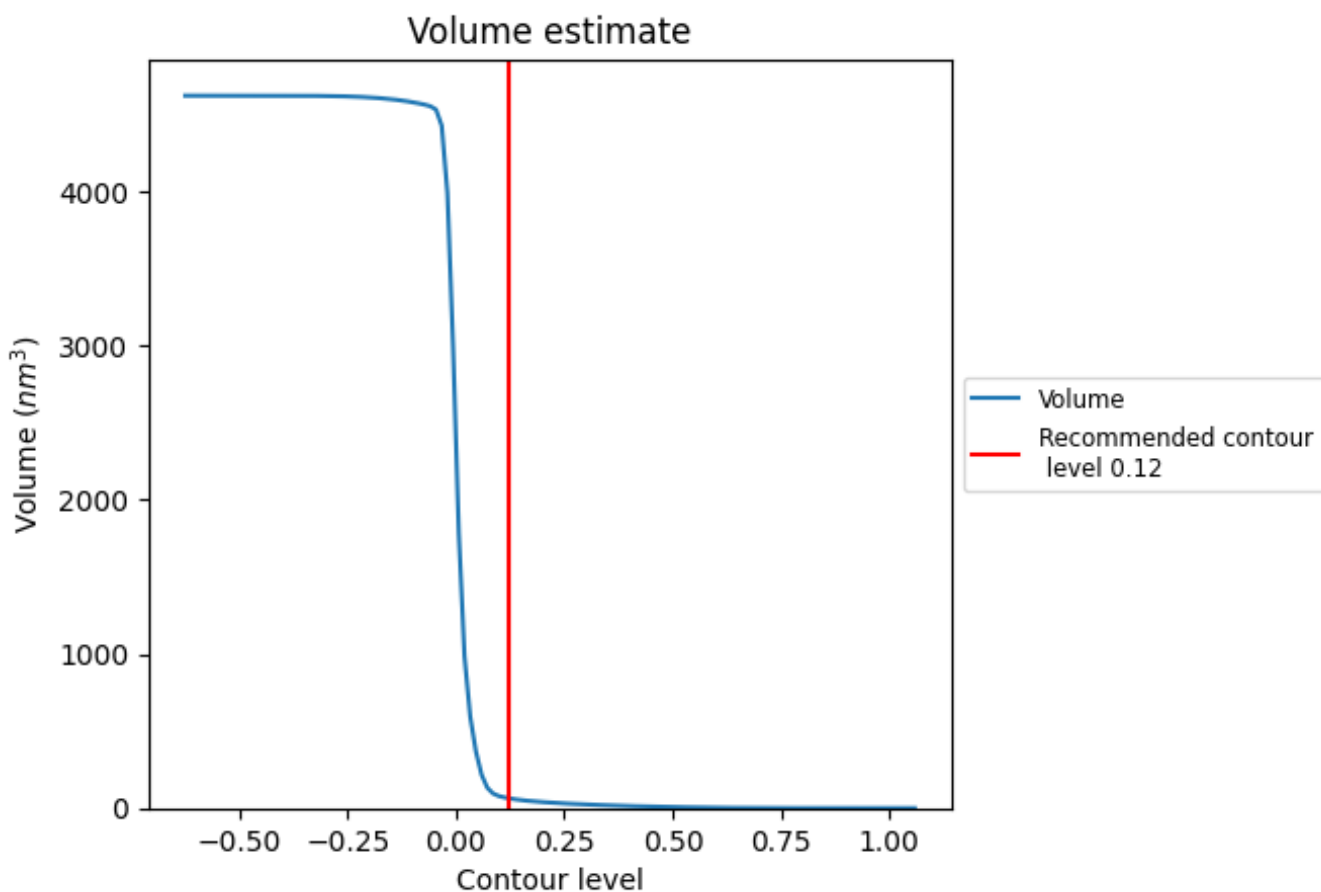
This section contains the results of statistical analysis of the map.

### 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.

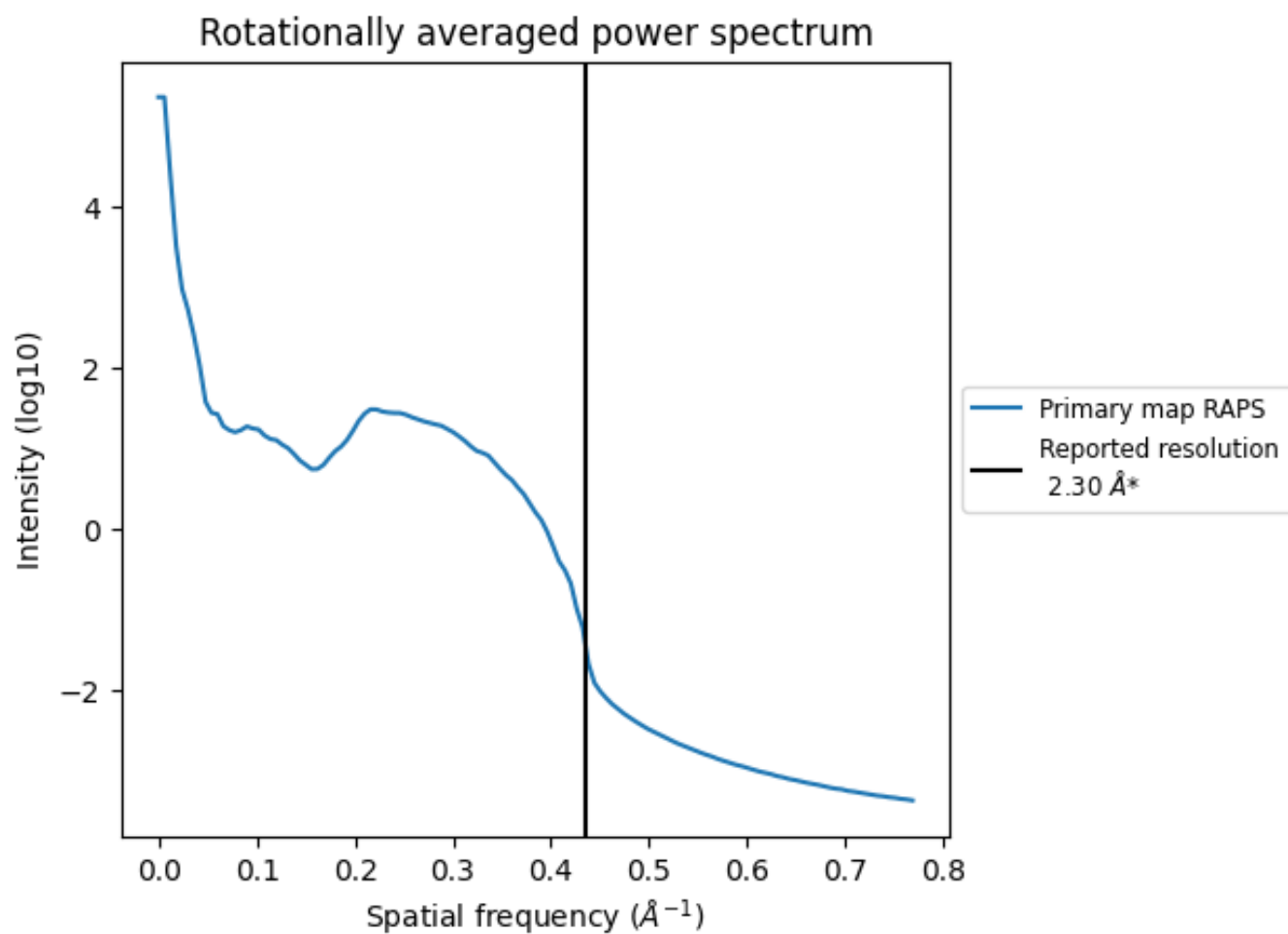
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is  $65 \text{ nm}^3$ ; this corresponds to an approximate mass of 59 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 0.435 Å<sup>-1</sup>

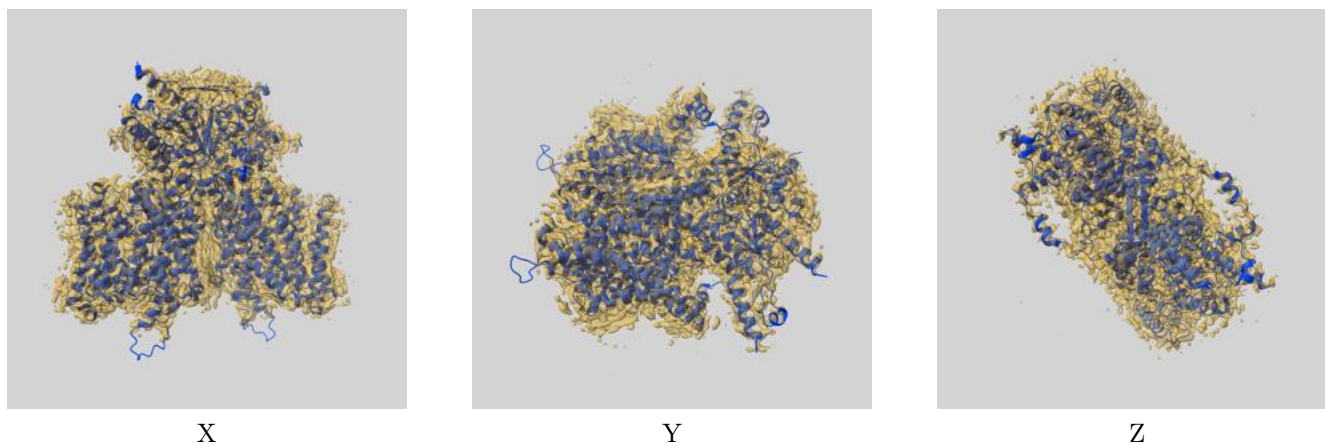
## 8 Fourier-Shell correlation

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-23329 and PDB model 7LGU. Per-residue inclusion information can be found in section 3 on page 14.

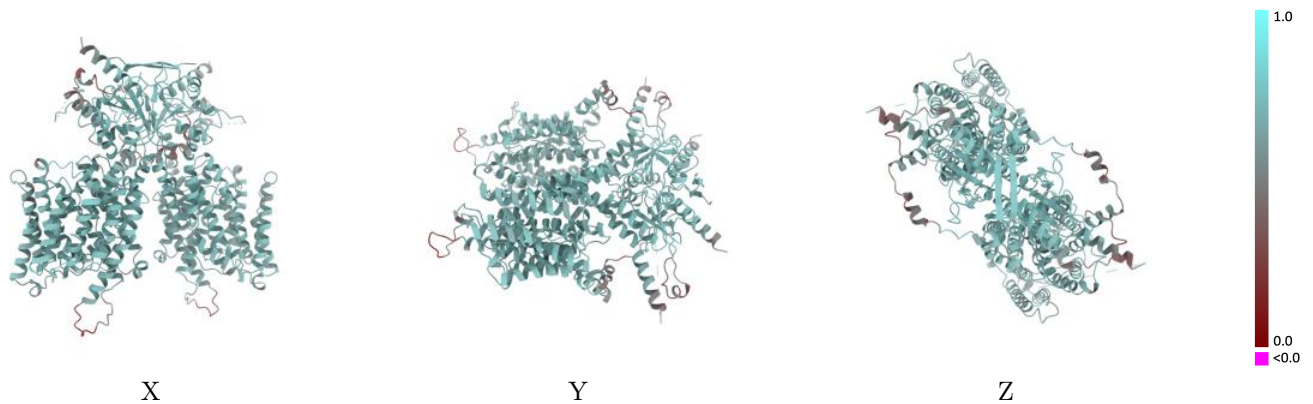
### 9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.12 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.

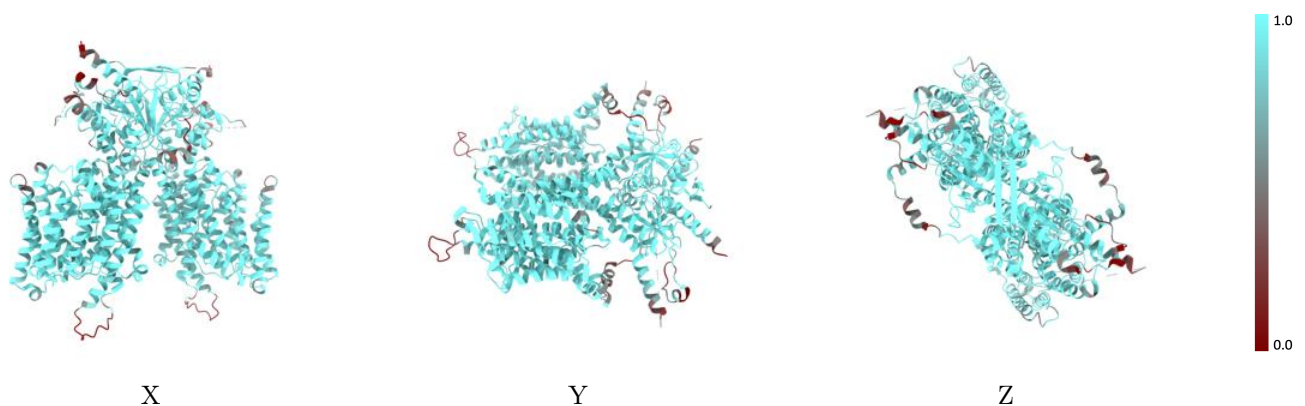


## 9.2 Q-score mapped to coordinate model [\(i\)](#)



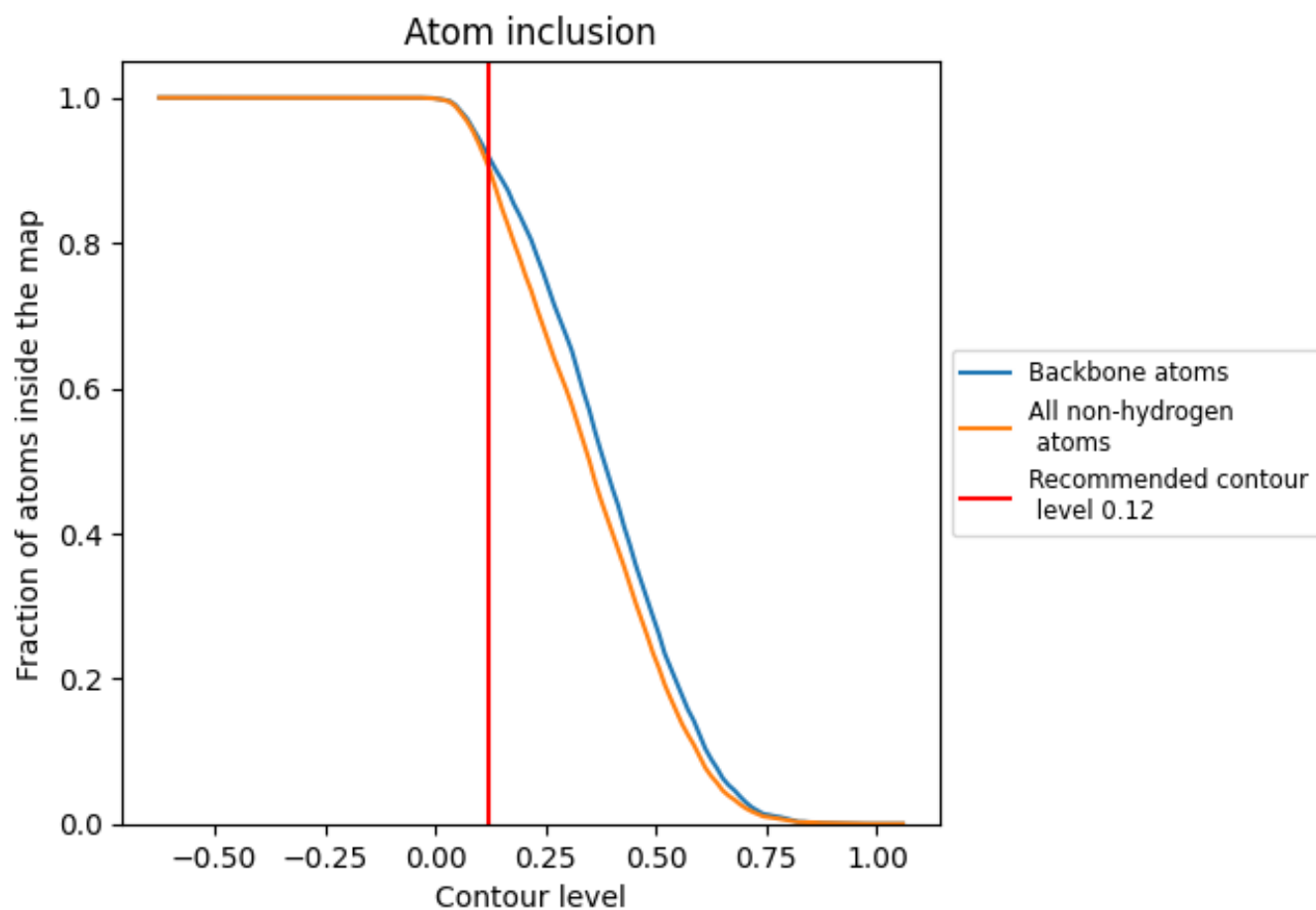
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.12).







## 9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 91% 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 (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9059	 0.6440
A	 0.9061	 0.6450
B	 0.9067	 0.6440

