



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 10:46 PM UTC

PDB ID : 6N07 / pdb_00006n07
EMDB ID : EMD-9312
Title : Structure of the HO BMC shell: BMC-TD focused map, open inner pore, compacted shell
Authors : Greber, B.J.; Sutter, M.; Kerfeld, C.A.
Deposited on : 2018-11-06
Resolution : 3.60 Å(reported)
Based on initial model : 5V74

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

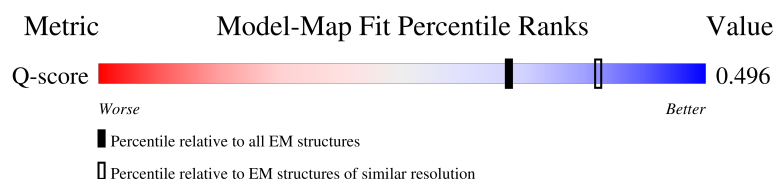
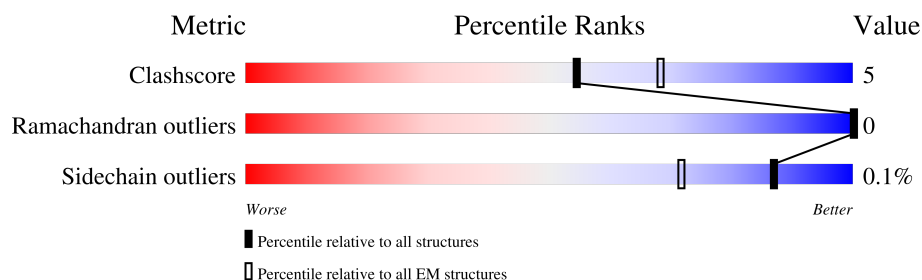
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



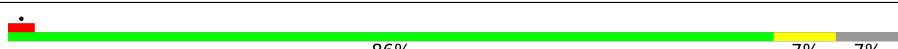
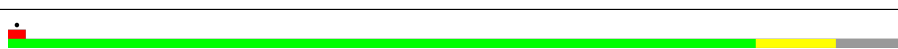
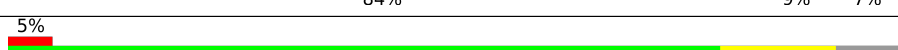
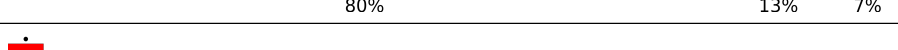
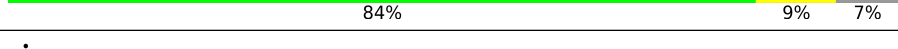
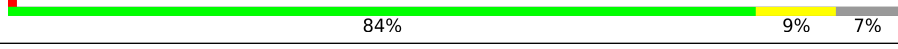
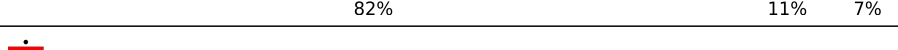
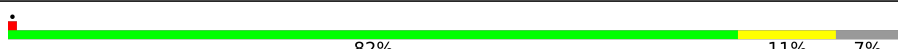
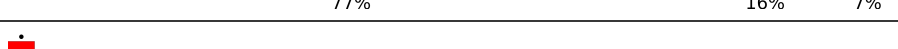
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	12797 (3.10 - 4.10)

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	205	
1	B	205	
1	C	205	
1	D	205	

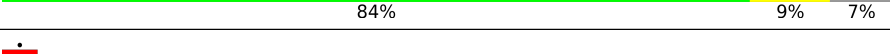
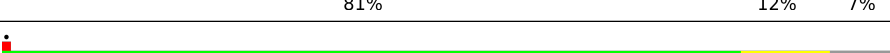
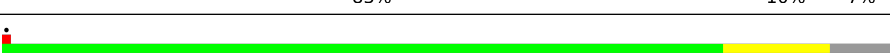
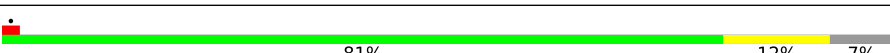
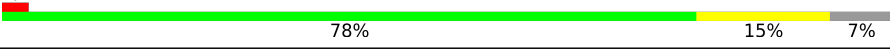
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Mol	Chain	Length	Quality of chain
1	E	205	
1	F	205	
2	GA	99	
2	GB	99	
2	GC	99	
2	GD	99	
2	GE	99	
2	GF	99	
2	HA	99	
2	HB	99	
2	HC	99	
2	HD	99	
2	HE	99	
2	HF	99	
2	IA	99	
2	IB	99	
2	IC	99	
2	ID	99	
2	IE	99	
2	IF	99	
2	JA	99	
2	JB	99	
2	JC	99	
2	JD	99	
2	JE	99	

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Mol	Chain	Length	Quality of chain
2	JF	99	 80%13%7%
2	KA	99	 81%12%7%
2	KB	99	 81%12%7%
2	KC	99	 75%18%7%
2	KD	99	 84%9%7%
2	KE	99	 81%12%7%
2	KF	99	 83%10%7%
2	LA	99	 81%12%7%
2	LB	99	 81%12%7%
2	LC	99	 83%10%7%
2	LD	99	 84%9%7%
2	LE	99	 78%15%7%
2	LF	99	 85%8%7%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 32833 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 Microcompartments protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	202	Total	C	N	O	S	0	0
			1525	971	265	286	3		
1	B	202	Total	C	N	O	S	0	0
			1525	971	265	286	3		
1	C	202	Total	C	N	O	S	0	0
			1525	971	265	286	3		
1	D	202	Total	C	N	O	S	0	0
			1525	971	265	286	3		
1	E	202	Total	C	N	O	S	0	0
			1525	971	265	286	3		
1	F	201	Total	C	N	O	S	0	0
			1520	968	264	285	3		

- Molecule 2 is a protein called Microcompartments protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	GA	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	GB	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	GC	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	GD	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	GE	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	GF	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	HA	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	HB	92	Total	C	N	O	S	0	0
			658	413	120	122	3		
2	HC	92	Total	C	N	O	S	0	0
			658	413	120	122	3		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	HD	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	HE	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	HF	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	IA	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	IB	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	IC	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	ID	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	IE	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	IF	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	JA	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	JB	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	JC	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	JD	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	JE	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	JF	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	KA	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	KB	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	KC	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	KD	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	KE	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	KF	92	Total 658	C 413	N 120	O 122	S 3	0	0

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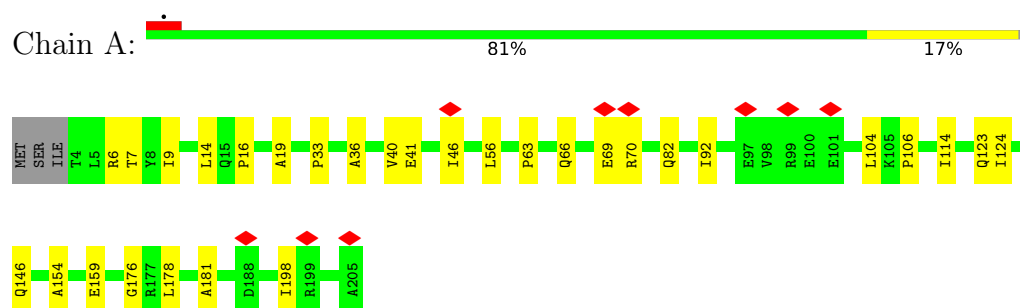
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Mol	Chain	Residues	Atoms					AltConf	Trace
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2	LB	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	LC	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	LD	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	LE	92	Total 658	C 413	N 120	O 122	S 3	0	0
2	LF	92	Total 658	C 413	N 120	O 122	S 3	0	0

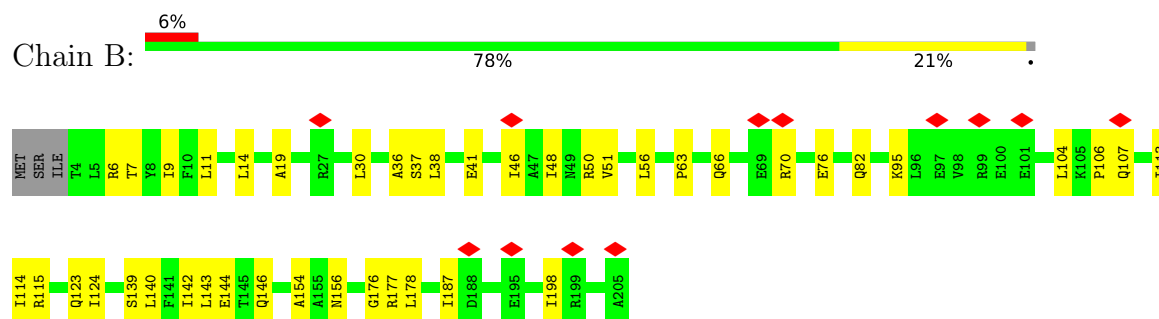
3 Residue-property plots [i](#)

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

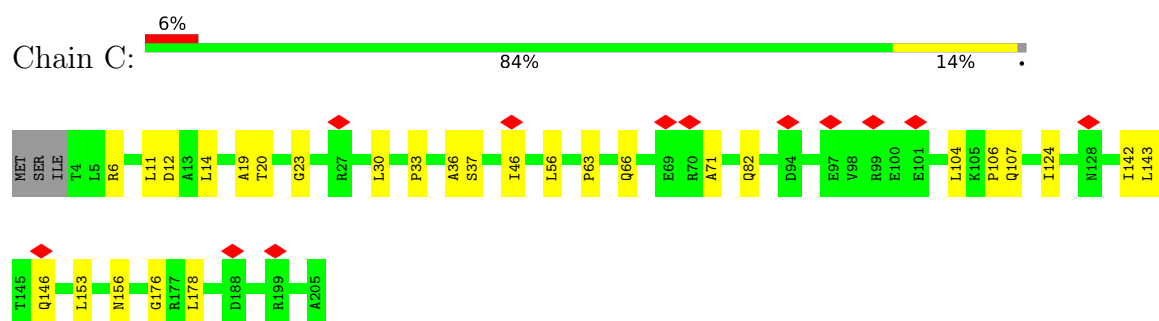
- Molecule 1: Microcompartments protein



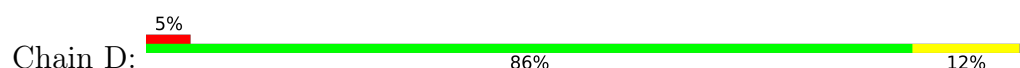
- Molecule 1: Microcompartments protein

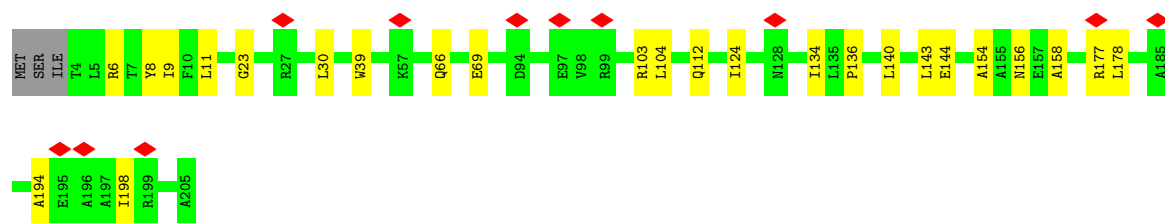


- Molecule 1: Microcompartments protein

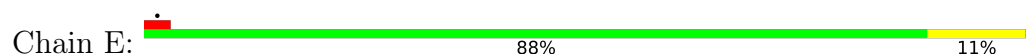


- Molecule 1: Microcompartments protein

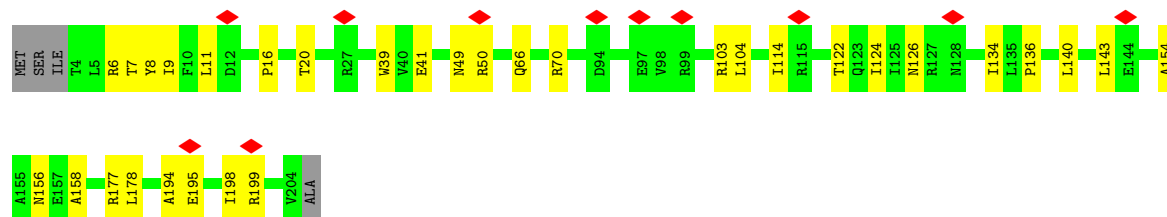
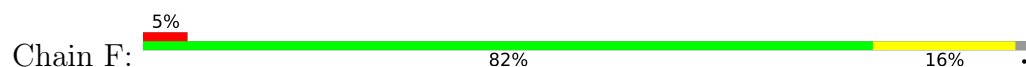




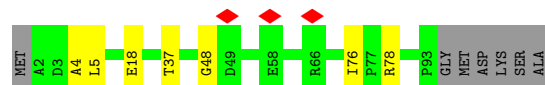
- Molecule 1: Microcompartments protein



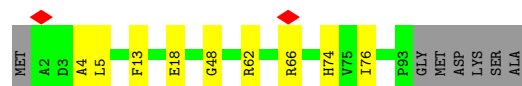
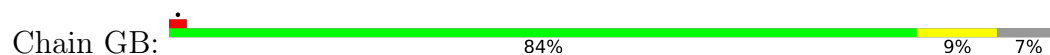
- Molecule 1: Microcompartments protein



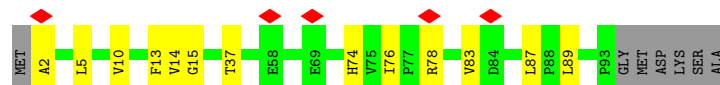
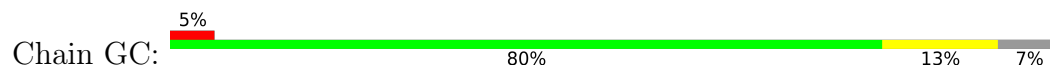
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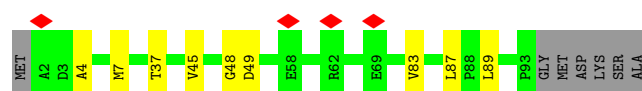
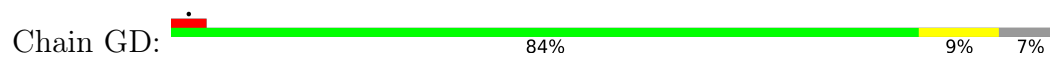
- Molecule 2: Microcompartments protein



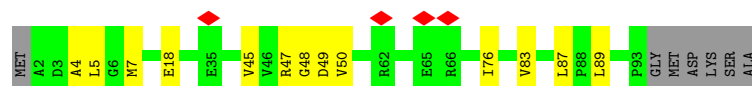
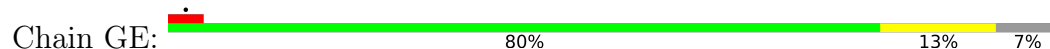
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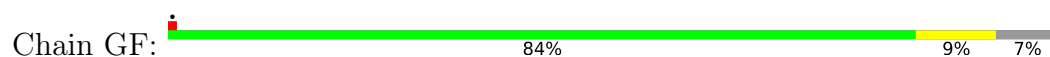
- Molecule 2: Microcompartments protein



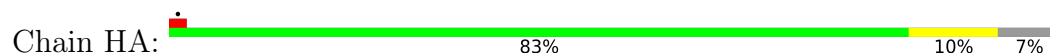
- Molecule 2: Microcompartments protein



- Molecule 2: Microcompartments protein



- Molecule 2: Microcompartments protein



- Molecule 2: Microcompartments protein



- Molecule 2: Microcompartments protein



- Molecule 2: Microcompartments protein



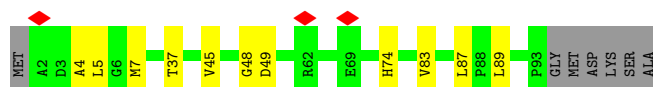
- Molecule 2: Microcompartments protein

Chain HE:  75% 17% 7%



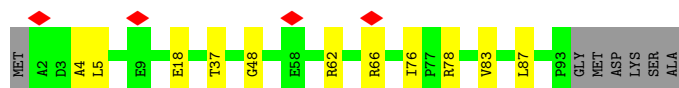
- Molecule 2: Microcompartments protein

Chain HF:  82% 11% 7%



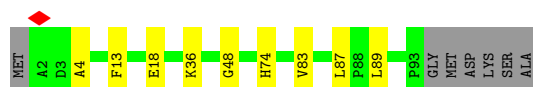
- Molecule 2: Microcompartments protein

Chain IA:  82% 11% 7%



- Molecule 2: Microcompartments protein

Chain IB:  84% 9% 7%



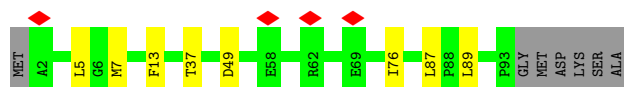
- Molecule 2: Microcompartments protein

Chain IC:  77% 16% 7%



- Molecule 2: Microcompartments protein

Chain ID:  85% 8% 7%

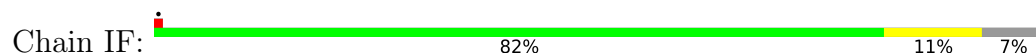


- Molecule 2: Microcompartments protein

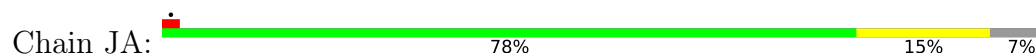
Chain IE:  73% 20% 7%



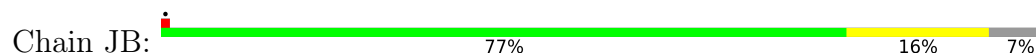
- Molecule 2: Microcompartments protein



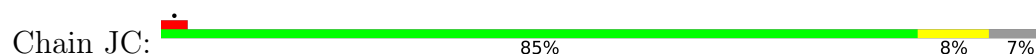
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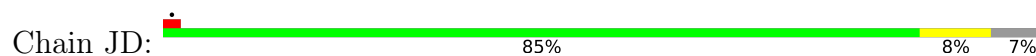
- Molecule 2: Microcompartments protein



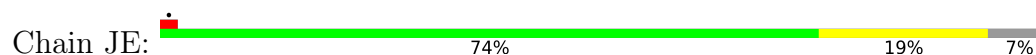
- Molecule 2: Microcompartments protein



- Molecule 2: Microcompartments protein

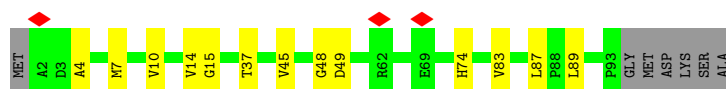


- Molecule 2: Microcompartments protein



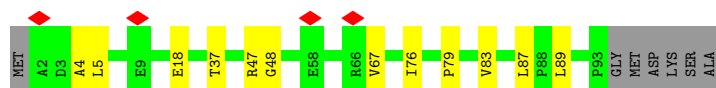
- Molecule 2: Microcompartments protein

Chain JF:  80% 13% 7%



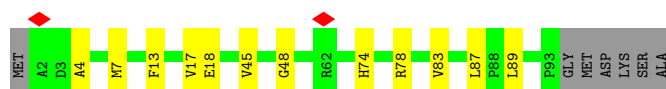
• Molecule 2: Microcompartments protein

Chain KA:  81% 12% 7%



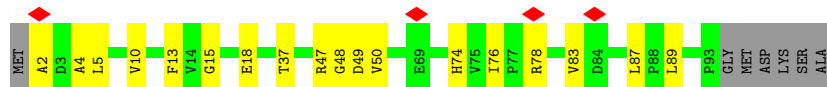
• Molecule 2: Microcompartments protein

Chain KB:  81% 12% 7%



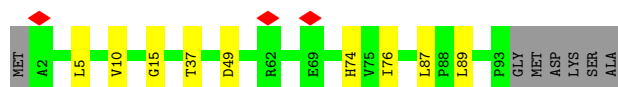
• Molecule 2: Microcompartments protein

Chain KC:  75% 18% 7%



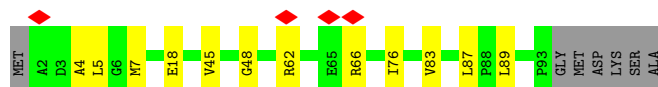
• Molecule 2: Microcompartments protein

Chain KD:  84% 9% 7%



• Molecule 2: Microcompartments protein

Chain KE:  81% 12% 7%



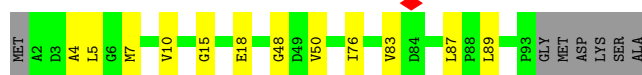
• Molecule 2: Microcompartments protein

Chain KF:  83% 10% 7%



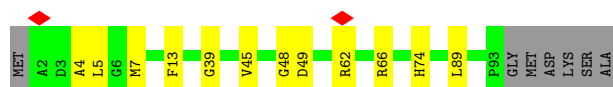
• Molecule 2: Microcompartments protein

Chain LA:  81% 12% 7%



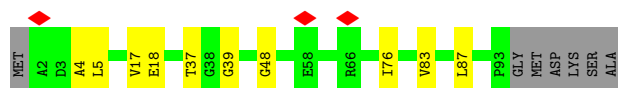
• Molecule 2: Microcompartments protein

Chain LB:  81% 12% 7%



• Molecule 2: Microcompartments protein

Chain LC:  83% 10% 7%



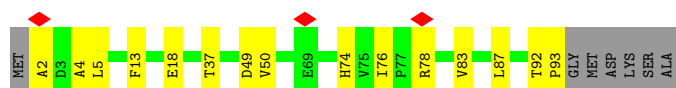
• Molecule 2: Microcompartments protein

Chain LD:  84% 9% 7%



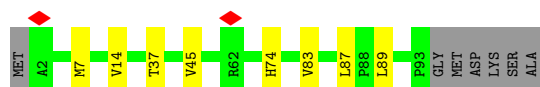
• Molecule 2: Microcompartments protein

Chain LE:  78% 15% 7%



• Molecule 2: Microcompartments protein

Chain LF:  85% 8% 7%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	83102	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; Initial CTF fitting using CTFFIND4, CTF correction applied within RELION	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	25	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	48543	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.128	Depositor
Minimum map value	-0.069	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.03	Depositor
Map size ($\text{\AA}$)	527.36, 527.36, 527.36	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.03, 1.03, 1.03	Depositor

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.28	0/1551	0.53	2/2111 (0.1%)
1	B	0.28	0/1551	0.54	0/2111
1	C	0.26	0/1551	0.50	1/2111 (0.0%)
1	D	0.24	0/1551	0.48	0/2111
1	E	0.24	0/1551	0.49	0/2111
1	F	0.24	0/1546	0.47	0/2104
2	GA	0.28	0/666	0.44	0/904
2	GB	0.27	0/666	0.43	0/904
2	GC	0.26	0/666	0.43	0/904
2	GD	0.26	0/666	0.45	0/904
2	GE	0.26	0/666	0.47	0/904
2	GF	0.26	0/666	0.43	0/904
2	HA	0.26	0/666	0.45	0/904
2	HB	0.26	0/666	0.45	0/904
2	HC	0.27	0/666	0.42	0/904
2	HD	0.27	0/666	0.46	0/904
2	HE	0.27	0/666	0.45	0/904
2	HF	0.27	0/666	0.47	0/904
2	IA	0.28	0/666	0.43	0/904
2	IB	0.27	0/666	0.45	0/904
2	IC	0.27	0/666	0.43	0/904
2	ID	0.26	0/666	0.44	0/904
2	IE	0.27	0/666	0.46	0/904
2	IF	0.27	0/666	0.49	0/904
2	JA	0.27	0/666	0.46	0/904
2	JB	0.26	0/666	0.43	0/904
2	JC	0.27	0/666	0.42	0/904
2	JD	0.28	0/666	0.45	0/904
2	JE	0.28	0/666	0.49	0/904
2	JF	0.27	0/666	0.47	0/904
2	KA	0.27	0/666	0.44	0/904
2	KB	0.28	0/666	0.45	0/904
2	KC	0.27	0/666	0.46	0/904
2	KD	0.26	0/666	0.48	0/904

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
2	KE	0.27	0/666	0.46	0/904
2	KF	0.28	0/666	0.48	0/904
2	LA	0.27	0/666	0.44	0/904
2	LB	0.26	0/666	0.45	0/904
2	LC	0.26	0/666	0.44	0/904
2	LD	0.26	0/666	0.45	0/904
2	LE	0.28	0/666	0.47	0/904
2	LF	0.27	0/666	0.47	0/904
All	All	0.27	0/33277	0.47	3/45203 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	HE	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	71	ALA	N-CA-C	-5.67	108.16	114.62
1	A	69	GLU	CA-C-N	5.40	131.85	121.54
1	A	69	GLU	C-N-CA	5.40	131.85	121.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	HE	49	ASP	Peptide

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	1525	0	1549	24	0
1	B	1525	0	1549	27	0
1	C	1525	0	1549	21	0
1	D	1525	0	1549	17	0
1	E	1525	0	1549	15	0
1	F	1520	0	1544	22	0
2	GA	658	0	678	5	0
2	GB	658	0	678	6	0
2	GC	658	0	678	9	0
2	GD	658	0	678	7	0
2	GE	658	0	678	9	0
2	GF	658	0	678	6	0
2	HA	658	0	678	6	0
2	HB	658	0	678	5	0
2	HC	658	0	678	5	0
2	HD	658	0	678	5	0
2	HE	658	0	678	12	0
2	HF	658	0	678	9	0
2	IA	658	0	678	8	0
2	IB	658	0	678	7	0
2	IC	658	0	678	13	0
2	ID	658	0	678	6	0
2	IE	658	0	678	13	0
2	IF	658	0	678	7	0
2	JA	658	0	678	12	0
2	JB	658	0	678	10	0
2	JC	658	0	678	5	0
2	JD	658	0	678	6	0
2	JE	658	0	678	14	0
2	JF	658	0	678	9	0
2	KA	658	0	678	9	0
2	KB	658	0	678	9	0
2	KC	658	0	678	13	0
2	KD	658	0	678	6	0
2	KE	658	0	678	7	0
2	KF	658	0	678	6	0
2	LA	658	0	678	9	0
2	LB	658	0	678	8	0
2	LC	658	0	678	7	0
2	LD	658	0	678	7	0
2	LE	658	0	678	11	0
2	LF	658	0	678	6	0
All	All	32833	0	33697	339	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (339) 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:C:106:PRO:HA	1:C:146:GLN:O	1.74	0.87
1:B:106:PRO:HA	1:B:146:GLN:O	1.78	0.84
2:IB:4:ALA:HB3	2:IB:48:GLY:O	1.83	0.76
2:LB:4:ALA:HB3	2:LB:48:GLY:O	1.91	0.70
1:C:11:LEU:HB2	1:C:37:SER:HB3	1.72	0.69
1:B:11:LEU:HB2	1:B:37:SER:HB3	1.75	0.68
1:A:143:LEU:HB3	1:A:178:LEU:HB3	1.76	0.67
2:JF:4:ALA:HB3	2:JF:48:GLY:O	1.95	0.66
2:HF:4:ALA:HB3	2:HF:48:GLY:O	1.95	0.66
2:KB:4:ALA:HB3	2:KB:48:GLY:O	1.97	0.64
1:D:66:GLN:HB3	1:F:124:ILE:HG21	1.78	0.64
2:LA:4:ALA:HB3	2:LA:48:GLY:O	1.99	0.62
2:GD:4:ALA:HB3	2:GD:48:GLY:O	1.98	0.62
1:A:66:GLN:HB3	1:B:124:ILE:HG21	1.82	0.61
1:B:11:LEU:HD22	1:B:156:ASN:HD22	1.67	0.60
1:B:66:GLN:HB3	1:C:124:ILE:HG21	1.82	0.60
1:C:11:LEU:HD22	1:C:156:ASN:HD22	1.66	0.60
1:D:11:LEU:HD22	1:D:156:ASN:HD22	1.68	0.59
1:B:38:LEU:O	1:B:76:GLU:HA	2.02	0.59
1:F:11:LEU:HD22	1:F:156:ASN:HD22	1.66	0.59
2:IA:5:LEU:HD23	2:IA:76:ILE:HD12	1.83	0.59
2:IF:78:ARG:HH21	2:JA:28:LYS:HD2	1.68	0.58
2:LE:83:VAL:HG13	2:LE:87:LEU:HD12	1.85	0.58
1:F:158:ALA:HB2	1:F:194:ALA:HB2	1.86	0.57
2:HA:7:MET:HG2	2:HA:45:VAL:HG22	1.86	0.57
1:E:124:ILE:HG21	1:F:66:GLN:HB3	1.86	0.57
2:JE:18:GLU:OE1	2:JF:74:HIS:NE2	2.38	0.57
1:D:69:GLU:O	1:F:177:ARG:NH2	2.38	0.56
2:HE:5:LEU:HD23	2:HE:76:ILE:HD12	1.87	0.56
2:JA:7:MET:HG2	2:JA:45:VAL:HG22	1.87	0.56
1:C:143:LEU:HB3	1:C:178:LEU:HB3	1.87	0.56
1:E:11:LEU:HD22	1:E:156:ASN:HD22	1.71	0.56
2:JB:4:ALA:HB3	2:JB:48:GLY:O	2.05	0.56
1:B:6:ARG:HA	1:B:104:LEU:HD13	1.86	0.56
2:KC:5:LEU:HD23	2:KC:76:ILE:HD12	1.87	0.56
2:KD:5:LEU:HD23	2:KD:76:ILE:HD12	1.88	0.56
2:KF:83:VAL:HG13	2:KF:87:LEU:HD12	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:9:ILE:HG22	1:F:39:TRP:HB2	1.88	0.56
2:JC:4:ALA:HB3	2:JC:48:GLY:O	2.06	0.56
2:LE:5:LEU:HD23	2:LE:76:ILE:HD12	1.88	0.56
2:KE:7:MET:HG2	2:KE:45:VAL:HG22	1.87	0.55
2:LC:4:ALA:HB3	2:LC:48:GLY:O	2.06	0.55
2:JC:83:VAL:HG13	2:JC:87:LEU:HD12	1.88	0.55
1:A:159:GLU:O	2:IA:78:ARG:NH2	2.40	0.55
2:JA:5:LEU:HD11	2:JA:89:LEU:HD12	1.87	0.55
2:JA:83:VAL:HG13	2:JA:87:LEU:HD12	1.88	0.55
1:B:50:ARG:HE	1:B:95:LYS:HD2	1.71	0.55
2:GC:5:LEU:HD23	2:GC:76:ILE:HD12	1.89	0.55
1:F:143:LEU:HB3	1:F:178:LEU:HB3	1.89	0.55
2:HE:4:ALA:HB3	2:HE:48:GLY:O	2.07	0.54
2:HF:7:MET:HG2	2:HF:45:VAL:HG22	1.87	0.54
2:KE:83:VAL:HG13	2:KE:87:LEU:HD12	1.90	0.54
1:A:142:ILE:HD12	1:C:46:ILE:HG12	1.90	0.54
2:KA:5:LEU:HD23	2:KA:76:ILE:HD12	1.88	0.54
1:A:106:PRO:HA	1:A:146:GLN:O	2.08	0.54
2:IE:7:MET:HG2	2:IE:45:VAL:HG22	1.89	0.54
2:GC:2:ALA:O	2:GC:78:ARG:NH1	2.41	0.54
2:LC:83:VAL:HG13	2:LC:87:LEU:HD12	1.89	0.54
1:B:46:ILE:HG12	1:C:142:ILE:HD12	1.90	0.54
2:KE:4:ALA:HB3	2:KE:48:GLY:O	2.08	0.54
1:A:56:LEU:HD21	1:A:63:PRO:HD3	1.91	0.53
2:GA:5:LEU:HD23	2:GA:76:ILE:HD12	1.90	0.53
2:JE:28:LYS:HB2	2:KB:78:ARG:HD2	1.90	0.53
2:KA:83:VAL:HG13	2:KA:87:LEU:HD12	1.91	0.53
2:KF:3:ASP:O	2:KF:47:ARG:NH1	2.41	0.53
2:IE:4:ALA:HB3	2:IE:48:GLY:O	2.09	0.53
2:LE:18:GLU:OE1	2:LF:74:HIS:NE2	2.40	0.53
1:A:46:ILE:HG12	1:B:142:ILE:HD12	1.89	0.53
2:HA:5:LEU:HD23	2:HA:76:ILE:HD12	1.90	0.53
2:HA:5:LEU:HD11	2:HA:89:LEU:HD12	1.91	0.53
2:JE:83:VAL:HG13	2:JE:87:LEU:HD12	1.90	0.53
2:JF:7:MET:HG2	2:JF:45:VAL:HG22	1.90	0.53
2:GB:4:ALA:HB3	2:GB:48:GLY:O	2.09	0.53
2:IB:18:GLU:OE1	2:IC:74:HIS:NE2	2.41	0.52
1:A:114:ILE:HD12	1:A:140:LEU:HD23	1.90	0.52
1:A:124:ILE:HG21	1:C:66:GLN:HB3	1.90	0.52
2:GC:83:VAL:HG13	2:GC:87:LEU:HD12	1.91	0.52
2:HC:83:VAL:HG13	2:HC:87:LEU:HD12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:HD:3:ASP:O	2:HD:47:ARG:NH1	2.43	0.52
2:HE:2:ALA:O	2:HE:78:ARG:NH1	2.42	0.52
2:LB:7:MET:HG2	2:LB:45:VAL:HG22	1.91	0.52
2:HC:18:GLU:OE1	2:HD:74:HIS:NE2	2.43	0.52
1:B:107:GLN:NE2	1:B:146:GLN:OE1	2.43	0.52
2:GB:18:GLU:OE1	2:GC:74:HIS:NE2	2.42	0.52
1:C:36:ALA:HB2	1:C:82:GLN:HG2	1.92	0.52
1:E:121:GLN:NE2	1:F:49:ASN:O	2.43	0.52
2:JF:83:VAL:HG13	2:JF:87:LEU:HD12	1.92	0.52
2:LF:7:MET:HG2	2:LF:45:VAL:HG22	1.92	0.52
1:D:112:GLN:HG3	1:E:50:ARG:HH11	1.74	0.51
2:JD:3:ASP:O	2:JD:47:ARG:NH1	2.43	0.51
1:B:139:SER:HB2	1:B:187:ILE:HG13	1.92	0.51
2:JA:5:LEU:HD23	2:JA:76:ILE:HD12	1.91	0.51
2:KE:18:GLU:OE1	2:KF:74:HIS:NE2	2.42	0.51
1:C:107:GLN:NE2	1:C:146:GLN:OE1	2.43	0.51
2:IE:62:ARG:O	2:IE:66:ARG:NH1	2.43	0.51
2:LC:18:GLU:OE1	2:LD:74:HIS:NE2	2.43	0.51
2:LD:18:GLU:OE1	2:LE:74:HIS:NE2	2.43	0.51
2:LF:87:LEU:HB3	2:LF:89:LEU:HD23	1.92	0.51
1:A:6:ARG:HA	1:A:104:LEU:HD13	1.92	0.51
2:IE:18:GLU:OE1	2:IF:74:HIS:NE2	2.44	0.51
1:A:14:LEU:HB3	1:A:19:ALA:HB2	1.93	0.51
2:LA:5:LEU:HD11	2:LA:89:LEU:HD12	1.91	0.51
2:LE:2:ALA:O	2:LE:78:ARG:NH1	2.43	0.51
1:B:114:ILE:HD12	1:B:140:LEU:HD23	1.92	0.51
2:IF:62:ARG:O	2:IF:66:ARG:NH1	2.44	0.51
2:GE:7:MET:HG2	2:GE:45:VAL:HG22	1.92	0.51
2:HC:4:ALA:HB3	2:HC:48:GLY:O	2.11	0.51
2:HE:13:PHE:HB2	2:HF:37:THR:HG21	1.92	0.51
1:B:7:THR:HB	1:B:41:GLU:HB3	1.93	0.50
2:KA:47:ARG:HH21	2:KA:89:LEU:HB3	1.76	0.50
2:LD:3:ASP:O	2:LD:47:ARG:NH1	2.44	0.50
2:HD:18:GLU:OE1	2:HE:74:HIS:NE2	2.44	0.50
2:KF:5:LEU:HD11	2:KF:89:LEU:HD12	1.94	0.50
1:A:7:THR:HB	1:A:41:GLU:HB3	1.94	0.50
1:D:124:ILE:HG21	1:E:66:GLN:HB3	1.94	0.50
2:HB:62:ARG:O	2:HB:66:ARG:NH1	2.43	0.50
2:GE:5:LEU:HD23	2:GE:76:ILE:HD12	1.93	0.50
2:JB:5:LEU:HD23	2:JB:76:ILE:HD12	1.94	0.50
2:JE:10:VAL:HG11	2:JE:15:GLY:HA3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:ILE:HD12	1:D:140:LEU:HD13	1.94	0.50
1:D:143:LEU:HB3	1:D:178:LEU:HB3	1.94	0.50
2:HE:83:VAL:HG13	2:HE:87:LEU:HD12	1.92	0.50
2:IB:13:PHE:HB2	2:IC:37:THR:HG21	1.93	0.50
2:IC:13:PHE:HB2	2:ID:37:THR:HG21	1.93	0.50
2:HE:18:GLU:OE1	2:HF:74:HIS:NE2	2.45	0.50
2:HF:49:ASP:OD1	2:HF:49:ASP:N	2.44	0.49
2:JF:49:ASP:OD1	2:JF:49:ASP:N	2.45	0.49
2:IC:83:VAL:HG13	2:IC:87:LEU:HD12	1.94	0.49
2:IC:8:ILE:O	2:IC:43:THR:HA	2.12	0.49
2:LE:13:PHE:HB2	2:LF:37:THR:HG21	1.94	0.49
1:B:36:ALA:HB2	1:B:82:GLN:HG2	1.93	0.49
2:JE:2:ALA:O	2:JE:78:ARG:NH1	2.45	0.49
2:KA:18:GLU:HG2	2:KA:67:VAL:HG21	1.95	0.49
2:KD:87:LEU:HB3	2:KD:89:LEU:HD23	1.94	0.49
2:LB:5:LEU:HD11	2:LB:89:LEU:HD12	1.95	0.49
1:B:144:GLU:HA	1:B:176:GLY:O	2.13	0.48
1:D:158:ALA:HB2	1:D:194:ALA:HB2	1.95	0.48
1:F:6:ARG:HA	1:F:104:LEU:HD13	1.95	0.48
1:C:107:GLN:HG2	1:C:146:GLN:HB3	1.95	0.48
2:GE:18:GLU:OE1	2:GF:74:HIS:NE2	2.45	0.48
2:IB:36:LYS:HE3	2:IC:37:THR:HA	1.95	0.48
2:JF:87:LEU:HB3	2:JF:89:LEU:HD23	1.95	0.48
1:F:8:TYR:O	1:F:103:ARG:NH1	2.42	0.48
2:GD:49:ASP:OD1	2:GD:49:ASP:N	2.45	0.48
1:B:56:LEU:HD21	1:B:63:PRO:HD3	1.96	0.48
2:JD:83:VAL:HG13	2:JD:87:LEU:HD12	1.95	0.48
1:D:8:TYR:O	1:D:103:ARG:NH1	2.43	0.48
2:KC:47:ARG:HH21	2:KC:89:LEU:HB3	1.78	0.48
2:LE:49:ASP:N	2:LE:49:ASP:OD1	2.46	0.48
1:A:36:ALA:HB2	1:A:82:GLN:HG2	1.96	0.47
1:E:112:GLN:HG3	1:F:50:ARG:HH11	1.79	0.47
2:KC:4:ALA:HB2	2:KC:50:VAL:HG22	1.96	0.47
2:LA:5:LEU:HD23	2:LA:76:ILE:HD12	1.96	0.47
2:HE:49:ASP:OD1	2:HE:49:ASP:N	2.47	0.47
2:LC:5:LEU:HD23	2:LC:76:ILE:HD12	1.96	0.47
2:IA:18:GLU:OE1	2:IB:74:HIS:NE2	2.48	0.47
2:IF:3:ASP:OD2	2:IF:91:ARG:NH2	2.40	0.47
2:JC:18:GLU:OE1	2:JD:74:HIS:NE2	2.46	0.47
2:KA:4:ALA:HB3	2:KA:48:GLY:O	2.14	0.47
1:A:40:VAL:HG21	1:A:92:ILE:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:114:ILE:HD12	1:E:140:LEU:HD23	1.97	0.47
2:GA:78:ARG:O	2:GF:25:LYS:NZ	2.40	0.47
1:A:123:GLN:HE21	1:D:30:LEU:HB3	1.79	0.47
2:GD:87:LEU:HB3	2:GD:89:LEU:HD23	1.95	0.47
2:GF:62:ARG:O	2:GF:66:ARG:NH1	2.48	0.47
2:IC:4:ALA:HB2	2:IC:50:VAL:HG22	1.95	0.47
2:KE:5:LEU:HD11	2:KE:89:LEU:HD12	1.95	0.47
2:HE:7:MET:HG2	2:HE:45:VAL:HG22	1.96	0.47
2:IE:28:LYS:HA	2:JB:78:ARG:HD2	1.97	0.47
1:E:69:GLU:OE2	1:F:70:ARG:NH1	2.42	0.47
2:IA:62:ARG:O	2:IA:66:ARG:NH1	2.48	0.46
2:LA:18:GLU:OE1	2:LB:74:HIS:NE2	2.48	0.46
2:JF:10:VAL:HG11	2:JF:15:GLY:HA3	1.97	0.46
1:E:154:ALA:HB2	1:E:198:ILE:HD11	1.97	0.46
2:HB:83:VAL:HG13	2:HB:87:LEU:HD12	1.97	0.46
2:IE:47:ARG:HH21	2:IE:89:LEU:HB3	1.80	0.46
2:JE:49:ASP:N	2:JE:49:ASP:OD1	2.48	0.46
2:KC:13:PHE:HB2	2:KD:37:THR:HG21	1.96	0.46
1:D:6:ARG:HA	1:D:104:LEU:HD13	1.96	0.46
1:E:144:GLU:HG2	1:E:177:ARG:HG2	1.98	0.46
2:GF:10:VAL:HG11	2:GF:15:GLY:HA3	1.98	0.46
2:HF:83:VAL:HG13	2:HF:87:LEU:HD12	1.95	0.46
2:LB:62:ARG:O	2:LB:66:ARG:NH1	2.49	0.46
1:C:144:GLU:HA	1:C:176:GLY:O	2.15	0.46
2:GB:62:ARG:O	2:GB:66:ARG:NH1	2.48	0.46
2:IA:4:ALA:HB3	2:IA:48:GLY:O	2.15	0.46
2:JD:13:PHE:HB2	2:JE:37:THR:HG21	1.98	0.46
2:JE:47:ARG:HH21	2:JE:89:LEU:HB3	1.81	0.46
2:KB:13:PHE:HB2	2:KC:37:THR:HG21	1.97	0.46
1:B:143:LEU:HB3	1:B:178:LEU:HB3	1.97	0.46
2:GB:5:LEU:HD23	2:GB:76:ILE:HD12	1.98	0.46
2:KE:62:ARG:O	2:KE:66:ARG:NH1	2.48	0.46
1:F:195:GLU:HB3	1:F:199:ARG:HH21	1.81	0.46
2:ID:87:LEU:HB3	2:ID:89:LEU:HD23	1.98	0.46
2:KA:37:THR:HG21	2:KF:13:PHE:HB2	1.97	0.46
2:LA:83:VAL:HG13	2:LA:87:LEU:HD12	1.97	0.46
2:LF:83:VAL:HG13	2:LF:87:LEU:HD12	1.96	0.46
2:GD:83:VAL:HG13	2:GD:87:LEU:HD12	1.97	0.46
2:HB:13:PHE:HB2	2:HC:37:THR:HG21	1.98	0.46
2:KA:47:ARG:HH12	2:KA:79:PRO:HG3	1.80	0.46
2:GA:4:ALA:HB3	2:GA:48:GLY:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:HC:19:ALA:HB2	2:HC:64:ALA:HB2	1.98	0.46
2:IC:5:LEU:HD23	2:IC:76:ILE:HD12	1.98	0.46
2:KF:10:VAL:HG11	2:KF:15:GLY:HA3	1.98	0.46
2:JC:19:ALA:HB2	2:JC:64:ALA:HB2	1.98	0.45
1:B:113:ILE:HG22	1:B:115:ARG:HG3	1.97	0.45
1:B:154:ALA:HB2	1:B:198:ILE:HD11	1.99	0.45
1:F:7:THR:HB	1:F:41:GLU:HB3	1.98	0.45
1:F:126:ASN:HB3	1:F:134:ILE:HG22	1.98	0.45
2:HB:3:ASP:O	2:HB:47:ARG:NH1	2.50	0.45
2:KB:7:MET:HG2	2:KB:45:VAL:HG22	1.99	0.45
2:LD:13:PHE:HB2	2:LE:37:THR:HG21	1.98	0.45
1:E:8:TYR:O	1:E:103:ARG:NH1	2.49	0.45
2:GB:13:PHE:HB2	2:GC:37:THR:HG21	1.99	0.45
2:IF:10:VAL:HG11	2:IF:15:GLY:HA3	1.99	0.45
1:C:11:LEU:HD21	1:C:153:LEU:HA	1.99	0.45
2:HA:83:VAL:HG13	2:HA:87:LEU:HD12	1.98	0.45
2:HA:4:ALA:HB3	2:HA:48:GLY:O	2.17	0.45
2:KC:83:VAL:HG13	2:KC:87:LEU:HD12	1.99	0.44
2:KB:87:LEU:HB3	2:KB:89:LEU:HD23	2.00	0.44
1:B:123:GLN:HE22	1:E:23:GLY:HA3	1.82	0.44
2:LB:13:PHE:HB2	2:LC:37:THR:HG21	1.98	0.44
2:LC:17:VAL:HG22	2:LD:87:LEU:HD13	1.99	0.44
1:A:144:GLU:HA	1:A:176:GLY:O	2.17	0.44
1:B:107:GLN:HG2	1:B:146:GLN:HB3	1.98	0.44
2:IE:4:ALA:HB2	2:IE:50:VAL:HG22	1.99	0.44
2:JD:17:VAL:HG22	2:JE:87:LEU:HD13	2.00	0.44
1:A:123:GLN:HE22	1:D:23:GLY:HA3	1.82	0.44
1:F:9:ILE:HD12	1:F:9:ILE:HA	1.91	0.44
2:GE:5:LEU:HD11	2:GE:89:LEU:HD12	1.99	0.44
2:IB:87:LEU:HB3	2:IB:89:LEU:HD23	1.99	0.44
2:JE:13:PHE:HB2	2:JF:37:THR:HG21	1.99	0.44
1:A:33:PRO:HG3	1:D:136:PRO:HG2	2.00	0.44
1:E:6:ARG:HA	1:E:104:LEU:HD13	2.00	0.44
2:JA:18:GLU:OE1	2:JB:74:HIS:NE2	2.51	0.44
2:JE:7:MET:HG2	2:JE:45:VAL:HG22	1.99	0.44
2:IA:37:THR:HG21	2:IF:13:PHE:HB2	2.00	0.44
2:IE:10:VAL:HG11	2:IE:15:GLY:HA3	2.00	0.44
2:JA:87:LEU:HB3	2:JA:89:LEU:HD23	2.00	0.44
2:LD:36:LYS:HE3	2:LE:37:THR:HA	1.99	0.44
1:A:16:PRO:HA	1:A:33:PRO:HB3	2.00	0.43
1:C:6:ARG:HA	1:C:104:LEU:HD13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:GC:13:PHE:HB2	2:GD:37:THR:HG21	2.01	0.43
2:GE:47:ARG:HH21	2:GE:89:LEU:HB3	1.83	0.43
2:HF:5:LEU:HD11	2:HF:89:LEU:HD12	2.00	0.43
2:KC:2:ALA:O	2:KC:78:ARG:NH1	2.51	0.43
2:IE:5:LEU:HD11	2:IE:89:LEU:HD12	2.00	0.43
2:JB:13:PHE:HB2	2:JC:37:THR:HG21	1.99	0.43
2:GE:4:ALA:HB3	2:GE:48:GLY:O	2.18	0.43
2:HF:87:LEU:HB3	2:HF:89:LEU:HD23	2.00	0.43
2:LE:4:ALA:HB2	2:LE:50:VAL:HG22	2.00	0.43
2:ID:49:ASP:OD1	2:ID:49:ASP:N	2.50	0.43
2:GD:7:MET:HG2	2:GD:45:VAL:HG22	1.99	0.43
2:IC:14:VAL:HG13	2:ID:7:MET:HE3	2.01	0.43
2:KB:18:GLU:OE1	2:KC:74:HIS:NE2	2.52	0.43
2:LA:4:ALA:HB2	2:LA:50:VAL:HG22	2.00	0.43
2:HD:13:PHE:HB2	2:HE:37:THR:HG21	1.99	0.43
2:JA:47:ARG:HH21	2:JA:89:LEU:HB3	1.82	0.43
2:JB:7:MET:HG2	2:JB:45:VAL:HG22	2.01	0.43
2:IB:83:VAL:HG13	2:IB:87:LEU:HD12	2.01	0.43
1:C:20:THR:HG23	1:F:126:ASN:ND2	2.34	0.43
1:D:154:ALA:HB2	1:D:198:ILE:HD11	2.00	0.43
2:GA:37:THR:HG21	2:GF:13:PHE:HB2	2.00	0.43
2:HE:10:VAL:HG11	2:HE:15:GLY:HA3	2.00	0.43
2:KC:4:ALA:HB3	2:KC:48:GLY:O	2.19	0.43
1:C:56:LEU:HD21	1:C:63:PRO:HD3	2.01	0.42
1:E:126:ASN:ND2	1:E:134:ILE:O	2.51	0.42
2:GE:83:VAL:HG13	2:GE:87:LEU:HD12	2.01	0.42
2:LA:87:LEU:HB3	2:LA:89:LEU:HD23	2.01	0.42
1:F:122:THR:O	1:F:126:ASN:ND2	2.51	0.42
2:IF:5:LEU:HD11	2:IF:89:LEU:HD12	2.01	0.42
2:JB:19:ALA:HB2	2:JB:64:ALA:HB2	2.01	0.42
1:F:16:PRO:O	1:F:20:THR:OG1	2.35	0.42
2:IA:83:VAL:HG13	2:IA:87:LEU:HD12	2.00	0.42
2:ID:5:LEU:HD23	2:ID:76:ILE:HD12	2.00	0.42
2:HA:18:GLU:OE1	2:HB:74:HIS:NE2	2.52	0.42
2:JA:10:VAL:HG11	2:JA:15:GLY:HA3	2.01	0.42
2:GE:4:ALA:HB2	2:GE:50:VAL:HG22	2.01	0.42
2:JB:45:VAL:HG11	2:JB:89:LEU:HD12	2.01	0.42
2:LA:7:MET:HE3	2:LF:14:VAL:HG13	2.01	0.42
2:GC:5:LEU:HD11	2:GC:89:LEU:HD12	2.00	0.42
2:JE:4:ALA:HB3	2:JE:48:GLY:O	2.19	0.42
1:A:139:SER:O	1:A:181:ALA:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:IA:5:LEU:HB3	2:IA:76:ILE:HB	2.02	0.42
2:IE:83:VAL:HG13	2:IE:87:LEU:HD12	2.01	0.42
2:LB:49:ASP:N	2:LB:49:ASP:OD1	2.53	0.42
1:C:12:ASP:OD1	1:C:12:ASP:N	2.53	0.42
2:IC:87:LEU:HB3	2:IC:89:LEU:HD23	2.01	0.42
2:KD:10:VAL:HG11	2:KD:15:GLY:HA3	2.02	0.42
1:A:123:GLN:HB3	1:D:30:LEU:HD13	2.02	0.41
1:B:14:LEU:HB3	1:B:19:ALA:HB2	2.01	0.41
2:KE:5:LEU:HD23	2:KE:76:ILE:HD12	2.02	0.41
1:B:9:ILE:HD12	1:B:9:ILE:HA	1.93	0.41
2:GA:18:GLU:OE1	2:GB:74:HIS:NE2	2.53	0.41
2:GC:10:VAL:HG11	2:GC:15:GLY:HA3	2.02	0.41
2:IC:49:ASP:OD1	2:IC:49:ASP:N	2.52	0.41
1:F:154:ALA:HB2	1:F:198:ILE:HD11	2.01	0.41
2:GE:49:ASP:OD1	2:GE:49:ASP:N	2.53	0.41
2:KA:18:GLU:OE1	2:KB:74:HIS:NE2	2.53	0.41
2:IE:8:ILE:O	2:IE:43:THR:HA	2.20	0.41
2:KD:49:ASP:OD1	2:KD:49:ASP:N	2.53	0.41
1:B:30:LEU:HD23	1:E:123:GLN:HB3	2.02	0.41
1:D:144:GLU:HG2	1:D:177:ARG:HG2	2.03	0.41
2:IC:5:LEU:HD11	2:IC:89:LEU:HD12	2.01	0.41
2:JE:4:ALA:HB2	2:JE:50:VAL:HG22	2.02	0.41
2:KC:18:GLU:OE1	2:KD:74:HIS:NE2	2.54	0.41
2:KC:49:ASP:OD1	2:KC:49:ASP:N	2.53	0.41
1:C:142:ILE:HA	1:C:178:LEU:O	2.21	0.41
2:IC:10:VAL:HG11	2:IC:15:GLY:HA3	2.02	0.41
2:JD:18:GLU:OE1	2:JE:74:HIS:NE2	2.53	0.41
2:KC:10:VAL:HG11	2:KC:15:GLY:HA3	2.03	0.41
1:A:9:ILE:HD12	1:A:9:ILE:HA	1.96	0.41
1:B:143:LEU:O	1:B:177:ARG:HA	2.21	0.41
1:F:114:ILE:HD12	1:F:140:LEU:HD23	2.03	0.41
2:JA:4:ALA:HB2	2:JA:50:VAL:HG22	2.03	0.41
1:C:23:GLY:HA2	1:C:30:LEU:HD22	2.02	0.41
2:HD:92:THR:HA	2:HD:93:PRO:HD3	1.94	0.41
2:HE:17:VAL:HG22	2:HF:87:LEU:HD13	2.03	0.41
2:ID:13:PHE:HB2	2:IE:37:THR:HG21	2.03	0.41
2:JA:7:MET:HE3	2:JF:14:VAL:HG13	2.03	0.41
2:JB:83:VAL:HG13	2:JB:87:LEU:HD12	2.03	0.41
2:LE:92:THR:HA	2:LE:93:PRO:HD3	1.95	0.41
1:C:14:LEU:HB3	1:C:19:ALA:HB2	2.02	0.41
2:GC:14:VAL:HG13	2:GD:7:MET:HE3	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:JA:4:ALA:HB3	2:JA:48:GLY:O	2.22	0.40
2:JB:92:THR:HA	2:JB:93:PRO:HD3	1.95	0.40
2:LA:10:VAL:HG11	2:LA:15:GLY:HA3	2.03	0.40
2:LB:39:GLY:HA2	2:LC:39:GLY:HA3	2.02	0.40
1:A:142:ILE:HA	1:A:178:LEU:O	2.20	0.40
1:A:154:ALA:HB2	1:A:198:ILE:HD11	2.04	0.40
1:C:33:PRO:HG2	1:F:136:PRO:HG2	2.03	0.40
2:LD:10:VAL:HG11	2:LD:15:GLY:HA3	2.03	0.40
2:GF:5:LEU:HD11	2:GF:89:LEU:HD12	2.04	0.40
2:IE:49:ASP:OD1	2:IE:49:ASP:N	2.54	0.40
2:KB:83:VAL:HG13	2:KB:87:LEU:HD12	2.03	0.40
1:B:48:ILE:HD12	1:B:51:VAL:HB	2.04	0.40
1:D:9:ILE:HG22	1:D:39:TRP:HB2	2.04	0.40
2:KB:17:VAL:HG22	2:KC:87:LEU:HD13	2.03	0.40
2:KA:5:LEU:HB3	2:KA:76:ILE:HB	2.04	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	200/205 (98%)	195 (98%)	5 (2%)	0	100	100
1	B	200/205 (98%)	192 (96%)	8 (4%)	0	100	100
1	C	200/205 (98%)	194 (97%)	6 (3%)	0	100	100
1	D	200/205 (98%)	191 (96%)	9 (4%)	0	100	100
1	E	200/205 (98%)	193 (96%)	7 (4%)	0	100	100
1	F	199/205 (97%)	191 (96%)	8 (4%)	0	100	100
2	GA	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	GB	90/99 (91%)	89 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	GC	90/99 (91%)	87 (97%)	3 (3%)	0	100	100
2	GD	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	GE	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	GF	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	HA	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	HB	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	HC	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	HD	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	HE	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	HF	90/99 (91%)	90 (100%)	0	0	100	100
2	IA	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	IB	90/99 (91%)	90 (100%)	0	0	100	100
2	IC	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	ID	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	IE	90/99 (91%)	90 (100%)	0	0	100	100
2	IF	90/99 (91%)	86 (96%)	4 (4%)	0	100	100
2	JA	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	JB	90/99 (91%)	87 (97%)	3 (3%)	0	100	100
2	JC	90/99 (91%)	90 (100%)	0	0	100	100
2	JD	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	JE	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	JF	90/99 (91%)	87 (97%)	3 (3%)	0	100	100
2	KA	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	KB	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	KC	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	KD	90/99 (91%)	87 (97%)	3 (3%)	0	100	100
2	KE	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	KF	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	LA	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	LB	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
2	LC	90/99 (91%)	88 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	LD	90/99 (91%)	89 (99%)	1 (1%)	0	100	100
2	LE	90/99 (91%)	86 (96%)	4 (4%)	0	100	100
2	LF	90/99 (91%)	88 (98%)	2 (2%)	0	100	100
All	All	4439/4794 (93%)	4339 (98%)	100 (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	154/157 (98%)	153 (99%)	1 (1%)	78	79
1	B	154/157 (98%)	153 (99%)	1 (1%)	78	79
1	C	154/157 (98%)	154 (100%)	0	100	100
1	D	154/157 (98%)	154 (100%)	0	100	100
1	E	154/157 (98%)	154 (100%)	0	100	100
1	F	154/157 (98%)	154 (100%)	0	100	100
2	GA	63/68 (93%)	63 (100%)	0	100	100
2	GB	63/68 (93%)	63 (100%)	0	100	100
2	GC	63/68 (93%)	63 (100%)	0	100	100
2	GD	63/68 (93%)	63 (100%)	0	100	100
2	GE	63/68 (93%)	63 (100%)	0	100	100
2	GF	63/68 (93%)	63 (100%)	0	100	100
2	HA	63/68 (93%)	63 (100%)	0	100	100
2	HB	63/68 (93%)	63 (100%)	0	100	100
2	HC	63/68 (93%)	63 (100%)	0	100	100
2	HD	63/68 (93%)	63 (100%)	0	100	100
2	HE	63/68 (93%)	63 (100%)	0	100	100
2	HF	63/68 (93%)	63 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	IA	63/68 (93%)	63 (100%)	0	100	100
2	IB	63/68 (93%)	63 (100%)	0	100	100
2	IC	63/68 (93%)	63 (100%)	0	100	100
2	ID	63/68 (93%)	63 (100%)	0	100	100
2	IE	63/68 (93%)	63 (100%)	0	100	100
2	IF	63/68 (93%)	63 (100%)	0	100	100
2	JA	63/68 (93%)	63 (100%)	0	100	100
2	JB	63/68 (93%)	63 (100%)	0	100	100
2	JC	63/68 (93%)	63 (100%)	0	100	100
2	JD	63/68 (93%)	63 (100%)	0	100	100
2	JE	63/68 (93%)	63 (100%)	0	100	100
2	JF	63/68 (93%)	63 (100%)	0	100	100
2	KA	63/68 (93%)	63 (100%)	0	100	100
2	KB	63/68 (93%)	63 (100%)	0	100	100
2	KC	63/68 (93%)	63 (100%)	0	100	100
2	KD	63/68 (93%)	63 (100%)	0	100	100
2	KE	63/68 (93%)	63 (100%)	0	100	100
2	KF	63/68 (93%)	63 (100%)	0	100	100
2	LA	63/68 (93%)	63 (100%)	0	100	100
2	LB	63/68 (93%)	63 (100%)	0	100	100
2	LC	63/68 (93%)	63 (100%)	0	100	100
2	LD	63/68 (93%)	63 (100%)	0	100	100
2	LE	63/68 (93%)	63 (100%)	0	100	100
2	LF	63/68 (93%)	63 (100%)	0	100	100
All	All	3192/3390 (94%)	3190 (100%)	2 (0%)	87	87

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ARG
1	B	70	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	HIS
1	A	123	GLN
1	A	130	GLN
1	B	79	HIS
1	B	123	GLN
1	B	130	GLN
1	B	156	ASN
1	C	79	HIS
1	C	123	GLN
1	C	130	GLN
1	C	156	ASN
1	D	156	ASN
1	E	79	HIS
1	E	156	ASN
1	F	66	GLN
1	F	112	GLN
1	F	126	ASN
1	F	130	GLN
1	F	156	ASN
2	GA	82	ASN
2	GB	82	ASN
2	GC	82	ASN
2	GD	82	ASN
2	GE	82	ASN
2	GF	82	ASN
2	HB	82	ASN
2	HC	82	ASN
2	HD	82	ASN
2	HE	82	ASN
2	HF	82	ASN
2	IA	82	ASN
2	IB	82	ASN
2	ID	82	ASN
2	IE	82	ASN
2	IF	82	ASN
2	JB	82	ASN
2	JC	82	ASN
2	JD	82	ASN
2	JE	82	ASN
2	JF	82	ASN
2	KA	82	ASN
2	KB	82	ASN
2	KC	82	ASN

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Mol	Chain	Res	Type
2	KD	82	ASN
2	KE	82	ASN
2	KF	82	ASN
2	LB	82	ASN
2	LC	82	ASN
2	LD	82	ASN
2	LE	82	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

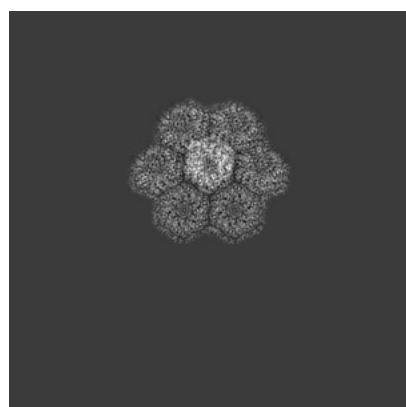
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9312. These allow visual inspection of the internal detail of the map and identification of artifacts.

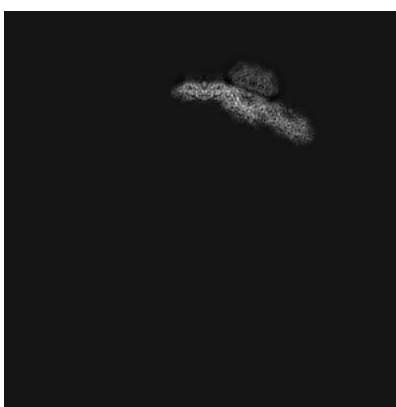
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

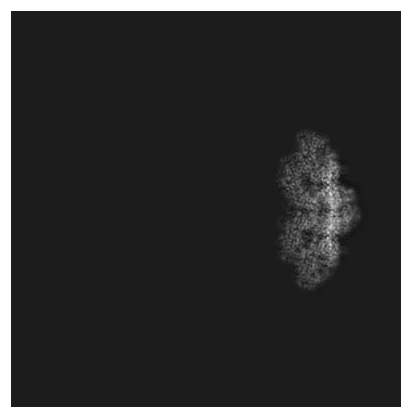
6.1.1 Primary map



X



Y

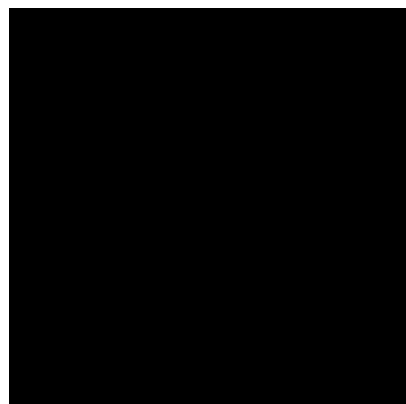


Z

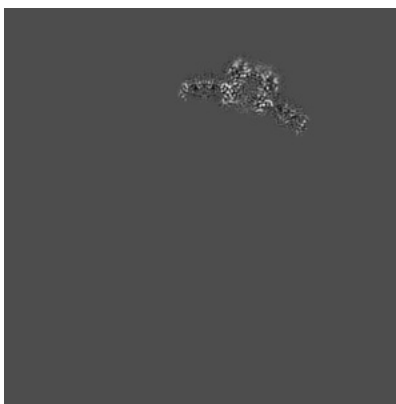
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 256



Y Index: 256

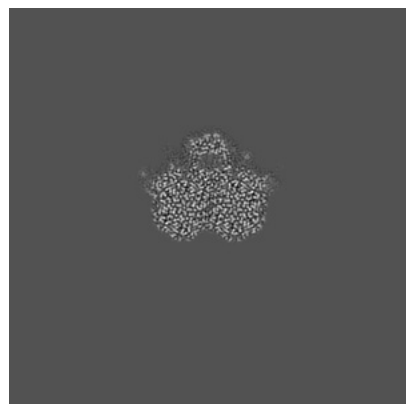


Z Index: 256

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

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 409



Y Index: 270

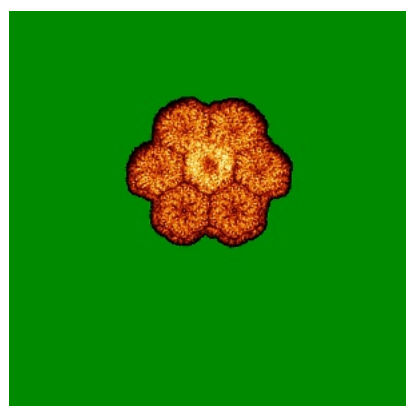


Z Index: 297

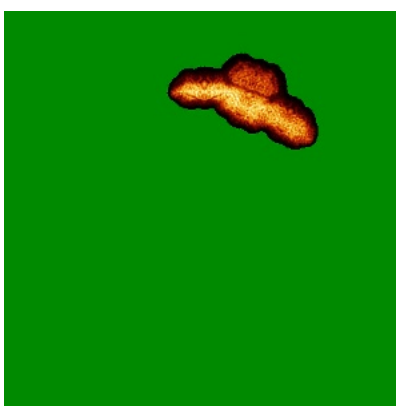
The images above show the largest variance slices of the map in three orthogonal directions.

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

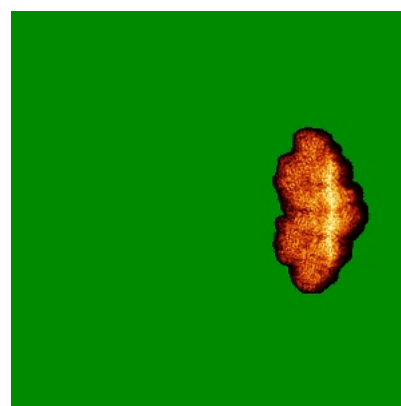
6.4.1 Primary map



X



Y

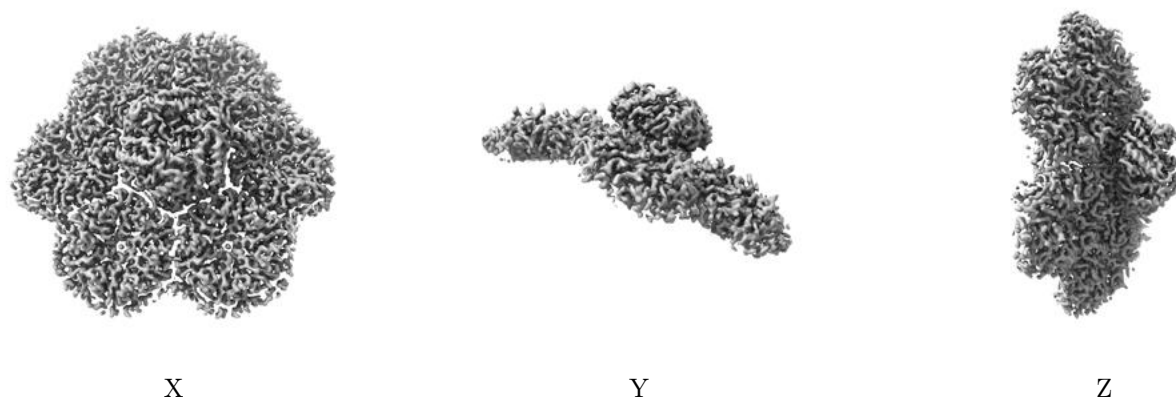


Z

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

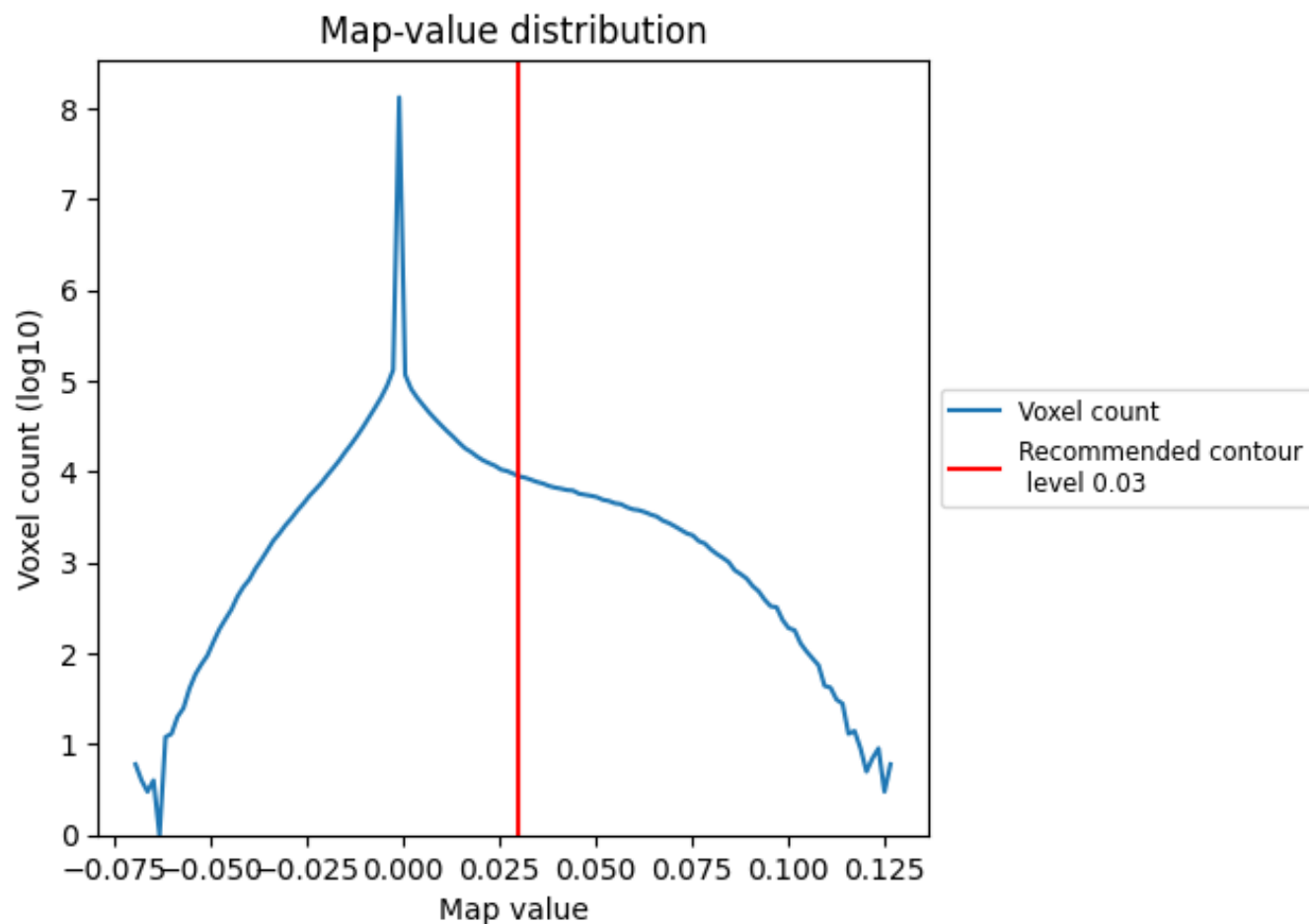
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

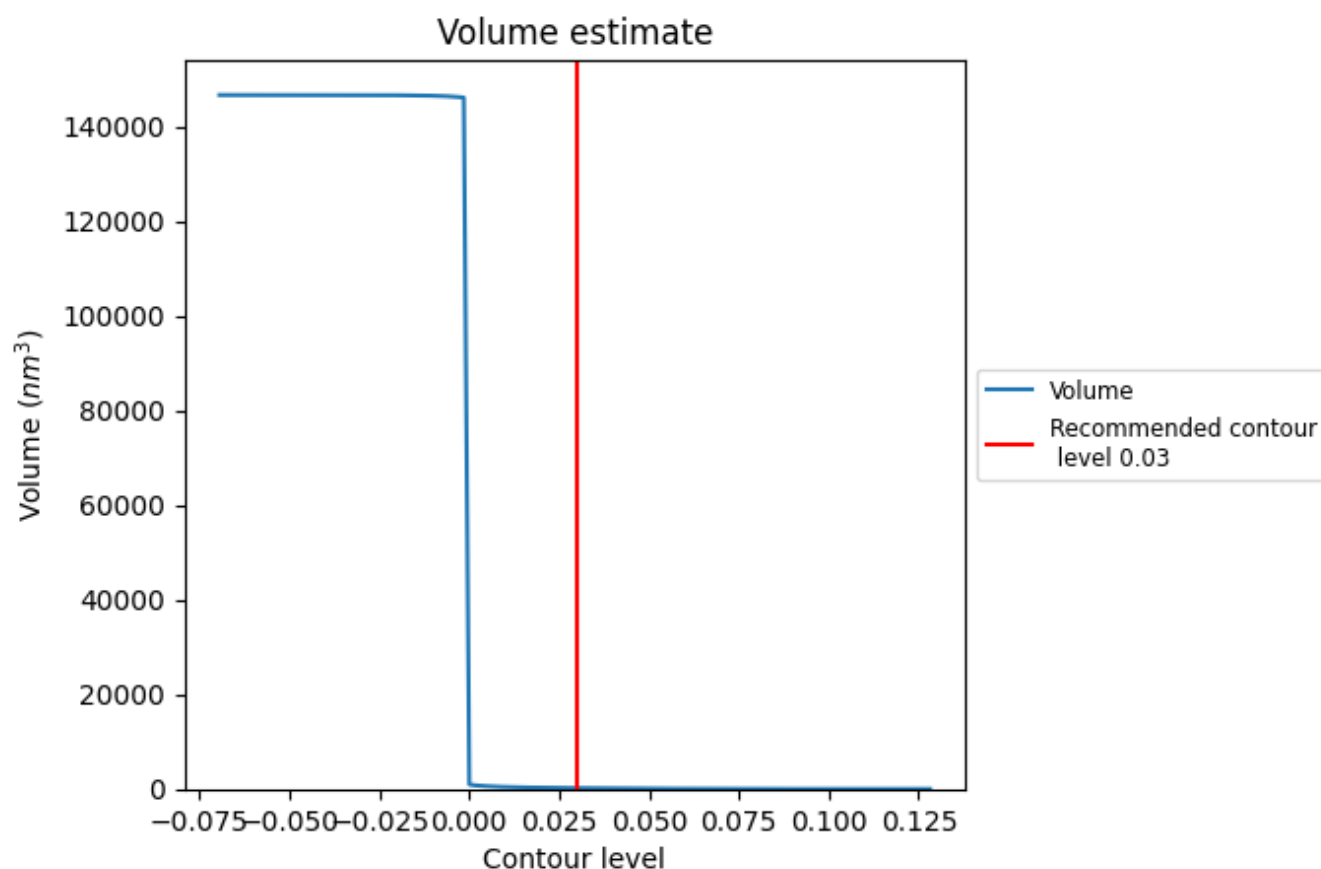
This section contains the results of statistical analysis of the map.

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.

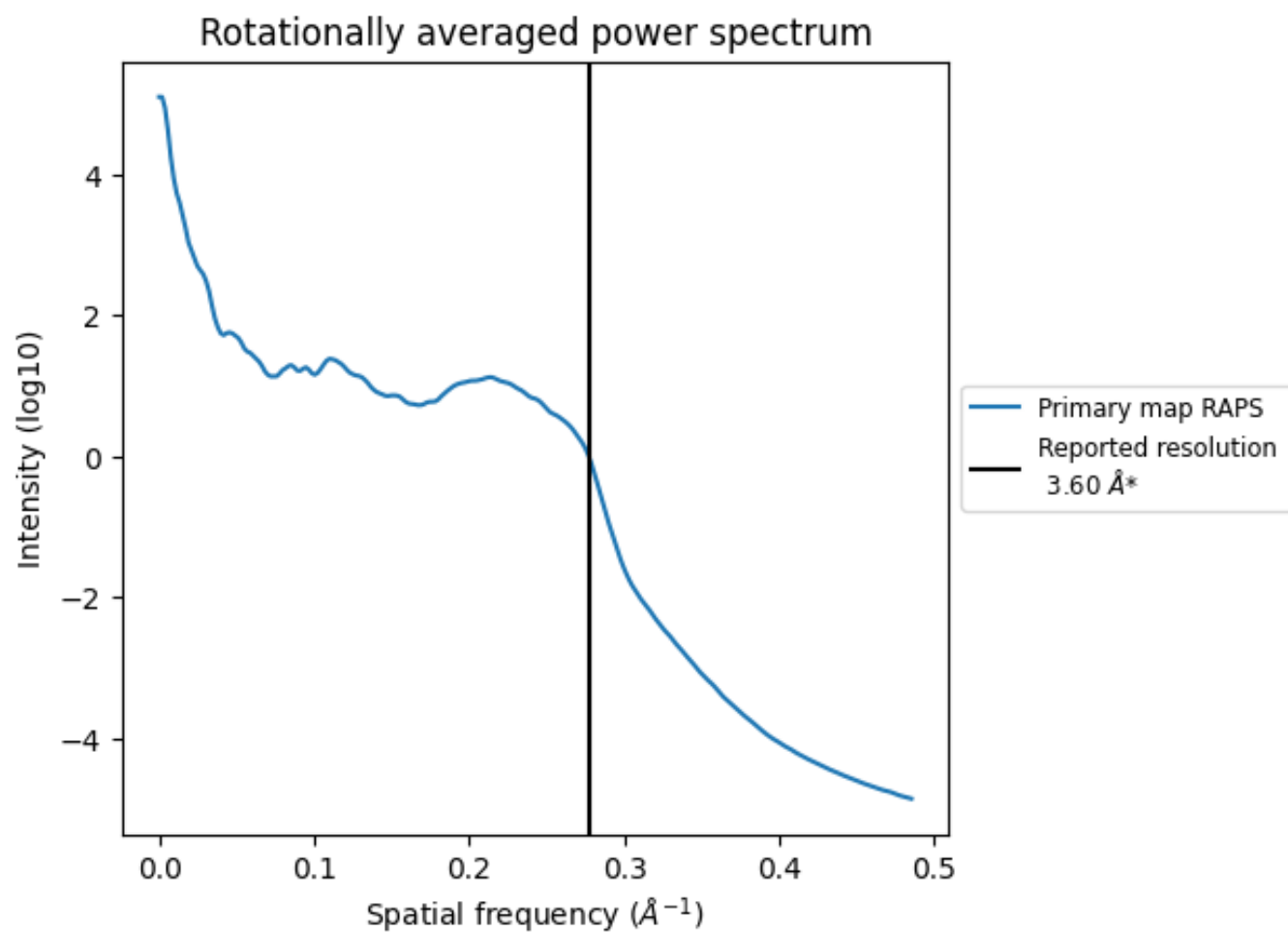
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 181 nm^3 ; this corresponds to an approximate mass of 163 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 0.278 Å⁻¹

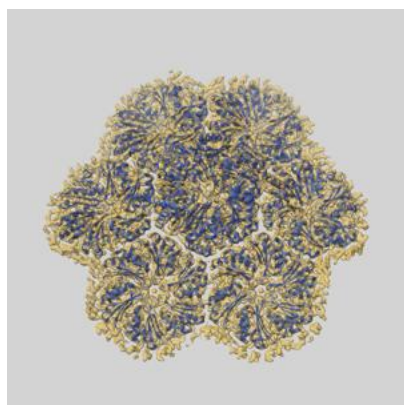
8 Fourier-Shell correlation ⓘ

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

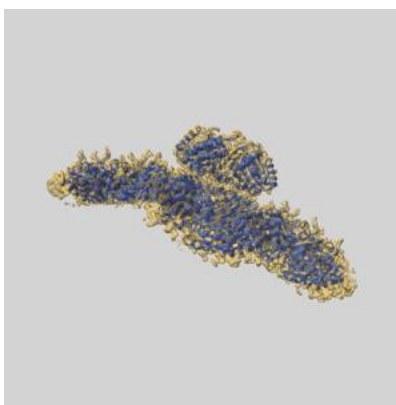
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-9312 and PDB model 6N07. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

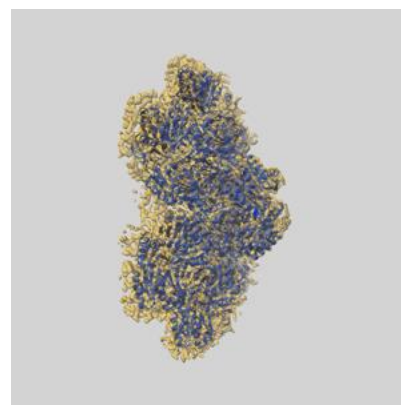
9.1 Map-model overlay [i](#)



X



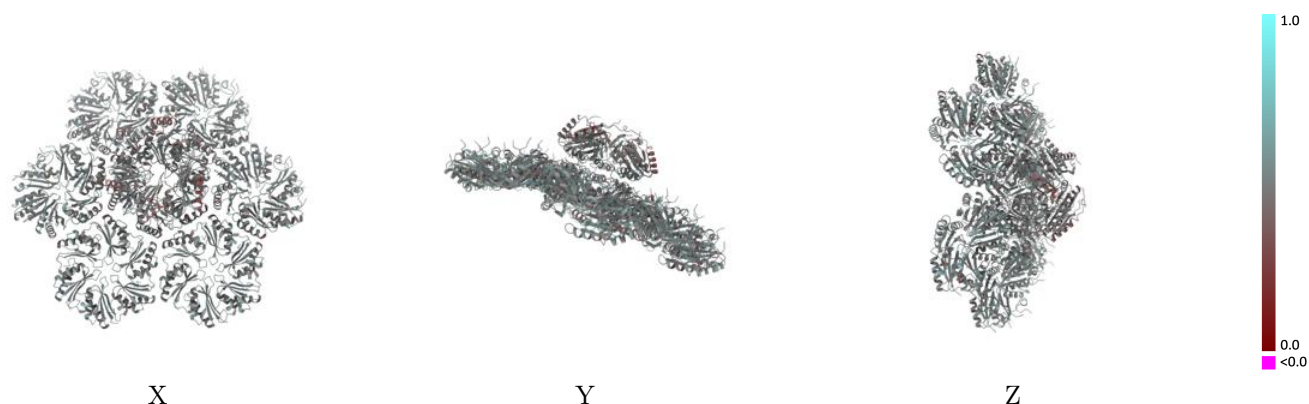
Y



Z

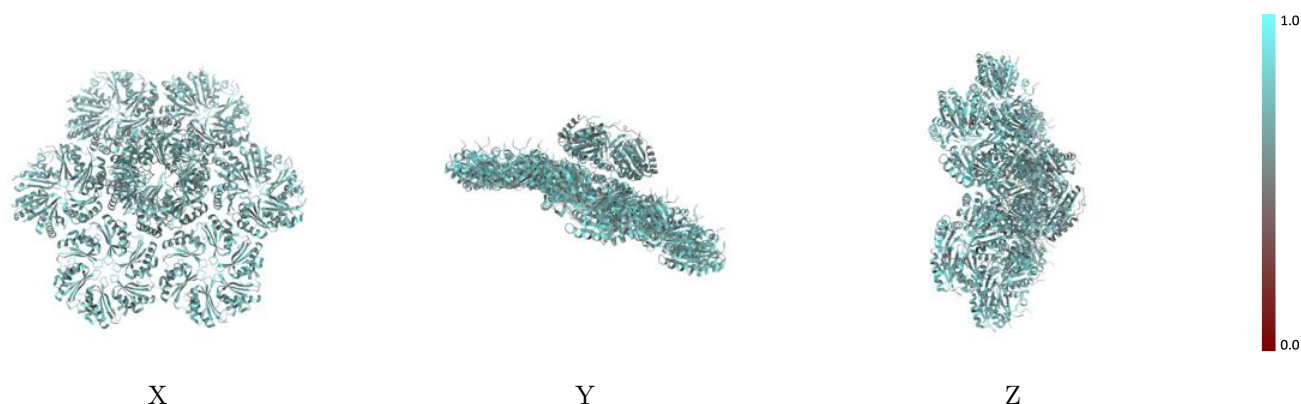
The images above show the 3D surface view of the map at the recommended contour level 0.03 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



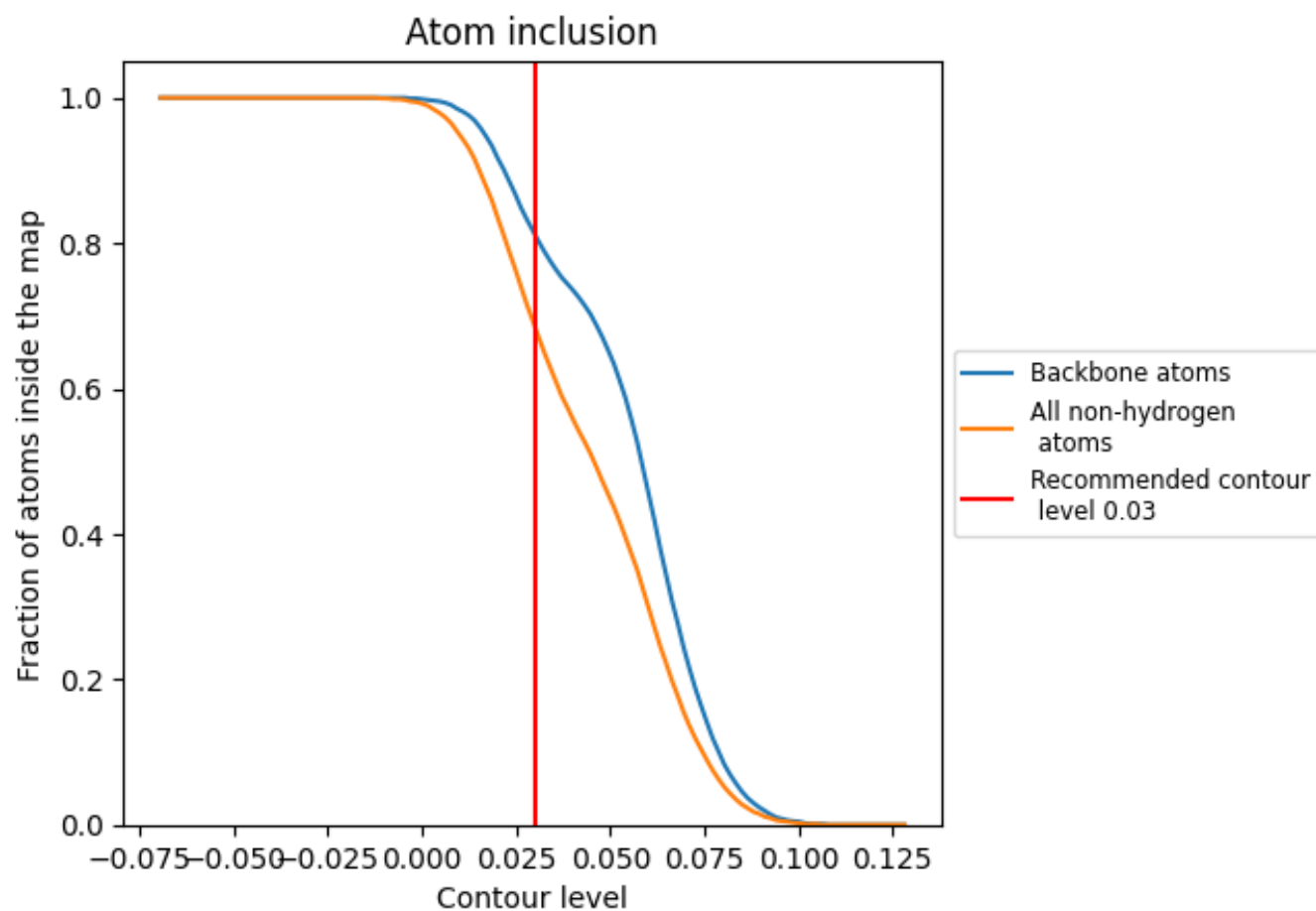
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.03).

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 68% 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 (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6850	 0.4960
A	 0.6530	 0.4870
B	 0.6550	 0.4840
C	 0.6530	 0.4840
D	 0.6550	 0.4680
E	 0.6520	 0.4670
F	 0.6520	 0.4710
GA	 0.7000	 0.5060
GB	 0.7030	 0.5040
GC	 0.6970	 0.5020
GD	 0.6970	 0.5000
GE	 0.6860	 0.5040
GF	 0.6860	 0.4980
HA	 0.6880	 0.5020
HB	 0.6880	 0.4930
HC	 0.6860	 0.5060
HD	 0.6930	 0.5060
HE	 0.7020	 0.5030
HF	 0.6940	 0.4990
IA	 0.7110	 0.5070
IB	 0.7070	 0.5090
IC	 0.6940	 0.5040
ID	 0.7080	 0.5010
IE	 0.6940	 0.5080
IF	 0.7000	 0.4990
JA	 0.6910	 0.5140
JB	 0.7020	 0.4970
JC	 0.6910	 0.5070
JD	 0.7050	 0.5080
JE	 0.6970	 0.5030
JF	 0.6940	 0.5000
KA	 0.6970	 0.5040
KB	 0.7020	 0.5040
KC	 0.6960	 0.5000
KD	 0.6970	 0.4970



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Chain	Atom inclusion	Q-score
KE	 0.7080	 0.5040
KF	 0.6990	 0.5020
LA	 0.7020	 0.5080
LB	 0.7070	 0.4970
LC	 0.6940	 0.5060
LD	 0.6970	 0.5030
LE	 0.6960	 0.5050
LF	 0.6990	 0.4960