



Full wwPDB EM Validation Report ⓘ

Mar 7, 2026 – 02:30 AM UTC

PDB ID : 9MDU / pdb_00009mdu
EMDB ID : EMD-48180
Title : Human CXCR4 tetramer
Authors : Zhang, Z.; Patel, D.J.; Junyu, X.
Deposited on : 2024-12-05
Resolution : 2.90 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

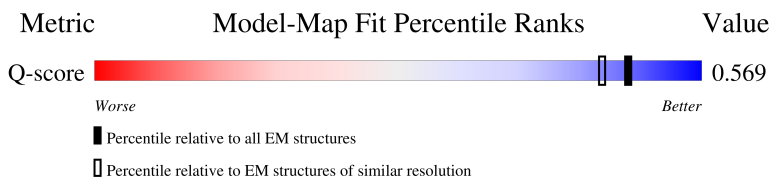
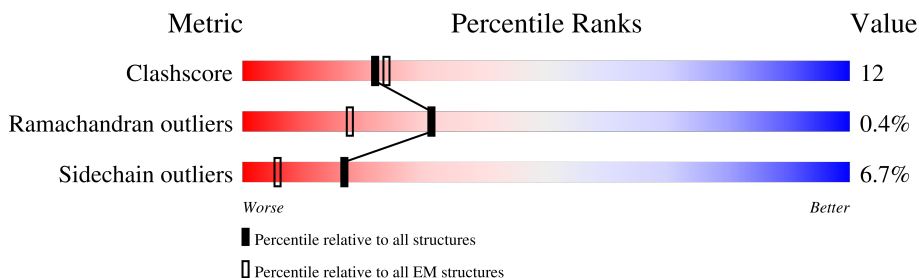
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.90 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	13054 (2.40 - 3.40)

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	360	 5% 57% 16% 22%
1	B	360	 5% 58% 16% 22%
1	C	360	 5% 55% 18% 22%
1	D	360	 5% 55% 19% 22%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9123 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 C-X-C chemokine receptor type 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	281	Total	C	N	O	S	0	0
			2282	1530	363	377	12		
1	A	281	Total	C	N	O	S	0	0
			2277	1526	362	377	12		
1	C	281	Total	C	N	O	S	0	0
			2282	1530	363	377	12		
1	D	281	Total	C	N	O	S	0	0
			2282	1530	363	377	12		

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	353	ASP	-	expression tag	UNP P61073
B	354	TYR	-	expression tag	UNP P61073
B	355	LYS	-	expression tag	UNP P61073
B	356	ASP	-	expression tag	UNP P61073
B	357	ASP	-	expression tag	UNP P61073
B	358	ASP	-	expression tag	UNP P61073
B	359	ASP	-	expression tag	UNP P61073
B	360	LYS	-	expression tag	UNP P61073
A	353	ASP	-	expression tag	UNP P61073
A	354	TYR	-	expression tag	UNP P61073
A	355	LYS	-	expression tag	UNP P61073
A	356	ASP	-	expression tag	UNP P61073
A	357	ASP	-	expression tag	UNP P61073
A	358	ASP	-	expression tag	UNP P61073
A	359	ASP	-	expression tag	UNP P61073
A	360	LYS	-	expression tag	UNP P61073
C	353	ASP	-	expression tag	UNP P61073
C	354	TYR	-	expression tag	UNP P61073
C	355	LYS	-	expression tag	UNP P61073
C	356	ASP	-	expression tag	UNP P61073
C	357	ASP	-	expression tag	UNP P61073
C	358	ASP	-	expression tag	UNP P61073

Continued on next page...

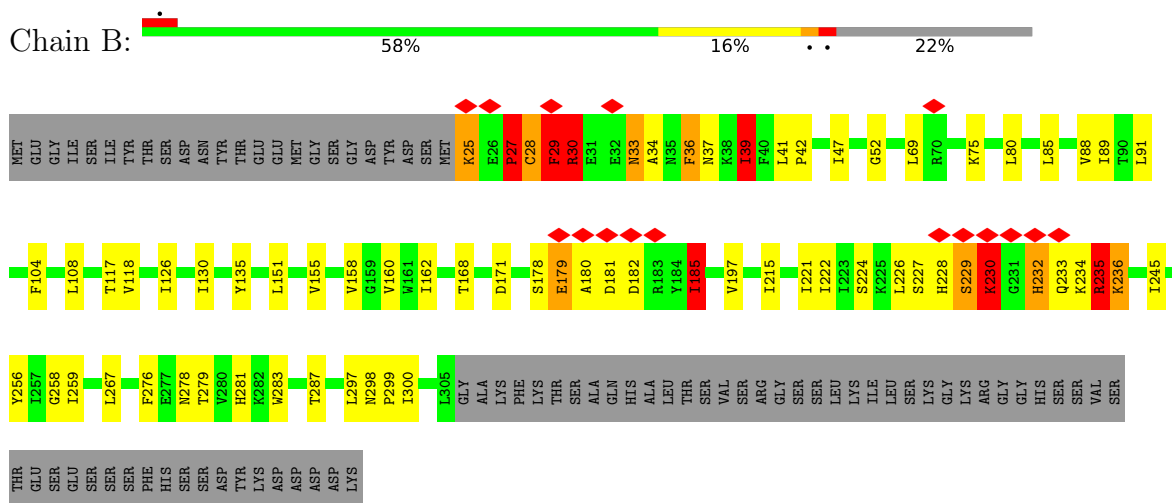
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	359	ASP	-	expression tag	UNP P61073
C	360	LYS	-	expression tag	UNP P61073
D	353	ASP	-	expression tag	UNP P61073
D	354	TYR	-	expression tag	UNP P61073
D	355	LYS	-	expression tag	UNP P61073
D	356	ASP	-	expression tag	UNP P61073
D	357	ASP	-	expression tag	UNP P61073
D	358	ASP	-	expression tag	UNP P61073
D	359	ASP	-	expression tag	UNP P61073
D	360	LYS	-	expression tag	UNP P61073

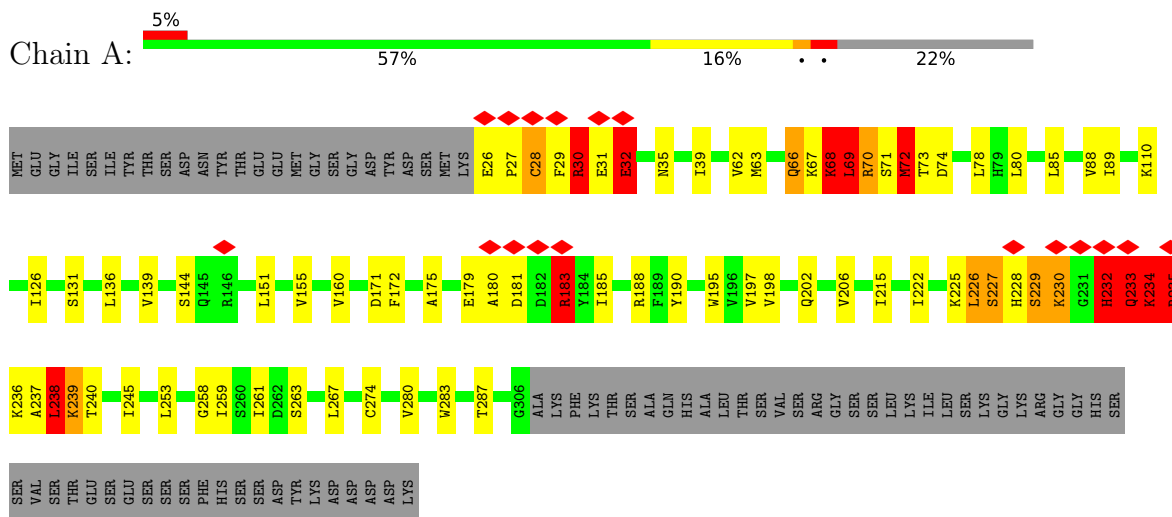
3 Residue-property plots

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

• Molecule 1: C-X-C chemokine receptor type 4

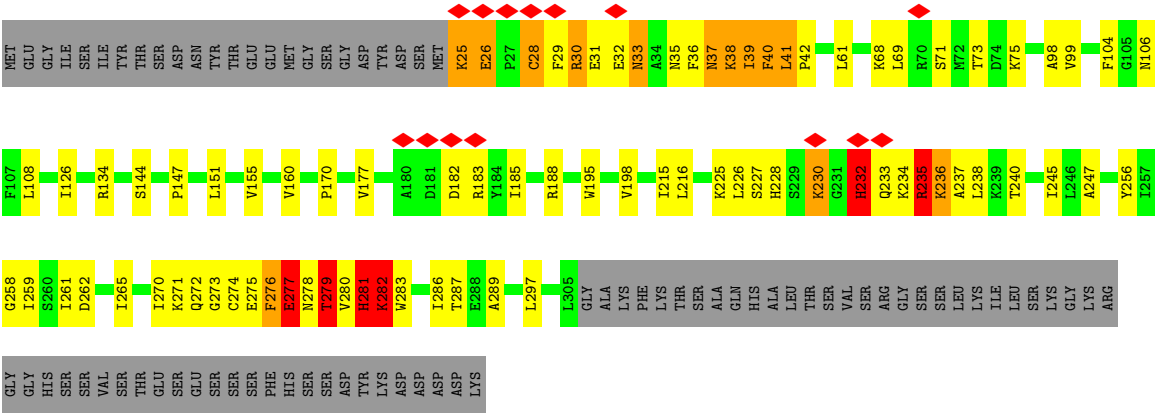


• Molecule 1: C-X-C chemokine receptor type 4

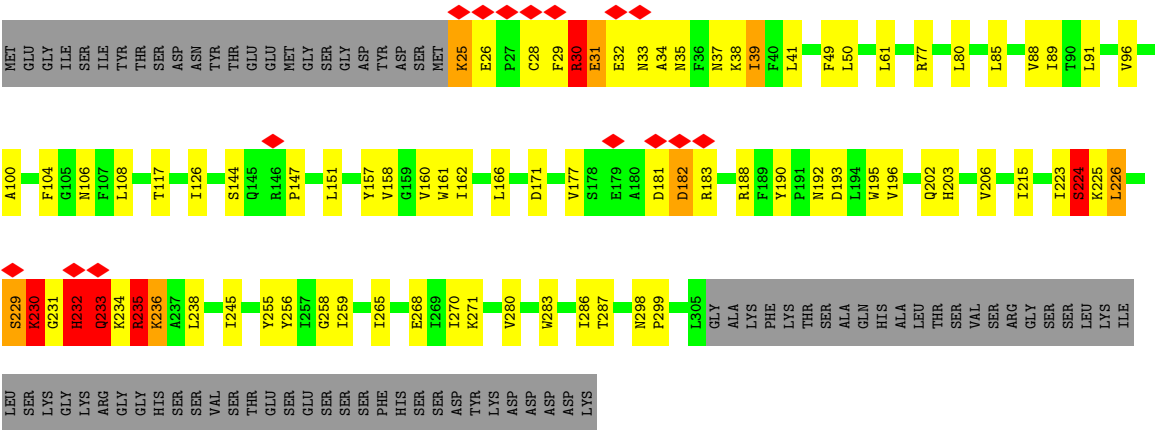


• Molecule 1: C-X-C chemokine receptor type 4





• Molecule 1: C-X-C chemokine receptor type 4



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	734989	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	59	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	2.901	Depositor
Minimum map value	-2.108	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.061	Depositor
Recommended contour level	0.28	Depositor
Map size (Å)	273.92, 273.92, 273.92	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07, 1.07, 1.07	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	0.59	1/2341 (0.0%)	1.13	29/3189 (0.9%)
1	B	0.53	2/2346 (0.1%)	1.31	24/3195 (0.8%)
1	C	0.55	1/2346 (0.0%)	1.13	28/3195 (0.9%)
1	D	0.50	0/2346	0.94	20/3195 (0.6%)
All	All	0.54	4/9379 (0.0%)	1.14	101/12774 (0.8%)

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
1	A	0	5
1	B	0	4
1	C	0	2
1	D	0	3
All	All	0	14

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	233	GLN	C-N	-18.69	1.09	1.34
1	B	230	LYS	C-N	-13.35	1.18	1.33
1	C	40	PHE	CA-C	-5.67	1.45	1.52
1	B	178	SER	CA-CB	-5.59	1.46	1.53

All (101) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	PRO	CB-CA-C	-22.29	74.78	111.56
1	B	28	CYS	N-CA-CB	19.84	144.01	110.49
1	B	228	HIS	CB-CA-C	-19.25	84.39	111.14
1	B	37	ASN	CB-CA-C	18.22	139.69	110.90

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	GLN	O-C-N	-17.77	98.95	122.59
1	A	66	GLN	N-CA-C	-17.25	80.62	108.73
1	B	230	LYS	CB-CA-C	16.80	143.86	110.42
1	B	37	ASN	N-CA-C	-16.46	93.37	111.14
1	C	29	PHE	CB-CA-C	-16.04	90.08	109.80
1	D	232	HIS	N-CA-C	15.97	135.14	111.34
1	B	229	SER	N-CA-CB	-14.70	93.88	110.35
1	C	38	LYS	N-CA-C	-14.49	89.03	110.48
1	D	235	ARG	CB-CA-C	14.26	134.42	110.74
1	B	27	PRO	CA-C-N	13.82	147.93	121.54
1	B	27	PRO	C-N-CA	13.82	147.93	121.54
1	A	234	LYS	CB-CA-C	13.58	136.80	110.46
1	A	239	LYS	N-CA-C	-13.14	82.81	110.80
1	D	231	GLY	N-CA-C	13.08	126.72	112.33
1	A	235	ARG	N-CA-CB	-12.74	90.92	110.42
1	C	272	GLN	N-CA-C	12.53	127.40	111.24
1	B	230	LYS	O-C-N	-12.19	106.38	122.59
1	B	29	PHE	N-CA-C	-12.01	94.70	111.52
1	B	185	ILE	N-CA-C	11.97	124.17	108.12
1	B	230	LYS	N-CA-C	-11.41	86.50	110.80
1	A	69	LEU	N-CA-C	-11.06	93.22	109.15
1	B	28	CYS	N-CA-C	-11.04	87.30	110.80
1	A	239	LYS	N-CA-CB	10.74	128.64	110.49
1	C	29	PHE	N-CA-C	10.32	125.91	110.64
1	A	234	LYS	N-CA-CB	-10.19	94.04	110.14
1	B	180	ALA	N-CA-C	10.06	125.09	108.48
1	D	223	ILE	CB-CA-C	9.90	127.53	111.29
1	C	272	GLN	CB-CA-C	-9.85	94.61	110.86
1	C	274	CYS	N-CA-C	-9.85	101.60	114.31
1	C	37	ASN	CB-CA-C	-9.49	93.08	110.63
1	A	229	SER	CB-CA-C	9.38	122.74	109.71
1	A	233	GLN	CA-C-N	9.32	133.99	120.38
1	A	233	GLN	C-N-CA	9.32	133.99	120.38
1	C	232	HIS	N-CA-C	-9.13	104.21	114.62
1	D	233	GLN	N-CA-CB	-9.06	96.04	109.82
1	C	278	ASN	N-CA-C	-9.04	100.89	114.64
1	D	229	SER	CA-C-N	8.98	138.70	121.54
1	D	229	SER	C-N-CA	8.98	138.70	121.54
1	D	230	LYS	N-CA-C	8.91	129.79	110.80
1	A	31	GLU	N-CA-CB	-8.86	99.29	110.98
1	B	39	ILE	N-CA-C	-8.85	104.85	113.53
1	C	226	LEU	N-CA-C	-8.80	101.23	112.93

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	277	GLU	N-CA-CB	-8.55	97.10	111.20
1	C	277	GLU	N-CA-C	-8.44	101.43	112.41
1	C	279	THR	N-CA-C	-8.34	103.24	113.41
1	B	229	SER	N-CA-C	-8.33	97.97	109.71
1	C	38	LYS	N-CA-CB	-8.25	97.75	110.29
1	A	229	SER	N-CA-C	-8.23	98.40	110.42
1	C	30	ARG	N-CA-CB	8.22	123.95	110.39
1	D	229	SER	CB-CA-C	8.11	125.96	109.67
1	A	35	ASN	N-CA-C	-8.05	105.44	114.62
1	B	30	ARG	CB-CA-C	7.96	126.25	110.42
1	D	224	SER	N-CA-CB	-7.95	98.84	110.53
1	A	68	LYS	N-CA-C	-7.87	102.99	112.59
1	A	235	ARG	N-CA-C	7.76	122.83	113.12
1	A	73	THR	N-CA-C	-7.55	103.01	112.90
1	C	30	ARG	N-CA-C	-7.55	93.71	107.75
1	A	32	GLU	N-CA-C	7.54	121.53	109.24
1	C	37	ASN	N-CA-C	7.40	120.22	111.71
1	C	41	LEU	N-CA-C	-7.35	102.63	112.55
1	A	233	GLN	N-CA-C	-6.99	95.91	110.80
1	C	38	LYS	CA-C-O	-6.92	113.51	121.47
1	A	72	MET	N-CA-C	-6.87	104.78	113.72
1	D	30	ARG	N-CA-C	-6.87	99.49	109.59
1	B	185	ILE	CB-CA-C	-6.71	104.04	111.35
1	D	232	HIS	CB-CA-C	-6.71	102.05	112.31
1	A	67	LYS	N-CA-CB	-6.68	99.20	110.49
1	B	228	HIS	N-CA-C	6.66	121.30	112.25
1	A	28	CYS	CB-CA-C	6.57	123.49	110.42
1	C	282	LYS	N-CA-C	-6.50	105.31	113.18
1	C	35	ASN	N-CA-C	-6.35	105.36	113.23
1	C	40	PHE	N-CA-CB	6.18	120.94	110.49
1	D	235	ARG	N-CA-C	-6.16	101.20	111.37
1	A	234	LYS	N-CA-C	-6.05	104.01	111.75
1	A	73	THR	CA-CB-OG1	-6.04	100.54	109.60
1	C	273	GLY	N-CA-C	-5.93	99.13	113.18
1	A	32	GLU	N-CA-CB	-5.90	100.82	110.43
1	C	281	HIS	N-CA-C	-5.88	106.75	114.04
1	A	238	LEU	N-CA-C	-5.81	98.42	110.80
1	B	36	PHE	CB-CA-C	-5.72	98.57	109.95
1	A	232	HIS	N-CA-C	-5.71	103.46	110.65
1	C	26	GLU	CB-CA-C	5.69	118.20	109.60
1	D	229	SER	N-CA-C	5.58	120.21	112.90
1	C	276	PHE	CA-CB-CG	-5.54	108.26	113.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	67	LYS	N-CA-C	-5.49	99.10	110.80
1	B	39	ILE	CB-CA-C	5.42	115.96	110.65
1	C	276	PHE	N-CA-C	-5.37	105.56	113.61
1	B	185	ILE	N-CA-CB	-5.33	105.76	111.46
1	D	226	LEU	N-CA-C	-5.27	106.24	113.30
1	D	236	LYS	CA-C-N	-5.23	110.82	121.18
1	D	236	LYS	C-N-CA	-5.23	110.82	121.18
1	C	28	CYS	N-CA-C	5.18	116.97	110.91
1	B	41	LEU	N-CA-C	-5.17	105.43	112.35
1	D	31	GLU	N-CA-C	-5.11	102.12	110.20
1	D	182	ASP	N-CA-C	-5.11	107.10	113.38
1	A	30	ARG	CB-CA-C	-5.06	102.46	110.81
1	D	30	ARG	CB-CA-C	5.05	117.84	110.26

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	ARG	Sidechain
1	A	233	GLN	Mainchain
1	A	235	ARG	Sidechain
1	A	30	ARG	Sidechain
1	A	70	ARG	Sidechain
1	B	230	LYS	Mainchain
1	B	235	ARG	Sidechain
1	B	27	PRO	Peptide
1	B	30	ARG	Sidechain
1	C	235	ARG	Peptide,Sidechain
1	D	229	SER	Peptide
1	D	235	ARG	Sidechain
1	D	30	ARG	Sidechain

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	2277	0	2340	60	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2282	0	2350	55	0
1	C	2282	0	2351	57	0
1	D	2282	0	2348	57	0
All	All	9123	0	9389	215	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 12.

All (215) 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:D:33:ASN:HD21	1:D:37:ASN:CB	1.18	1.53
1:D:33:ASN:ND2	1:D:37:ASN:HB2	1.14	1.43
1:B:27:PRO:O	1:B:27:PRO:CG	1.66	1.22
1:A:181:ASP:OD2	1:A:183:ARG:HG3	1.38	1.19
1:A:30:ARG:O	1:A:30:ARG:HG3	1.42	1.09
1:A:62:VAL:O	1:A:66:GLN:O	1.77	1.03
1:B:229:SER:O	1:B:232:HIS:CE1	2.13	1.01
1:A:181:ASP:OD2	1:A:183:ARG:CG	2.11	0.97
1:B:27:PRO:C	1:B:29:PHE:H	1.62	0.91
1:B:27:PRO:O	1:B:27:PRO:HG2	1.09	0.91
1:A:232:HIS:O	1:A:233:GLN:C	2.14	0.90
1:D:33:ASN:HD21	1:D:37:ASN:HB3	1.40	0.85
1:A:226:LEU:HD11	1:A:234:LYS:O	1.78	0.84
1:C:36:PHE:O	1:C:38:LYS:O	1.97	0.83
1:B:27:PRO:C	1:B:29:PHE:N	2.36	0.83
1:A:181:ASP:CG	1:A:183:ARG:HG3	2.02	0.82
1:B:27:PRO:O	1:B:29:PHE:N	2.10	0.81
1:C:236:LYS:HZ1	1:C:240:THR:HG23	1.46	0.81
1:B:229:SER:C	1:B:232:HIS:CE1	2.60	0.78
1:B:229:SER:O	1:B:232:HIS:ND1	2.15	0.78
1:B:39:ILE:HD11	1:A:258:GLY:HA2	1.67	0.77
1:C:236:LYS:NZ	1:C:240:THR:HG23	1.99	0.77
1:C:275:GLU:O	1:C:279:THR:OG1	2.03	0.77
1:C:277:GLU:O	1:C:277:GLU:HG3	1.85	0.77
1:C:28:CYS:SG	1:C:28:CYS:O	2.44	0.76
1:A:233:GLN:O	1:A:234:LYS:C	2.23	0.75
1:B:185:ILE:O	1:B:185:ILE:HG12	1.89	0.73
1:C:277:GLU:HA	1:C:280:VAL:HG12	1.69	0.73
1:A:233:GLN:O	1:A:235:ARG:N	2.23	0.72
1:A:63:MET:HB3	1:A:69:LEU:HD13	1.72	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:PHE:HB3	1:C:108:LEU:HD23	1.72	0.71
1:D:33:ASN:CG	1:D:37:ASN:HB2	2.13	0.70
1:A:232:HIS:O	1:A:233:GLN:O	2.08	0.70
1:A:228:HIS:O	1:A:229:SER:C	2.36	0.69
1:B:233:GLN:OE1	1:B:235:ARG:HD2	1.94	0.67
1:A:151:LEU:HA	1:A:155:VAL:HB	1.76	0.66
1:C:42:PRO:HG3	1:C:98:ALA:HB3	1.78	0.65
1:C:232:HIS:HA	1:C:235:ARG:HD3	1.77	0.65
1:B:104:PHE:HB3	1:B:108:LEU:HD23	1.77	0.65
1:D:104:PHE:HB3	1:D:108:LEU:HD23	1.79	0.64
1:A:30:ARG:O	1:A:30:ARG:CG	2.29	0.63
1:C:69:LEU:HD12	1:C:75:LYS:HG2	1.80	0.62
1:B:197:VAL:HG11	1:B:267:LEU:HD13	1.81	0.62
1:B:28:CYS:SG	1:B:30:ARG:C	2.84	0.61
1:C:36:PHE:C	1:C:38:LYS:O	2.43	0.61
1:D:33:ASN:ND2	1:D:37:ASN:CB	2.00	0.61
1:C:30:ARG:HD3	1:C:281:HIS:HD2	1.65	0.61
1:A:235:ARG:HG3	1:A:239:LYS:HB2	1.83	0.60
1:D:230:LYS:O	1:D:230:LYS:CG	2.44	0.60
1:B:39:ILE:HD12	1:A:261:ILE:HD12	1.83	0.59
1:C:39:ILE:HD13	1:D:280:VAL:HG22	1.84	0.59
1:A:226:LEU:CD1	1:A:234:LYS:O	2.51	0.59
1:B:224:SER:HB3	1:D:61:LEU:HD21	1.85	0.59
1:D:235:ARG:O	1:D:238:LEU:HB3	2.02	0.59
1:A:63:MET:HB3	1:A:69:LEU:CD1	2.33	0.58
1:B:258:GLY:HA2	1:D:39:ILE:HD11	1.86	0.58
1:D:230:LYS:O	1:D:230:LYS:CD	2.52	0.58
1:B:126:ILE:O	1:B:130:ILE:HG13	2.04	0.58
1:A:283:TRP:O	1:A:287:THR:OG1	2.18	0.57
1:B:151:LEU:HA	1:B:155:VAL:HB	1.85	0.57
1:A:69:LEU:HD12	1:A:74:ASP:HB3	1.86	0.57
1:D:232:HIS:HA	1:D:235:ARG:HE	1.70	0.56
1:C:28:CYS:SG	1:C:277:GLU:HG2	2.45	0.56
1:A:27:PRO:O	1:A:28:CYS:C	2.47	0.56
1:C:236:LYS:C	1:C:238:LEU:N	2.62	0.56
1:C:41:LEU:HB2	1:C:42:PRO:HD3	1.87	0.56
1:D:35:ASN:HA	1:D:38:LYS:HD2	1.87	0.56
1:A:85:LEU:O	1:A:89:ILE:HG12	2.06	0.55
1:B:28:CYS:SG	1:B:30:ARG:O	2.63	0.55
1:D:190:TYR:HD2	1:D:196:VAL:HG22	1.71	0.55
1:D:215:ILE:HG21	1:D:245:ILE:HG21	1.89	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:ILE:HG21	1:B:245:ILE:HG21	1.89	0.54
1:C:38:LYS:HG3	1:C:99:VAL:HA	1.90	0.54
1:D:77:ARG:NH2	1:D:298:ASN:O	2.39	0.54
1:D:162:ILE:HG23	1:D:166:LEU:HD13	1.89	0.54
1:B:39:ILE:HD13	1:A:280:VAL:HG22	1.89	0.54
1:C:41:LEU:HD21	1:C:286:ILE:HD13	1.89	0.54
1:A:227:SER:O	1:A:229:SER:O	2.26	0.53
1:B:28:CYS:HB3	1:B:278:ASN:HD21	1.74	0.53
1:B:85:LEU:O	1:B:89:ILE:HG12	2.07	0.53
1:A:215:ILE:HG21	1:A:245:ILE:HG21	1.90	0.53
1:A:139:VAL:HG22	1:A:225:LYS:HG3	1.92	0.53
1:D:33:ASN:O	1:D:34:ALA:C	2.51	0.52
1:A:39:ILE:HD12	1:C:261:ILE:HD12	1.91	0.52
1:A:188:ARG:NE	1:A:190:TYR:OH	2.38	0.52
1:B:47:ILE:HG12	1:A:253:LEU:HD23	1.91	0.52
1:C:151:LEU:HA	1:C:155:VAL:HB	1.92	0.52
1:D:85:LEU:O	1:D:89:ILE:HG12	2.09	0.52
1:D:234:LYS:O	1:D:235:ARG:C	2.53	0.52
1:C:39:ILE:HD11	1:D:258:GLY:HA2	1.92	0.52
1:D:232:HIS:CD2	1:D:236:LYS:HZ2	2.27	0.52
1:C:170:PRO:HB2	1:C:188:ARG:HH11	1.75	0.52
1:D:188:ARG:NE	1:D:190:TYR:OH	2.42	0.51
1:B:117:THR:OG1	1:B:171:ASP:OD2	2.15	0.51
1:B:234:LYS:HA	1:B:236:LYS:HD2	1.92	0.51
1:A:69:LEU:HD21	1:A:78:LEU:HD22	1.92	0.51
1:A:171:ASP:OD1	1:A:188:ARG:NH1	2.45	0.50
1:B:69:LEU:HD13	1:B:75:LYS:HG2	1.93	0.50
1:D:96:VAL:O	1:D:100:ALA:N	2.42	0.50
1:A:32:GLU:H	1:A:32:GLU:CD	2.16	0.50
1:A:226:LEU:HD21	1:A:234:LYS:HA	1.93	0.50
1:A:195:TRP:HA	1:A:198:VAL:HG12	1.94	0.50
1:C:37:ASN:O	1:C:41:LEU:HG	2.12	0.50
1:D:203:HIS:HE1	1:D:255:TYR:HE2	1.59	0.50
1:B:88:VAL:HG12	1:B:91:LEU:HD12	1.94	0.49
1:C:126:ILE:HG13	1:C:160:VAL:HG11	1.93	0.49
1:C:215:ILE:HG21	1:C:245:ILE:HG21	1.94	0.49
1:D:202:GLN:O	1:D:206:VAL:HG12	2.12	0.49
1:A:27:PRO:C	1:A:29:PHE:N	2.70	0.48
1:B:80:LEU:HG	1:B:126:ILE:HG21	1.94	0.48
1:D:232:HIS:CD2	1:D:236:LYS:NZ	2.82	0.48
1:A:126:ILE:HG13	1:A:160:VAL:HG11	1.94	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:ILE:HG13	1:D:160:VAL:HG11	1.96	0.48
1:B:126:ILE:HG13	1:B:160:VAL:HG11	1.94	0.48
1:A:226:LEU:HD11	1:A:234:LYS:C	2.38	0.47
1:D:25:LYS:HB2	1:D:25:LYS:HE2	1.45	0.47
1:C:276:PHE:O	1:C:276:PHE:CG	2.67	0.47
1:A:237:ALA:O	1:A:240:THR:HG22	2.14	0.47
1:D:232:HIS:O	1:D:235:ARG:HB3	2.14	0.47
1:A:197:VAL:HG11	1:A:267:LEU:HD21	1.97	0.47
1:C:144:SER:HA	1:C:147:PRO:HD2	1.96	0.46
1:D:230:LYS:O	1:D:230:LYS:HD2	2.15	0.46
1:A:28:CYS:HB2	1:A:274:CYS:HB3	1.96	0.46
1:C:61:LEU:HD21	1:D:224:SER:HB2	1.96	0.46
1:D:226:LEU:HD23	1:D:226:LEU:HA	1.78	0.46
1:A:230:LYS:H	1:A:230:LYS:HG2	1.52	0.46
1:D:41:LEU:HD21	1:D:286:ILE:HD13	1.98	0.46
1:B:230:LYS:HB2	1:B:232:HIS:NE2	2.31	0.46
1:C:236:LYS:HZ2	1:C:236:LYS:HG2	1.36	0.46
1:A:68:LYS:CE	1:C:227:SER:O	2.64	0.45
1:A:179:GLU:O	1:A:181:ASP:O	2.34	0.45
1:D:117:THR:OG1	1:D:171:ASP:OD2	2.28	0.45
1:D:108:LEU:HD12	1:D:108:LEU:HA	1.83	0.45
1:C:42:PRO:HD3	1:C:98:ALA:HB1	1.99	0.45
1:D:49:PHE:HD2	1:D:50:LEU:HD22	1.81	0.45
1:D:144:SER:OG	1:D:144:SER:O	2.34	0.45
1:D:147:PRO:O	1:D:151:LEU:HB2	2.17	0.45
1:B:118:VAL:HG23	1:B:168:THR:HG21	1.99	0.45
1:B:283:TRP:O	1:B:287:THR:OG1	2.21	0.45
1:C:106:ASN:HA	1:C:177:VAL:HG23	1.98	0.45
1:D:29:PHE:CE1	1:D:30:ARG:O	2.70	0.45
1:A:72:MET:HE2	1:A:72:MET:HB2	1.59	0.45
1:C:39:ILE:O	1:C:39:ILE:HG23	2.17	0.45
1:B:276:PHE:HA	1:B:279:THR:HG22	1.97	0.45
1:B:85:LEU:HA	1:B:88:VAL:HG22	1.99	0.45
1:C:71:SER:C	1:C:73:THR:H	2.24	0.45
1:B:28:CYS:SG	1:B:278:ASN:ND2	2.90	0.45
1:B:29:PHE:C	1:B:30:ARG:O	2.60	0.45
1:C:68:LYS:HA	1:C:68:LYS:HD3	1.68	0.44
1:C:282:LYS:HE2	1:C:282:LYS:HB3	1.54	0.44
1:B:233:GLN:CD	1:B:235:ARG:HD2	2.41	0.44
1:A:110:LYS:NZ	1:A:172:PHE:O	2.42	0.44
1:C:25:LYS:HE2	1:C:25:LYS:HB2	1.42	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:283:TRP:O	1:D:287:THR:OG1	2.30	0.44
1:A:39:ILE:HD11	1:C:258:GLY:HA2	1.99	0.44
1:A:202:GLN:O	1:A:206:VAL:HG12	2.17	0.44
1:D:192:ASN:HB3	1:D:195:TRP:HD1	1.82	0.44
1:A:136:LEU:O	1:A:144:SER:OG	2.35	0.44
1:D:106:ASN:HA	1:D:177:VAL:HG23	2.00	0.44
1:B:52:GLY:HA3	1:B:88:VAL:HG11	1.99	0.44
1:D:158:VAL:HA	1:D:162:ILE:HD12	1.99	0.43
1:C:265:ILE:HD12	1:C:280:VAL:HG11	2.00	0.43
1:C:134:ARG:HD3	1:C:238:LEU:HD12	2.01	0.43
1:A:68:LYS:HD2	1:C:227:SER:O	2.19	0.43
1:C:33:ASN:HD22	1:C:33:ASN:HA	1.53	0.43
1:D:190:TYR:CD2	1:D:196:VAL:HG22	2.53	0.43
1:D:265:ILE:HG13	1:D:270:ILE:HD12	1.99	0.43
1:A:85:LEU:HA	1:A:88:VAL:HG22	2.00	0.43
1:B:36:PHE:O	1:B:36:PHE:CG	2.71	0.43
1:C:236:LYS:O	1:C:238:LEU:N	2.45	0.43
1:C:195:TRP:HA	1:C:198:VAL:HG22	2.01	0.42
1:C:262:ASP:OD1	1:C:281:HIS:HE1	2.02	0.42
1:D:268:GLU:O	1:D:271:LYS:NZ	2.46	0.42
1:C:236:LYS:C	1:C:238:LEU:H	2.23	0.42
1:B:298:ASN:HB3	1:B:299:PRO:HD3	2.00	0.42
1:D:298:ASN:HB3	1:D:299:PRO:HD3	2.01	0.42
1:A:222:ILE:HD13	1:A:238:LEU:HD11	2.02	0.42
1:B:33:ASN:O	1:B:34:ALA:C	2.63	0.42
1:A:179:GLU:O	1:A:180:ALA:C	2.61	0.42
1:D:88:VAL:HG12	1:D:91:LEU:HD12	2.02	0.42
1:D:157:TYR:HA	1:D:161:TRP:HD1	1.84	0.42
1:D:85:LEU:HA	1:D:88:VAL:HG22	2.01	0.42
1:C:41:LEU:HD22	1:C:289:ALA:HB2	2.01	0.42
1:C:236:LYS:NZ	1:C:237:ALA:H	2.18	0.42
1:A:259:ILE:O	1:A:263:SER:OG	2.37	0.42
1:C:247:ALA:HB1	1:C:297:LEU:HD22	2.02	0.42
1:C:270:ILE:O	1:C:271:LYS:C	2.61	0.42
1:B:25:LYS:HD3	1:B:25:LYS:HA	1.79	0.41
1:B:179:GLU:H	1:B:179:GLU:HG2	1.57	0.41
1:C:216:LEU:HD23	1:C:216:LEU:HA	1.90	0.41
1:D:256:TYR:HA	1:D:259:ILE:HG12	2.01	0.41
1:A:175:ALA:HA	1:A:188:ARG:HA	2.02	0.41
1:B:28:CYS:O	1:B:30:ARG:O	2.39	0.41
1:B:69:LEU:HD12	1:B:69:LEU:O	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:ILE:HD13	1:C:280:VAL:HG23	2.02	0.41
1:D:181:ASP:HB3	1:D:182:ASP:H	1.66	0.41
1:D:80:LEU:HG	1:D:126:ILE:HG21	2.03	0.41
1:B:256:TYR:HA	1:B:259:ILE:HG12	2.02	0.41
1:A:183:ARG:HG3	1:A:183:ARG:H	1.42	0.41
1:B:158:VAL:HG12	1:B:162:ILE:HD12	2.03	0.41
1:A:80:LEU:HG	1:A:126:ILE:HG21	2.03	0.41
1:C:283:TRP:O	1:C:287:THR:OG1	2.35	0.41
1:B:135:TYR:HE1	1:B:221:ILE:HG21	1.86	0.41
1:A:131:SER:OG	1:A:215:ILE:HG12	2.21	0.40
1:B:28:CYS:O	1:B:29:PHE:C	2.65	0.40
1:A:69:LEU:HD12	1:A:69:LEU:HA	1.60	0.40
1:B:230:LYS:HD2	1:B:230:LYS:HA	1.73	0.40
1:B:297:LEU:HA	1:B:300:ILE:HB	2.03	0.40
1:C:230:LYS:H	1:C:230:LYS:HG2	1.65	0.40
1:C:256:TYR:HA	1:C:259:ILE:HG12	2.02	0.40
1:B:30:ARG:CZ	1:B:281:HIS:HB3	2.52	0.40
1:B:222:ILE:O	1:B:226:LEU:HB2	2.21	0.40
1:C:236:LYS:HZ3	1:C:240:THR:HG23	1.83	0.40
1:D:192:ASN:OD1	1:D:193:ASP:N	2.54	0.40
1:D:233:GLN:H	1:D:233:GLN:HG3	1.07	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	279/360 (78%)	253 (91%)	25 (9%)	1 (0%)	30	59
1	B	279/360 (78%)	253 (91%)	24 (9%)	2 (1%)	18	48
1	C	279/360 (78%)	256 (92%)	22 (8%)	1 (0%)	30	59
1	D	279/360 (78%)	262 (94%)	17 (6%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1116/1440 (78%)	1024 (92%)	88 (8%)	4 (0%)	31 59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	GLN
1	B	230	LYS
1	C	40	PHE
1	B	27	PRO

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	250/319 (78%)	231 (92%)	19 (8%)	12 36
1	B	251/319 (79%)	238 (95%)	13 (5%)	21 52
1	C	251/319 (79%)	230 (92%)	21 (8%)	10 31
1	D	251/319 (79%)	237 (94%)	14 (6%)	19 50
All	All	1003/1276 (79%)	936 (93%)	67 (7%)	17 43

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	25	LYS
1	B	29	PHE
1	B	33	ASN
1	B	39	ILE
1	B	42	PRO
1	B	179	GLU
1	B	181	ASP
1	B	182	ASP
1	B	185	ILE
1	B	227	SER
1	B	232	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	235	ARG
1	B	236	LYS
1	A	26	GLU
1	A	30	ARG
1	A	32	GLU
1	A	68	LYS
1	A	69	LEU
1	A	70	ARG
1	A	71	SER
1	A	72	MET
1	A	183	ARG
1	A	185	ILE
1	A	226	LEU
1	A	227	SER
1	A	230	LYS
1	A	232	HIS
1	A	233	GLN
1	A	234	LYS
1	A	235	ARG
1	A	236	LYS
1	A	238	LEU
1	C	25	LYS
1	C	26	GLU
1	C	31	GLU
1	C	32	GLU
1	C	33	ASN
1	C	39	ILE
1	C	182	ASP
1	C	183	ARG
1	C	185	ILE
1	C	225	LYS
1	C	228	HIS
1	C	230	LYS
1	C	232	HIS
1	C	233	GLN
1	C	234	LYS
1	C	235	ARG
1	C	236	LYS
1	C	277	GLU
1	C	279	THR
1	C	281	HIS
1	C	282	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	25	LYS
1	D	26	GLU
1	D	28	CYS
1	D	30	ARG
1	D	31	GLU
1	D	32	GLU
1	D	39	ILE
1	D	183	ARG
1	D	224	SER
1	D	225	LYS
1	D	230	LYS
1	D	232	HIS
1	D	233	GLN
1	D	235	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	33	ASN
1	B	145	GLN
1	B	232	HIS
1	B	278	ASN
1	A	33	ASN
1	A	228	HIS
1	A	272	GLN
1	A	281	HIS
1	C	33	ASN
1	C	35	ASN
1	C	145	GLN
1	C	192	ASN
1	C	233	GLN
1	C	281	HIS
1	D	33	ASN
1	D	37	ASN
1	D	119	ASN
1	D	233	GLN
1	D	272	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	230:LYS	C	231:GLY	N	1.17
1	A	233:GLN	C	234:LYS	N	1.09

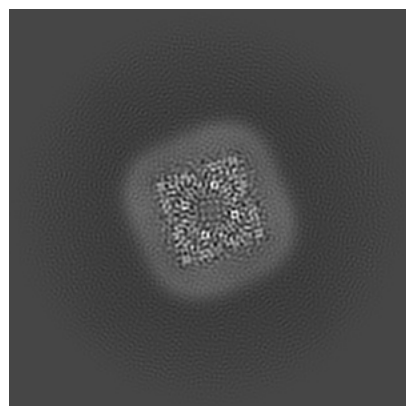
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-48180. These allow visual inspection of the internal detail of the map and identification of artifacts.

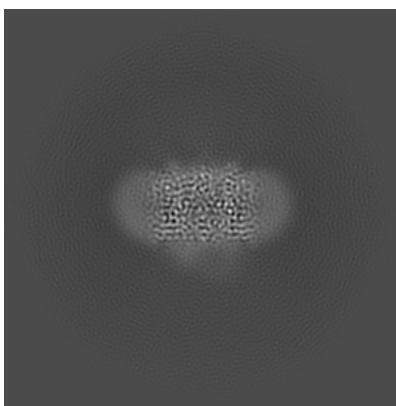
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

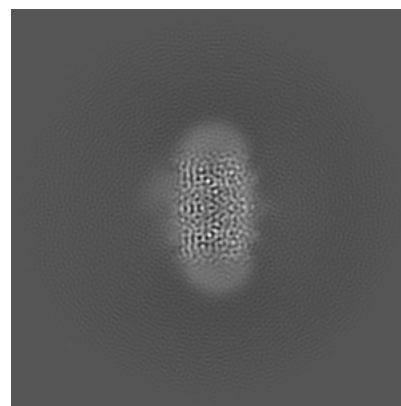
6.1.1 Primary map



X

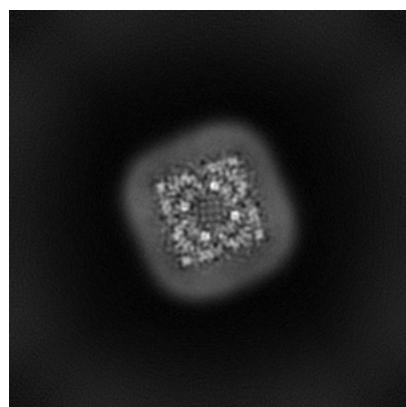


Y

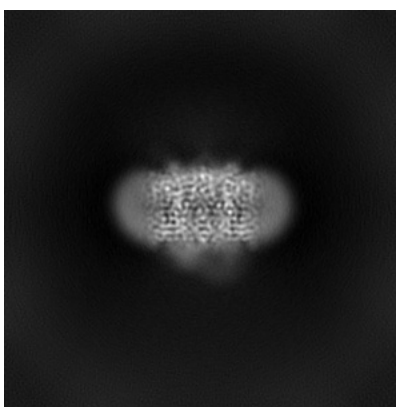


Z

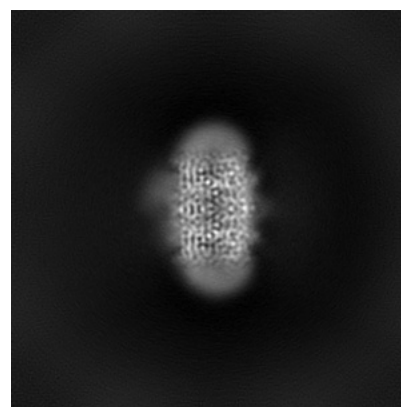
6.1.2 Raw map



X



Y

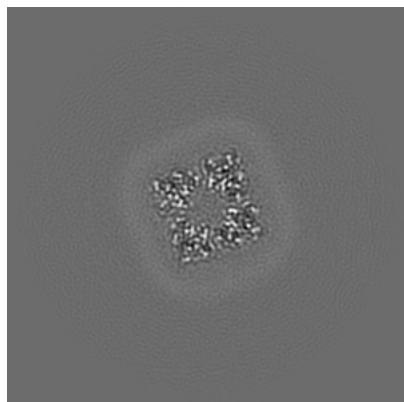


Z

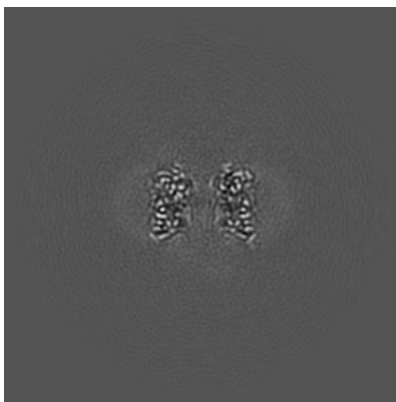
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

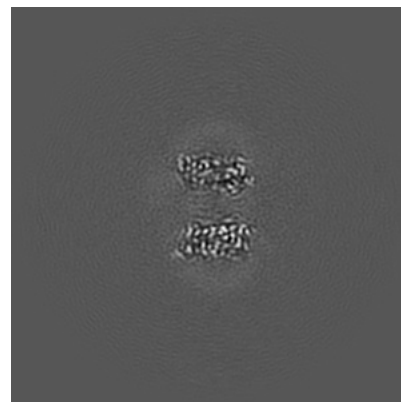
6.2.1 Primary map



X Index: 128

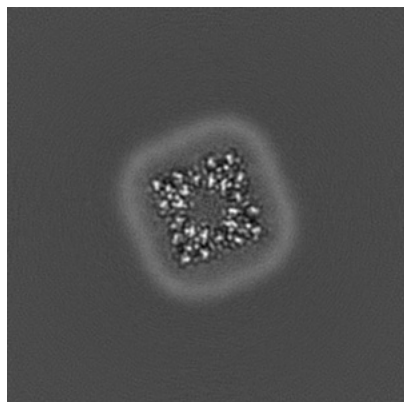


Y Index: 128

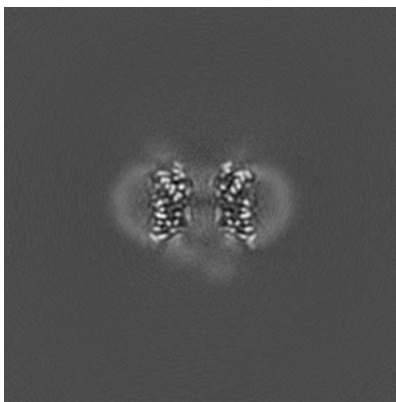


Z Index: 128

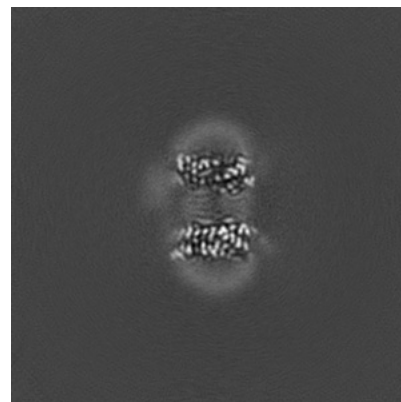
6.2.2 Raw map



X Index: 128



Y Index: 128

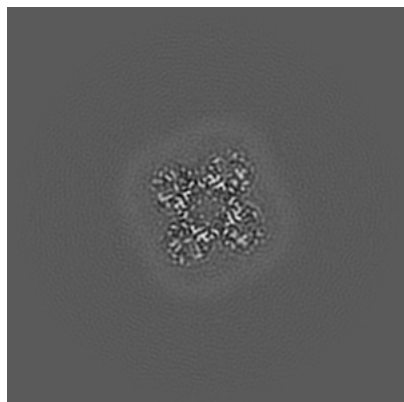


Z Index: 128

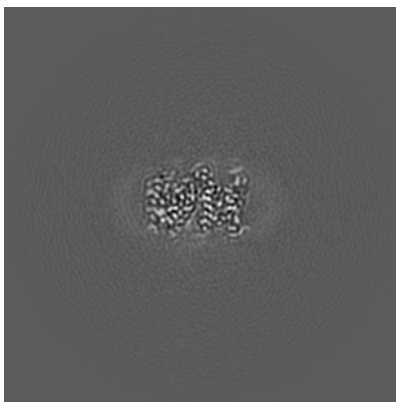
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

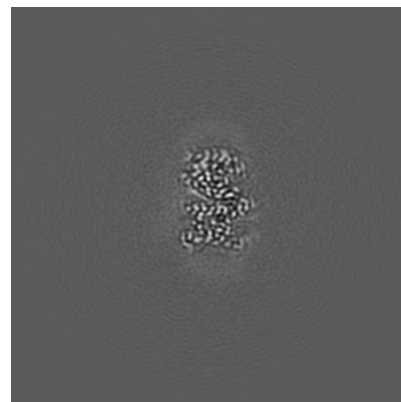
6.3.1 Primary map



X Index: 137

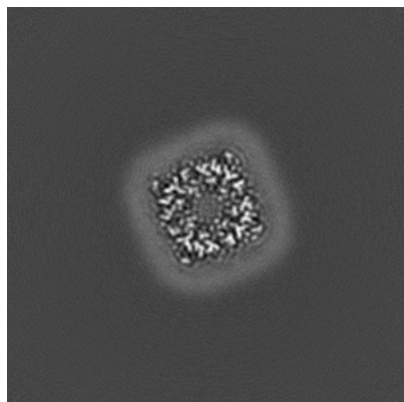


Y Index: 111

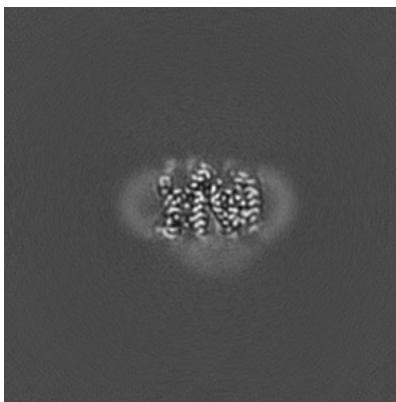


Z Index: 110

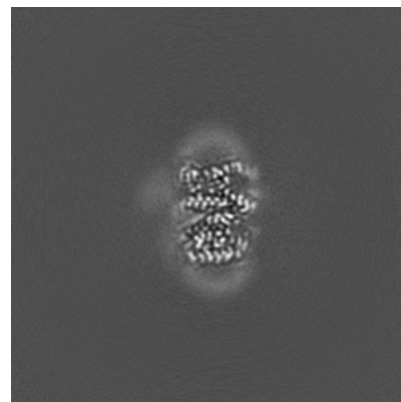
6.3.2 Raw map



X Index: 117



Y Index: 145

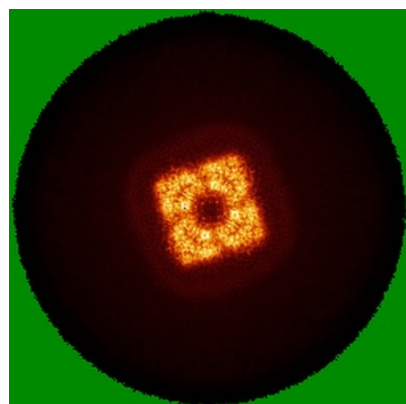


Z Index: 143

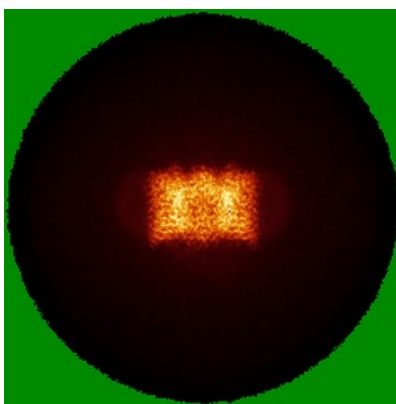
The images above show the largest variance slices of the map in three orthogonal directions.

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

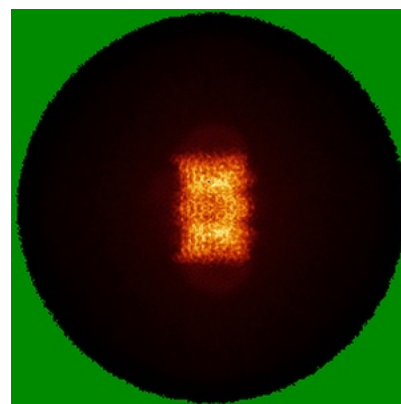
6.4.1 Primary map



X

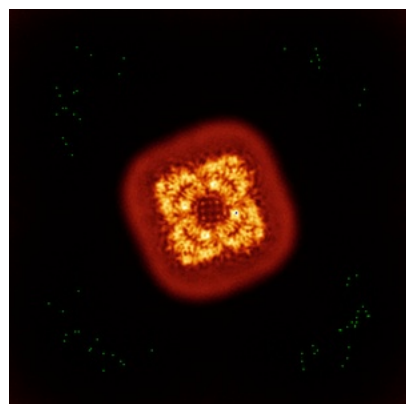


Y

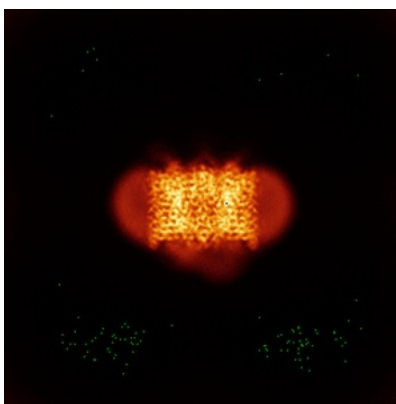


Z

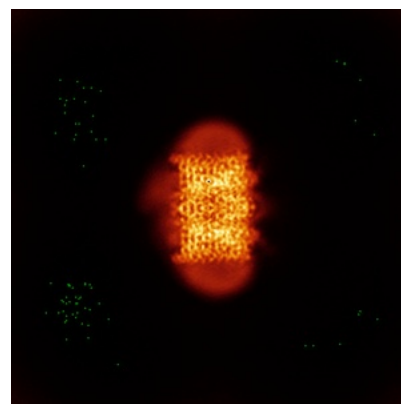
6.4.2 Raw map



X



Y

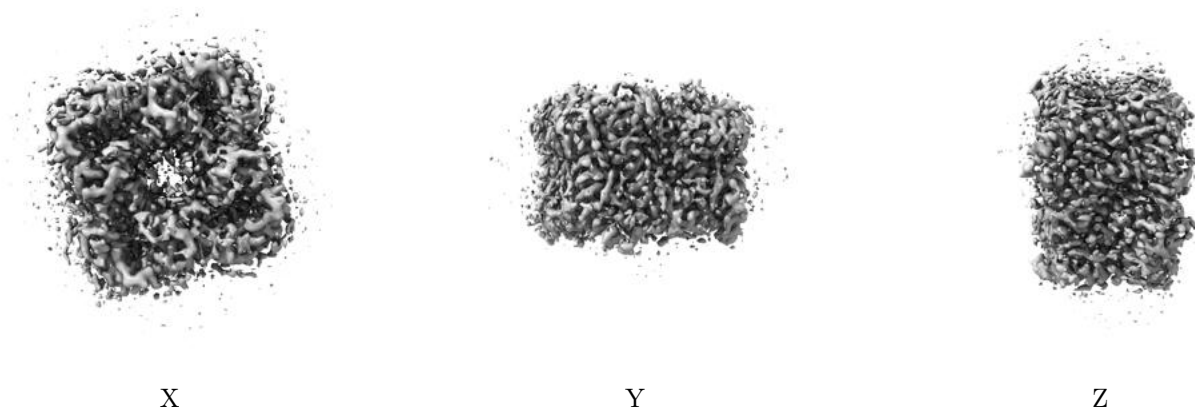


Z

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.

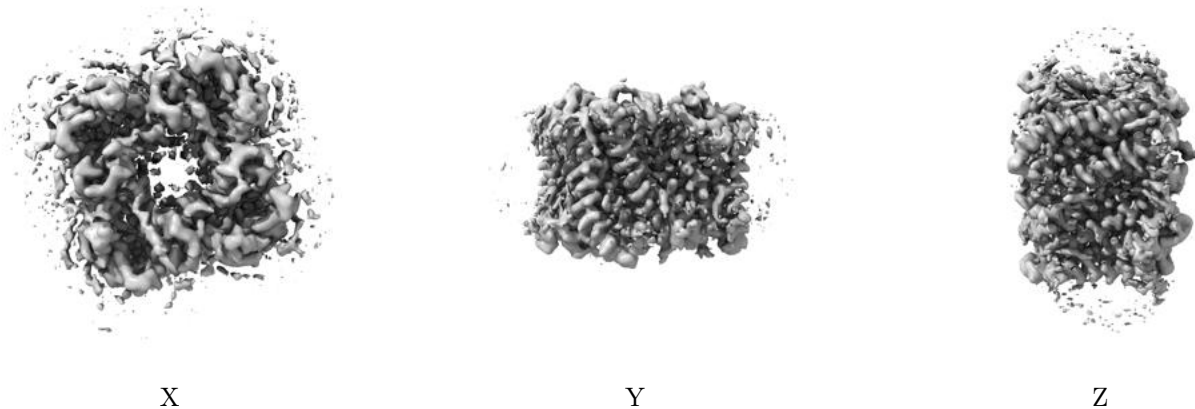
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.28. 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.

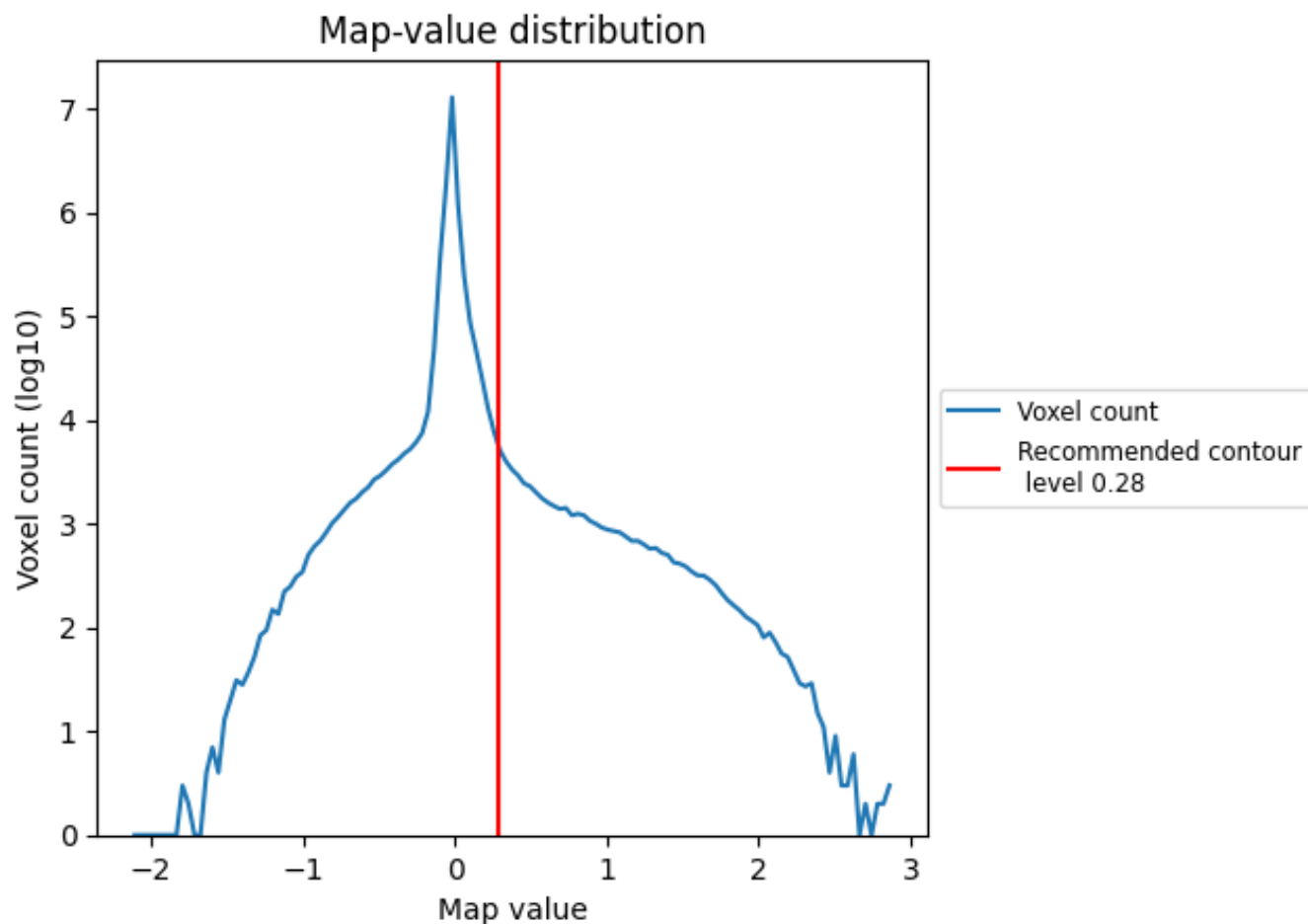
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

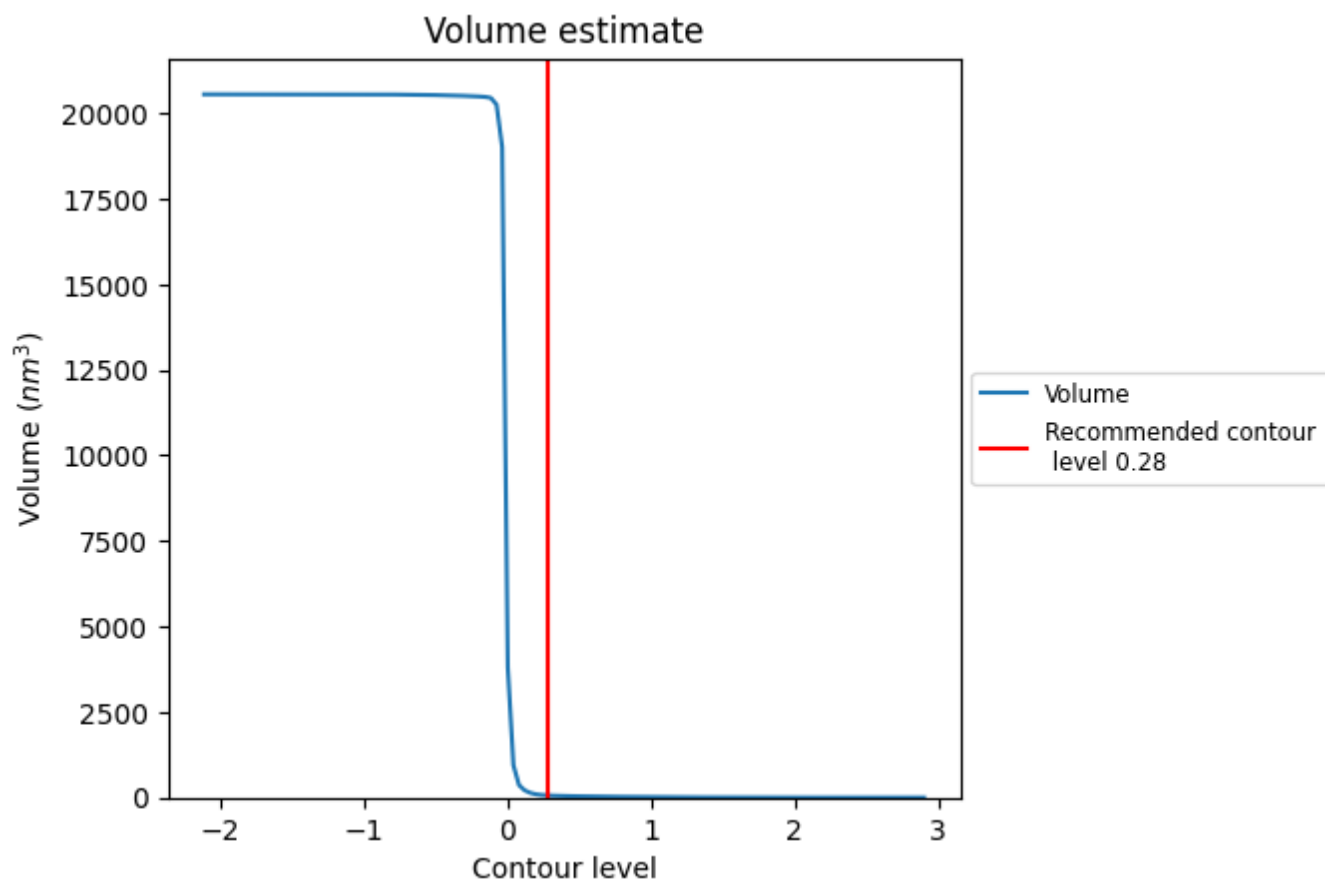
This section contains the results of statistical analysis of the map.

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.

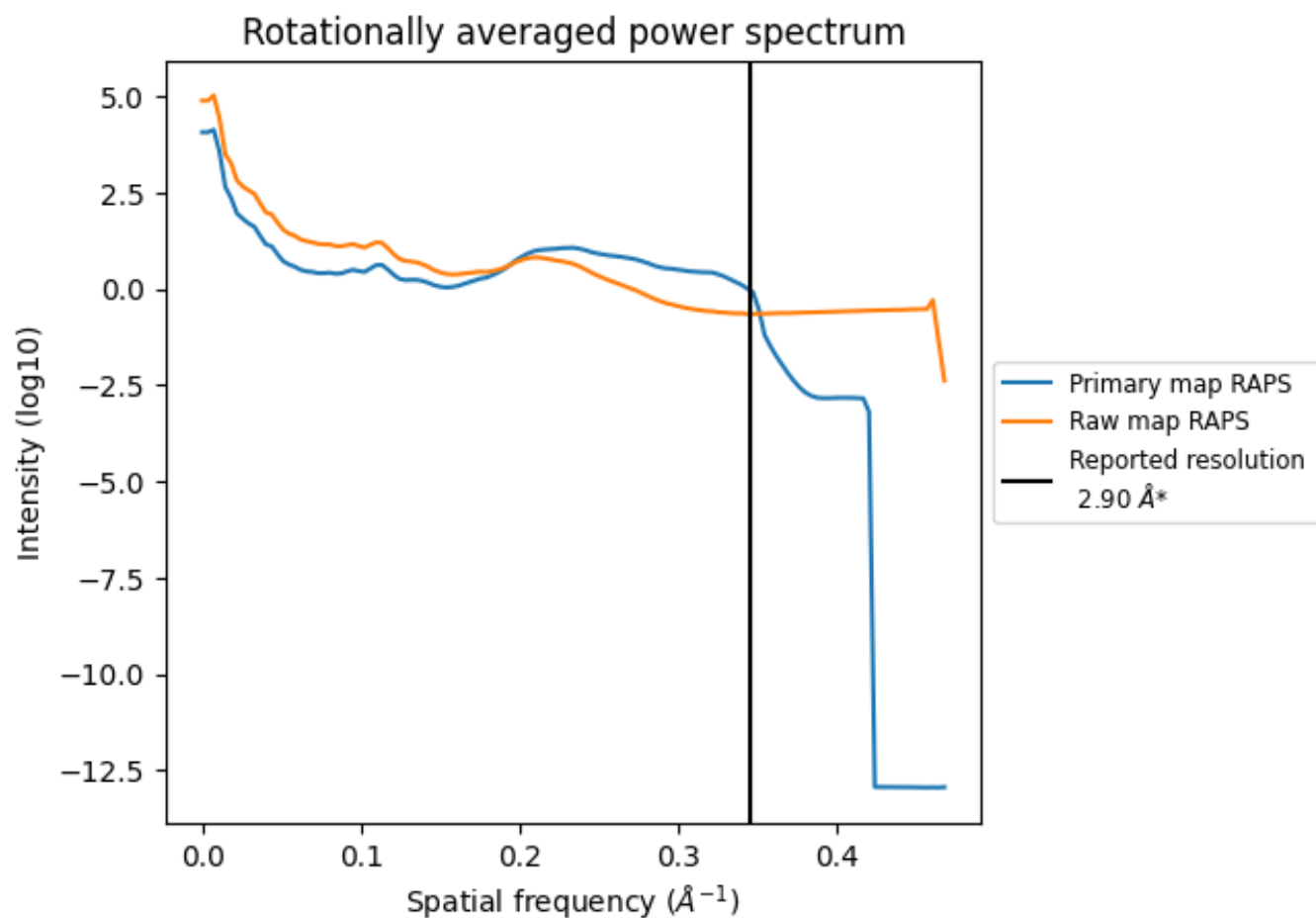
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 63 nm^3 ; this corresponds to an approximate mass of 57 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 ⓘ

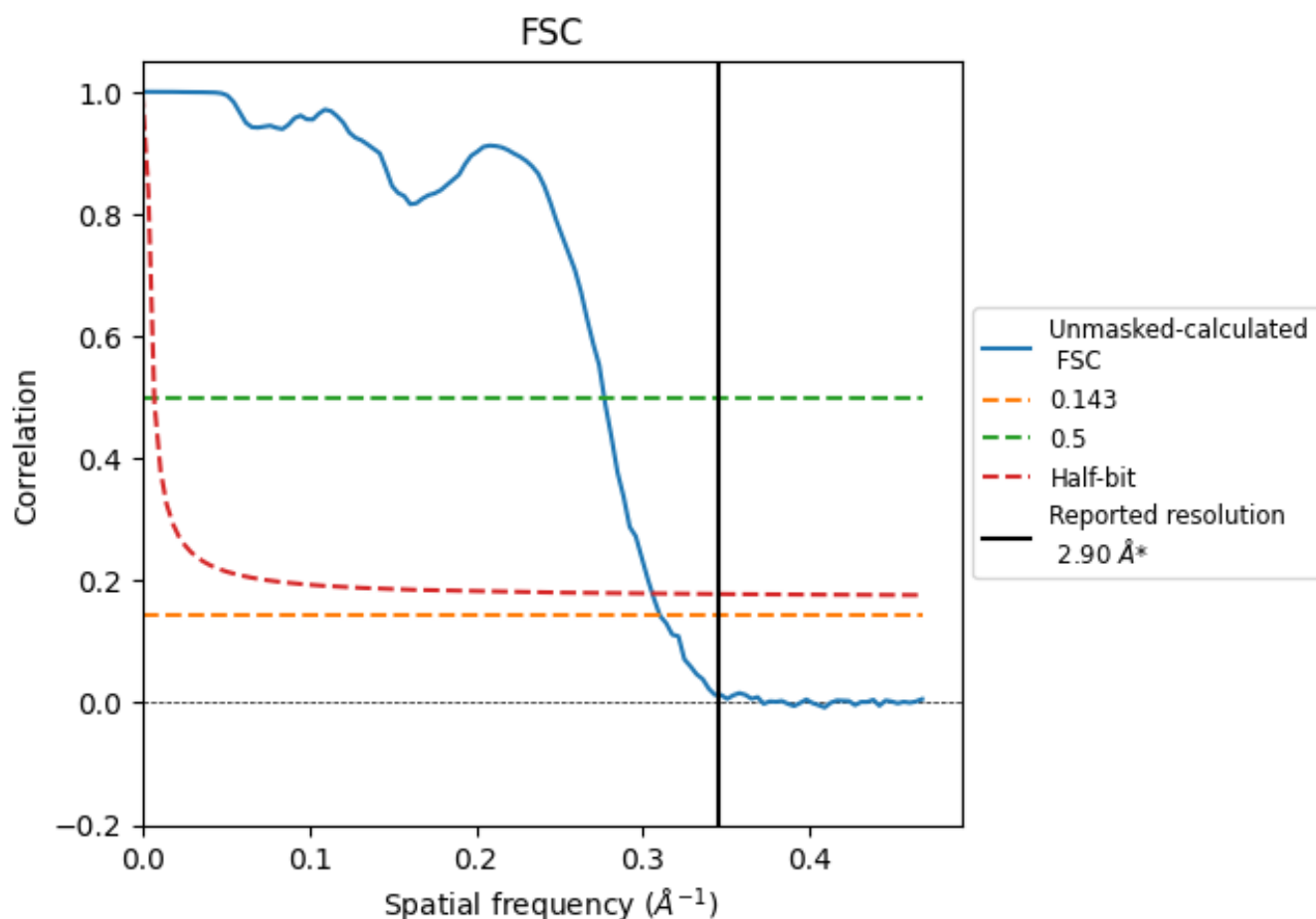


*Reported resolution corresponds to spatial frequency of 0.345 Å⁻¹

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 0.345 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

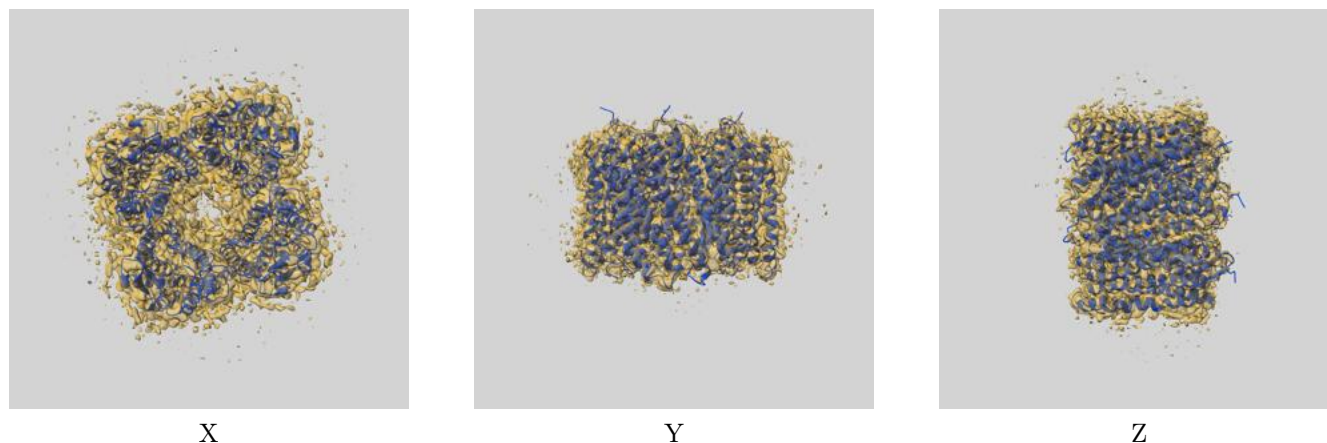
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.22	3.61	3.27

*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 3.22 differs from the reported value 2.9 by more than 10 %

9 Map-model fit [i](#)

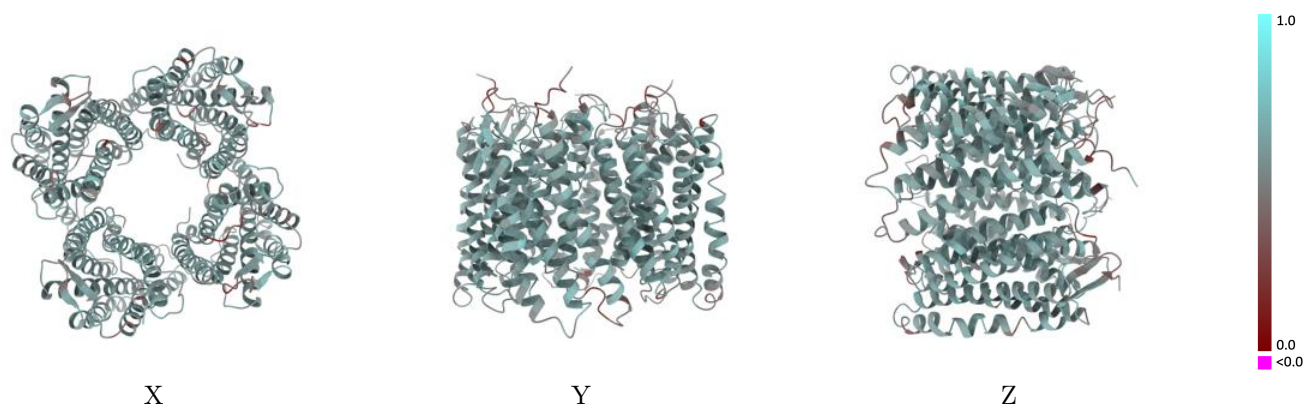
This section contains information regarding the fit between EMDB map EMD-48180 and PDB model 9MDU. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



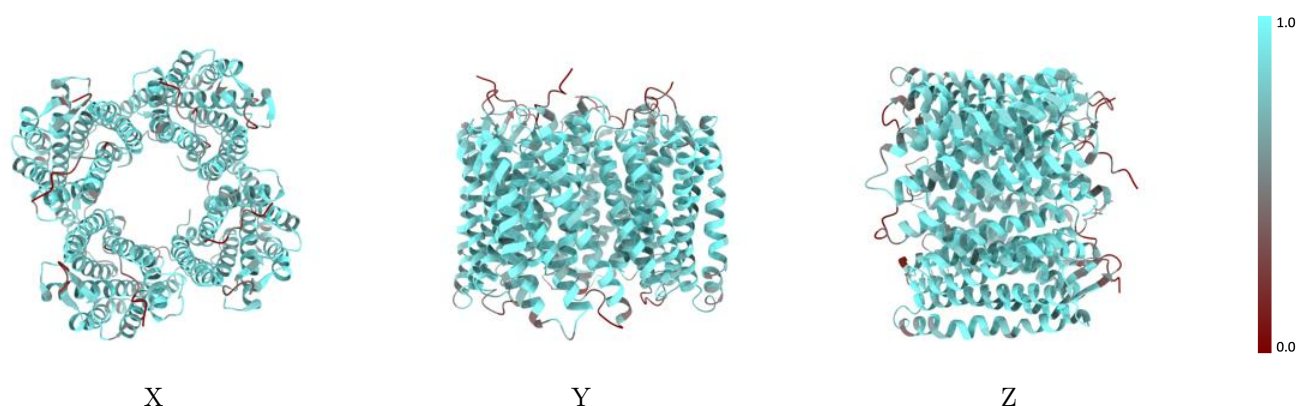
The images above show the 3D surface view of the map at the recommended contour level 0.28 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



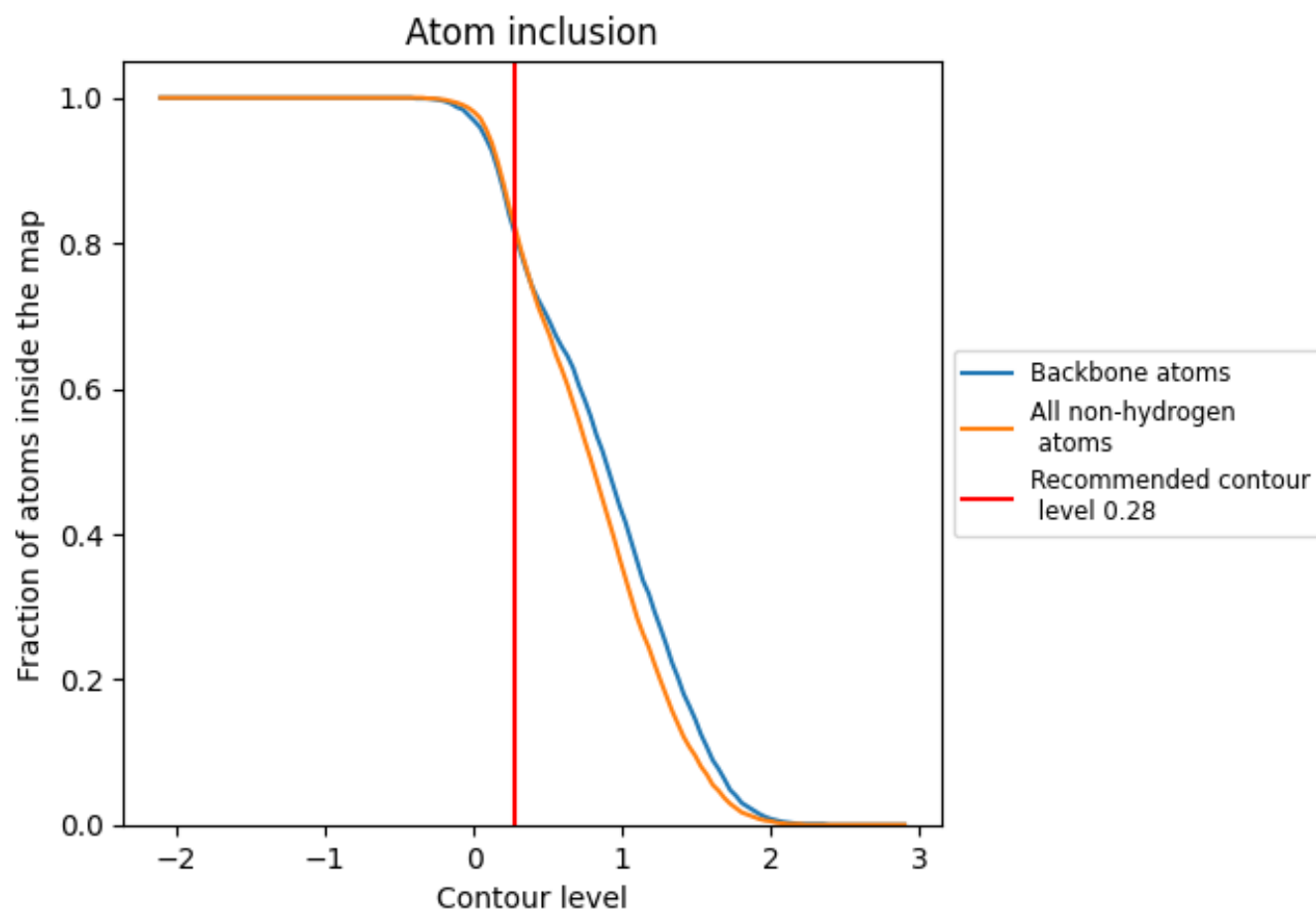
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.28).

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 82% 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 (0.28) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8210	0.5690
A	0.8210	0.5660
B	0.8150	0.5650
C	0.8190	0.5750
D	0.8280	0.5700

