



Full wwPDB EM Validation Report ⓘ

May 2, 2026 – 08:14 AM EDT

PDB ID : 9OSI / pdb_00009osi
EMDB ID : EMD-70807
Title : The intact LBD state of GluK2/K5 with alpha-amino-3-hydroxy-5-methyl-4-i
soxazolepropionic acid (AMPA)
Authors : Khanra, N.K.; Meyerson, J.R.
Deposited on : 2025-05-24
Resolution : 3.95 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

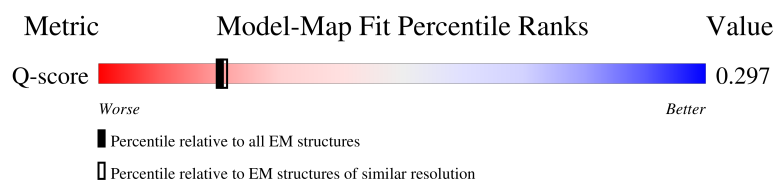
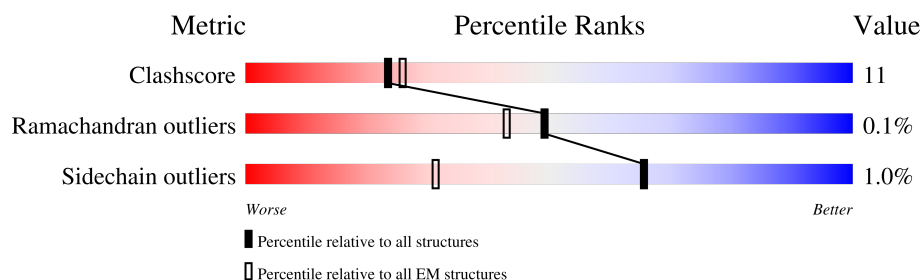
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.95 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	7707 (3.45 - 4.45)

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	1101	
1	C	1101	
2	B	942	
2	D	942	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 22601 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 Glutamate receptor ionotropic, kainate 5, Green fluorescent protein chimera.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	747	Total	C	N	O	S	0	0
			5589	3574	944	1044	27		
1	C	747	Total	C	N	O	S	0	0
			5592	3575	944	1046	27		

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	559	VAL	CYS	conflict	UNP Q63273
A	578	SER	CYS	conflict	UNP Q63273
A	619	ILE	CYS	conflict	UNP Q63273
A	813	ALA	CYS	conflict	UNP Q63273
A	828	TYR	-	linker	UNP Q63273
A	829	LYS	-	linker	UNP Q63273
A	830	SER	-	linker	UNP Q63273
A	831	ARG	-	linker	UNP Q63273
A	832	ALA	-	linker	UNP Q63273
A	833	GLU	-	linker	UNP Q63273
A	834	ALA	-	linker	UNP Q63273
A	835	LYS	-	linker	UNP Q63273
A	836	ARG	-	linker	UNP Q63273
A	837	MET	-	linker	UNP Q63273
A	838	LYS	-	linker	UNP Q63273
A	839	GLY	-	linker	UNP Q63273
A	840	LEU	-	linker	UNP Q63273
A	841	VAL	-	linker	UNP Q63273
A	842	PRO	-	linker	UNP Q63273
A	843	ARG	-	linker	UNP Q63273
A	844	GLY	-	linker	UNP Q63273
A	845	SER	-	linker	UNP Q63273
A	846	ALA	-	linker	UNP Q63273
A	847	ALA	-	linker	UNP Q63273
A	848	ALA	-	linker	UNP Q63273

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Chain	Residue	Modelled	Actual	Comment	Reference
A	849	ALA	-	linker	UNP Q63273
A	850	MET	-	linker	UNP Q63273
A	851	VAL	-	linker	UNP Q63273
A	914	LEU	PHE	conflict	UNP P42212
A	915	THR	SER	conflict	UNP P42212
A	1056	LYS	ALA	conflict	UNP P42212
A	1081	LEU	HIS	conflict	UNP P42212
A	1089	SER	-	expression tag	UNP P42212
A	1090	GLY	-	expression tag	UNP P42212
A	1091	LEU	-	expression tag	UNP P42212
A	1092	ARG	-	expression tag	UNP P42212
A	1093	THR	-	expression tag	UNP P42212
A	1094	GLU	-	expression tag	UNP P42212
A	1095	THR	-	expression tag	UNP P42212
A	1096	SER	-	expression tag	UNP P42212
A	1097	GLN	-	expression tag	UNP P42212
A	1098	VAL	-	expression tag	UNP P42212
A	1099	ALA	-	expression tag	UNP P42212
A	1100	PRO	-	expression tag	UNP P42212
A	1101	ALA	-	expression tag	UNP P42212
C	559	VAL	CYS	conflict	UNP Q63273
C	578	SER	CYS	conflict	UNP Q63273
C	619	ILE	CYS	conflict	UNP Q63273
C	813	ALA	CYS	conflict	UNP Q63273
C	828	TYR	-	linker	UNP Q63273
C	829	LYS	-	linker	UNP Q63273
C	830	SER	-	linker	UNP Q63273
C	831	ARG	-	linker	UNP Q63273
C	832	ALA	-	linker	UNP Q63273
C	833	GLU	-	linker	UNP Q63273
C	834	ALA	-	linker	UNP Q63273
C	835	LYS	-	linker	UNP Q63273
C	836	ARG	-	linker	UNP Q63273
C	837	MET	-	linker	UNP Q63273
C	838	LYS	-	linker	UNP Q63273
C	839	GLY	-	linker	UNP Q63273
C	840	LEU	-	linker	UNP Q63273
C	841	VAL	-	linker	UNP Q63273
C	842	PRO	-	linker	UNP Q63273
C	843	ARG	-	linker	UNP Q63273
C	844	GLY	-	linker	UNP Q63273
C	845	SER	-	linker	UNP Q63273

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Chain	Residue	Modelled	Actual	Comment	Reference
C	846	ALA	-	linker	UNP Q63273
C	847	ALA	-	linker	UNP Q63273
C	848	ALA	-	linker	UNP Q63273
C	849	ALA	-	linker	UNP Q63273
C	850	MET	-	linker	UNP Q63273
C	851	VAL	-	linker	UNP Q63273
C	914	LEU	PHE	conflict	UNP P42212
C	915	THR	SER	conflict	UNP P42212
C	1056	LYS	ALA	conflict	UNP P42212
C	1081	LEU	HIS	conflict	UNP P42212
C	1089	SER	-	expression tag	UNP P42212
C	1090	GLY	-	expression tag	UNP P42212
C	1091	LEU	-	expression tag	UNP P42212
C	1092	ARG	-	expression tag	UNP P42212
C	1093	THR	-	expression tag	UNP P42212
C	1094	GLU	-	expression tag	UNP P42212
C	1095	THR	-	expression tag	UNP P42212
C	1096	SER	-	expression tag	UNP P42212
C	1097	GLN	-	expression tag	UNP P42212
C	1098	VAL	-	expression tag	UNP P42212
C	1099	ALA	-	expression tag	UNP P42212
C	1100	PRO	-	expression tag	UNP P42212
C	1101	ALA	-	expression tag	UNP P42212

- Molecule 2 is a protein called Glutamate receptor ionotropic, kainate 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	755	Total	C	N	O	S	0	0
			5716	3680	946	1066	24		
2	D	755	Total	C	N	O	S	0	0
			5704	3669	945	1066	24		

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	567	VAL	ILE	conflict	UNP P42260
B	576	VAL	CYS	conflict	UNP P42260
B	595	SER	CYS	conflict	UNP P42260
B	909	SER	-	expression tag	UNP P42260
B	910	GLY	-	expression tag	UNP P42260
B	911	LEU	-	expression tag	UNP P42260
B	912	ARG	-	expression tag	UNP P42260

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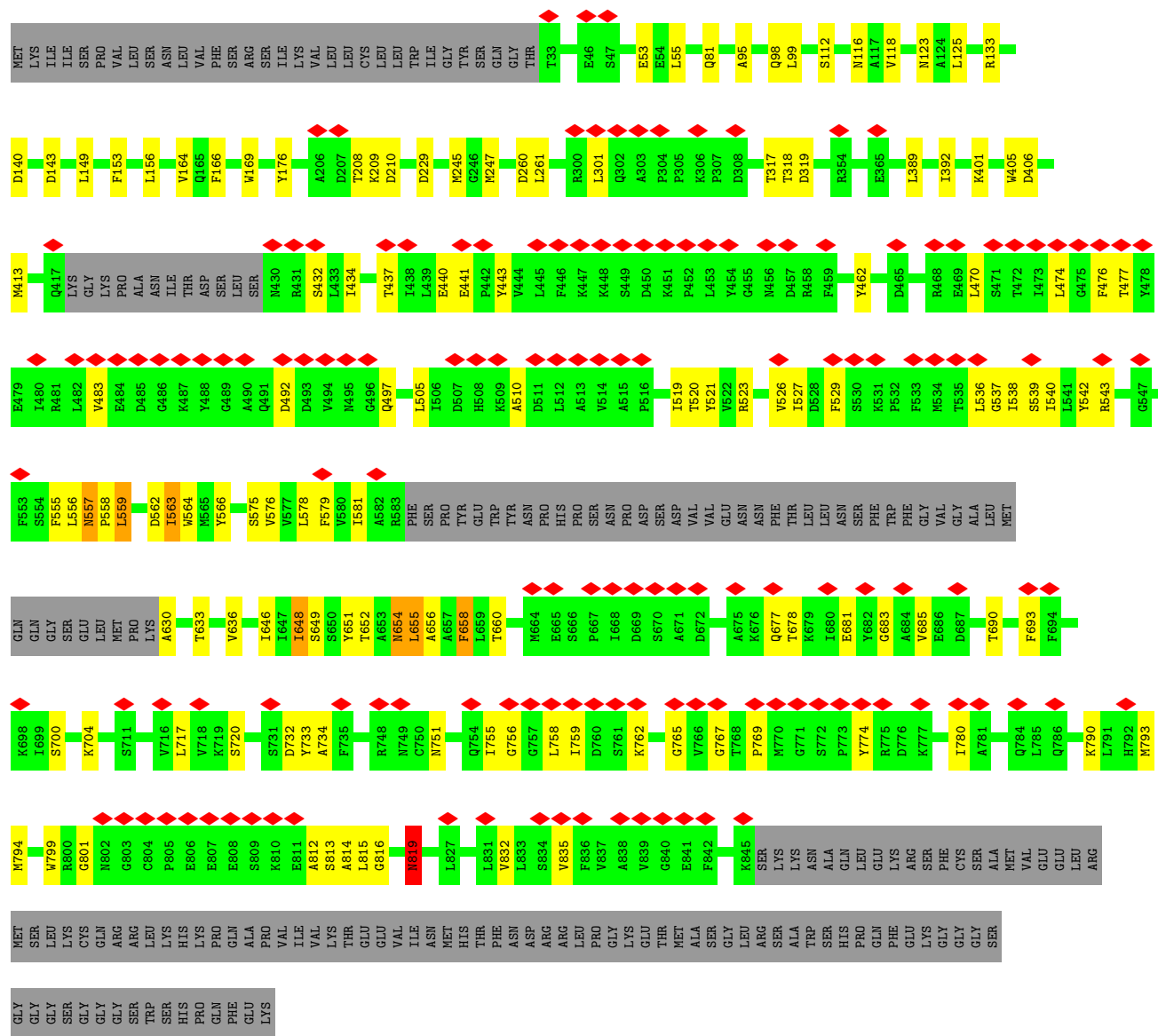
Chain	Residue	Modelled	Actual	Comment	Reference
B	913	SER	-	expression tag	UNP P42260
B	914	ALA	-	expression tag	UNP P42260
B	915	TRP	-	expression tag	UNP P42260
B	916	SER	-	expression tag	UNP P42260
B	917	HIS	-	expression tag	UNP P42260
B	918	PRO	-	expression tag	UNP P42260
B	919	GLN	-	expression tag	UNP P42260
B	920	PHE	-	expression tag	UNP P42260
B	921	GLU	-	expression tag	UNP P42260
B	922	LYS	-	expression tag	UNP P42260
B	923	GLY	-	expression tag	UNP P42260
B	924	GLY	-	expression tag	UNP P42260
B	925	GLY	-	expression tag	UNP P42260
B	926	SER	-	expression tag	UNP P42260
B	927	GLY	-	expression tag	UNP P42260
B	928	GLY	-	expression tag	UNP P42260
B	929	GLY	-	expression tag	UNP P42260
B	930	SER	-	expression tag	UNP P42260
B	931	GLY	-	expression tag	UNP P42260
B	932	GLY	-	expression tag	UNP P42260
B	933	GLY	-	expression tag	UNP P42260
B	934	SER	-	expression tag	UNP P42260
B	935	TRP	-	expression tag	UNP P42260
B	936	SER	-	expression tag	UNP P42260
B	937	HIS	-	expression tag	UNP P42260
B	938	PRO	-	expression tag	UNP P42260
B	939	GLN	-	expression tag	UNP P42260
B	940	PHE	-	expression tag	UNP P42260
B	941	GLU	-	expression tag	UNP P42260
B	942	LYS	-	expression tag	UNP P42260
D	567	VAL	ILE	conflict	UNP P42260
D	576	VAL	CYS	conflict	UNP P42260
D	595	SER	CYS	conflict	UNP P42260
D	909	SER	-	expression tag	UNP P42260
D	910	GLY	-	expression tag	UNP P42260
D	911	LEU	-	expression tag	UNP P42260
D	912	ARG	-	expression tag	UNP P42260
D	913	SER	-	expression tag	UNP P42260
D	914	ALA	-	expression tag	UNP P42260
D	915	TRP	-	expression tag	UNP P42260
D	916	SER	-	expression tag	UNP P42260
D	917	HIS	-	expression tag	UNP P42260

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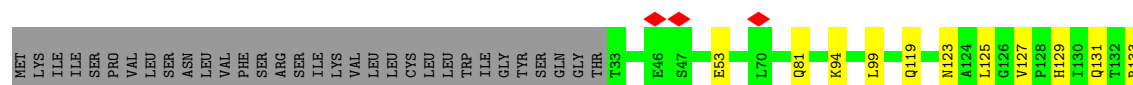
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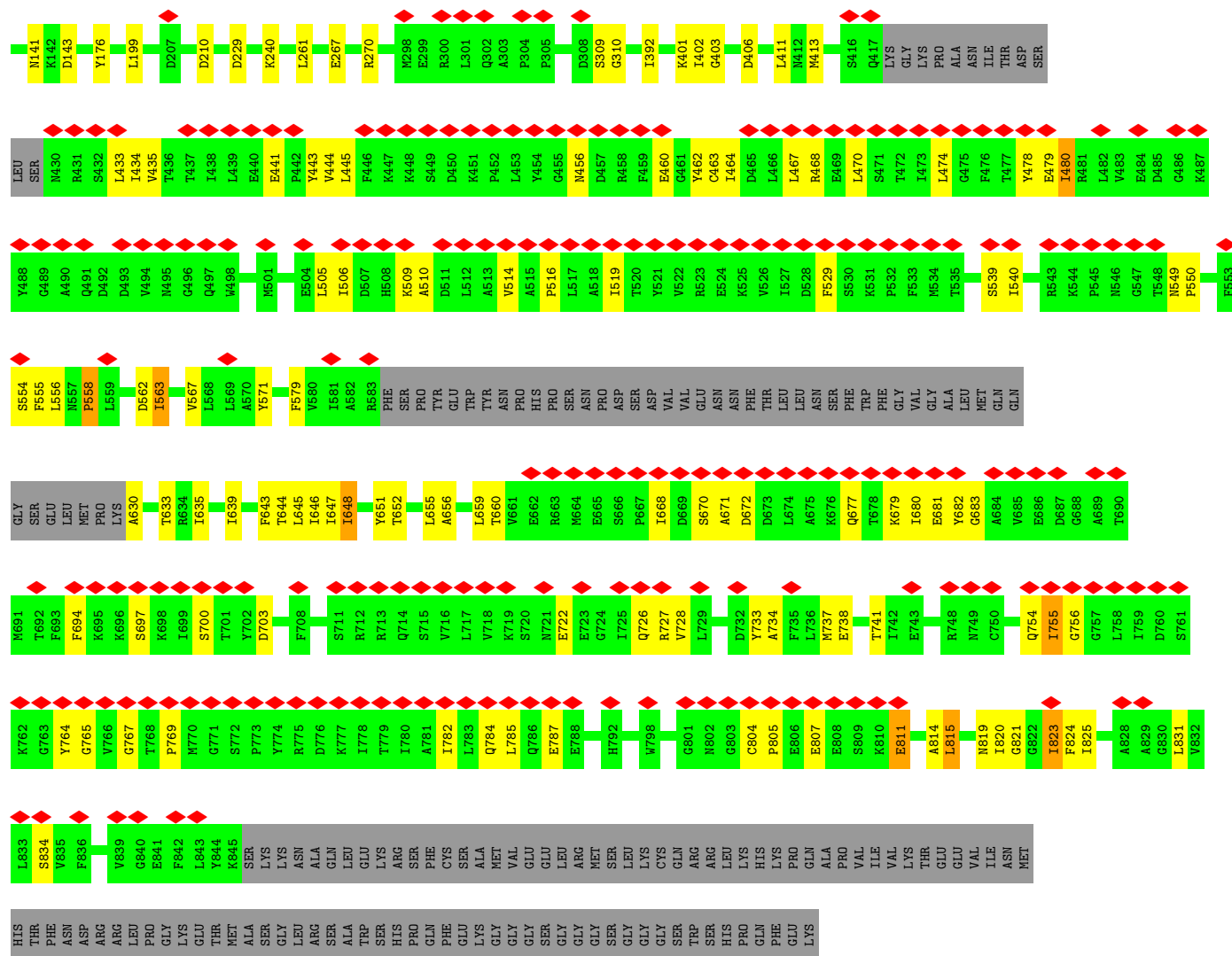
Chain	Residue	Modelled	Actual	Comment	Reference
D	918	PRO	-	expression tag	UNP P42260
D	919	GLN	-	expression tag	UNP P42260
D	920	PHE	-	expression tag	UNP P42260
D	921	GLU	-	expression tag	UNP P42260
D	922	LYS	-	expression tag	UNP P42260
D	923	GLY	-	expression tag	UNP P42260
D	924	GLY	-	expression tag	UNP P42260
D	925	GLY	-	expression tag	UNP P42260
D	926	SER	-	expression tag	UNP P42260
D	927	GLY	-	expression tag	UNP P42260
D	928	GLY	-	expression tag	UNP P42260
D	929	GLY	-	expression tag	UNP P42260
D	930	SER	-	expression tag	UNP P42260
D	931	GLY	-	expression tag	UNP P42260
D	932	GLY	-	expression tag	UNP P42260
D	933	GLY	-	expression tag	UNP P42260
D	934	SER	-	expression tag	UNP P42260
D	935	TRP	-	expression tag	UNP P42260
D	936	SER	-	expression tag	UNP P42260
D	937	HIS	-	expression tag	UNP P42260
D	938	PRO	-	expression tag	UNP P42260
D	939	GLN	-	expression tag	UNP P42260
D	940	PHE	-	expression tag	UNP P42260
D	941	GLU	-	expression tag	UNP P42260
D	942	LYS	-	expression tag	UNP P42260

- Molecule 2: Glutamate receptor ionotropic, kainate 2



- Molecule 2: Glutamate receptor ionotropic, kainate 2





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	189632	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51.28	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.718	Depositor
Minimum map value	-1.686	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.042	Depositor
Recommended contour level	0.38	Depositor
Map size (\AA)	450.52798, 450.52798, 450.52798	wwPDB
Map dimensions	416, 416, 416	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.083, 1.083, 1.083	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.52	8/5696 (0.1%)	0.71	8/7757 (0.1%)
1	C	0.46	0/5699	0.69	7/7761 (0.1%)
2	B	0.44	2/5839 (0.0%)	0.69	8/7949 (0.1%)
2	D	0.42	1/5825 (0.0%)	0.68	5/7929 (0.1%)
All	All	0.46	11/23059 (0.0%)	0.69	28/31396 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	636	THR	CA-C	-8.69	1.40	1.52
1	A	633	SER	CA-C	-8.19	1.41	1.52
1	A	634	SER	CA-C	-7.38	1.43	1.52
1	A	636	THR	C-O	-7.23	1.15	1.24
2	B	655	LEU	CA-C	-7.20	1.43	1.52
1	A	633	SER	C-O	-6.96	1.14	1.24
1	A	634	SER	C-O	-6.61	1.16	1.24
1	A	536	PHE	N-CA	-6.56	1.37	1.46
2	D	648	ILE	C-O	-6.24	1.16	1.24
1	A	532	LYS	CA-C	-5.35	1.47	1.53
2	B	654	ASN	C-O	-5.08	1.17	1.24

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	819	ASN	N-CA-C	-13.31	97.50	113.88
2	D	563	ILE	N-CA-C	9.49	120.30	110.62
2	D	811	GLU	N-CA-C	9.44	123.80	108.41
2	B	563	ILE	N-CA-C	-8.08	100.95	111.05
2	D	562	ASP	N-CA-C	-7.89	102.76	111.36
1	A	431	ARG	N-CA-C	-7.45	100.38	110.36
1	A	633	SER	N-CA-C	-7.10	104.09	112.89
2	B	655	LEU	N-CA-C	-6.65	104.11	111.36
1	A	533	PRO	N-CA-C	-6.49	99.11	112.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	802	GLU	N-CA-C	-6.37	105.16	113.12
1	A	635	TYR	N-CA-C	6.37	117.88	111.07
1	A	534	GLY	N-CA-C	-6.27	102.25	111.03
2	B	557	ASN	N-CA-C	6.00	123.07	109.81
1	C	787	ARG	CA-C-N	-5.89	115.92	123.16
1	C	787	ARG	C-N-CA	-5.89	115.92	123.16
1	A	732	LEU	N-CA-C	-5.79	106.83	114.31
1	A	431	ARG	CA-C-N	5.65	126.91	119.84
1	A	431	ARG	C-N-CA	5.65	126.91	119.84
2	D	549	ASN	N-CA-C	-5.62	102.56	109.65
1	C	702	VAL	N-CA-C	-5.37	100.33	108.17
1	C	799	LEU	N-CA-C	5.35	116.86	109.15
1	C	634	SER	CA-CB-OG	-5.31	100.48	111.10
2	D	807	GLU	N-CA-C	5.31	116.80	109.15
2	B	819	ASN	CB-CA-C	-5.19	100.62	109.03
2	B	557	ASN	CA-C-N	-5.17	113.37	119.84
2	B	557	ASN	C-N-CA	-5.17	113.37	119.84
2	B	658	PHE	N-CA-CB	5.16	117.55	110.07
1	C	701	PHE	N-CA-C	5.16	116.82	108.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5589	0	5395	129	0
1	C	5592	0	5397	158	0
2	B	5716	0	5514	132	0
2	D	5704	0	5497	124	0
All	All	22601	0	21803	490	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 11.

All (490) 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)
2:B:559:LEU:CD1	2:B:564:TRP:HE1	1.03	1.62
2:B:563:ILE:HG21	1:C:804:ILE:CD1	1.27	1.58
1:C:629:LEU:HD22	2:D:648:ILE:CD1	1.33	1.58
2:B:559:LEU:HD12	2:B:564:TRP:NE1	1.19	1.43
2:B:563:ILE:CG2	1:C:804:ILE:HD11	1.49	1.41
1:A:539:LEU:HD21	1:A:547:TRP:CE3	1.57	1.39
2:B:563:ILE:CG2	1:C:804:ILE:CD1	2.01	1.38
1:A:539:LEU:CD2	1:A:547:TRP:CZ3	2.10	1.35
1:A:539:LEU:HD21	1:A:547:TRP:CD2	1.63	1.32
1:A:539:LEU:HD21	1:A:547:TRP:CZ3	1.67	1.29
2:D:815:LEU:HD13	2:D:815:LEU:O	1.12	1.28
1:A:536:PHE:CG	1:A:538:PHE:HE2	1.53	1.27
2:B:559:LEU:CD1	2:B:564:TRP:NE1	1.73	1.27
1:A:539:LEU:CD2	1:A:547:TRP:CE3	2.16	1.26
1:C:670:HIS:N	1:C:702:VAL:O	1.68	1.25
1:C:669:ILE:HA	1:C:702:VAL:CG1	1.68	1.24
2:B:556:LEU:HD23	2:B:556:LEU:O	1.38	1.20
1:C:668:THR:O	1:C:702:VAL:HB	1.04	1.18
1:C:669:ILE:CA	1:C:702:VAL:HG12	1.74	1.17
1:C:546:VAL:HG11	2:D:820:ILE:HD11	1.17	1.16
1:C:668:THR:O	1:C:702:VAL:CB	1.93	1.16
2:B:654:ASN:ND2	1:C:799:LEU:HD13	1.61	1.16
1:A:539:LEU:HD21	1:A:547:TRP:CE2	1.80	1.16
1:A:536:PHE:CG	1:A:538:PHE:CE2	2.36	1.12
1:C:629:LEU:CD2	2:D:648:ILE:HD13	1.79	1.12
2:B:559:LEU:HD13	2:B:564:TRP:NE1	1.64	1.11
1:C:629:LEU:HD22	2:D:648:ILE:HD11	1.28	1.10
1:A:539:LEU:HD21	1:A:547:TRP:CH2	1.85	1.10
1:C:629:LEU:CD2	2:D:648:ILE:CD1	2.28	1.10
2:B:654:ASN:HD21	1:C:799:LEU:CD1	1.64	1.09
1:A:539:LEU:HD23	1:A:547:TRP:CZ3	1.88	1.08
2:D:815:LEU:HD13	2:D:815:LEU:C	1.80	1.06
2:B:816:GLY:N	2:B:819:ASN:OD1	1.87	1.05
1:A:539:LEU:CD2	1:A:547:TRP:CH2	2.38	1.04
2:B:654:ASN:HD21	1:C:799:LEU:HD13	0.93	1.04
2:D:815:LEU:O	2:D:815:LEU:CD1	2.04	1.04
1:C:629:LEU:HD22	2:D:648:ILE:HD13	1.08	1.03
2:B:563:ILE:HD11	1:C:801:MET:H	1.20	1.03
1:C:546:VAL:CG1	2:D:820:ILE:HD11	1.88	1.02
1:A:539:LEU:HD21	1:A:547:TRP:CZ2	1.95	1.01
2:B:559:LEU:HD12	2:B:564:TRP:CE2	1.95	1.01
1:A:536:PHE:CD2	1:A:538:PHE:HE2	1.77	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:563:ILE:HG21	1:C:804:ILE:HD13	1.39	1.00
1:A:536:PHE:CE2	1:A:806:GLY:HA2	1.96	0.99
2:D:815:LEU:CD1	2:D:819:ASN:HB3	1.92	0.99
1:C:702:VAL:HG11	1:C:719:PHE:HE1	1.29	0.98
2:B:563:ILE:HD13	1:C:804:ILE:CD1	1.93	0.96
2:B:563:ILE:HG21	1:C:804:ILE:HD12	1.44	0.95
2:D:815:LEU:HD11	2:D:819:ASN:HB3	1.45	0.95
2:B:563:ILE:HG21	1:C:804:ILE:HD11	1.12	0.95
2:B:563:ILE:HG23	1:C:804:ILE:HD11	1.51	0.93
2:D:555:PHE:CE1	2:D:556:LEU:HG	2.04	0.93
2:B:563:ILE:CG2	1:C:804:ILE:HD12	1.96	0.92
1:A:539:LEU:CD2	1:A:547:TRP:CD2	2.48	0.90
1:A:644:THR:O	1:A:648:MET:HB2	1.70	0.90
1:C:535:TYR:HB2	1:C:806:GLY:HA2	1.53	0.90
1:C:546:VAL:HG11	2:D:820:ILE:CD1	2.02	0.90
1:A:541:PRO:HG3	1:A:635:TYR:CE1	2.07	0.89
2:B:563:ILE:HD11	1:C:801:MET:N	1.88	0.88
1:A:538:PHE:CE2	1:A:806:GLY:HA3	2.09	0.88
1:A:538:PHE:CD2	1:A:806:GLY:HA3	2.09	0.87
1:C:804:ILE:O	1:C:807:ILE:HG12	1.74	0.87
2:B:658:PHE:HE1	1:C:647:ARG:HH12	1.18	0.86
1:A:536:PHE:CD1	1:A:538:PHE:CE2	2.65	0.85
2:D:555:PHE:CZ	2:D:556:LEU:HG	2.12	0.84
1:A:541:PRO:CG	1:A:635:TYR:CE1	2.60	0.83
2:B:576:VAL:HA	2:B:579:PHE:HD2	1.44	0.83
1:A:541:PRO:CG	1:A:635:TYR:HE1	1.92	0.83
2:B:556:LEU:O	2:B:556:LEU:CD2	2.24	0.82
1:A:536:PHE:CE2	1:A:806:GLY:CA	2.63	0.82
2:D:558:PRO:HB2	2:D:651:TYR:CE1	2.14	0.82
1:A:630:ILE:O	1:A:634:SER:HB2	1.80	0.81
1:A:721:LEU:HD13	1:A:725:MET:HG3	1.64	0.80
2:B:563:ILE:CG1	1:C:804:ILE:HD11	2.13	0.78
1:C:415:LYS:CB	1:C:459:ARG:O	2.32	0.78
2:D:402:ILE:HG13	2:D:413:MET:HE1	1.63	0.78
1:C:643:LEU:HD23	1:C:646:GLN:OE1	1.85	0.77
2:B:575:SER:O	2:B:579:PHE:CD2	2.37	0.77
1:A:536:PHE:CD2	1:A:538:PHE:CE2	2.67	0.76
1:A:631:ILE:O	1:A:634:SER:HB3	1.86	0.76
2:B:563:ILE:HD13	1:C:804:ILE:HD11	1.65	0.76
1:A:539:LEU:HD22	1:A:547:TRP:CE3	2.19	0.75
1:A:541:PRO:HG2	1:A:635:TYR:HE1	1.51	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:821:GLY:O	2:D:824:PHE:N	2.20	0.75
2:D:470:LEU:O	2:D:474:LEU:HB2	1.87	0.74
1:A:561:LEU:O	1:A:565:ALA:HB2	1.86	0.74
2:B:563:ILE:HG23	1:C:804:ILE:CD1	2.06	0.74
2:D:555:PHE:CZ	2:D:556:LEU:CD1	2.70	0.73
2:B:816:GLY:O	2:B:819:ASN:HB2	1.89	0.73
1:A:721:LEU:HD13	1:A:725:MET:CG	2.18	0.73
2:B:559:LEU:HD12	2:B:564:TRP:HE1	0.57	0.73
2:B:562:ASP:O	2:B:566:TYR:N	2.14	0.73
2:B:434:ILE:O	2:B:510:ALA:HB1	1.89	0.73
2:B:559:LEU:HD13	2:B:564:TRP:CD1	2.24	0.72
1:C:416:THR:HA	1:C:461:ARG:HB2	1.71	0.72
1:C:415:LYS:CB	1:C:460:PHE:HD1	2.03	0.71
1:C:702:VAL:HG21	1:C:717:TYR:OH	1.90	0.71
1:C:541:PRO:O	2:D:814:ALA:HA	1.91	0.71
2:B:563:ILE:CB	1:C:804:ILE:HD11	2.21	0.70
1:A:539:LEU:CD2	1:A:547:TRP:CZ2	2.70	0.70
1:A:536:PHE:CD1	1:A:538:PHE:CZ	2.80	0.70
2:D:815:LEU:C	2:D:815:LEU:CD1	2.55	0.70
1:C:669:ILE:HA	1:C:702:VAL:HG12	0.79	0.69
2:D:677:GLN:HE21	2:D:679:LYS:H	1.40	0.69
2:B:563:ILE:CD1	1:C:804:ILE:HD11	2.22	0.69
1:C:661:GLN:HE21	1:C:664:ILE:HG12	1.57	0.69
1:C:702:VAL:HG11	1:C:719:PHE:CE1	2.20	0.69
1:C:430:ARG:HA	1:C:443:PHE:HA	1.74	0.69
1:A:539:LEU:HD23	1:A:547:TRP:CH2	2.16	0.68
1:C:415:LYS:HA	1:C:460:PHE:CD1	2.28	0.68
2:B:559:LEU:CD1	2:B:564:TRP:CE2	2.65	0.68
2:D:815:LEU:HD13	2:D:819:ASN:HB3	1.76	0.68
2:B:558:PRO:HG2	2:B:655:LEU:HD23	1.75	0.68
1:A:521:ILE:H	1:A:744:ASP:HB3	1.59	0.67
2:B:529:PHE:HA	2:B:767:GLY:HA2	1.77	0.67
1:C:151:ARG:HD2	1:C:391:TRP:HE1	1.59	0.67
2:B:654:ASN:OD1	1:C:643:LEU:HD12	1.94	0.66
2:B:563:ILE:HD13	1:C:804:ILE:HD13	1.75	0.66
1:C:541:PRO:C	2:D:814:ALA:HA	2.20	0.66
2:B:649:SER:HG	1:C:636:THR:HG1	1.37	0.66
1:A:539:LEU:C	1:A:541:PRO:HD2	2.20	0.65
2:D:555:PHE:CZ	2:D:556:LEU:CG	2.80	0.65
1:C:654:SER:H	1:C:657:ASP:HB3	1.61	0.65
1:C:536:PHE:H	1:C:806:GLY:HA3	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:PHE:HZ	1:A:809:VAL:HG11	1.62	0.64
2:B:558:PRO:O	2:B:559:LEU:HG	1.97	0.64
1:A:539:LEU:HD11	1:A:547:TRP:CE2	2.31	0.64
1:A:542:PHE:CD2	1:A:547:TRP:HZ3	2.16	0.64
2:B:526:VAL:HG13	2:B:527:ILE:HG23	1.79	0.63
1:A:539:LEU:CD2	1:A:547:TRP:CE2	2.71	0.63
1:A:541:PRO:HG2	1:A:635:TYR:CE1	2.30	0.63
1:C:669:ILE:HD12	1:C:703:LYS:O	1.98	0.62
1:C:644:THR:O	1:C:648:MET:HB2	2.00	0.62
1:C:807:ILE:O	1:C:810:VAL:N	2.32	0.62
1:A:151:ARG:HD2	1:A:391:TRP:HE1	1.64	0.62
2:D:648:ILE:HG12	2:D:648:ILE:O	2.00	0.62
2:D:555:PHE:CE1	2:D:556:LEU:CG	2.81	0.62
2:D:529:PHE:HA	2:D:767:GLY:HA2	1.81	0.61
2:B:559:LEU:CG	2:B:564:TRP:HE1	2.01	0.61
2:B:813:SER:OG	2:B:814:ALA:N	2.30	0.61
2:D:411:LEU:HG	2:D:413:MET:HG2	1.82	0.61
1:C:193:LEU:HD23	1:C:202:LEU:HD21	1.83	0.61
1:C:712:VAL:HA	1:C:717:TYR:HB3	1.82	0.61
2:D:434:ILE:HB	2:D:510:ALA:HA	1.83	0.60
2:B:558:PRO:HB3	2:B:654:ASN:HB3	1.83	0.60
1:C:669:ILE:HD13	1:C:702:VAL:CG1	2.31	0.60
1:C:415:LYS:CA	1:C:460:PHE:HD1	2.13	0.60
1:C:643:LEU:CD2	1:C:646:GLN:OE1	2.49	0.60
1:C:668:THR:O	1:C:702:VAL:CG1	2.50	0.60
1:C:415:LYS:CB	1:C:460:PHE:HA	2.31	0.60
2:B:99:LEU:HD13	2:B:125:LEU:HD12	1.85	0.59
1:A:542:PHE:HD2	1:A:547:TRP:CZ3	2.20	0.59
1:C:670:HIS:H	1:C:702:VAL:C	1.93	0.59
2:D:99:LEU:HD13	2:D:125:LEU:HD12	1.85	0.58
1:A:635:TYR:O	1:A:638:ASN:N	2.32	0.58
2:B:536:LEU:N	2:B:762:LYS:O	2.33	0.58
1:A:536:PHE:HE2	1:A:806:GLY:HA2	1.60	0.58
1:C:565:ALA:O	1:C:566:ARG:NH2	2.37	0.57
2:D:558:PRO:HG2	2:D:655:LEU:HD13	1.87	0.57
2:B:557:ASN:O	2:B:559:LEU:N	2.37	0.57
1:C:779:LYS:O	1:C:783:TRP:HB2	2.03	0.57
1:A:397:LEU:C	1:A:397:LEU:HD12	2.29	0.57
1:C:801:MET:O	1:C:801:MET:HG2	2.04	0.57
1:C:629:LEU:CD2	2:D:648:ILE:HD11	2.19	0.57
2:B:576:VAL:HA	2:B:579:PHE:CD2	2.32	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:415:LYS:HA	1:C:460:PHE:HD1	1.70	0.56
2:D:682:TYR:HA	2:D:734:ALA:HB3	1.87	0.56
2:D:119:GLN:HE22	2:D:141:ASN:HD21	1.53	0.56
2:D:555:PHE:CD1	2:D:556:LEU:HG	2.39	0.56
2:D:722:GLU:HG3	2:D:726:GLN:HE21	1.70	0.56
1:C:543:SER:N	2:D:814:ALA:HB1	2.20	0.56
2:D:648:ILE:O	2:D:648:ILE:CG1	2.54	0.56
2:B:685:VAL:HG23	2:B:720:SER:HA	1.88	0.56
1:A:151:ARG:NH1	1:A:395:ARG:O	2.39	0.55
2:D:53:GLU:OE2	2:D:133:ARG:NH2	2.38	0.55
2:D:683:GLY:N	2:D:734:ALA:O	2.38	0.55
1:C:534:GLY:O	1:C:537:SER:OG	2.22	0.55
2:D:456:ASN:ND2	2:D:478:TYR:O	2.39	0.55
2:D:558:PRO:HB2	2:D:651:TYR:CD1	2.41	0.55
1:A:630:ILE:O	1:A:634:SER:CB	2.53	0.55
2:D:668:ILE:HG22	2:D:755:ILE:HG12	1.87	0.55
1:A:633:SER:HB2	2:B:652:THR:OG1	2.05	0.55
2:B:578:LEU:HD21	2:B:636:VAL:HG11	1.88	0.55
1:C:681:SER:O	1:C:682:ARG:NH2	2.39	0.55
1:A:487:GLU:HA	1:A:490:ASN:HB2	1.87	0.54
2:B:558:PRO:CB	2:B:654:ASN:HB3	2.37	0.54
1:A:536:PHE:CB	1:A:538:PHE:CE2	2.88	0.54
2:B:437:THR:O	2:B:483:VAL:N	2.38	0.54
2:B:575:SER:O	2:B:579:PHE:HD2	1.90	0.54
1:C:151:ARG:NH1	1:C:395:ARG:O	2.41	0.54
1:A:807:ILE:HD11	2:D:643:PHE:HA	1.90	0.54
2:D:540:ILE:HB	2:D:755:ILE:HG23	1.87	0.54
1:A:536:PHE:CD2	1:A:806:GLY:CA	2.91	0.54
1:A:138:LEU:HB3	1:A:306:LEU:HD22	1.88	0.54
1:C:670:HIS:CA	1:C:702:VAL:O	2.53	0.54
1:A:536:PHE:CD2	1:A:806:GLY:HA3	2.43	0.54
1:A:656:ASP:OD1	1:A:656:ASP:N	2.40	0.54
1:C:629:LEU:HD13	2:D:648:ILE:HD12	1.89	0.54
1:A:676:THR:O	1:A:680:ASN:ND2	2.41	0.54
1:C:702:VAL:CG2	1:C:717:TYR:OH	2.56	0.54
2:B:732:ASP:OD1	2:B:732:ASP:N	2.40	0.54
1:A:633:SER:HB2	2:B:652:THR:HG1	1.73	0.53
1:C:774:ARG:HA	1:C:777:ILE:HD12	1.90	0.53
1:A:684:GLN:HA	1:A:687:GLN:HB2	1.89	0.53
1:C:726:ASN:ND2	1:C:738:GLN:OE1	2.41	0.53
2:D:728:VAL:HG23	2:D:733:TYR:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:53:GLU:OE2	2:B:133:ARG:NH2	2.41	0.53
2:B:559:LEU:HD12	2:B:564:TRP:CZ2	2.41	0.53
1:A:535:TYR:CE1	1:A:805:GLY:HA3	2.44	0.53
1:A:652:VAL:HA	1:A:739:ILE:HD11	1.91	0.53
1:C:669:ILE:CD1	1:C:703:LYS:O	2.57	0.53
2:D:119:GLN:NE2	2:D:143:ASP:OD2	2.42	0.53
2:D:433:LEU:O	2:D:479:GLU:HB3	2.09	0.53
2:B:790:LYS:O	2:B:794:MET:CB	2.57	0.53
2:D:670:SER:OG	2:D:671:ALA:N	2.41	0.53
1:A:419:VAL:HB	1:A:464:LEU:HD23	1.90	0.52
2:D:672:ASP:OD1	2:D:672:ASP:N	2.41	0.52
2:D:435:VAL:HB	2:D:480:ILE:HA	1.91	0.52
1:A:488:LEU:HD11	1:A:753:PRO:HD3	1.92	0.52
1:C:449:ASP:OD2	1:C:782:TRP:NE1	2.36	0.52
1:C:496:ALA:HB3	1:C:751:GLY:HA3	1.91	0.52
1:C:31:ASP:OD2	1:C:99:SER:OG	2.28	0.52
2:B:562:ASP:C	2:B:564:TRP:N	2.62	0.52
2:D:668:ILE:HD11	2:D:680:ILE:HD13	1.92	0.52
2:B:432:SER:O	2:B:432:SER:OG	2.22	0.52
2:D:659:LEU:HD11	2:D:815:LEU:HD23	1.90	0.52
2:D:694:PHE:O	2:D:697:SER:OG	2.28	0.52
2:B:505:LEU:HD22	2:B:769:PRO:HD3	1.90	0.52
1:C:669:ILE:HD13	1:C:702:VAL:HG13	1.92	0.52
2:B:520:THR:HG23	2:B:523:ARG:H	1.75	0.52
1:A:633:SER:CB	2:B:652:THR:OG1	2.57	0.51
1:A:666:TYR:HA	1:A:718:ALA:HB3	1.92	0.51
2:D:467:LEU:HB3	2:D:514:VAL:HG11	1.91	0.51
1:A:539:LEU:HD11	1:A:547:TRP:NE1	2.25	0.51
2:B:389:LEU:HB2	2:B:405:TRP:HB3	1.92	0.51
1:C:502:ILE:HD12	1:C:515:PRO:HG3	1.92	0.51
2:D:815:LEU:HD11	2:D:819:ASN:CB	2.31	0.51
1:A:514:LYS:H	1:A:763:THR:HG23	1.76	0.51
1:C:165:CYS:HB2	1:C:192:MET:HG2	1.93	0.51
1:C:803:ASN:C	1:C:805:GLY:H	2.18	0.51
1:C:803:ASN:C	1:C:805:GLY:N	2.69	0.51
2:B:559:LEU:O	1:C:799:LEU:O	2.28	0.51
1:C:670:HIS:CB	1:C:702:VAL:O	2.59	0.51
2:B:520:THR:OG1	2:B:521:TYR:N	2.44	0.51
2:B:816:GLY:O	2:B:819:ASN:CB	2.59	0.51
2:D:267:GLU:OE2	2:D:270:ARG:NH2	2.44	0.51
2:D:392:ILE:HA	2:D:401:LYS:HA	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:LEU:O	1:C:187:THR:OG1	2.30	0.50
1:C:669:ILE:CA	1:C:702:VAL:CG1	2.55	0.50
1:C:643:LEU:HD23	1:C:643:LEU:N	2.26	0.50
1:C:681:SER:O	1:C:687:GLN:NE2	2.45	0.50
2:D:700:SER:HA	2:D:703:ASP:HB2	1.94	0.50
1:C:259:GLU:O	1:C:382:ARG:NH1	2.39	0.50
1:C:631:ILE:O	1:C:634:SER:OG	2.30	0.50
1:A:536:PHE:CD1	1:A:538:PHE:HE2	2.04	0.50
2:B:210:ASP:N	2:B:210:ASP:OD1	2.43	0.50
2:B:492:ASP:N	2:B:497:GLN:O	2.42	0.50
2:D:819:ASN:C	2:D:821:GLY:N	2.68	0.50
1:A:469:ASP:N	1:A:469:ASP:OD1	2.36	0.50
1:C:733:ASN:HB3	1:C:736:LEU:HD23	1.92	0.50
2:B:649:SER:OG	1:C:636:THR:OG1	2.10	0.50
2:B:681:GLU:HG2	2:B:733:TYR:HA	1.92	0.49
1:A:48:ARG:HD2	1:A:64:GLU:HA	1.94	0.49
1:A:205:GLU:O	1:A:209:ASP:HB2	2.12	0.49
1:A:195:ASP:OD1	1:A:195:ASP:N	2.45	0.49
2:D:434:ILE:HG13	2:D:509:LYS:O	2.12	0.49
2:D:464:ILE:HD13	2:D:467:LEU:HD21	1.95	0.49
1:C:523:ILE:HB	1:C:739:ILE:HB	1.92	0.49
1:C:543:SER:O	1:C:543:SER:OG	2.30	0.49
2:B:406:ASP:OD1	2:B:406:ASP:N	2.42	0.49
1:C:420:THR:OG1	1:C:495:LEU:O	2.30	0.49
1:C:203:LEU:HB3	1:C:235:MET:HE1	1.95	0.49
2:B:123:ASN:ND2	2:B:143:ASP:OD1	2.46	0.49
2:D:406:ASP:OD1	2:D:406:ASP:N	2.46	0.49
2:D:754:GLN:NE2	2:D:756:GLY:O	2.36	0.49
1:A:203:LEU:HB3	1:A:235:MET:HE1	1.94	0.48
2:D:819:ASN:C	2:D:821:GLY:H	2.20	0.48
2:B:539:SER:OG	2:B:756:GLY:O	2.31	0.48
1:C:205:GLU:O	1:C:209:ASP:HB2	2.13	0.48
2:D:402:ILE:CG1	2:D:413:MET:HE1	2.37	0.48
2:D:811:GLU:O	2:D:811:GLU:HG3	2.13	0.48
1:A:496:ALA:HB3	1:A:751:GLY:HA3	1.95	0.48
2:D:505:LEU:HD22	2:D:769:PRO:HD3	1.94	0.48
1:C:691:ASN:O	1:C:695:SER:OG	2.30	0.48
1:A:474:ALA:N	1:A:482:THR:OG1	2.47	0.48
1:A:654:SER:OG	1:A:656:ASP:OD1	2.32	0.48
2:D:831:LEU:O	2:D:834:SER:OG	2.29	0.48
2:B:519:ILE:HD11	2:B:765:GLY:HA3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:651:TYR:CE2	2:B:655:LEU:HD11	2.49	0.48
1:C:92:SER:OG	1:C:116:PRO:O	2.29	0.48
2:B:540:ILE:HB	2:B:755:ILE:HB	1.95	0.48
1:C:494:ASP:OD1	1:C:494:ASP:N	2.45	0.48
1:C:618:ARG:NH1	2:D:571:TYR:OH	2.46	0.48
1:A:769:LEU:HA	1:A:772:ASN:HD22	1.78	0.47
2:B:559:LEU:CD1	2:B:564:TRP:CD1	2.76	0.47
1:C:474:ALA:H	1:C:482:THR:H	1.60	0.47
1:C:523:ILE:HG22	1:C:739:ILE:HD13	1.95	0.47
2:D:554:SER:O	2:D:555:PHE:C	2.56	0.47
1:A:128:LEU:HD23	1:A:130:TYR:H	1.79	0.47
1:A:522:SER:HB2	1:A:726:ASN:HD22	1.80	0.47
2:B:646:ILE:HG21	1:C:807:ILE:HD13	1.96	0.47
1:A:27:ALA:HB2	1:A:90:VAL:HG11	1.97	0.47
1:C:422:ILE:HG21	1:C:472:TYR:HE1	1.80	0.47
1:C:557:VAL:HA	1:C:560:VAL:HG22	1.96	0.47
2:D:443:TYR:HA	2:D:462:TYR:HB3	1.97	0.47
2:D:644:THR:O	2:D:648:ILE:HG22	2.15	0.47
2:B:537:GLY:HA3	2:B:758:LEU:HB3	1.96	0.47
2:B:543:ARG:HE	2:B:751:ASN:HD21	1.63	0.47
2:B:656:ALA:O	2:B:660:THR:OG1	2.26	0.47
1:A:542:PHE:CD2	1:A:547:TRP:CZ3	2.98	0.47
1:C:52:ASN:HD21	1:C:63:VAL:H	1.63	0.47
2:B:557:ASN:HB2	2:B:558:PRO:HD3	1.97	0.46
2:D:555:PHE:CZ	2:D:556:LEU:HD12	2.48	0.46
1:A:381:SER:OG	1:A:383:GLN:O	2.32	0.46
2:B:700:SER:O	2:B:704:LYS:N	2.37	0.46
1:C:524:LEU:HB2	1:C:719:PHE:HB3	1.96	0.46
1:C:653:GLU:HB2	1:C:739:ILE:HG23	1.97	0.46
1:C:799:LEU:N	1:C:799:LEU:HD12	2.29	0.46
2:D:555:PHE:CZ	2:D:556:LEU:HD11	2.50	0.46
1:C:26:MET:HE1	1:C:311:VAL:HG23	1.97	0.46
2:D:821:GLY:O	2:D:824:PHE:HB2	2.15	0.46
1:A:182:LEU:O	1:A:187:THR:OG1	2.32	0.46
1:A:674:THR:OG1	1:A:675:MET:N	2.48	0.46
2:B:562:ASP:O	2:B:566:TYR:HB2	2.15	0.46
1:C:769:LEU:HA	1:C:772:ASN:HD22	1.80	0.46
1:A:520:GLY:O	1:A:723:SER:N	2.48	0.46
1:A:526:ARG:HA	1:A:736:LEU:HD13	1.97	0.46
1:A:621:SER:OG	1:A:622:GLY:N	2.47	0.46
2:B:153:PHE:HA	2:B:156:LEU:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:681:GLU:HB2	2:D:733:TYR:HA	1.97	0.46
2:B:166:PHE:CE2	2:B:413:MET:HE3	2.50	0.46
1:C:806:GLY:O	1:C:809:VAL:HB	2.16	0.46
1:A:670:HIS:N	1:A:702:VAL:O	2.43	0.46
2:D:456:ASN:HA	2:D:480:ILE:HD13	1.98	0.46
1:A:419:VAL:HG13	1:A:495:LEU:HB3	1.96	0.46
1:A:424:GLU:O	1:A:428:VAL:N	2.44	0.46
1:A:531:ARG:HH12	2:B:812:ALA:HB2	1.81	0.46
1:A:659:ALA:O	1:A:662:THR:OG1	2.34	0.46
2:B:477:THR:O	2:B:477:THR:OG1	2.32	0.46
2:D:630:ALA:O	2:D:633:THR:OG1	2.33	0.46
2:D:738:GLU:O	2:D:741:THR:OG1	2.27	0.46
2:B:434:ILE:HB	2:B:510:ALA:HA	1.98	0.45
2:B:780:ILE:HG23	1:C:504:ALA:HB1	1.98	0.45
2:D:645:LEU:HD13	2:D:645:LEU:HA	1.79	0.45
1:A:26:MET:HE1	1:A:311:VAL:HG23	1.98	0.45
1:A:467:VAL:HA	1:A:484:MET:HE3	1.98	0.45
1:A:724:THR:O	1:A:728:TYR:N	2.48	0.45
2:B:208:THR:OG1	2:B:209:LYS:N	2.49	0.45
1:C:183:ILE:HB	2:D:199:LEU:HB3	1.97	0.45
1:A:665:GLU:HB3	1:A:717:TYR:HA	1.99	0.45
1:A:630:ILE:O	1:A:634:SER:N	2.48	0.45
1:C:480:SER:OG	1:C:481:TRP:N	2.50	0.45
2:D:441:GLU:HG3	2:D:445:LEU:HD21	1.99	0.45
1:C:153:LEU:HD21	1:C:213:THR:HG21	1.99	0.45
1:A:631:ILE:HD13	1:A:631:ILE:HA	1.68	0.45
1:C:153:LEU:HD11	1:C:215:ILE:HD11	1.98	0.45
1:A:803:ASN:OD1	1:A:803:ASN:N	2.48	0.45
2:B:651:TYR:CZ	2:B:655:LEU:HD11	2.52	0.45
1:C:539:LEU:HB2	1:C:547:TRP:HE1	1.81	0.45
2:D:99:LEU:HD21	2:D:127:VAL:HG21	1.99	0.45
2:D:464:ILE:HG22	2:D:468:ARG:HH11	1.82	0.45
2:D:647:ILE:HD13	2:D:647:ILE:HA	1.74	0.45
2:B:799:TRP:O	2:B:801:GLY:N	2.50	0.45
2:D:443:TYR:O	2:D:463:CYS:N	2.48	0.45
1:A:469:ASP:OD2	1:A:483:GLY:N	2.50	0.44
2:B:164:VAL:HG13	2:B:169:TRP:HB2	1.99	0.44
1:A:543:SER:O	1:A:543:SER:OG	2.35	0.44
2:D:123:ASN:ND2	2:D:143:ASP:OD1	2.50	0.44
2:B:392:ILE:HA	2:B:401:LYS:HA	1.99	0.44
1:C:24:LEU:HD11	1:C:318:VAL:HG11	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:419:VAL:HA	1:C:495:LEU:HB3	1.99	0.44
1:A:628:THR:O	1:A:632:ILE:HG12	2.18	0.44
1:C:198:ASP:OD1	1:C:200:THR:OG1	2.35	0.44
2:D:681:GLU:OE1	2:D:727:ARG:NH2	2.51	0.44
2:D:555:PHE:CE2	2:D:556:LEU:HG	2.51	0.44
2:B:440:GLU:HG2	2:B:443:TYR:HD2	1.84	0.43
1:C:415:LYS:CB	1:C:460:PHE:CD1	2.94	0.43
2:B:140:ASP:OD1	2:B:140:ASP:N	2.43	0.43
1:C:670:HIS:CB	1:C:702:VAL:H	2.31	0.43
2:D:210:ASP:OD1	2:D:210:ASP:N	2.51	0.43
2:B:95:ALA:HB2	2:B:118:VAL:HG23	2.00	0.43
2:B:319:ASP:N	2:B:319:ASP:OD1	2.48	0.43
2:D:445:LEU:N	2:D:460:GLU:O	2.42	0.43
2:B:413:MET:HE2	2:B:413:MET:HB2	1.88	0.43
1:C:426:PRO:HG2	1:C:783:TRP:HD1	1.83	0.43
1:A:454:LEU:HD12	1:A:457:LEU:HB2	2.01	0.43
1:A:517:MET:SD	1:A:517:MET:N	2.92	0.43
1:A:536:PHE:CD1	1:A:538:PHE:HZ	2.35	0.43
1:A:638:ASN:HD21	2:B:815:LEU:HD13	1.83	0.43
1:C:526:ARG:HH21	1:C:713:LEU:HD13	1.84	0.43
1:A:645:VAL:HA	1:A:648:MET:HE2	2.01	0.43
1:A:814:GLY:HA3	2:D:639:ILE:HG13	2.01	0.43
2:B:81:GLN:HG2	2:B:98:GLN:HE21	1.84	0.43
1:C:246:THR:HG22	1:C:248:ASP:H	1.83	0.43
2:D:444:VAL:HG22	2:D:464:ILE:HG12	2.01	0.43
1:A:48:ARG:HG3	1:A:63:VAL:HG13	2.01	0.42
2:D:129:HIS:CD2	2:D:131:GLN:HE21	2.36	0.42
2:B:542:TYR:HB3	2:B:755:ILE:HD13	2.01	0.42
2:D:176:TYR:HA	2:D:229:ASP:HB3	2.01	0.42
2:D:635:ILE:HD13	2:D:635:ILE:HA	1.85	0.42
1:A:700:VAL:O	1:A:717:TYR:OH	2.34	0.42
1:C:99:SER:HB3	1:C:102:SER:HB3	2.01	0.42
1:C:138:LEU:HB3	1:C:306:LEU:HD22	2.01	0.42
2:D:506:ILE:HD13	2:D:506:ILE:HA	1.93	0.42
1:A:153:LEU:HD11	1:A:215:ILE:HD11	2.01	0.42
1:A:671:ALA:H	1:A:675:MET:HG3	1.84	0.42
2:B:317:THR:OG1	2:B:319:ASP:OD1	2.35	0.42
1:C:194:ASP:OD1	1:C:194:ASP:N	2.49	0.42
2:D:403:GLY:HA3	2:D:413:MET:SD	2.58	0.42
2:D:456:ASN:HD21	2:D:479:GLU:HA	1.82	0.42
2:D:558:PRO:CG	2:D:655:LEU:HD13	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:737:MET:SD	2:D:741:THR:OG1	2.77	0.42
2:D:782:ILE:HA	2:D:785:LEU:HD12	2.01	0.42
2:B:443:TYR:HA	2:B:462:TYR:HB3	2.01	0.42
2:B:683:GLY:N	2:B:734:ALA:O	2.52	0.42
1:C:521:ILE:HB	1:C:743:LEU:HB2	2.01	0.42
1:A:487:GLU:HG3	1:A:492:LYS:HB2	2.01	0.42
1:A:732:LEU:HD12	1:A:732:LEU:HA	1.91	0.42
2:B:578:LEU:HD23	2:B:581:ILE:HD12	2.01	0.42
2:B:654:ASN:ND2	1:C:799:LEU:CD1	2.44	0.42
1:C:519:LEU:HD12	1:C:746:LYS:HB2	2.01	0.42
1:C:687:GLN:HA	1:C:690:TRP:HB3	2.01	0.42
2:B:470:LEU:HG	2:B:476:PHE:HE2	1.85	0.42
2:B:815:LEU:HD12	2:B:815:LEU:HA	1.84	0.42
1:C:36:CYS:HB3	1:C:292:CYS:HB3	1.43	0.42
2:D:519:ILE:HD11	2:D:765:GLY:HA3	2.02	0.42
2:D:563:ILE:O	2:D:567:VAL:HG23	2.19	0.42
1:A:356:ASP:OD1	1:A:356:ASP:N	2.51	0.42
2:B:245:MET:HB3	2:B:247:MET:HG3	2.02	0.42
1:A:536:PHE:CE2	1:A:809:VAL:HG21	2.54	0.42
1:A:775:LEU:O	1:A:779:LYS:N	2.52	0.42
2:D:656:ALA:O	2:D:660:THR:OG1	2.26	0.42
2:D:784:GLN:HA	2:D:787:GLU:HB2	2.01	0.42
2:D:555:PHE:CG	2:D:556:LEU:N	2.87	0.41
2:B:55:LEU:HD13	2:B:301:LEU:HB3	2.01	0.41
2:B:112:SER:O	2:B:116:ASN:ND2	2.49	0.41
2:B:176:TYR:HA	2:B:229:ASP:HB3	2.02	0.41
2:B:557:ASN:CB	2:B:558:PRO:HD3	2.49	0.41
1:C:629:LEU:HD23	1:C:629:LEU:HA	1.81	0.41
1:C:683:TYR:HB2	1:C:686:TYR:CG	2.55	0.41
1:C:804:ILE:O	1:C:807:ILE:CG1	2.56	0.41
2:D:261:LEU:HB3	2:D:392:ILE:HD11	2.01	0.41
2:B:474:LEU:HD23	2:B:474:LEU:HA	1.91	0.41
1:A:681:SER:HB3	1:A:686:TYR:HB2	2.02	0.41
2:B:563:ILE:HD13	1:C:804:ILE:CG1	2.47	0.41
1:C:633:SER:OG	2:D:652:THR:OG1	2.14	0.41
2:D:646:ILE:HD13	2:D:646:ILE:HA	1.86	0.41
2:D:655:LEU:HA	2:D:655:LEU:HD12	1.75	0.41
2:D:804:CYS:HA	2:D:805:PRO:HD3	1.91	0.41
2:B:630:ALA:O	2:B:633:THR:OG1	2.32	0.41
2:B:793:MET:HE2	2:B:793:MET:HB3	1.78	0.41
1:C:668:THR:C	1:C:702:VAL:CG1	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:309:SER:OG	2:D:310:GLY:N	2.53	0.41
2:D:579:PHE:HD1	2:D:579:PHE:HA	1.79	0.41
1:A:654:SER:H	1:A:657:ASP:HB3	1.85	0.41
1:A:721:LEU:HD13	1:A:725:MET:HG2	1.99	0.41
1:C:415:LYS:CA	1:C:460:PHE:CD1	2.93	0.41
1:C:780:ARG:NH1	1:C:785:GLY:O	2.42	0.41
1:C:426:PRO:HD3	1:C:728:TYR:CZ	2.55	0.41
1:A:713:LEU:HD21	1:A:736:LEU:HD21	2.03	0.41
2:B:690:THR:HA	2:B:693:PHE:HD2	1.84	0.41
1:C:520:GLY:HA3	1:C:742:LEU:HD22	2.02	0.41
1:C:560:VAL:HG21	2:D:831:LEU:HD11	2.02	0.41
2:D:539:SER:OG	2:D:756:GLY:O	2.38	0.41
2:D:823:ILE:HD12	2:D:823:ILE:HA	1.80	0.41
1:A:526:ARG:NH2	1:A:713:LEU:O	2.54	0.41
1:A:540:ASP:N	1:A:541:PRO:HD2	2.35	0.41
2:B:677:GLN:OE1	2:B:678:THR:N	2.53	0.41
2:B:832:VAL:HA	2:B:835:VAL:HG22	2.02	0.41
1:C:499:ALA:HB1	1:C:748:TYR:CZ	2.56	0.41
1:C:555:LEU:HD12	1:C:555:LEU:HA	1.92	0.41
2:D:81:GLN:HE21	2:D:94:LYS:HD2	1.85	0.41
2:D:516:PRO:HB3	2:D:764:TYR:CE2	2.56	0.41
1:C:113:LYS:HD3	1:C:113:LYS:HA	1.93	0.40
2:D:240:LYS:NZ	2:D:309:SER:O	2.52	0.40
1:A:721:LEU:HA	1:A:721:LEU:HD23	1.88	0.40
2:B:260:ASP:OD1	2:B:318:THR:OG1	2.33	0.40
2:B:556:LEU:HD11	2:B:564:TRP:CZ3	2.56	0.40
1:C:120:VAL:HG21	1:C:307:MET:HE2	2.02	0.40
1:A:669:ILE:HA	1:A:702:VAL:HG22	2.02	0.40
1:A:712:VAL:HA	1:A:717:TYR:HB3	2.03	0.40
2:B:648:ILE:HD13	2:B:648:ILE:HA	1.63	0.40
1:A:27:ALA:HB3	1:A:93:VAL:HG23	2.04	0.40
1:A:502:ILE:HD11	1:A:749:GLY:HA3	2.02	0.40
1:A:524:LEU:HB3	1:A:726:ASN:ND2	2.36	0.40
2:B:261:LEU:HB3	2:B:392:ILE:HD11	2.03	0.40
2:B:538:ILE:HB	2:B:759:ILE:HD12	2.04	0.40
2:B:717:LEU:HD23	2:B:717:LEU:HA	1.90	0.40
1:C:153:LEU:HD23	1:C:153:LEU:HA	1.92	0.40
1:C:227:LEU:HD23	1:C:227:LEU:HA	1.94	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	741/1101 (67%)	699 (94%)	42 (6%)	0	100	100
1	C	741/1101 (67%)	683 (92%)	57 (8%)	1 (0%)	48	81
2	B	749/942 (80%)	710 (95%)	38 (5%)	1 (0%)	48	81
2	D	749/942 (80%)	702 (94%)	46 (6%)	1 (0%)	48	81
All	All	2980/4086 (73%)	2794 (94%)	183 (6%)	3 (0%)	49	81

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	550	PRO
2	B	559	LEU
1	C	431	ARG

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	568/950 (60%)	562 (99%)	6 (1%)	65	74
1	C	569/950 (60%)	564 (99%)	5 (1%)	70	76
2	B	587/817 (72%)	581 (99%)	6 (1%)	68	76
2	D	585/817 (72%)	579 (99%)	6 (1%)	68	76
All	All	2309/3534 (65%)	2286 (99%)	23 (1%)	65	76

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	THR
1	A	539	LEU
1	A	542	PHE
1	A	619	ILE
1	A	644	THR
1	A	799	LEU
2	B	149	LEU
2	B	441	GLU
2	B	555	PHE
2	B	648	ILE
2	B	774	TYR
2	B	819	ASN
1	C	397	LEU
1	C	419	VAL
1	C	542	PHE
1	C	643	LEU
1	C	702	VAL
2	D	480	ILE
2	D	558	PRO
2	D	755	ILE
2	D	815	LEU
2	D	823	ILE
2	D	825	ILE

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

Mol	Chain	Res	Type
1	A	312	HIS
1	A	348	ASN
1	A	638	ASN
1	A	691	ASN
1	A	726	ASN
1	A	772	ASN
2	B	67	ASN
2	B	81	GLN
2	B	98	GLN
2	B	119	GLN
2	B	123	ASN
2	B	129	HIS
2	B	203	GLN
2	B	288	GLN
2	B	412	ASN
2	B	721	ASN

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Mol	Chain	Res	Type
1	C	129	GLN
1	C	285	ASN
1	C	291	ASN
1	C	312	HIS
1	C	322	ASN
1	C	342	HIS
1	C	348	ASN
1	C	385	HIS
1	C	490	ASN
1	C	661	GLN
1	C	663	ASN
1	C	680	ASN
1	C	726	ASN
1	C	735	ASN
1	C	738	GLN
1	C	772	ASN
2	D	81	GLN
2	D	119	GLN
2	D	123	ASN
2	D	129	HIS
2	D	136	HIS
2	D	203	GLN
2	D	328	HIS
2	D	412	ASN
2	D	417	GLN
2	D	557	ASN
2	D	677	GLN
2	D	726	GLN
2	D	747	GLN

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.

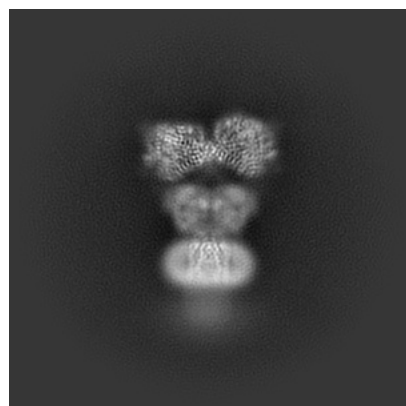
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-70807. These allow visual inspection of the internal detail of the map and identification of artifacts.

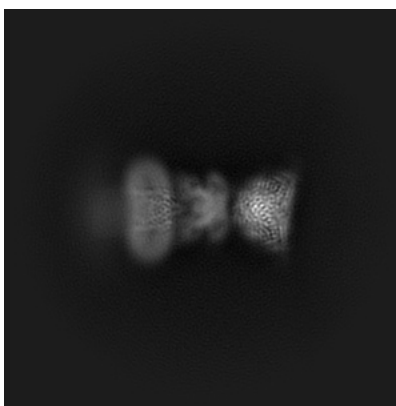
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

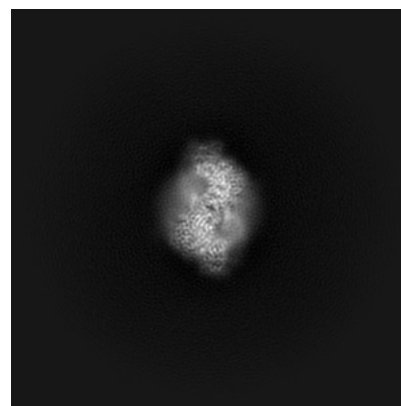
6.1.1 Primary map



X

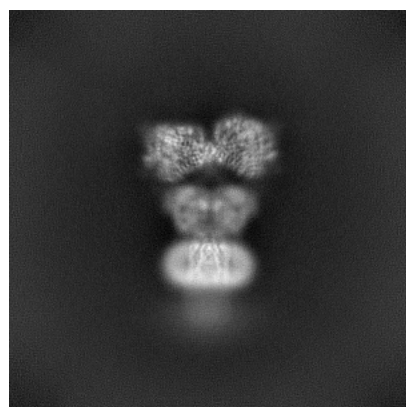


Y

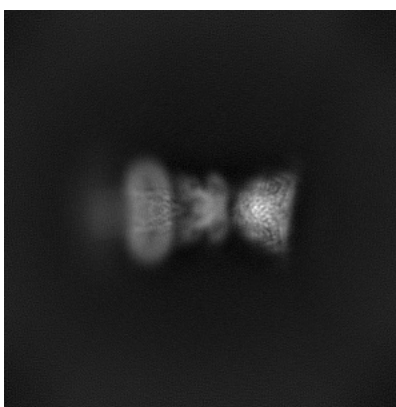


Z

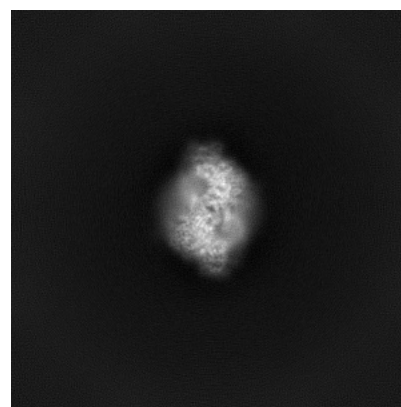
6.1.2 Raw map



X



Y

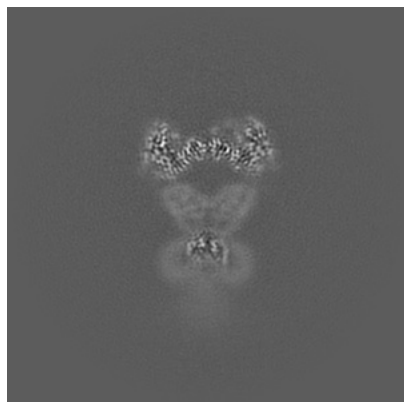


Z

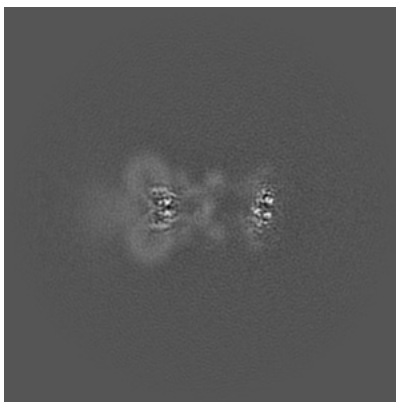
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

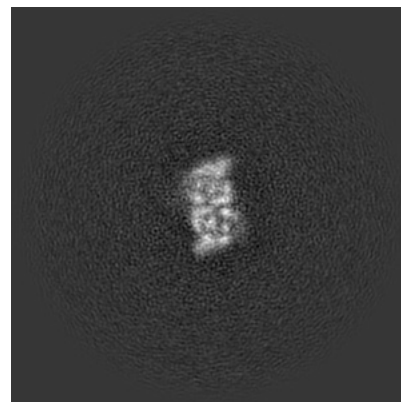
6.2.1 Primary map



X Index: 208

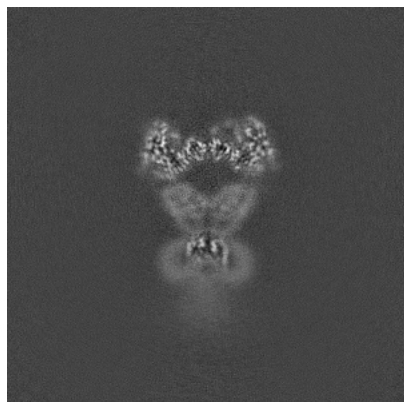


Y Index: 208

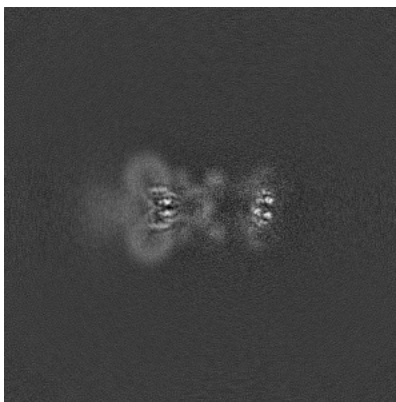


Z Index: 208

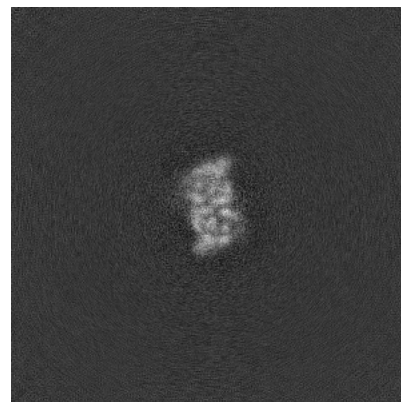
6.2.2 Raw map



X Index: 208



Y Index: 208

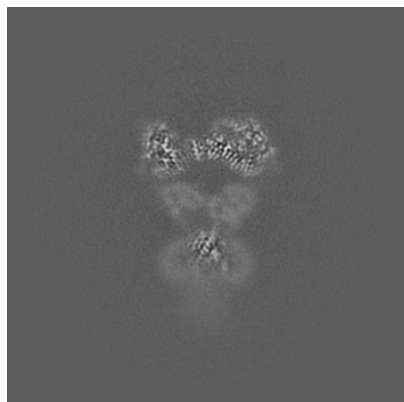


Z Index: 208

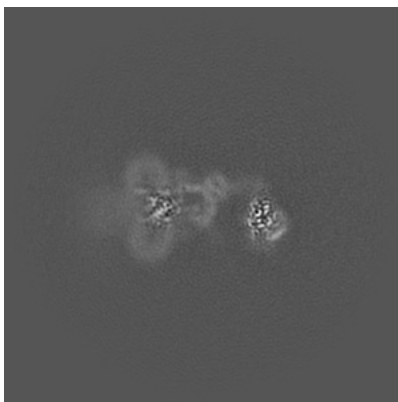
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

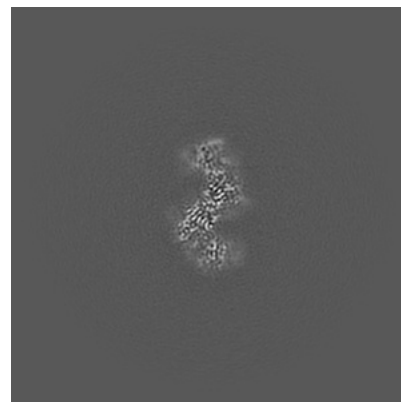
6.3.1 Primary map



X Index: 212

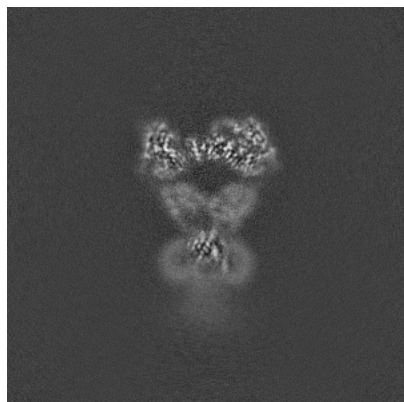


Y Index: 200

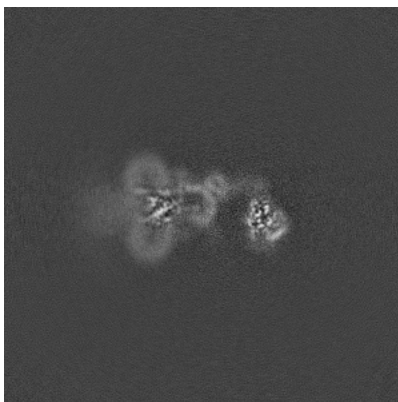


Z Index: 267

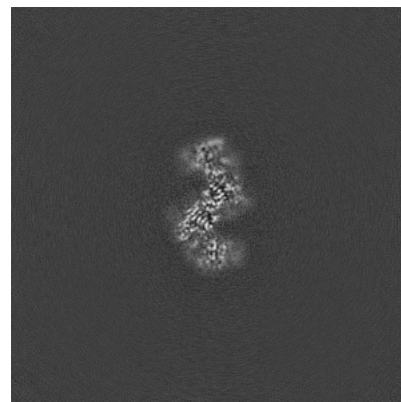
6.3.2 Raw map



X Index: 211



Y Index: 200



Z Index: 267

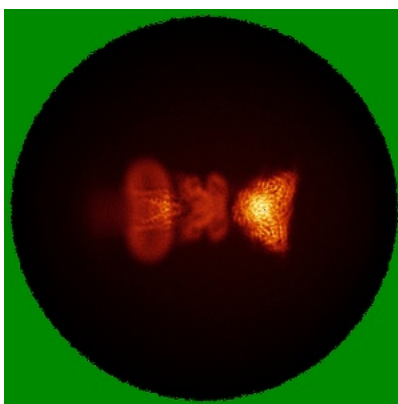
The images above show the largest variance slices of the map in three orthogonal directions.

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

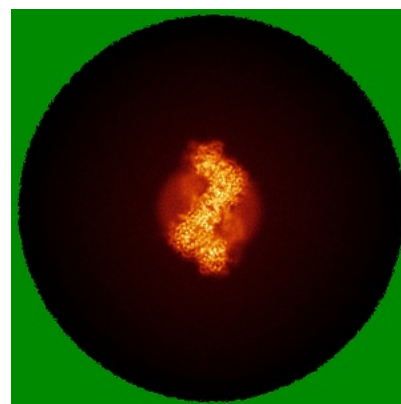
6.4.1 Primary map



X

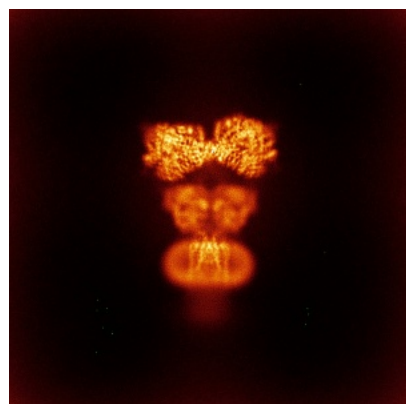


Y

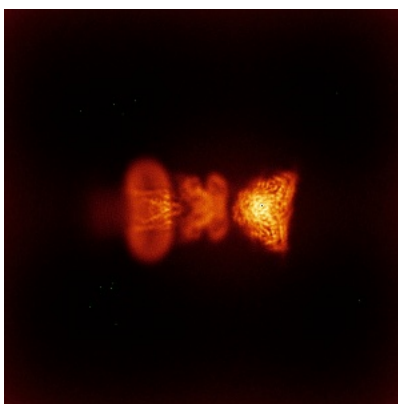


Z

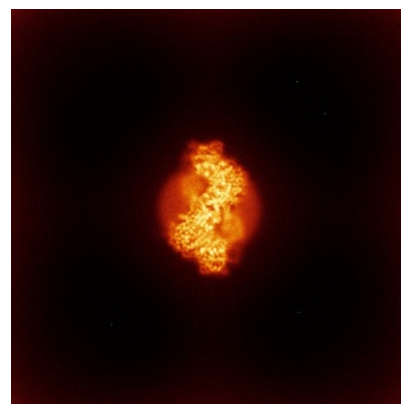
6.4.2 Raw map



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.

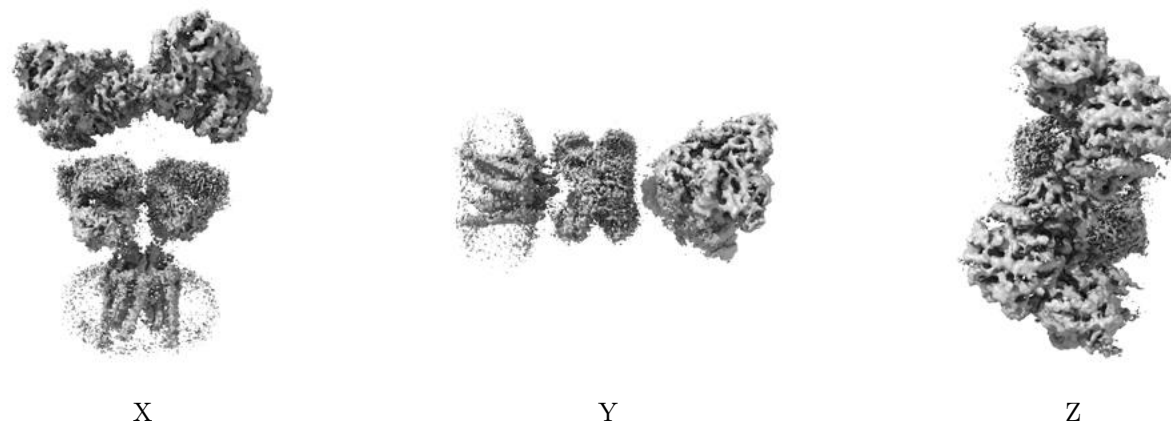
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.38. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

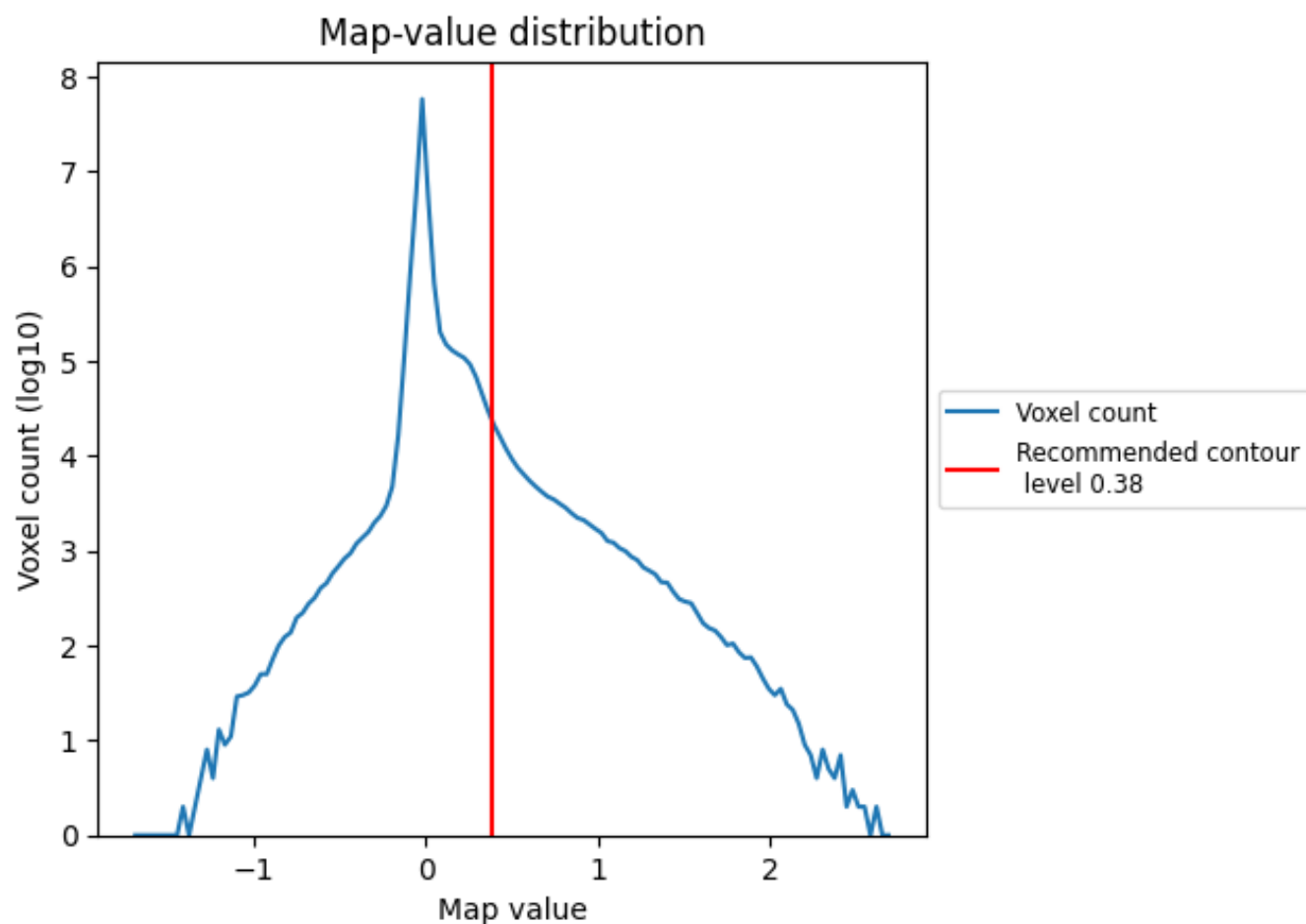
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

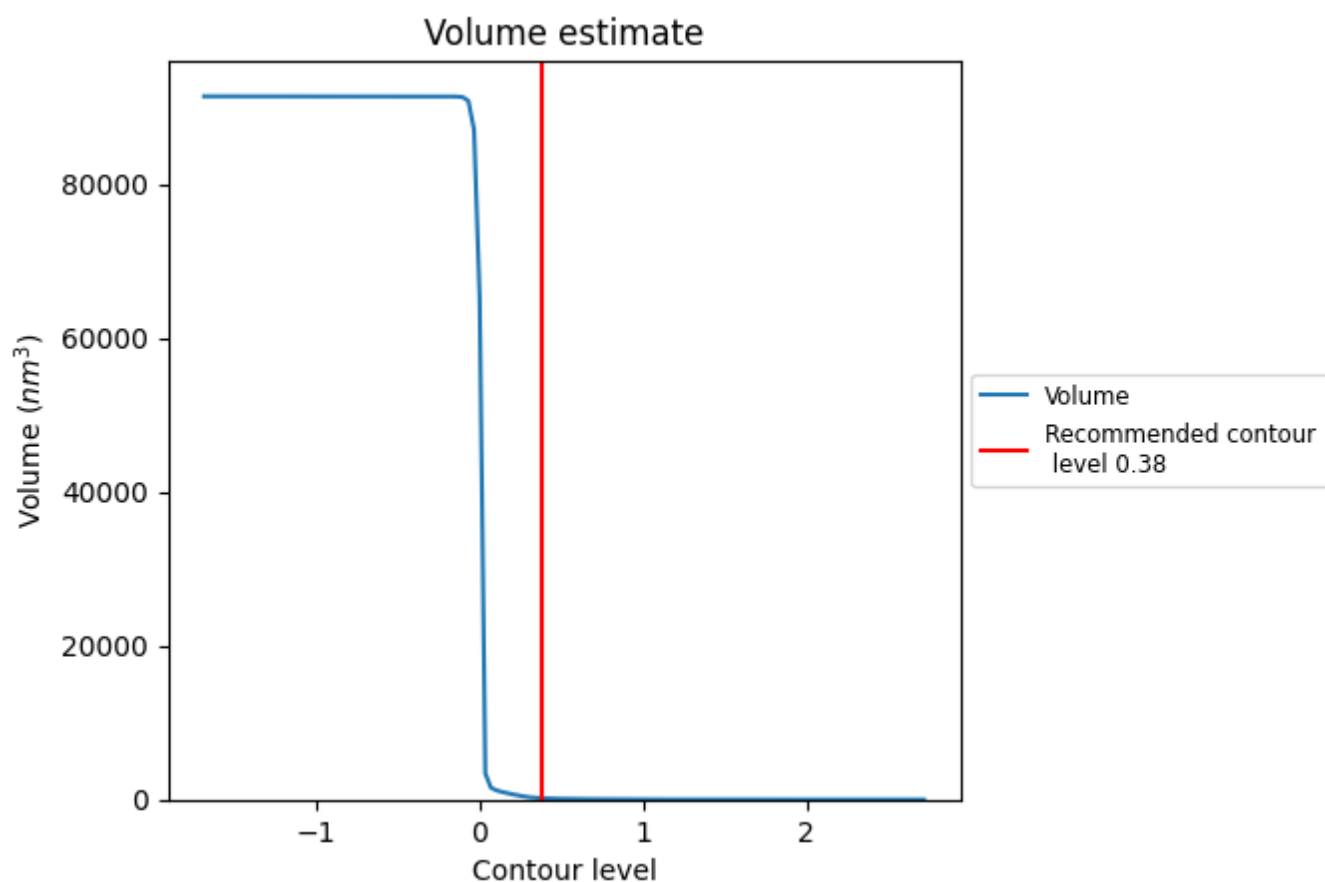
This section contains the results of statistical analysis of the map.

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.

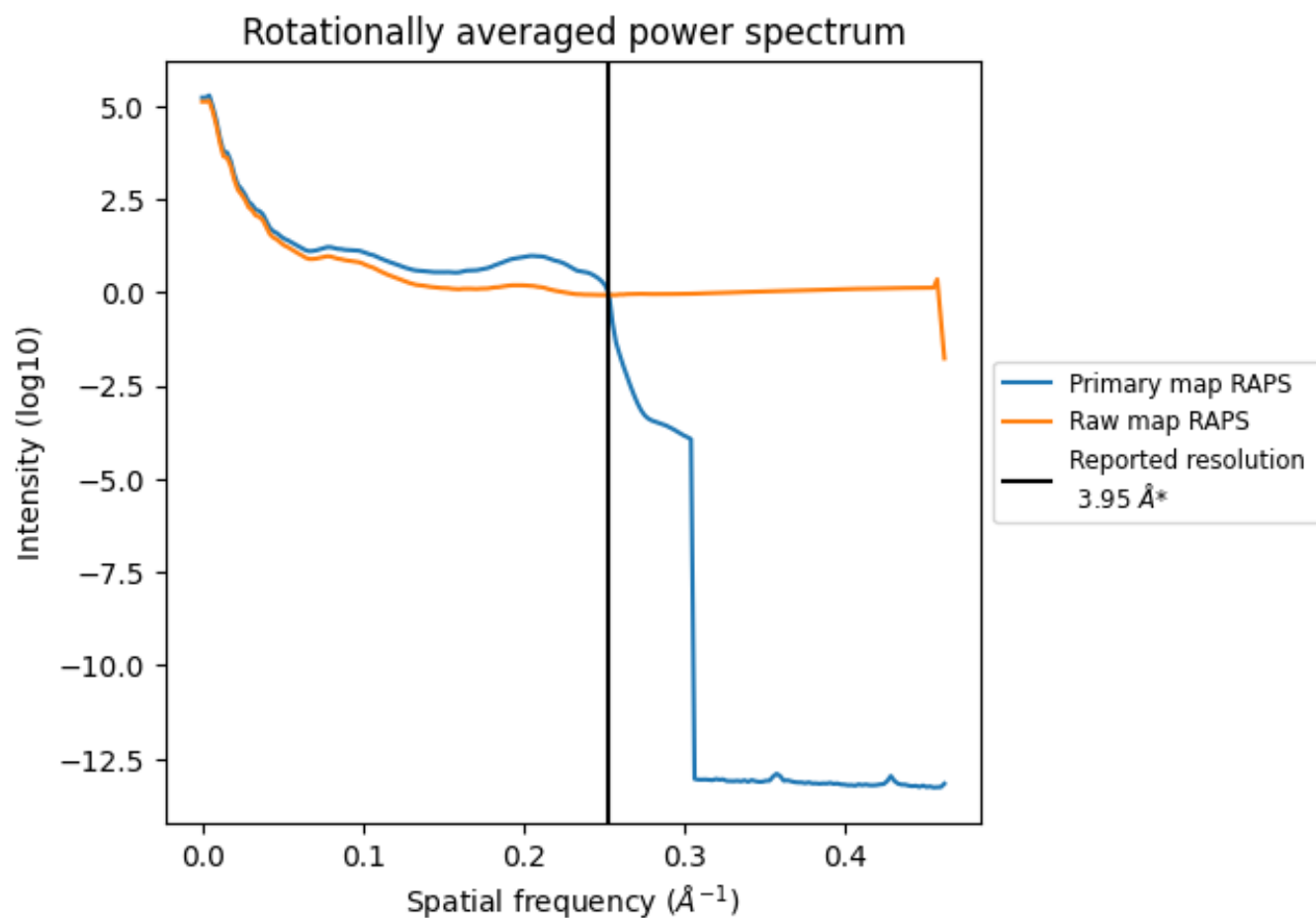
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 172 nm³; this corresponds to an approximate mass of 155 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 ⓘ

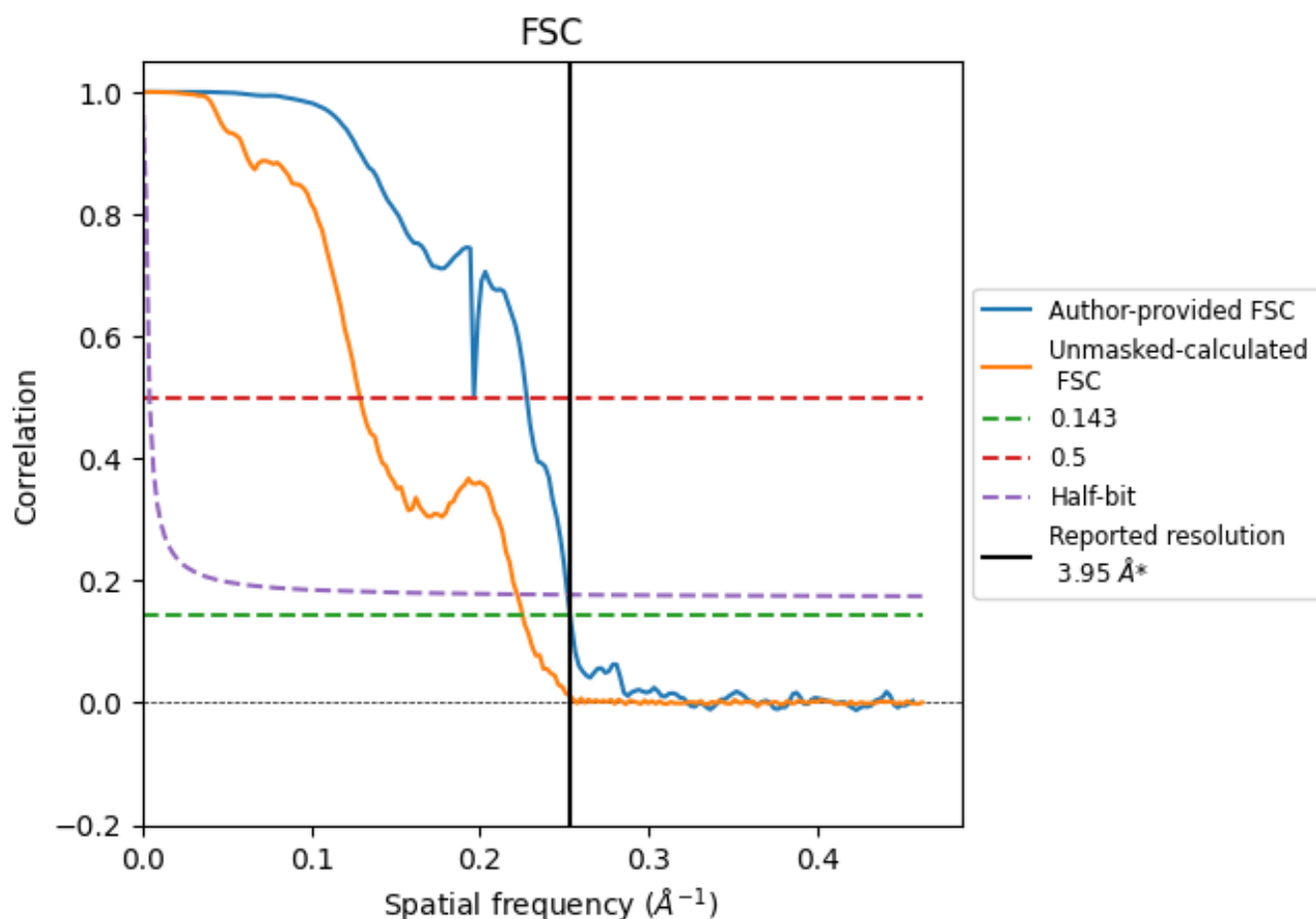


*Reported resolution corresponds to spatial frequency of 0.253 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.253 \AA^{-1}

8.2 Resolution estimates [i](#)

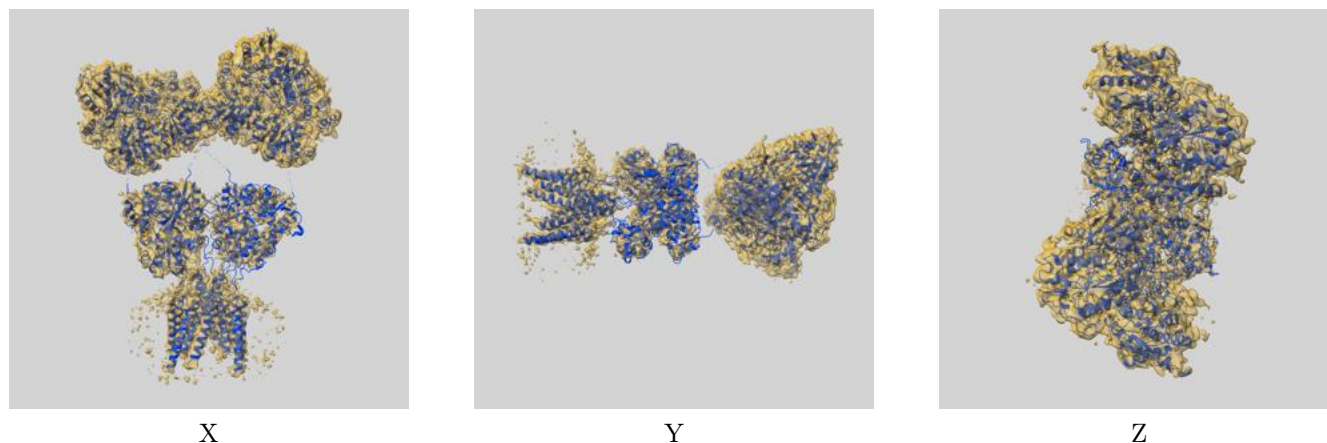
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.95	-	-
Author-provided FSC curve	3.95	4.39	3.97
Unmasked-calculated*	4.44	7.76	4.51

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.44 differs from the reported value 3.95 by more than 10 %

9 Map-model fit [i](#)

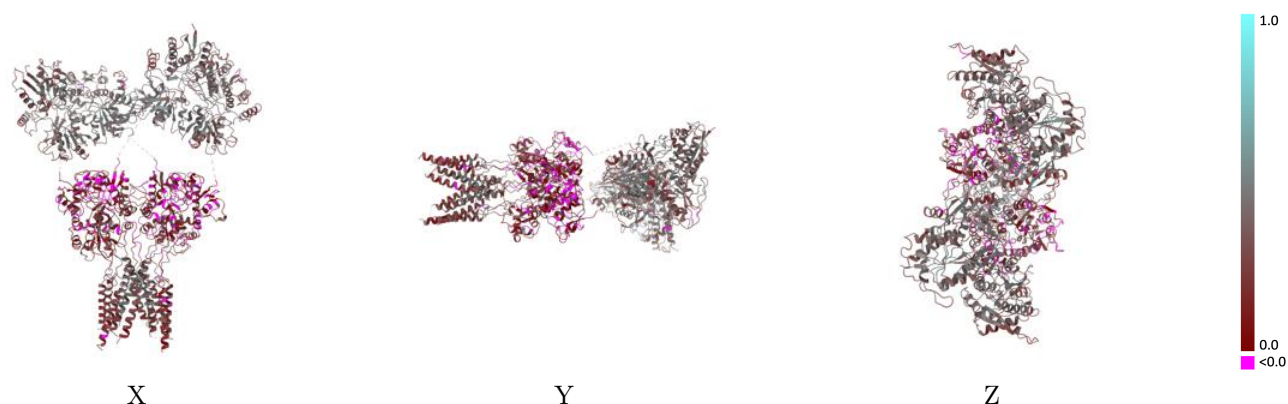
This section contains information regarding the fit between EMDB map EMD-70807 and PDB model 9OSI. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



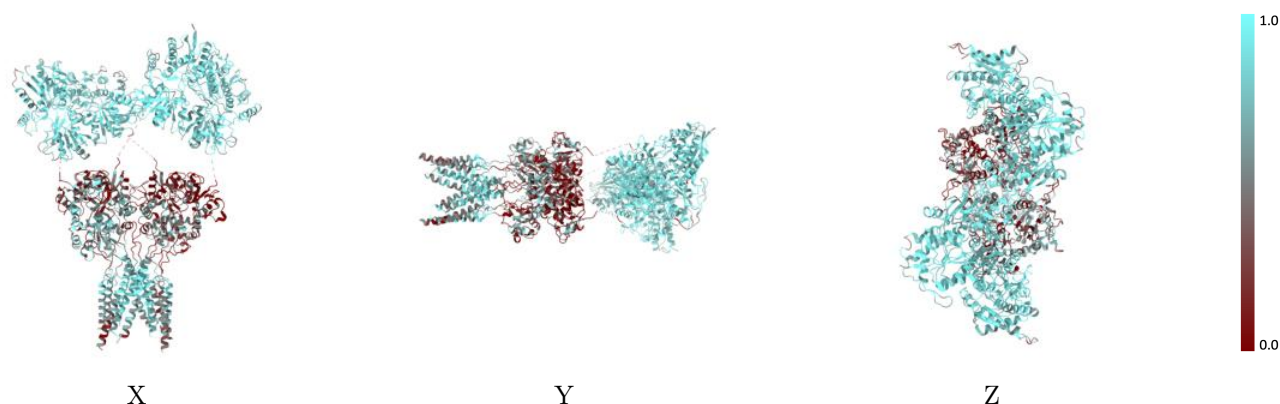
The images above show the 3D surface view of the map at the recommended contour level 0.38 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



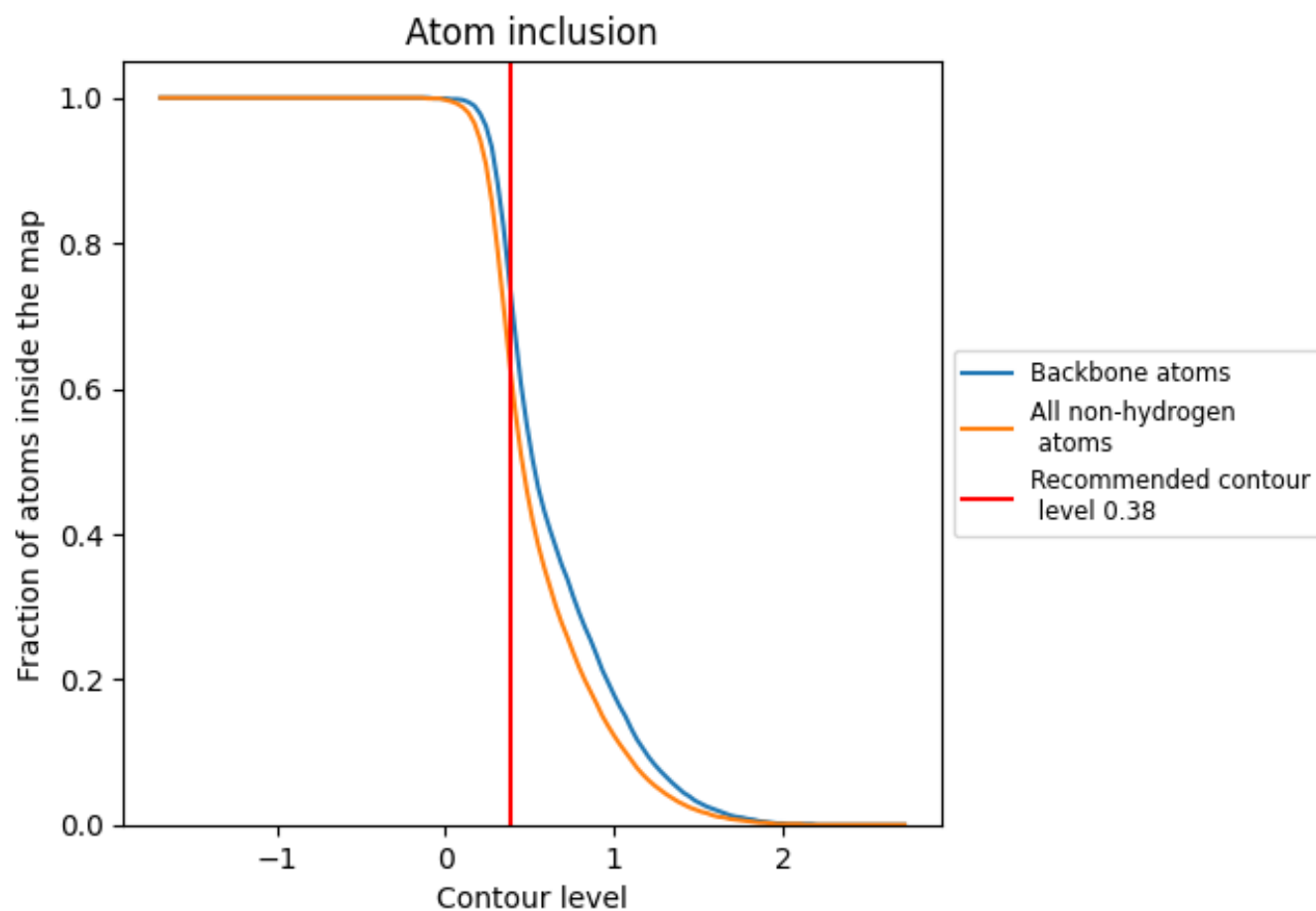
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.38).

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 64% 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 (0.38) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6360	<div></div> 0.2970
A	<div></div> 0.6020	<div></div> 0.2730
B	<div></div> 0.6770	<div></div> 0.3170
C	<div></div> 0.6490	<div></div> 0.2920
D	<div></div> 0.6160	<div></div> 0.3050

