



Full wwPDB EM Validation Report ⓘ

May 18, 2024 – 09:35 PM EDT

PDB ID : 7LGW
EMDB ID : EMD-23331
Title : Structure of human Prestin in nanodisc in the presence of NaCl
Authors : Ge, J.; Gouaux, E.
Deposited on : 2021-01-21
Resolution : 2.70 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

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.dev92
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

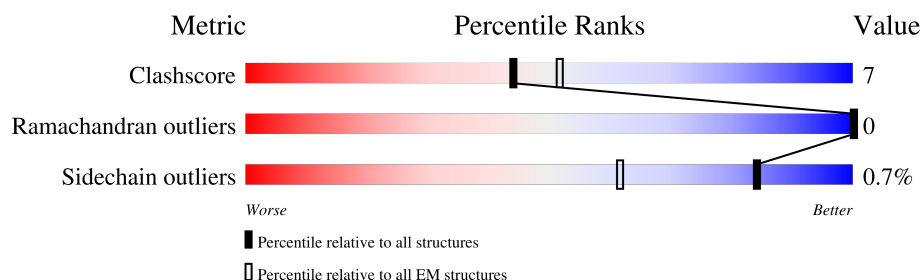
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.70 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 ≥ 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	<div> <div>5%</div> <div>76%</div> <div>14%</div> <div>9%</div> </div>
1	B	750	<div> <div>5%</div> <div>76%</div> <div>15%</div> <div>9%</div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 11390 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
1	A	680	Total	C	N	O	S	0	0
			5226	3431	845	926	24		
1	B	680	Total	C	N	O	S	0	0
			5226	3431	845	926	24		

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).

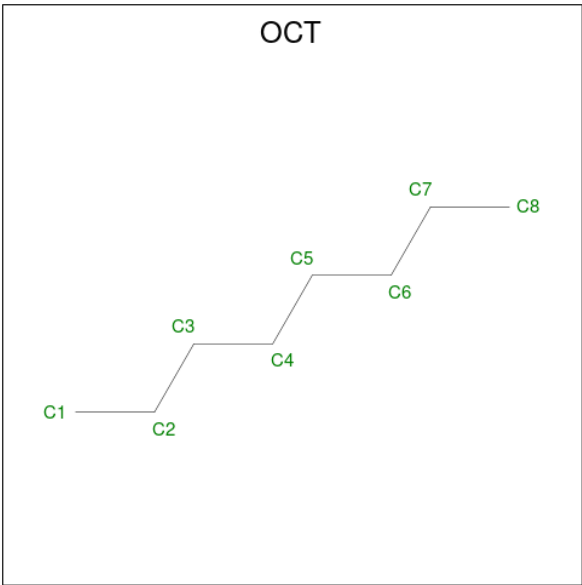
Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total	Cl	0
			1	1	
2	B	1	Total	Cl	0
			1	1	

- Molecule 3 is CHOLESTEROL (three-letter code: CLR) (formula: C₂₇H₄₆O).



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: C₈H₁₈).



Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	C	0
			8	8	
4	A	1	Total	C	0
			8	8	

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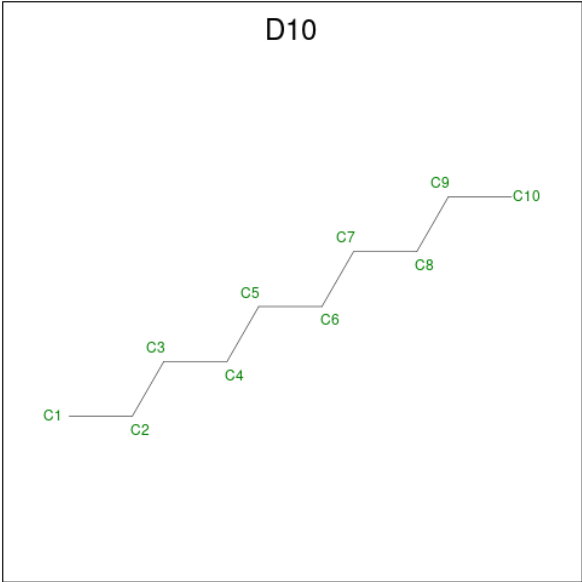
Mol	Chain	Residues	Atoms	AltConf
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total C 8 8	0
4	A	1	Total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

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Mol	Chain	Residues	Atoms	AltConf
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0
4	B	1	Total C 8 8	0

- Molecule 5 is DECANE (three-letter code: D10) (formula: C₁₀H₂₂).



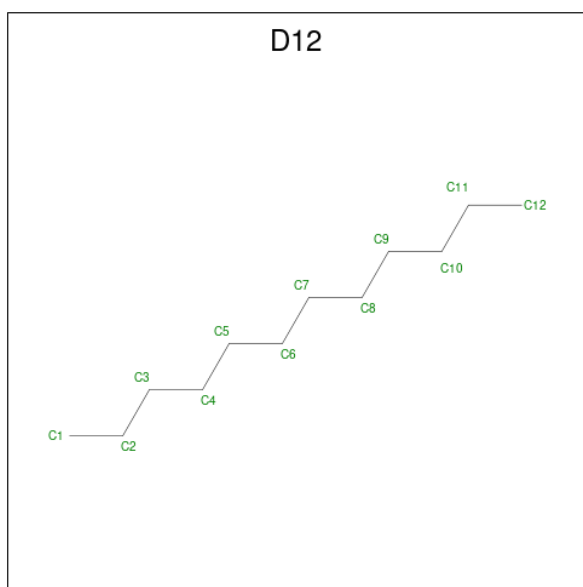
Mol	Chain	Residues	Atoms	AltConf
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0

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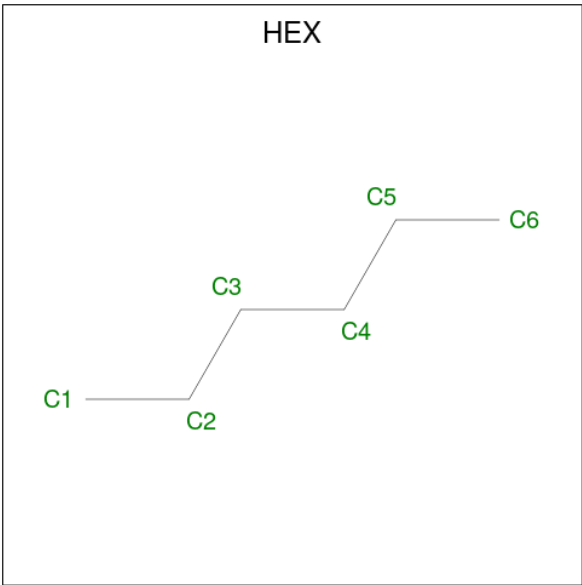
Mol	Chain	Residues	Atoms	AltConf
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	A	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0
5	B	1	Total C 10 10	0

- Molecule 6 is DODECANE (three-letter code: D12) (formula: C₁₂H₂₆).



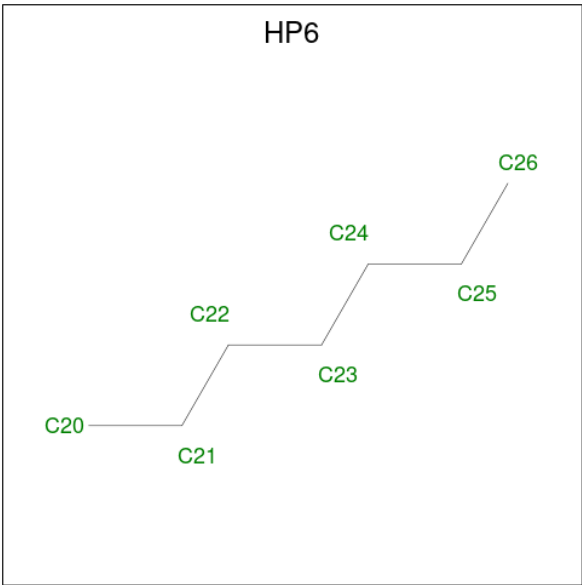
Mol	Chain	Residues	Atoms	AltConf
6	A	1	Total C 12 12	0
6	A	1	Total C 12 12	0
6	A	1	Total C 12 12	0
6	A	1	Total C 12 12	0
6	A	1	Total C 12 12	0
6	B	1	Total C 12 12	0
6	B	1	Total C 12 12	0
6	B	1	Total C 12 12	0
6	B	1	Total C 12 12	0
6	B	1	Total C 12 12	0

- Molecule 7 is HEXANE (three-letter code: HEX) (formula: C₆H₁₄).



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

- Molecule 8 is HEPTANE (three-letter code: HP6) (formula: C₇H₁₆).



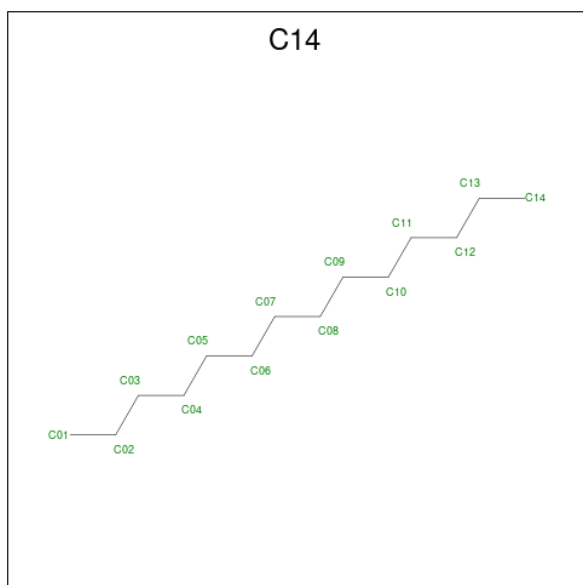
Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total	C	0
			7	7	
8	A	1	Total	C	0
			7	7	

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Mol	Chain	Residues	Atoms	AltConf
8	B	1	Total C 7 7	0
8	B	1	Total C 7 7	0

- Molecule 9 is TETRADECANE (three-letter code: C14) (formula: C₁₄H₃₀).



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

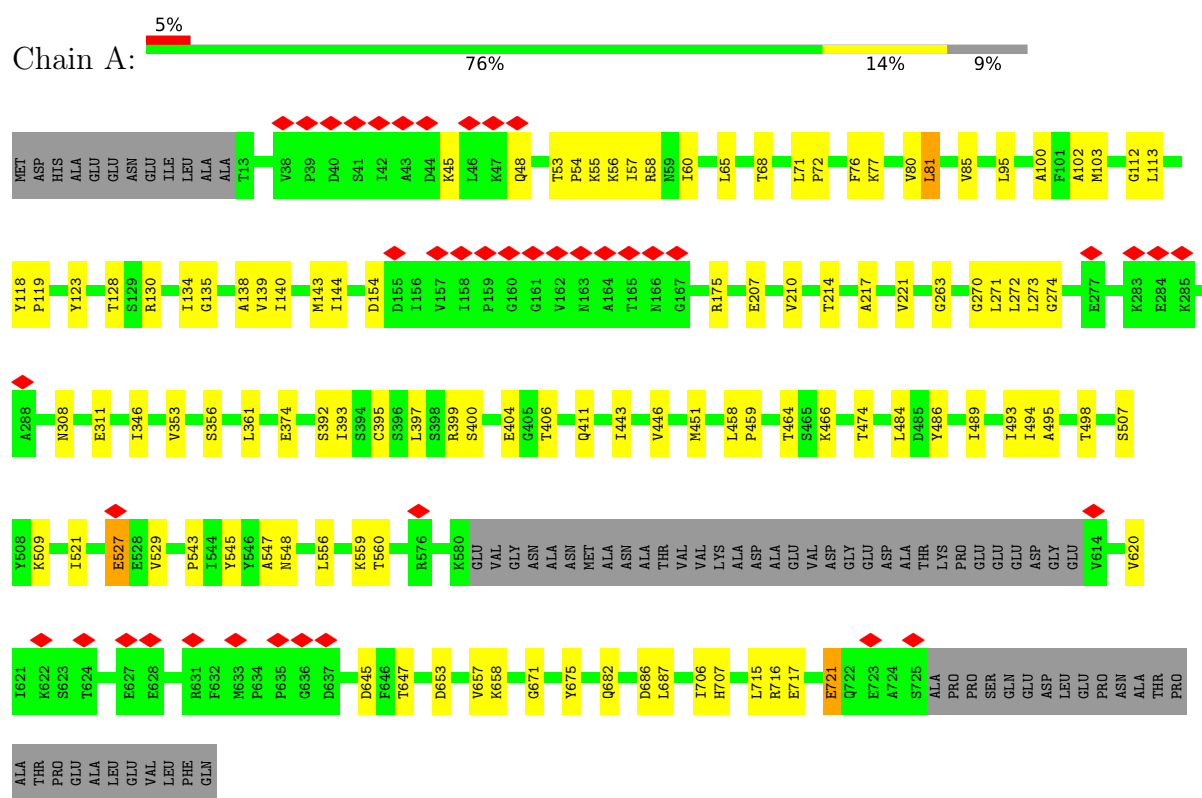
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	AltConf
10	A	15	Total O 15 15	0
10	B	15	Total O 15 15	0

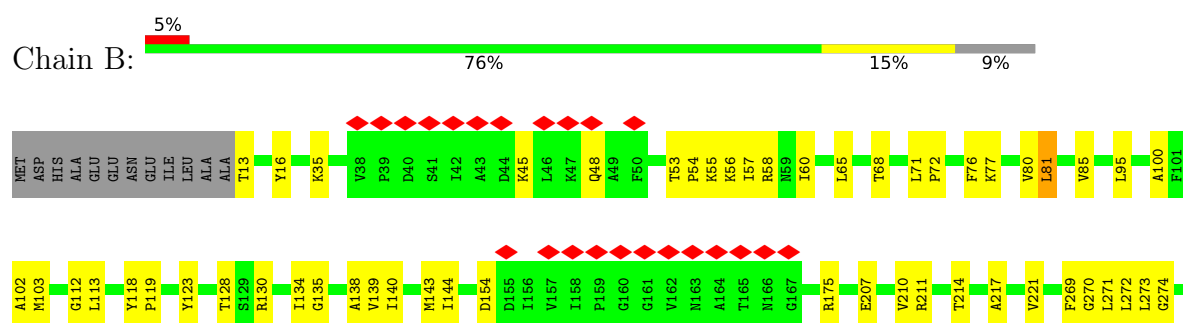
3 Residue-property plots

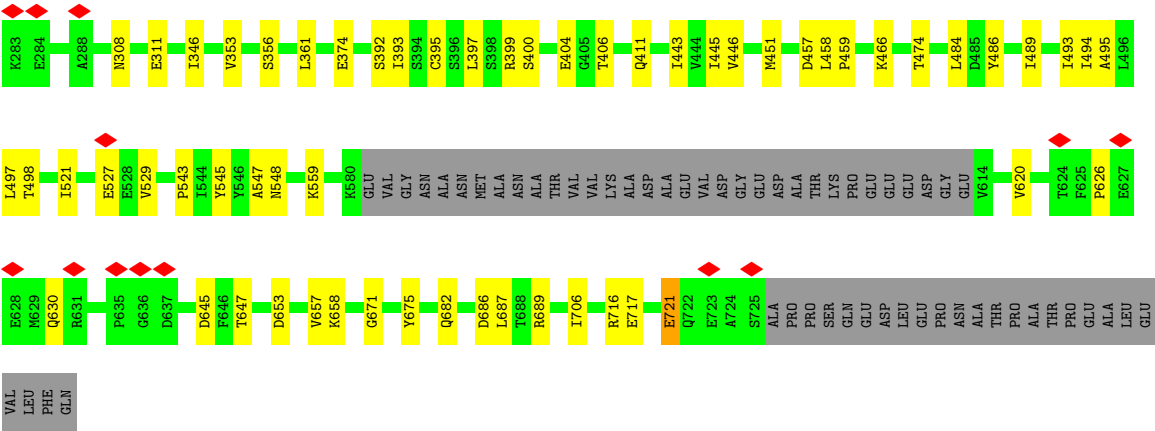
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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	155000	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	0.860	Depositor
Minimum map value	-0.443	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.037	Depositor
Recommended contour level	0.09	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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.27	0/5344	0.42	0/7266
1	B	0.27	0/5344	0.42	0/7266
All	All	0.27	0/10688	0.42	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 [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	5226	0	5384	74	0
1	B	5226	0	5384	75	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	28	0	46	7	0
3	B	28	0	46	11	0
4	A	152	0	342	15	0
4	B	152	0	342	9	0
5	A	170	0	374	6	0
5	B	160	0	352	10	0
6	A	60	0	130	2	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	4	0
8	B	14	0	32	0	0
9	A	28	0	60	1	0
9	B	28	0	60	1	0
10	A	15	0	0	0	0
10	B	15	0	0	0	0
All	All	11390	0	12742	166	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 7.

All (166) 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:B:497:LEU:CD1	3:B:802:CLR:H261	2.06	0.86
3:B:802:CLR:H152	5:B:807:D10:H62	1.64	0.78
1:B:497:LEU:HD12	3:B:802:CLR:H261	1.66	0.78
1:A:119:PRO:HB3	1:A:393:ILE:HD13	1.71	0.73
1:B:119:PRO:HB3	1:B:393:ILE:HD13	1.71	0.73
3:B:802:CLR:H161	5:B:807:D10:H82	1.73	0.70
4:A:847:OCT:H21	4:A:848:OCT:H52	1.78	0.66
6:A:823:D12:H71	4:A:824:OCT:H82	1.78	0.66
1:A:138:ALA:HB3	1:A:221:VAL:HG22	1.78	0.65
1:A:214:THR:HB	1:A:451:MET:HE1	1.79	0.65
1:B:138:ALA:HB3	1:B:221:VAL:HG22	1.79	0.65
6:B:823:D12:H71	4:B:824:OCT:H82	1.78	0.64
1:B:497:LEU:HD12	3:B:802:CLR:C26	2.27	0.64
1:A:308:ASN:ND2	1:A:311:GLU:OE1	2.32	0.63
1:B:308:ASN:ND2	1:B:311:GLU:OE1	2.32	0.63
3:B:802:CLR:C15	5:B:807:D10:H62	2.28	0.63
1:A:495:ALA:O	1:A:498:THR:OG1	2.18	0.61
1:A:493:ILE:HG21	3:A:802:CLR:H241	1.83	0.61
1:B:272:LEU:HB3	1:B:446:VAL:HG11	1.83	0.61
1:A:272:LEU:HB3	1:A:446:VAL:HG11	1.82	0.61
1:B:77:LYS:HE3	4:B:844:OCT:H31	1.83	0.61
1:B:214:THR:HB	1:B:451:MET:HE1	1.83	0.60
1:B:102:ALA:HB2	1:B:112:GLY:HA3	1.84	0.60
1:B:210:VAL:HG13	1:B:356:SER:HB3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:495:ALA:O	1:B:498:THR:OG1	2.18	0.60
6:A:815:D12:H52	5:A:816:D10:H81	1.85	0.59
6:B:815:D12:H52	5:B:816:D10:H81	1.85	0.59
1:A:210:VAL:HG13	1:A:356:SER:HB3	1.83	0.59
1:A:102:ALA:HB2	1:A:112:GLY:HA3	1.84	0.59
1:A:494:ILE:HG12	3:A:802:CLR:H263	1.85	0.59
1:A:77:LYS:HE3	4:A:844:OCT:H31	1.83	0.59
4:A:837:OCT:H71	8:A:838:HP6:H221	1.84	0.59
1:A:80:VAL:HG13	1:A:81:LEU:HD23	1.85	0.58
1:A:545:TYR:HA	1:A:653:ASP:OD1	2.03	0.58
1:A:207:GLU:OE1	1:B:658:LYS:NZ	2.36	0.58
1:B:545:TYR:HA	1:B:653:ASP:OD1	2.04	0.58
1:B:80:VAL:HG13	1:B:81:LEU:HD23	1.85	0.58
1:B:100:ALA:HA	1:B:103:MET:HE2	1.86	0.58
3:B:802:CLR:C16	5:B:807:D10:H82	2.33	0.58
1:A:395:CYS:HG	1:A:400:SER:HG	1.51	0.57
1:B:486:TYR:CE1	3:B:802:CLR:H12	2.40	0.57
1:A:100:ALA:HA	1:A:103:MET:HE2	1.87	0.57
1:A:274:GLY:HA2	4:A:847:OCT:H72	1.87	0.56
1:B:657:VAL:HG22	1:B:687:LEU:HD23	1.87	0.56
1:A:543:PRO:O	1:A:548:ASN:ND2	2.38	0.56
1:B:274:GLY:HA2	4:B:847:OCT:H72	1.87	0.56
1:B:543:PRO:O	1:B:548:ASN:ND2	2.38	0.56
1:A:493:ILE:CG2	3:A:802:CLR:H241	2.36	0.56
1:B:497:LEU:HD11	3:B:802:CLR:H261	1.87	0.55
4:B:805:OCT:H52	5:B:816:D10:H41	1.89	0.55
1:A:657:VAL:HG22	1:A:687:LEU:HD23	1.88	0.55
1:A:489:ILE:HB	5:A:812:D10:H62	1.89	0.55
4:A:805:OCT:H52	5:A:816:D10:H41	1.89	0.54
1:B:489:ILE:HB	5:B:812:D10:H62	1.89	0.54
1:B:645:ASP:OD1	1:B:647:THR:OG1	2.23	0.54
4:A:847:OCT:H32	4:A:848:OCT:H31	1.91	0.53
1:B:270:GLY:HA2	4:B:847:OCT:H13	1.91	0.53
1:B:154:ASP:OD1	1:B:175:ARG:NH2	2.38	0.52
1:A:645:ASP:OD1	1:A:647:THR:OG1	2.23	0.52
1:A:270:GLY:HA2	4:A:847:OCT:H13	1.91	0.52
1:B:645:ASP:OD2	1:B:706:ILE:HD11	2.10	0.51
1:A:645:ASP:OD2	1:A:706:ILE:HD11	2.10	0.51
1:B:65:LEU:O	1:B:68:THR:OG1	2.26	0.51
1:A:715:LEU:HD21	1:B:13:THR:HG22	1.91	0.51
1:A:53:THR:HG23	1:A:56:LYS:H	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:LEU:O	1:A:68:THR:OG1	2.26	0.51
1:A:527:GLU:HB2	1:B:35:LYS:HE3	1.93	0.50
1:B:53:THR:HG23	1:B:56:LYS:H	1.76	0.50
4:A:837:OCT:H51	8:A:838:HP6:H241	1.92	0.50
1:B:54:PRO:HA	1:B:57:ILE:HG12	1.94	0.50
1:B:395:CYS:HG	1:B:400:SER:HG	1.58	0.50
1:A:54:PRO:HA	1:A:57:ILE:HG12	1.94	0.50
1:A:154:ASP:OD1	1:A:175:ARG:NH2	2.38	0.49
1:A:263:GLY:HA3	8:A:838:HP6:H211	1.94	0.49
1:B:620:VAL:HG13	1:B:671:GLY:HA2	1.95	0.49
1:A:620:VAL:HG13	1:A:671:GLY:HA2	1.95	0.49
1:B:139:VAL:HG11	1:B:484:LEU:HD12	1.95	0.49
1:A:717:GLU:O	1:A:721:GLU:HG2	2.13	0.48
1:B:139:VAL:O	1:B:143:MET:HG3	2.14	0.48
1:B:717:GLU:O	1:B:721:GLU:HG2	2.13	0.48
1:A:55:LYS:HE3	1:A:58:ARG:HH21	1.79	0.48
1:A:139:VAL:HG11	1:A:484:LEU:HD12	1.95	0.48
3:A:802:CLR:H151	5:A:807:D10:H62	1.95	0.47
1:B:55:LYS:HE3	1:B:58:ARG:HH21	1.79	0.47
1:A:494:ILE:CG1	3:A:802:CLR:H263	2.44	0.47
1:B:144:ILE:HD11	1:B:392:SER:HB3	1.96	0.47
1:B:466:LYS:HB2	9:B:830:C14:H021	1.97	0.47
1:B:76:PHE:HB3	4:B:844:OCT:H22	1.97	0.47
4:B:805:OCT:H21	4:B:806:OCT:H32	1.97	0.47
1:A:139:VAL:O	1:A:143:MET:HG3	2.14	0.47
4:A:805:OCT:H21	4:A:806:OCT:H32	1.97	0.47
1:B:272:LEU:HD11	1:B:443:ILE:HG12	1.97	0.46
1:A:95:LEU:HD12	1:A:113:LEU:HD21	1.97	0.46
1:A:144:ILE:HD11	1:A:392:SER:HB3	1.96	0.46
1:A:272:LEU:HD11	1:A:443:ILE:HG12	1.97	0.46
1:B:486:TYR:CZ	3:B:802:CLR:H12	2.51	0.46
1:A:85:VAL:HG12	1:A:406:THR:HG22	1.98	0.46
1:A:458:LEU:HB2	1:A:459:PRO:HD3	1.98	0.46
1:A:658:LYS:NZ	1:B:207:GLU:OE1	2.45	0.46
1:B:95:LEU:HD12	1:B:113:LEU:HD21	1.97	0.46
1:A:466:LYS:HB2	9:A:830:C14:H021	1.96	0.46
1:A:76:PHE:HB3	4:A:844:OCT:H22	1.97	0.46
1:B:395:CYS:SG	1:B:400:SER:OG	2.66	0.46
6:B:811:D12:H72	6:B:811:D12:H41	1.80	0.46
1:A:682:GLN:NE2	1:A:686:ASP:OD2	2.49	0.45
1:B:474:THR:HG21	1:B:494:ILE:HG21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:675:TYR:OH	1:B:716:ARG:NH1	2.49	0.45
5:B:834:D10:H62	5:B:834:D10:H92	1.75	0.45
1:B:85:VAL:HG12	1:B:406:THR:HG22	1.98	0.45
1:A:135:GLY:HA3	1:A:399:ARG:HH12	1.82	0.45
1:B:682:GLN:NE2	1:B:686:ASP:OD2	2.50	0.45
1:B:128:THR:OG1	1:B:374:GLU:OE2	2.32	0.45
1:B:135:GLY:HA3	1:B:399:ARG:HH12	1.82	0.45
1:A:675:TYR:OH	1:A:716:ARG:NH1	2.49	0.45
1:B:458:LEU:HB2	1:B:459:PRO:HD3	1.98	0.45
1:A:486:TYR:CE1	3:A:802:CLR:H12	2.52	0.44
1:A:707:HIS:HB3	1:B:16:TYR:HE2	1.81	0.44
1:B:521:ILE:HB	1:B:529:VAL:HG11	1.99	0.44
1:B:217:ALA:O	1:B:221:VAL:HG23	2.18	0.44
1:A:140:ILE:HG23	1:A:346:ILE:HG23	2.00	0.44
1:A:217:ALA:O	1:A:221:VAL:HG23	2.18	0.44
1:A:474:THR:HG21	1:A:494:ILE:HG21	1.99	0.44
1:B:71:LEU:HB3	1:B:72:PRO:HD3	2.00	0.44
1:B:140:ILE:HG23	1:B:346:ILE:HG23	2.00	0.43
1:B:273:LEU:HD23	1:B:273:LEU:HA	1.87	0.43
1:A:556:LEU:O	1:A:560:THR:OG1	2.29	0.43
1:B:559:LYS:HE2	1:B:559:LYS:HB3	1.81	0.43
1:A:119:PRO:HG2	1:A:393:ILE:HG21	2.00	0.43
1:B:56:LYS:O	1:B:60:ILE:HG12	2.18	0.43
1:B:119:PRO:HG2	1:B:393:ILE:HG21	2.00	0.43
1:A:521:ILE:HB	1:A:529:VAL:HG11	1.99	0.43
1:A:56:LYS:O	1:A:60:ILE:HG12	2.18	0.43
1:A:71:LEU:HB3	1:A:72:PRO:HD3	2.00	0.43
1:A:464:THR:O	1:B:689:ARG:HD2	2.19	0.43
1:B:361:LEU:HD22	1:B:404:GLU:HG3	2.00	0.43
4:A:847:OCT:H12	4:A:848:OCT:H13	2.00	0.43
1:A:361:LEU:HD22	1:A:404:GLU:HG3	2.00	0.43
1:A:123:TYR:OH	1:A:374:GLU:OE1	2.31	0.42
1:A:128:THR:OG1	1:A:374:GLU:OE2	2.32	0.42
1:B:211:ARG:NE	1:B:457:ASP:OD2	2.27	0.42
4:A:837:OCT:H32	8:A:838:HP6:H241	2.02	0.42
1:B:545:TYR:CE2	1:B:547:ALA:HB3	2.54	0.42
1:A:71:LEU:HD23	4:A:841:OCT:H51	2.02	0.41
1:A:353:VAL:HG13	1:A:397:LEU:HD21	2.02	0.41
1:B:353:VAL:HG13	1:B:397:LEU:HD21	2.02	0.41
1:B:493:ILE:HG21	3:B:802:CLR:H25	2.02	0.41
1:A:130:ARG:HA	1:A:411:GLN:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:545:TYR:CE2	1:A:547:ALA:HB3	2.54	0.41
1:B:81:LEU:HD11	4:B:846:OCT:H31	2.03	0.41
1:A:81:LEU:HD11	4:A:846:OCT:H31	2.03	0.41
1:A:118:TYR:HE1	5:A:822:D10:H81	1.86	0.41
1:A:273:LEU:HD23	1:A:273:LEU:HA	1.88	0.41
1:A:559:LYS:HE2	1:A:559:LYS:HB3	1.81	0.41
1:A:45:LYS:O	1:A:48:GLN:HG3	2.21	0.41
1:A:123:TYR:CG	1:A:134:ILE:HG13	2.56	0.41
1:B:123:TYR:CG	1:B:134:ILE:HG13	2.56	0.41
1:B:130:ARG:HA	1:B:411:GLN:HG2	2.02	0.41
1:B:269:PHE:HE2	1:B:445:ILE:HG22	1.86	0.41
5:B:834:D10:H22	5:B:834:D10:H51	1.88	0.41
1:A:493:ILE:HG21	3:A:802:CLR:C24	2.49	0.41
5:A:834:D10:H62	5:A:834:D10:H92	1.75	0.41
1:B:45:LYS:O	1:B:48:GLN:HG3	2.20	0.40
1:B:118:TYR:HE1	5:B:822:D10:H81	1.86	0.40
1:B:626:PRO:O	1:B:630:GLN:N	2.52	0.40
1:B:71:LEU:HD23	4:B:841:OCT:H51	2.02	0.40
1:A:507:SER:HB3	1:A:509:LYS:NZ	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	676/750 (90%)	660 (98%)	16 (2%)	0	100	100
1	B	676/750 (90%)	660 (98%)	16 (2%)	0	100	100
All	All	1352/1500 (90%)	1320 (98%)	32 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	567/631 (90%)	563 (99%)	4 (1%)	84	94
1	B	567/631 (90%)	563 (99%)	4 (1%)	84	94
All	All	1134/1262 (90%)	1126 (99%)	8 (1%)	84	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	LEU
1	A	271	LEU
1	A	527	GLU
1	A	721	GLU
1	B	81	LEU
1	B	271	LEU
1	B	527	GLU
1	B	721	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	650	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 95 ligands modelled in this entry, 2 are monoatomic - leaving 93 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	A	808	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	A	842	-	7,7,7	0.29	0	6,6,6	0.74	0
5	D10	A	818	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	B	844	-	7,7,7	0.30	0	6,6,6	0.71	0
5	D10	A	836	-	9,9,9	0.29	0	8,8,8	0.81	0
6	D12	B	810	-	11,11,11	0.29	0	10,10,10	0.83	0
5	D10	A	832	-	9,9,9	0.28	0	8,8,8	0.82	0
4	OCT	B	845	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	B	841	-	7,7,7	0.29	0	6,6,6	0.73	0
5	D10	A	828	-	9,9,9	0.30	0	8,8,8	0.79	0
5	D10	A	835	-	9,9,9	0.29	0	8,8,8	0.79	0
8	HP6	B	826	-	6,6,6	0.29	0	5,5,5	0.68	0
5	D10	B	828	-	9,9,9	0.30	0	8,8,8	0.78	0
5	D10	B	832	-	9,9,9	0.28	0	8,8,8	0.82	0
5	D10	A	840	-	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	0.80	0
4	OCT	A	844	-	7,7,7	0.30	0	6,6,6	0.72	0
6	D12	B	838	-	11,11,11	0.29	0	10,10,10	0.85	0
6	D12	A	839	-	11,11,11	0.29	0	10,10,10	0.84	0
8	HP6	B	837	-	6,6,6	0.31	0	5,5,5	0.68	0
3	CLR	B	802	-	31,31,31	0.26	0	48,48,48	0.57	1 (2%)
4	OCT	B	808	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	A	803	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	B	843	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	827	-	7,7,7	0.30	0	6,6,6	0.74	0
3	CLR	A	802	-	31,31,31	0.28	0	48,48,48	0.39	0
5	D10	B	813	-	9,9,9	0.29	0	8,8,8	0.80	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	OCT	A	809	-	7,7,7	0.29	0	6,6,6	0.77	0
5	D10	A	831	-	9,9,9	0.28	0	8,8,8	0.82	0
6	D12	B	815	-	11,11,11	0.28	0	10,10,10	0.85	0
9	C14	A	830	-	13,13,13	0.27	0	12,12,12	0.87	0
4	OCT	B	847	-	7,7,7	0.29	0	6,6,6	0.72	0
9	C14	B	829	-	13,13,13	0.28	0	12,12,12	0.88	0
4	OCT	B	824	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	A	847	-	7,7,7	0.30	0	6,6,6	0.73	0
6	D12	B	823	-	11,11,11	0.29	0	10,10,10	0.85	0
4	OCT	B	836	-	7,7,7	0.30	0	6,6,6	0.71	0
4	OCT	A	848	-	7,7,7	0.31	0	6,6,6	0.70	0
5	D10	B	822	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	B	809	-	7,7,7	0.28	0	6,6,6	0.77	0
5	D10	B	835	-	9,9,9	0.30	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	B	805	-	7,7,7	0.29	0	6,6,6	0.74	0
5	D10	B	839	-	9,9,9	0.29	0	8,8,8	0.78	0
5	D10	A	833	-	9,9,9	0.29	0	8,8,8	0.82	0
4	OCT	A	843	-	7,7,7	0.29	0	6,6,6	0.73	0
4	OCT	B	803	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	A	813	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	B	842	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	A	804	-	7,7,7	0.29	0	6,6,6	0.73	0
7	HEX	B	825	-	5,5,5	0.30	0	4,4,4	0.58	0
5	D10	A	820	-	9,9,9	0.29	0	8,8,8	0.80	0
4	OCT	A	819	-	7,7,7	0.30	0	6,6,6	0.73	0
6	D12	A	823	-	11,11,11	0.29	0	10,10,10	0.84	0
4	OCT	A	845	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	A	816	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	A	814	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	B	812	-	9,9,9	0.30	0	8,8,8	0.79	0
8	HP6	A	826	-	6,6,6	0.30	0	5,5,5	0.69	0
9	C14	A	829	-	13,13,13	0.28	0	12,12,12	0.88	0
5	D10	B	831	-	9,9,9	0.28	0	8,8,8	0.82	0
7	HEX	A	825	-	5,5,5	0.31	0	4,4,4	0.58	0
5	D10	B	816	-	9,9,9	0.29	0	8,8,8	0.80	0
4	OCT	A	841	-	7,7,7	0.29	0	6,6,6	0.72	0
9	C14	B	830	-	13,13,13	0.27	0	12,12,12	0.88	0
5	D10	B	833	-	9,9,9	0.29	0	8,8,8	0.82	0
4	OCT	B	840	-	7,7,7	0.30	0	6,6,6	0.72	0
6	D12	A	811	-	11,11,11	0.29	0	10,10,10	0.86	0
5	D10	B	820	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	A	807	-	9,9,9	0.29	0	8,8,8	0.80	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	HP6	A	838	-	6,6,6	0.29	0	5,5,5	0.65	0
4	OCT	B	819	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	A	822	-	9,9,9	0.29	0	8,8,8	0.80	0
6	D12	B	811	-	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	0.75	0
5	D10	A	834	-	9,9,9	0.29	0	8,8,8	0.80	0
4	OCT	A	805	-	7,7,7	0.30	0	6,6,6	0.74	0
4	OCT	B	817	-	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	0.73	0
5	D10	A	812	-	9,9,9	0.29	0	8,8,8	0.79	0
4	OCT	A	824	-	7,7,7	0.30	0	6,6,6	0.73	0
5	D10	B	807	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	A	827	-	7,7,7	0.29	0	6,6,6	0.74	0
4	OCT	A	837	-	7,7,7	0.30	0	6,6,6	0.73	0
4	OCT	B	806	-	7,7,7	0.30	0	6,6,6	0.74	0
5	D10	B	818	-	9,9,9	0.29	0	8,8,8	0.81	0
4	OCT	B	804	-	7,7,7	0.29	0	6,6,6	0.73	0
4	OCT	A	817	-	7,7,7	0.30	0	6,6,6	0.73	0
6	D12	A	815	-	11,11,11	0.28	0	10,10,10	0.85	0
4	OCT	A	846	-	7,7,7	0.30	0	6,6,6	0.73	0
6	D12	A	810	-	11,11,11	0.29	0	10,10,10	0.83	0
5	D10	A	821	-	9,9,9	0.29	0	8,8,8	0.80	0
5	D10	B	814	-	9,9,9	0.30	0	8,8,8	0.80	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	A	808	-	-	0/5/5/5	-
4	OCT	A	842	-	-	0/5/5/5	-
5	D10	A	818	-	-	1/7/7/7	-
4	OCT	B	844	-	-	0/5/5/5	-
5	D10	A	836	-	-	0/7/7/7	-
6	D12	B	810	-	-	2/9/9/9	-
5	D10	A	832	-	-	0/7/7/7	-
4	OCT	B	845	-	-	0/5/5/5	-
4	OCT	B	841	-	-	0/5/5/5	-
5	D10	A	828	-	-	0/7/7/7	-
5	D10	A	835	-	-	0/7/7/7	-
8	HP6	B	826	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	D10	B	828	-	-	0/7/7/7	-
5	D10	B	832	-	-	0/7/7/7	-
5	D10	A	840	-	-	0/7/7/7	-
5	D10	B	834	-	-	0/7/7/7	-
4	OCT	A	844	-	-	0/5/5/5	-
6	D12	B	838	-	-	1/9/9/9	-
6	D12	A	839	-	-	1/9/9/9	-
8	HP6	B	837	-	-	2/4/4/4	-
3	CLR	B	802	-	-	0/10/68/68	0/4/4/4
4	OCT	B	808	-	-	0/5/5/5	-
4	OCT	A	803	-	-	0/5/5/5	-
4	OCT	B	843	-	-	0/5/5/5	-
4	OCT	B	827	-	-	0/5/5/5	-
3	CLR	A	802	-	-	0/10/68/68	0/4/4/4
5	D10	B	813	-	-	0/7/7/7	-
4	OCT	A	809	-	-	0/5/5/5	-
5	D10	A	831	-	-	1/7/7/7	-
6	D12	B	815	-	-	0/9/9/9	-
9	C14	A	830	-	-	0/11/11/11	-
4	OCT	B	847	-	-	1/5/5/5	-
9	C14	B	829	-	-	0/11/11/11	-
4	OCT	B	824	-	-	0/5/5/5	-
4	OCT	A	847	-	-	1/5/5/5	-
6	D12	B	823	-	-	0/9/9/9	-
4	OCT	B	836	-	-	0/5/5/5	-
4	OCT	A	848	-	-	1/5/5/5	-
5	D10	B	822	-	-	0/7/7/7	-
4	OCT	B	809	-	-	0/5/5/5	-
5	D10	B	835	-	-	0/7/7/7	-
5	D10	B	821	-	-	0/7/7/7	-
4	OCT	B	805	-	-	0/5/5/5	-
5	D10	B	839	-	-	0/7/7/7	-
5	D10	A	833	-	-	0/7/7/7	-
4	OCT	A	843	-	-	0/5/5/5	-
4	OCT	B	803	-	-	0/5/5/5	-
5	D10	A	813	-	-	0/7/7/7	-
4	OCT	B	842	-	-	0/5/5/5	-
4	OCT	A	804	-	-	1/5/5/5	-
7	HEX	B	825	-	-	0/3/3/3	-
5	D10	A	820	-	-	0/7/7/7	-
4	OCT	A	819	-	-	0/5/5/5	-
6	D12	A	823	-	-	0/9/9/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	OCT	A	845	-	-	0/5/5/5	-
5	D10	A	816	-	-	0/7/7/7	-
5	D10	A	814	-	-	0/7/7/7	-
5	D10	B	812	-	-	0/7/7/7	-
8	HP6	A	826	-	-	0/4/4/4	-
9	C14	A	829	-	-	0/11/11/11	-
5	D10	B	831	-	-	1/7/7/7	-
7	HEX	A	825	-	-	0/3/3/3	-
5	D10	B	816	-	-	0/7/7/7	-
4	OCT	A	841	-	-	0/5/5/5	-
9	C14	B	830	-	-	0/11/11/11	-
5	D10	B	833	-	-	0/7/7/7	-
4	OCT	B	840	-	-	0/5/5/5	-
6	D12	A	811	-	-	2/9/9/9	-
5	D10	B	820	-	-	0/7/7/7	-
5	D10	A	807	-	-	0/7/7/7	-
8	HP6	A	838	-	-	3/4/4/4	-
4	OCT	B	819	-	-	0/5/5/5	-
5	D10	A	822	-	-	0/7/7/7	-
6	D12	B	811	-	-	2/9/9/9	-
4	OCT	A	806	-	-	0/5/5/5	-
5	D10	A	834	-	-	0/7/7/7	-
4	OCT	A	805	-	-	0/5/5/5	-
4	OCT	B	817	-	-	0/5/5/5	-
4	OCT	B	846	-	-	0/5/5/5	-
5	D10	A	812	-	-	0/7/7/7	-
4	OCT	A	824	-	-	0/5/5/5	-
5	D10	B	807	-	-	0/7/7/7	-
4	OCT	A	827	-	-	0/5/5/5	-
4	OCT	A	837	-	-	2/5/5/5	-
4	OCT	B	806	-	-	0/5/5/5	-
5	D10	B	818	-	-	1/7/7/7	-
4	OCT	B	804	-	-	1/5/5/5	-
4	OCT	A	817	-	-	0/5/5/5	-
6	D12	A	815	-	-	0/9/9/9	-
4	OCT	A	846	-	-	0/5/5/5	-
6	D12	A	810	-	-	2/9/9/9	-
5	D10	A	821	-	-	0/7/7/7	-
5	D10	B	814	-	-	0/7/7/7	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	802	CLR	C16-C17-C13	-2.46	100.88	103.84

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	810	D12	C4-C5-C6-C7
6	B	810	D12	C4-C5-C6-C7
6	B	810	D12	C3-C4-C5-C6
6	A	810	D12	C3-C4-C5-C6
6	A	811	D12	C7-C8-C9-C10
6	B	811	D12	C7-C8-C9-C10
4	A	837	OCT	C4-C5-C6-C7
8	B	837	HP6	C21-C22-C23-C24
8	A	838	HP6	C21-C22-C23-C24
6	B	838	D12	C11-C10-C9-C8
6	A	839	D12	C11-C10-C9-C8
5	A	831	D10	C4-C5-C6-C7
5	B	831	D10	C4-C5-C6-C7
8	A	838	HP6	C20-C21-C22-C23
4	A	804	OCT	C4-C5-C6-C7
4	B	804	OCT	C4-C5-C6-C7
4	A	848	OCT	C3-C4-C5-C6
6	B	811	D12	C3-C4-C5-C6
6	A	811	D12	C3-C4-C5-C6
8	A	838	HP6	C22-C23-C24-C25
4	A	847	OCT	C2-C3-C4-C5
4	B	847	OCT	C2-C3-C4-C5
5	A	818	D10	C6-C7-C8-C9
5	B	818	D10	C6-C7-C8-C9
4	A	837	OCT	C5-C6-C7-C8
8	B	837	HP6	C20-C21-C22-C23

There are no ring outliers.

36 monomers are involved in 55 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	844	OCT	2	0
4	B	841	OCT	1	0
5	B	834	D10	2	0

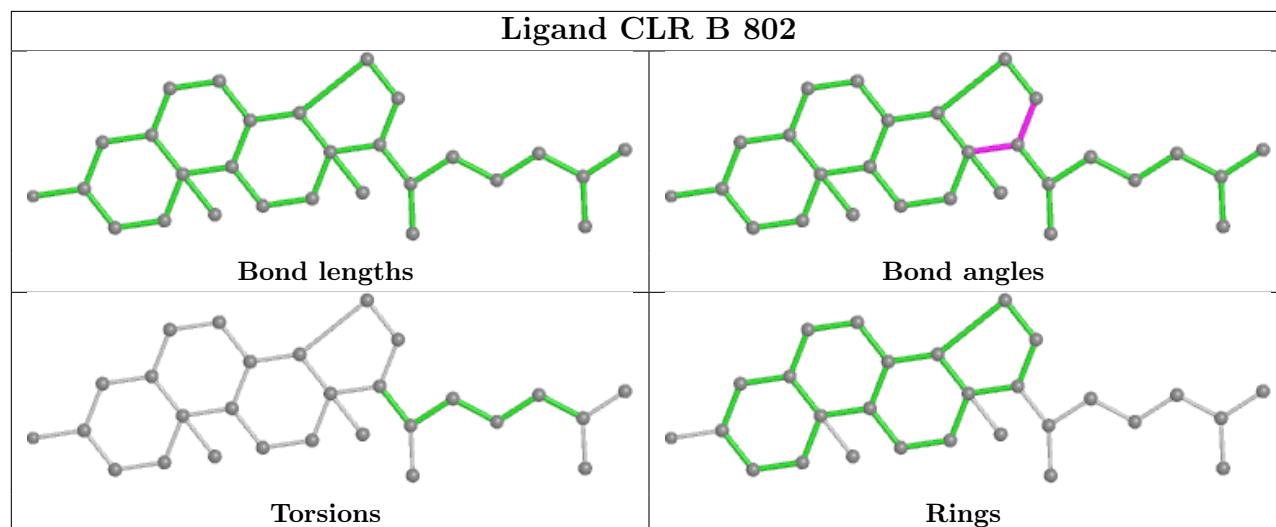
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	844	OCT	2	0
3	B	802	CLR	11	0
3	A	802	CLR	7	0
6	B	815	D12	1	0
9	A	830	C14	1	0
4	B	847	OCT	2	0
4	B	824	OCT	1	0
4	A	847	OCT	5	0
6	B	823	D12	1	0
4	A	848	OCT	3	0
5	B	822	D10	1	0
4	B	805	OCT	2	0
6	A	823	D12	1	0
5	A	816	D10	2	0
5	B	812	D10	1	0
5	B	816	D10	2	0
4	A	841	OCT	1	0
9	B	830	C14	1	0
5	A	807	D10	1	0
8	A	838	HP6	4	0
5	A	822	D10	1	0
6	B	811	D12	1	0
4	A	806	OCT	1	0
5	A	834	D10	1	0
4	A	805	OCT	2	0
4	B	846	OCT	1	0
5	A	812	D10	1	0
4	A	824	OCT	1	0
5	B	807	D10	4	0
4	A	837	OCT	3	0
4	B	806	OCT	1	0
6	A	815	D12	1	0
4	A	846	OCT	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

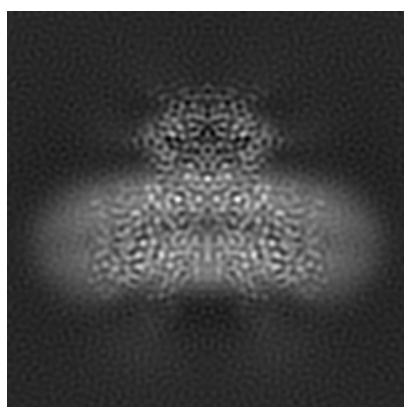
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23331. These allow visual inspection of the internal detail of the map and identification of artifacts.

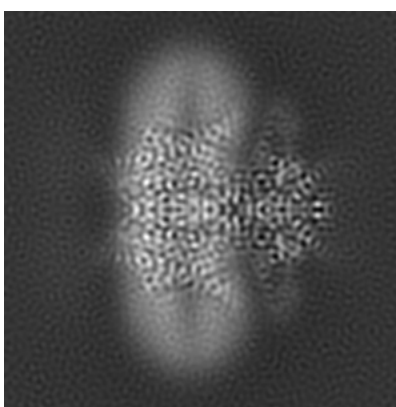
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

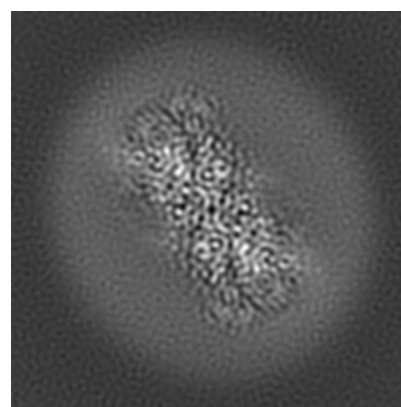
6.1.1 Primary map



X



Y

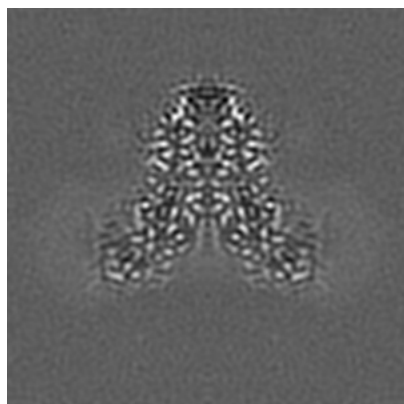


Z

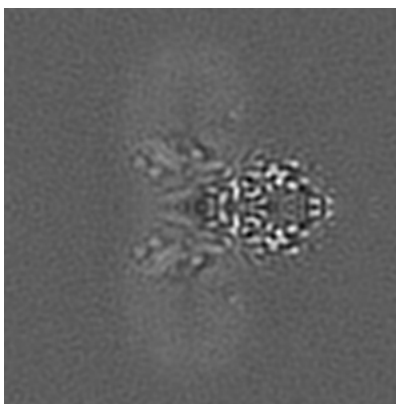
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

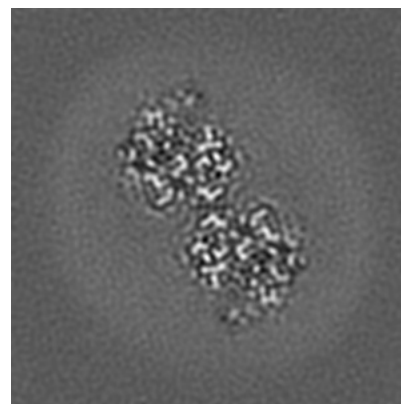
6.2.1 Primary map



X Index: 128



Y Index: 128

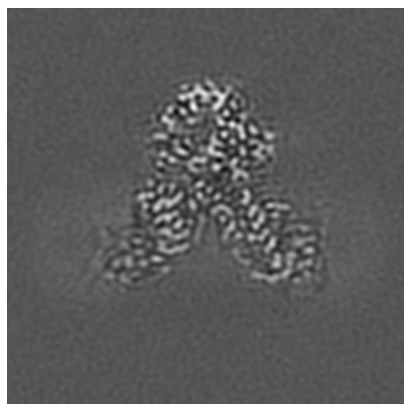


Z Index: 128

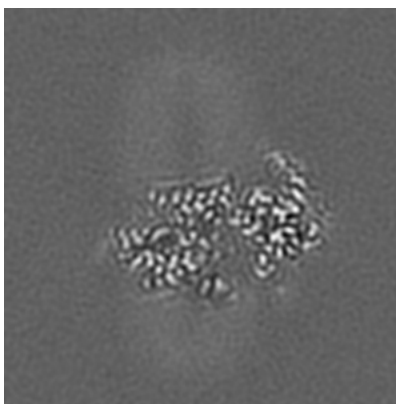
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

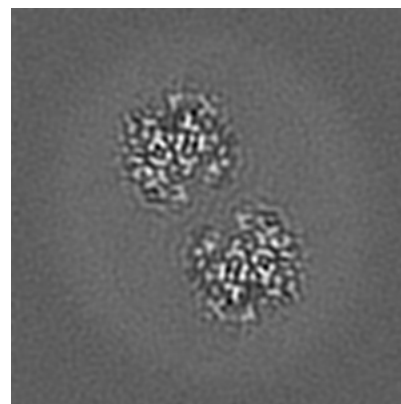
6.3.1 Primary map



X Index: 125



Y Index: 150

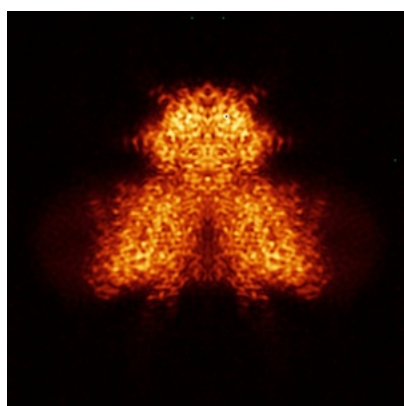


Z Index: 99

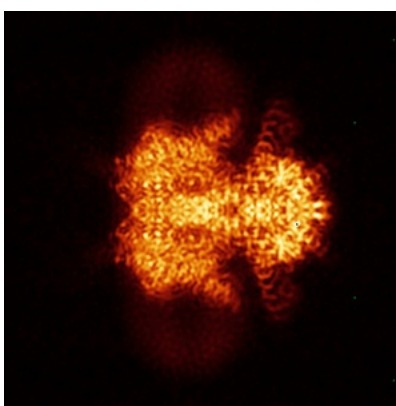
The images above show the largest variance slices of the map in three orthogonal directions.

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

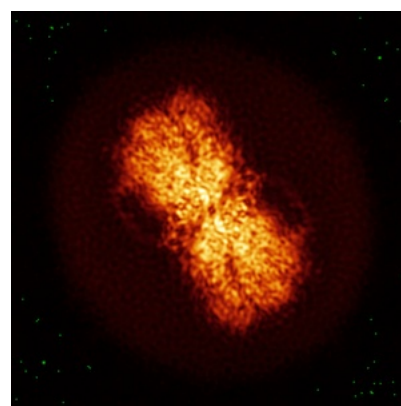
6.4.1 Primary map



X



Y

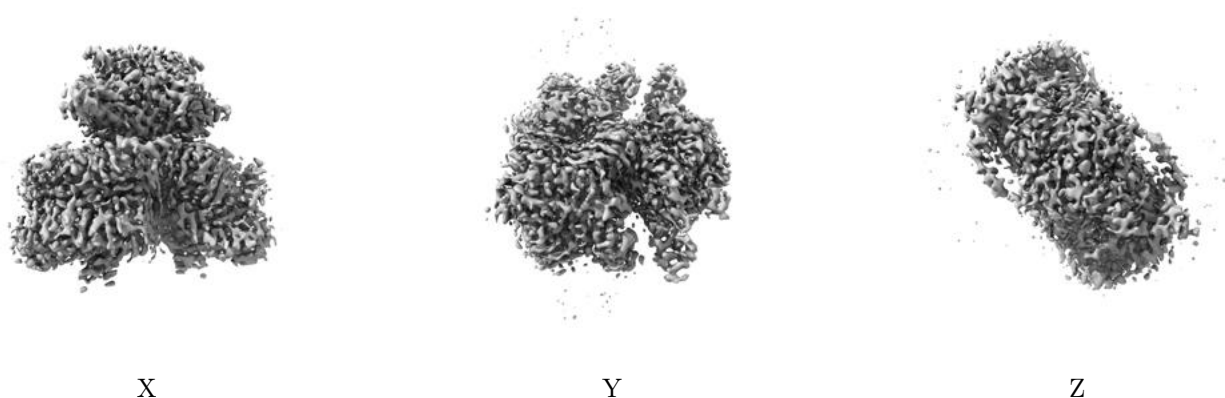


Z

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

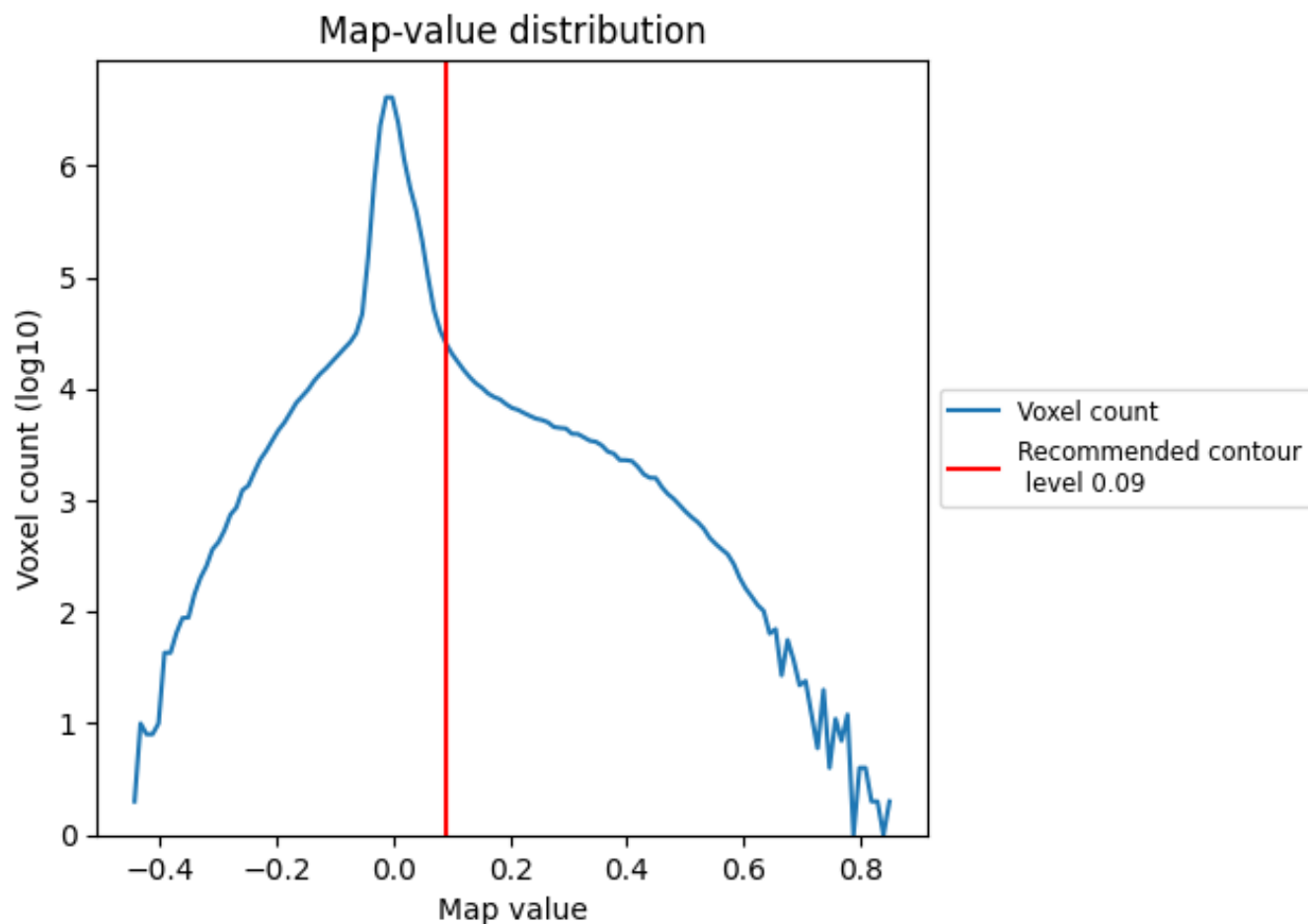
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

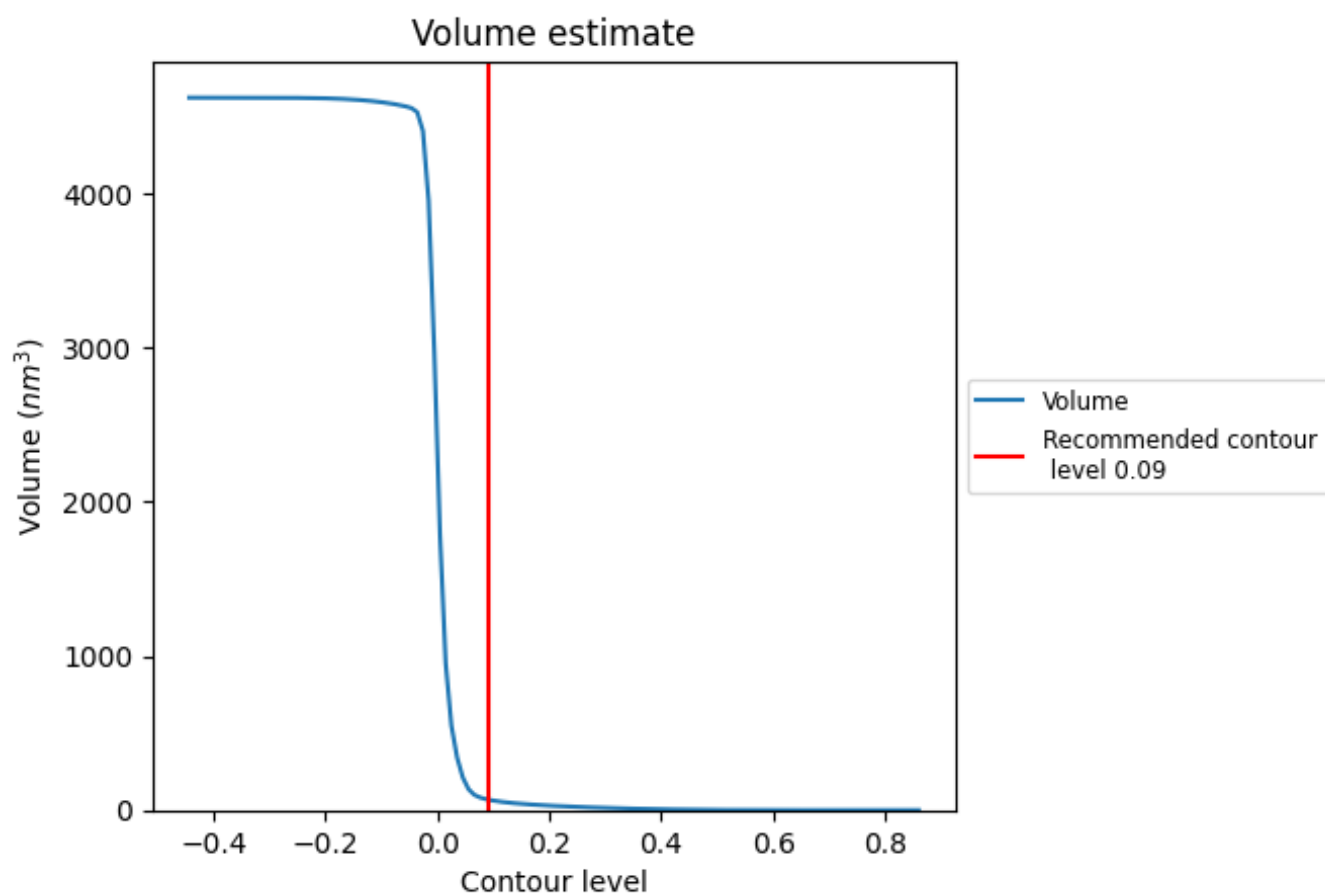
This section contains the results of statistical analysis of the map.

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.

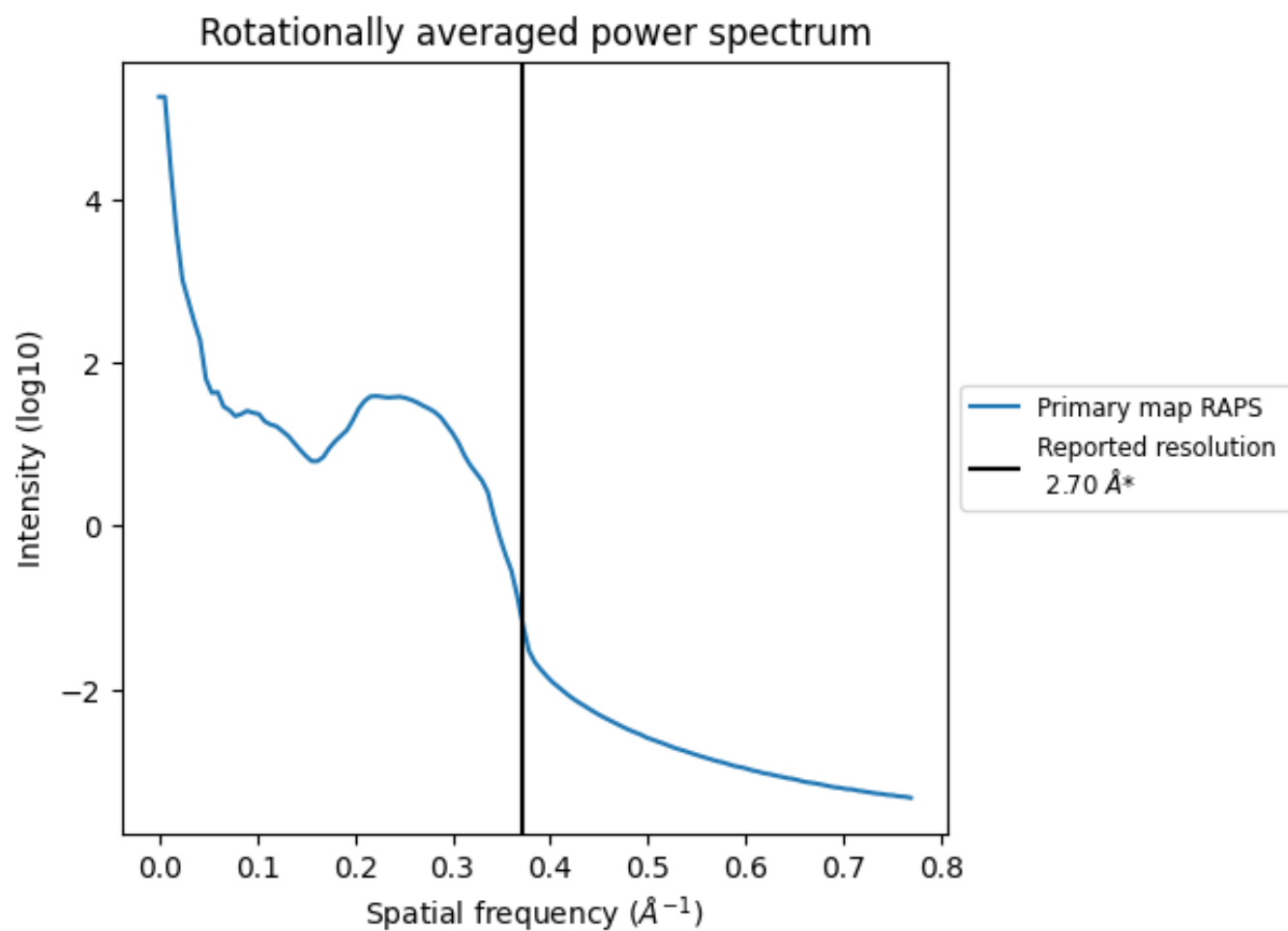
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 69 nm^3 ; this corresponds to an approximate mass of 62 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 0.370 Å⁻¹

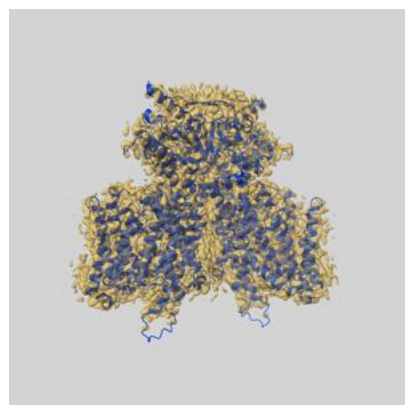
8 Fourier-Shell correlation

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

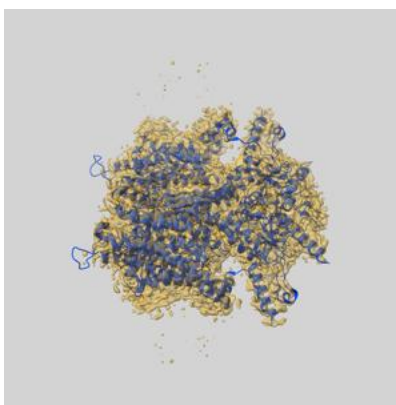
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23331 and PDB model 7LGW. Per-residue inclusion information can be found in section [3](#) on page [12](#).

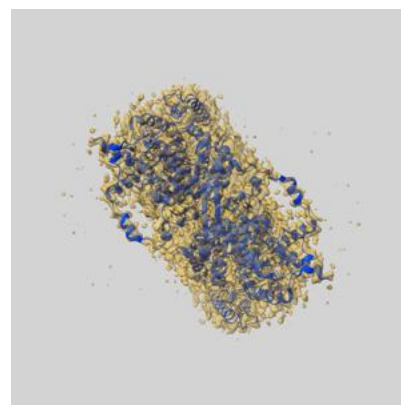
9.1 Map-model overlay [i](#)



X



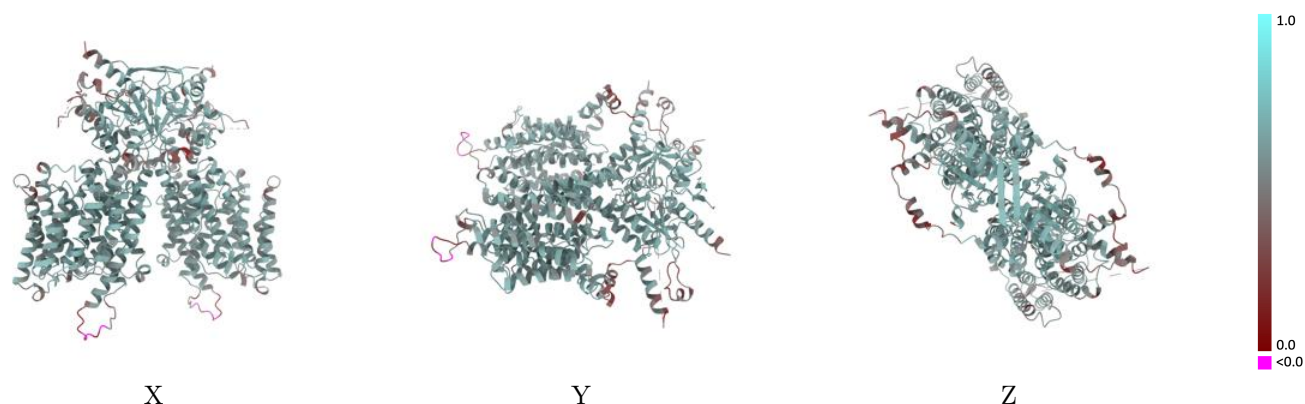
Y



Z

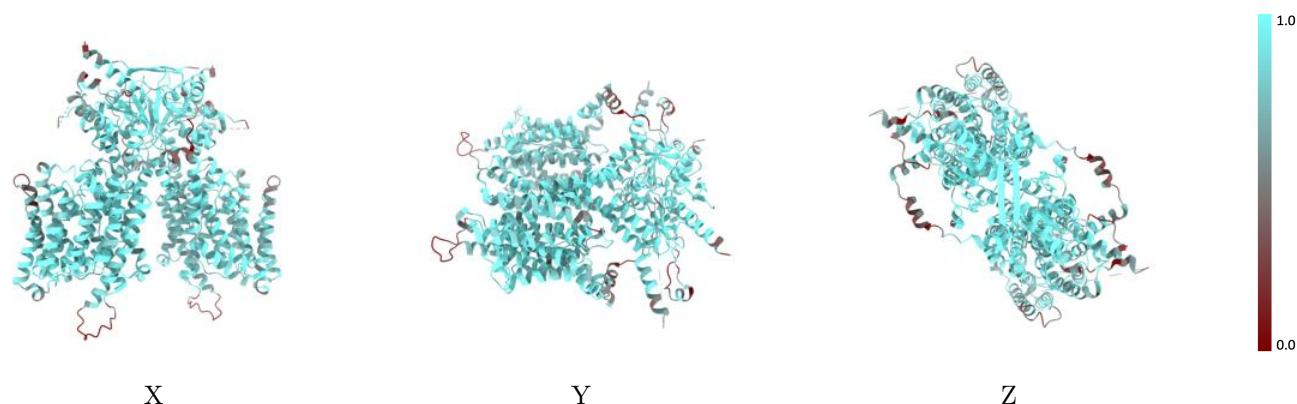
The images above show the 3D surface view of the map at the recommended contour level 0.09 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



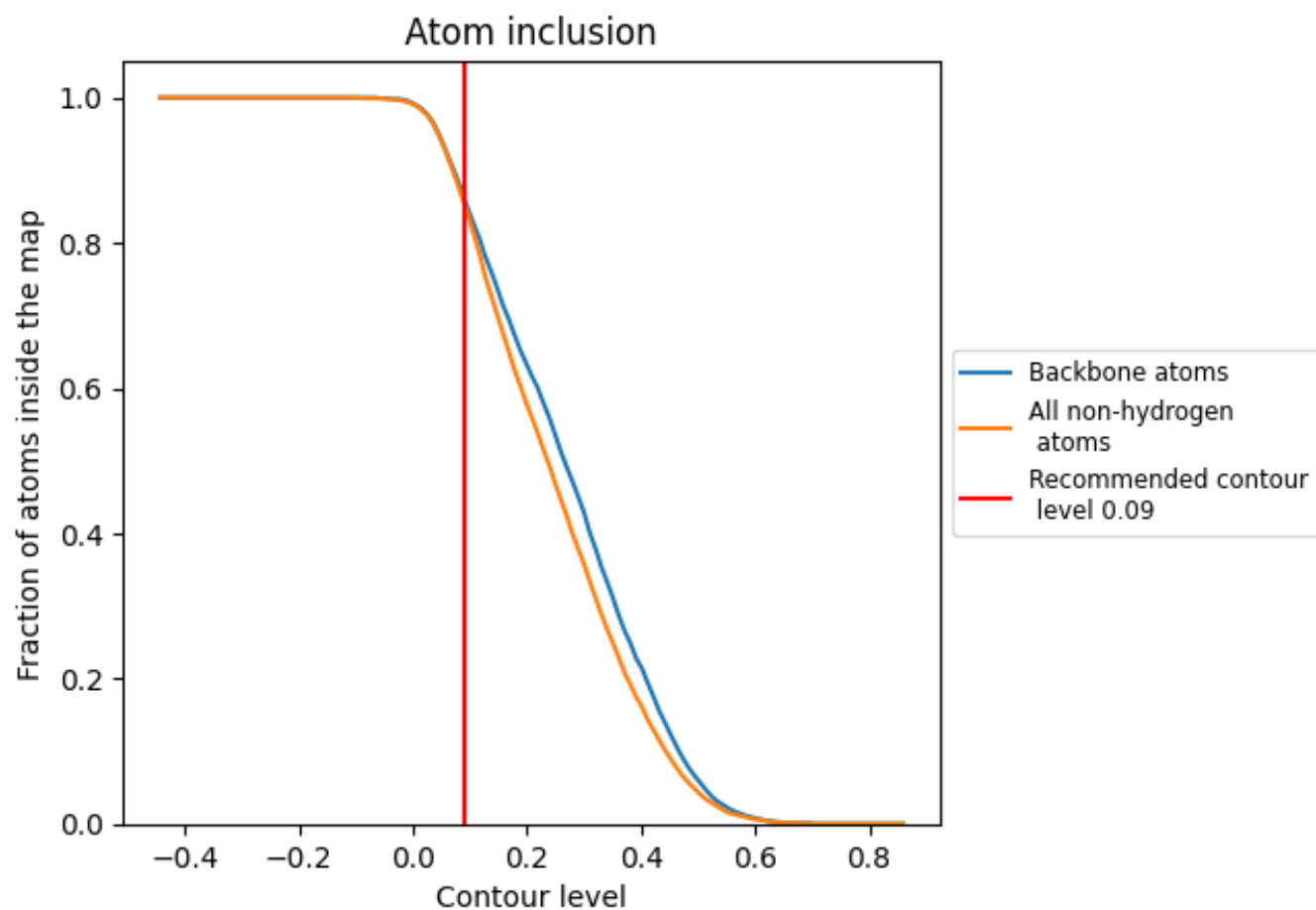
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.09).

9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 85% 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 (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.8540	<div><div></div></div> 0.5670
A	<div><div></div></div> 0.8540	<div><div></div></div> 0.5650
B	<div><div></div></div> 0.8540	<div><div></div></div> 0.5680

