



## Full wwPDB EM Validation Report ⓘ

Apr 8, 2026 – 06:12 PM UTC

PDB ID : 9ZN5 / pdb\_00009zn5  
EMDB ID : EMD-74435  
Title : Hybrid model of a dimer of BrxC-BrxB fusion complexed with PglZ from the Acinetobacter BREX system  
Authors : Doyle, L.A.; Stoddard, B.L.; Kaiser, B.; Kaiser, A.  
Deposited on : 2025-12-12  
Resolution : 6.94 Å(reported)  
Based on initial model : .

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.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

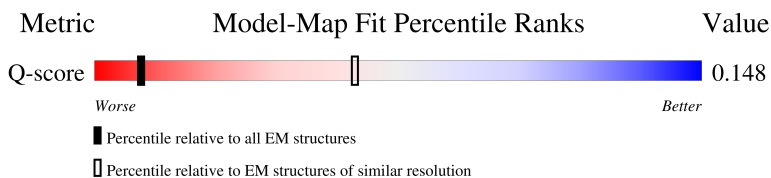
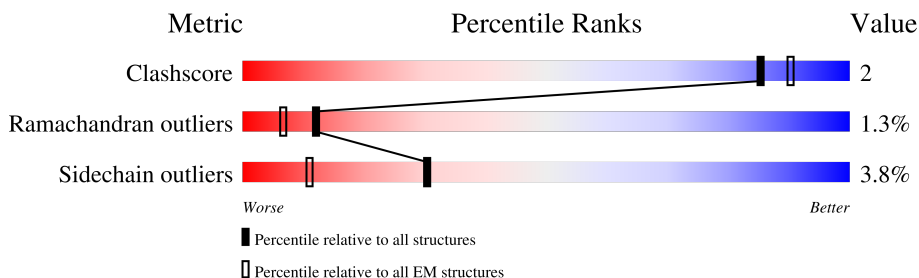
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 6.94 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	454 ( 6.45 - 7.41 )

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	754	
1	C	754	
2	B	879	
2	D	879	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 51990 atoms, of which 25836 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 BrxC-BrxB.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	754	Total	C	H	N	O	S	0	0
			11950	3837	5921	1031	1143	18		
1	C	754	Total	C	H	N	O	S	0	0
			11950	3837	5921	1031	1143	18		

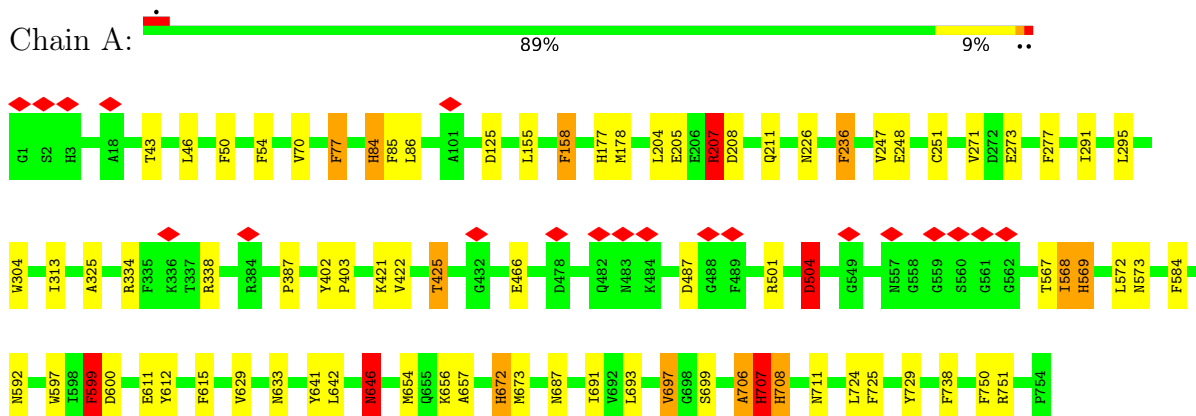
- Molecule 2 is a protein called PglZ.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	879	Total	C	H	N	O	S	0	0
			14045	4480	6997	1196	1350	22		
2	D	879	Total	C	H	N	O	S	0	0
			14045	4480	6997	1196	1350	22		

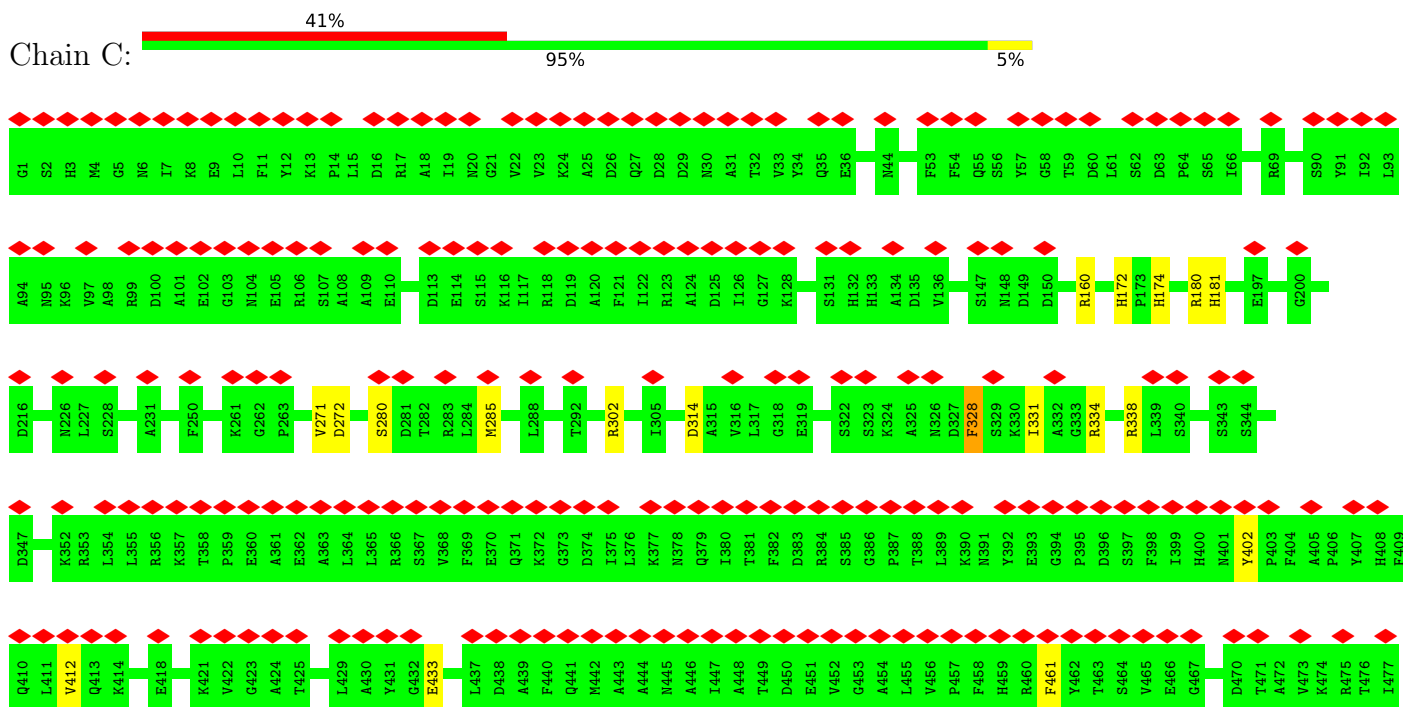
### 3 Residue-property plots

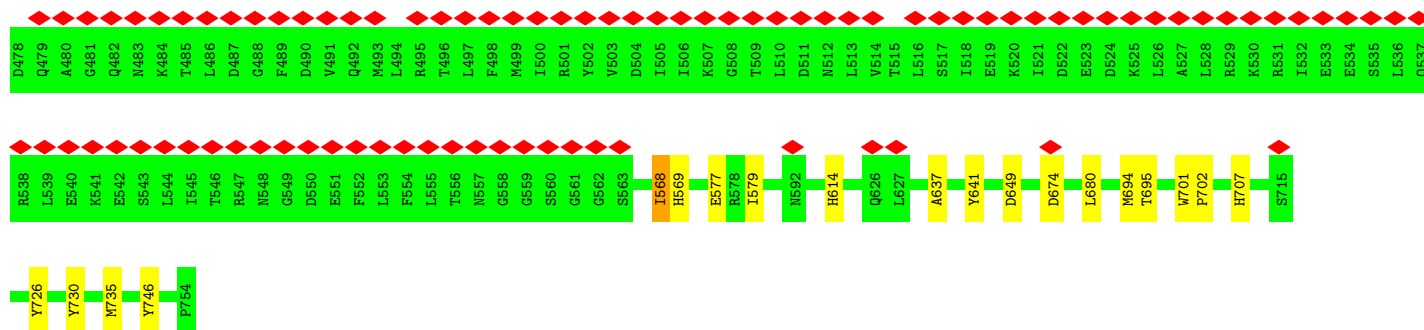
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: BrxC-BrxB

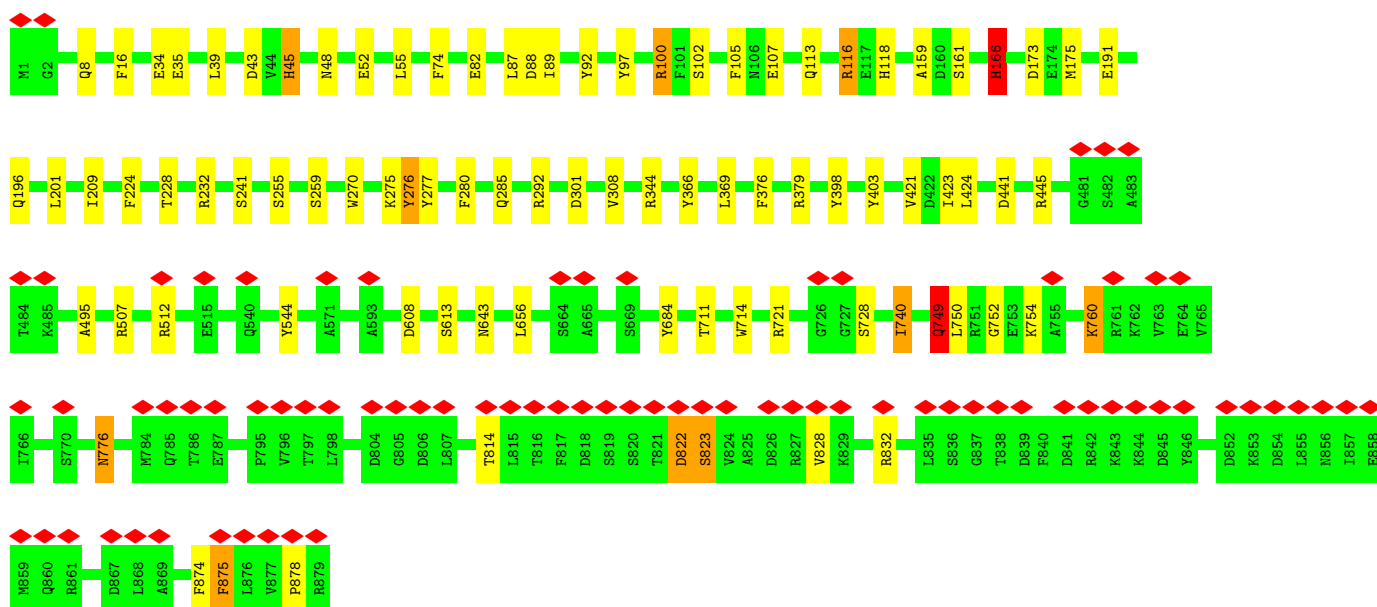
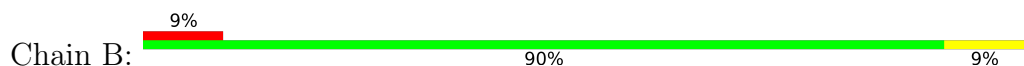


#### • Molecule 1: BrxC-BrxB

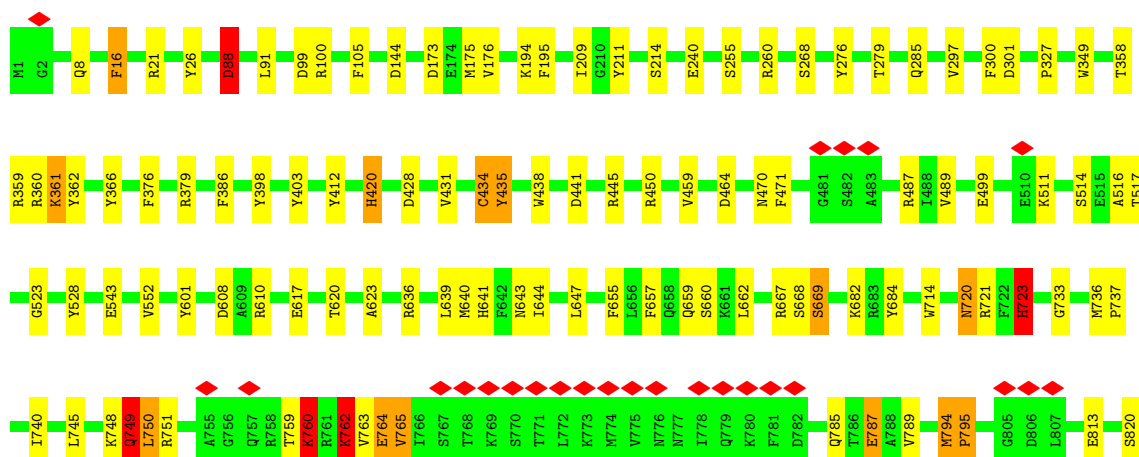
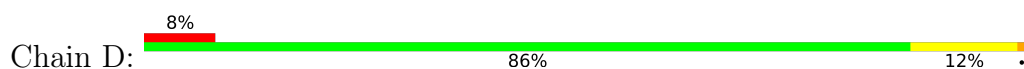




• Molecule 2: PglZ



• Molecule 2: PglZ



L835	S836	G837	T838	D839	F840	D841	R842	K843	K844	D845	Y846	F847	L848	V849	L850	K851	D852	K853	D854	L855	N856	I857	E858	M859	Q860	R861	Y862	K863	V864	T865	I866	D867	L868	A869	F870	T871	D872	D873	F874	F875	L876	V877	P878	R879
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## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	195504	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.183	Depositor
Minimum map value	-1.092	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.045	Depositor
Recommended contour level	0.3	Depositor
Map size ( $\text{\AA}$ )	448.8, 448.8, 448.8	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	2.244, 2.244, 2.244	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	0.78	0/6157	1.41	19/8310 (0.2%)
1	C	0.79	0/6157	1.31	8/8310 (0.1%)
2	B	1.55	3/7185 (0.0%)	1.43	34/9709 (0.4%)
2	D	0.88	6/7185 (0.1%)	1.42	39/9709 (0.4%)
All	All	1.07	9/26684 (0.0%)	1.40	100/36038 (0.3%)

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	C	0	6
2	B	0	10
2	D	0	8
All	All	0	29

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	823	SER	N-CA	112.95	2.85	1.46
2	D	875	PHE	CG-CD1	18.65	1.78	1.38
2	D	875	PHE	CG-CD2	18.33	1.77	1.38
2	D	875	PHE	CE1-CZ	13.05	1.77	1.38
2	D	875	PHE	CD2-CE2	12.86	1.77	1.38
2	D	875	PHE	CD1-CE1	12.71	1.76	1.38
2	D	875	PHE	CE2-CZ	12.69	1.76	1.38
2	B	823	SER	CA-CB	11.47	1.68	1.53
2	B	823	SER	CA-C	7.03	1.61	1.52

All (100) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	822	ASP	CA-C-N	16.89	145.24	122.77
2	B	822	ASP	C-N-CA	16.89	145.24	122.77
2	B	823	SER	N-CA-C	12.95	129.83	108.73
2	D	794	MET	N-CA-CB	-11.48	97.74	110.71
2	D	762	LYS	CA-C-N	10.80	141.14	121.70
2	D	762	LYS	C-N-CA	10.80	141.14	121.70
2	B	749	GLN	N-CA-C	9.77	125.04	112.34
2	D	762	LYS	N-CA-C	9.60	128.37	114.16
2	D	762	LYS	O-C-N	-9.56	112.00	121.85
2	B	823	SER	N-CA-CB	9.42	125.77	110.42
2	B	823	SER	CB-CA-C	-8.93	95.73	110.19
2	D	16	PHE	CA-CB-CG	8.72	122.52	113.80
2	D	760	LYS	N-CA-C	8.46	128.82	110.80
2	D	514	SER	N-CA-C	8.15	121.64	108.76
2	D	794	MET	CA-C-N	8.10	129.96	119.84
2	D	794	MET	C-N-CA	8.10	129.96	119.84
1	A	599	PHE	CA-CB-CG	-8.01	105.79	113.80
2	D	88	ASP	CA-CB-CG	7.93	120.53	112.60
2	B	874	PHE	CA-CB-CG	7.86	121.66	113.80
2	B	209	ILE	CB-CA-C	7.58	118.20	110.70
1	A	158	PHE	CA-CB-CG	7.03	120.83	113.80
2	B	776	ASN	CA-CB-CG	7.01	119.61	112.60
2	B	760	LYS	N-CA-C	6.96	125.63	110.80
2	D	874	PHE	CA-CB-CG	6.68	120.48	113.80
2	D	386	PHE	CA-CB-CG	6.54	120.34	113.80
1	A	84	HIS	CA-CB-CG	6.48	120.28	113.80
2	D	668	SER	CA-C-N	6.47	133.89	121.54
2	D	668	SER	C-N-CA	6.47	133.89	121.54
2	B	823	SER	CA-C-O	6.46	127.27	120.36
1	A	504	ASP	CA-CB-CG	6.41	119.01	112.60
2	B	48	ASN	N-CA-C	-6.38	98.78	108.67
1	A	751	ARG	NE-CZ-NH2	6.36	124.92	119.20
1	A	207	ARG	NE-CZ-NH2	6.25	124.83	119.20
2	D	723	HIS	CB-CG-CD2	-6.19	123.15	131.20
1	A	646	ASN	CA-CB-CG	-6.09	106.51	112.60
2	B	173	ASP	N-CA-C	6.08	118.69	111.33
1	A	50	PHE	CA-CB-CG	6.08	119.88	113.80
2	B	16	PHE	CA-CB-CG	6.07	119.86	113.80
1	A	573	ASN	CA-CB-CG	-5.97	106.63	112.60
2	B	721	ARG	NE-CZ-NH2	5.95	124.56	119.20
2	D	794	MET	N-CA-C	5.92	117.29	108.76
2	D	749	GLN	N-CA-C	5.86	123.28	110.80
2	D	748	LYS	CA-C-N	5.83	132.67	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	748	LYS	C-N-CA	5.83	132.67	121.54
1	A	672	HIS	CB-CG-CD2	-5.74	123.74	131.20
1	C	180	ARG	NE-CZ-NH2	5.70	124.33	119.20
1	A	177	HIS	CB-CG-CD2	-5.68	123.82	131.20
2	D	435	TYR	CA-C-N	5.66	128.13	120.38
2	D	435	TYR	C-N-CA	5.66	128.13	120.38
1	A	569	HIS	CA-CB-CG	5.64	119.44	113.80
2	D	450	ARG	NE-CZ-NH2	5.63	124.27	119.20
2	B	750	LEU	N-CA-C	5.62	119.64	112.34
2	B	166	HIS	CB-CG-CD2	-5.59	123.93	131.20
1	A	84	HIS	CB-CG-CD2	-5.56	123.97	131.20
2	B	344	ARG	NE-CZ-NH2	5.52	124.17	119.20
2	B	100	ARG	NE-CZ-NH2	5.48	124.13	119.20
2	B	507	ARG	NE-CZ-NH2	5.47	124.12	119.20
2	B	159	ALA	N-CA-C	5.46	117.74	110.53
2	D	379	ARG	NE-CZ-NH2	5.45	124.11	119.20
1	A	208	ASP	CA-CB-CG	5.39	117.99	112.60
2	D	543	GLU	N-CA-C	5.39	115.65	108.38
2	D	105	PHE	CA-CB-CG	5.37	119.17	113.80
1	A	592	ASN	CA-CB-CG	5.37	117.97	112.60
2	B	292	ARG	NE-CZ-NH2	5.37	124.03	119.20
2	D	723	HIS	CA-CB-CG	5.36	119.16	113.80
1	C	328	PHE	CA-CB-CG	5.35	119.15	113.80
2	B	423	ILE	CA-C-N	5.34	131.74	121.54
2	B	423	ILE	C-N-CA	5.34	131.74	121.54
1	C	334	ARG	NE-CZ-NH2	5.29	123.97	119.20
1	A	277	PHE	CA-CB-CG	-5.22	108.58	113.80
2	B	34	GLU	CA-C-N	5.22	127.27	120.28
2	B	34	GLU	C-N-CA	5.22	127.27	120.28
2	D	785	GLN	CA-C-N	5.21	128.64	120.82
2	D	785	GLN	C-N-CA	5.21	128.64	120.82
2	B	740	ILE	N-CA-C	5.21	118.20	113.20
2	D	523	GLY	CA-C-N	5.19	129.31	122.51
2	D	523	GLY	C-N-CA	5.19	129.31	122.51
2	B	832	ARG	NE-CZ-NH2	5.18	123.87	119.20
1	A	707	HIS	CB-CG-CD2	-5.18	124.46	131.20
1	C	302	ARG	NE-CZ-NH2	5.18	123.86	119.20
2	D	514	SER	CA-C-N	5.17	130.71	122.59
2	D	514	SER	C-N-CA	5.17	130.71	122.59
1	C	172	HIS	CB-CG-CD2	-5.15	124.50	131.20
2	D	360	ARG	NE-CZ-NH2	5.15	123.84	119.20
2	B	255	SER	N-CA-C	5.15	119.64	112.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	795	PRO	N-CA-C	5.14	123.06	112.47
1	A	77	PHE	N-CA-C	5.12	121.72	110.80
2	D	260	ARG	NE-CZ-NH2	5.12	123.81	119.20
1	A	633	ASN	CA-CB-CG	-5.12	107.48	112.60
2	D	428	ASP	CA-CB-CG	5.11	117.71	112.60
2	B	118	HIS	CB-CG-CD2	-5.09	124.58	131.20
2	D	144	ASP	CA-CB-CG	5.07	117.67	112.60
2	B	512	ARG	NE-CZ-NH2	5.03	123.73	119.20
2	B	776	ASN	OD1-CG-ND2	-5.02	117.58	122.60
2	B	379	ARG	NE-CZ-NH2	5.00	123.70	119.20
1	C	614	HIS	CB-CG-CD2	-5.00	124.69	131.20
2	D	842	ARG	NE-CZ-NH2	5.00	123.70	119.20
2	B	684	TYR	N-CA-C	5.00	115.91	108.60
1	C	272	ASP	CA-C-N	5.00	129.47	122.36
1	C	272	ASP	C-N-CA	5.00	129.47	122.36

There are no chirality outliers.

All (29) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	207	ARG	Sidechain
1	A	501	ARG	Sidechain
1	A	599	PHE	Sidechain
1	A	641	TYR	Sidechain
1	A	729	TYR	Sidechain
2	B	100	ARG	Sidechain
2	B	116	ARG	Sidechain
2	B	166	HIS	Sidechain
2	B	232	ARG	Sidechain
2	B	276	TYR	Sidechain
2	B	277	TYR	Sidechain
2	B	45	HIS	Sidechain
2	B	544	TYR	Sidechain
2	B	875	PHE	Peptide
2	B	97	TYR	Sidechain
1	C	160	ARG	Sidechain
1	C	402	TYR	Sidechain
1	C	641	TYR	Sidechain
1	C	726	TYR	Sidechain
1	C	730	TYR	Sidechain
1	C	746	TYR	Sidechain
2	D	21	ARG	Sidechain

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Mol	Chain	Res	Type	Group
2	D	211	TYR	Sidechain
2	D	26	TYR	Sidechain
2	D	276	TYR	Sidechain
2	D	366	TYR	Sidechain
2	D	601	TYR	Sidechain
2	D	636	ARG	Sidechain
2	D	846	TYR	Sidechain

## 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6029	5921	5921	28	0
1	C	6029	5921	5921	8	0
2	B	7048	6997	6997	44	0
2	D	7048	6997	6997	62	0
All	All	26154	25836	25836	115	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 2.

All (115) 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:D:875:PHE:CE2	2:D:875:PHE:CZ	1.76	1.69
2:D:875:PHE:CG	2:D:875:PHE:CD1	1.78	1.66
2:D:875:PHE:CD1	2:D:875:PHE:CE1	1.76	1.64
2:D:875:PHE:CE2	2:D:875:PHE:CD2	1.77	1.62
2:D:875:PHE:CG	2:D:875:PHE:CD2	1.77	1.62
2:D:875:PHE:CZ	2:D:875:PHE:CE1	1.77	1.61
2:B:823:SER:HA	2:D:875:PHE:CZ	1.48	1.45
2:B:823:SER:HA	2:D:875:PHE:CE1	1.53	1.42
2:B:823:SER:CA	2:D:875:PHE:CE1	2.26	1.17
2:B:823:SER:CA	2:D:875:PHE:CZ	2.26	1.16
2:B:823:SER:CA	2:D:875:PHE:CD1	2.29	1.16
2:B:823:SER:CA	2:D:875:PHE:CE2	2.28	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:823:SER:CA	2:D:875:PHE:CD2	2.30	1.14
2:B:823:SER:CA	2:D:875:PHE:CG	2.31	1.14
2:B:823:SER:N	2:D:875:PHE:CD1	2.24	1.05
2:B:823:SER:N	2:D:875:PHE:CD2	2.24	1.05
2:B:823:SER:N	2:D:875:PHE:CG	2.24	1.04
2:B:823:SER:N	2:D:875:PHE:CE2	2.26	1.03
2:B:823:SER:N	2:D:875:PHE:CE1	2.28	1.02
2:B:823:SER:N	2:D:875:PHE:CZ	2.28	1.00
2:B:823:SER:CB	2:D:875:PHE:CD1	2.55	0.90
2:B:823:SER:C	2:D:875:PHE:CE2	2.57	0.81
2:B:823:SER:CB	2:D:875:PHE:CG	2.69	0.76
2:B:823:SER:HA	2:D:875:PHE:CE2	2.22	0.75
2:B:823:SER:HB2	2:D:875:PHE:CD1	2.24	0.70
2:B:823:SER:HA	2:D:875:PHE:CD1	2.29	0.68
2:D:760:LYS:HA	2:D:794:MET:HA	1.76	0.67
2:B:823:SER:HB3	2:D:875:PHE:CG	2.30	0.67
2:B:823:SER:C	2:D:875:PHE:CD2	2.77	0.63
1:A:211:GLN:H	1:A:211:GLN:CD	2.06	0.63
2:B:822:ASP:C	2:D:875:PHE:CZ	2.79	0.60
2:D:639:LEU:HB3	2:D:750:LEU:HA	1.86	0.58
2:B:822:ASP:C	2:D:875:PHE:CE1	2.82	0.57
1:C:328:PHE:HA	1:C:331:ILE:HD12	1.86	0.56
1:A:697:VAL:HG11	1:A:725:PHE:O	2.07	0.54
2:B:749:GLN:HA	2:B:752:GLY:C	2.37	0.50
2:D:516:ALA:HB1	2:D:745:LEU:HD11	1.94	0.50
2:D:763:VAL:HG23	2:D:794:MET:C	2.36	0.50
2:D:194:LYS:HE3	2:D:195:PHE:CZ	2.48	0.48
1:A:707:HIS:CE1	1:A:708:HIS:CE1	3.02	0.48
2:D:88:ASP:HA	2:D:91:LEU:HD12	1.95	0.48
1:C:701:TRP:CD2	1:C:702:PRO:HA	2.49	0.48
2:B:8:GLN:OE1	2:B:39:LEU:HD11	2.13	0.48
1:C:412:VAL:HG21	1:C:461:PHE:CD1	2.47	0.48
2:B:270:TRP:CZ2	2:B:280:PHE:HA	2.50	0.47
2:D:327:PRO:HA	2:D:376:PHE:CZ	2.50	0.47
1:A:597:TRP:CE3	1:A:750:PHE:HB3	2.50	0.46
2:B:161:SER:HB3	2:B:166:HIS:CD2	2.50	0.46
2:B:89:ILE:HA	2:B:92:TYR:CE2	2.51	0.46
1:A:572:LEU:HD22	1:A:612:TYR:CG	2.51	0.46
2:D:762:LYS:H	2:D:794:MET:HG2	1.81	0.46
2:B:275:LYS:HE2	2:B:276:TYR:CE1	2.51	0.46
1:C:637:ALA:HB1	1:C:680:LEU:CD2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:412:TYR:CZ	2:D:431:VAL:HG21	2.51	0.45
1:A:54:PHE:CD1	1:A:125:ASP:HB3	2.52	0.45
2:B:823:SER:O	2:D:875:PHE:CD2	2.69	0.45
2:B:102:SER:HA	2:B:116:ARG:HD3	1.99	0.45
2:D:349:TRP:HA	2:D:359:ARG:HE	1.81	0.45
2:D:489:VAL:HB	2:D:647:LEU:HD13	1.99	0.45
1:A:697:VAL:HG22	1:A:724:LEU:HD11	1.98	0.45
1:A:248:GLU:H	1:A:248:GLU:CD	2.24	0.44
1:A:642:LEU:O	1:A:646:ASN:HA	2.17	0.44
1:A:568:ILE:HG23	1:A:569:HIS:H	1.82	0.44
1:C:707:HIS:CE1	1:C:735:MET:SD	3.10	0.44
2:B:822:ASP:O	2:D:875:PHE:CZ	2.71	0.44
2:D:763:VAL:HG13	2:D:765:VAL:HB	2.00	0.44
2:B:398:TYR:CE2	2:B:403:TYR:HA	2.53	0.43
2:D:720:ASN:C	2:D:721:ARG:HG3	2.43	0.43
2:D:657:PHE:CE1	2:D:659:GLN:HA	2.54	0.43
2:B:35:GLU:H	2:B:35:GLU:CD	2.27	0.43
2:D:750:LEU:HD22	2:D:750:LEU:N	2.34	0.43
1:A:236:PHE:CD1	1:A:236:PHE:C	2.97	0.43
1:A:672:HIS:CG	1:A:673:MET:N	2.87	0.43
1:A:725:PHE:CD1	1:A:725:PHE:N	2.86	0.43
1:C:568:ILE:HG23	1:C:569:HIS:H	1.84	0.43
1:A:70:VAL:HA	1:A:304:TRP:CE3	2.54	0.43
1:A:155:LEU:HA	1:A:158:PHE:CD2	2.54	0.43
1:A:85:PHE:CD1	1:A:85:PHE:C	2.95	0.42
1:A:706:ALA:O	1:A:707:HIS:C	2.62	0.42
2:D:620:THR:O	2:D:623:ALA:HB3	2.18	0.42
1:A:251:CYS:SG	1:A:295:LEU:HD21	2.59	0.42
2:B:270:TRP:CH2	2:B:280:PHE:HA	2.54	0.42
2:B:105:PHE:CZ	2:B:113:GLN:HA	2.54	0.42
2:B:270:TRP:CE3	2:B:280:PHE:CD2	3.08	0.42
2:B:366:TYR:CD2	2:B:369:LEU:HD12	2.55	0.42
1:A:247:VAL:HG13	1:A:291:ILE:HG23	2.01	0.42
1:C:285:MET:HG3	1:C:331:ILE:HD13	2.02	0.42
2:D:759:THR:O	2:D:794:MET:HG3	2.20	0.42
1:A:313:ILE:HG23	1:A:338:ARG:NH2	2.35	0.42
1:A:402:TYR:CD1	1:A:403:PRO:HA	2.55	0.41
2:D:398:TYR:CZ	2:D:403:TYR:HA	2.56	0.41
2:D:434:CYS:HA	2:D:438:TRP:HB3	2.03	0.41
1:C:314:ASP:CG	1:C:338:ARG:HH22	2.29	0.41
2:D:640:MET:HE1	2:D:787:GLU:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:412:TYR:CE1	2:D:431:VAL:HG21	2.55	0.41
1:A:629:VAL:HG22	1:A:691:ILE:HD13	2.02	0.41
2:D:610:ARG:HD2	2:D:623:ALA:HB2	2.03	0.41
2:D:669:SER:HA	2:D:723:HIS:CD2	2.55	0.41
2:D:684:TYR:CD1	2:D:733:GLY:HA3	2.55	0.41
1:A:43:THR:H	1:A:46:LEU:HB2	1.86	0.41
1:A:673:MET:N	1:A:673:MET:HE2	2.36	0.41
2:D:736:MET:HB3	2:D:737:PRO:HD2	2.03	0.41
1:A:656:LYS:HE2	2:B:52:GLU:HA	2.02	0.41
2:B:241:SER:HB3	2:B:308:VAL:HA	2.03	0.41
2:D:487:ARG:HB2	2:D:644:ILE:HG23	2.02	0.41
1:A:85:PHE:CG	1:A:86:LEU:N	2.88	0.41
2:D:297:VAL:HA	2:D:300:PHE:CD2	2.56	0.41
2:D:641:HIS:HA	2:D:820:SER:HB2	2.02	0.41
1:A:654:MET:HA	1:A:657:ALA:HB3	2.02	0.41
2:B:45:HIS:CD2	2:B:45:HIS:N	2.88	0.40
2:B:403:TYR:CD1	2:B:403:TYR:C	3.00	0.40
2:D:662:LEU:C	2:D:667:ARG:HE	2.29	0.40
1:A:599:PHE:CG	1:A:600:ASP:N	2.90	0.40
2:D:361:LYS:HE2	2:D:362:TYR:CE2	2.56	0.40
2:B:823:SER:CA	2:B:823:SER:N	2.85	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	752/754 (100%)	665 (88%)	71 (9%)	16 (2%)	5	30
1	C	752/754 (100%)	701 (93%)	48 (6%)	3 (0%)	30	67
2	B	877/879 (100%)	782 (89%)	85 (10%)	10 (1%)	11	46
2	D	877/879 (100%)	777 (89%)	87 (10%)	13 (2%)	8	40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	3258/3266 (100%)	2925 (90%)	291 (9%)	42 (1%)	12	42

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	77	PHE
1	A	568	ILE
1	A	646	ASN
1	A	697	VAL
1	A	706	ALA
1	A	707	HIS
2	B	87	LEU
2	B	445	ARG
2	D	660	SER
2	D	669	SER
2	D	749	GLN
2	D	760	LYS
2	D	795	PRO
1	A	425	THR
1	A	699	SER
2	D	682	LYS
1	A	273	GLU
1	A	325	ALA
1	A	466	GLU
2	B	43	ASP
2	B	760	LYS
2	D	255	SER
2	D	435	TYR
2	D	789	VAL
1	A	421	LYS
1	A	504	ASP
2	B	259	SER
2	B	613	SER
2	D	420	HIS
2	D	445	ARG
2	D	751	ARG
2	D	764	GLU
2	B	495	ALA
2	B	643	ASN
1	C	433	GLU
1	A	207	ARG
1	A	567	THR

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Mol	Chain	Res	Type
1	C	280	SER
2	B	878	PRO
1	A	387	PRO
2	B	828	VAL
1	C	568	ILE

### 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	648/648 (100%)	627 (97%)	21 (3%)	34	56
1	C	648/648 (100%)	639 (99%)	9 (1%)	59	72
2	B	760/760 (100%)	732 (96%)	28 (4%)	30	51
2	D	760/760 (100%)	712 (94%)	48 (6%)	16	37
All	All	2816/2816 (100%)	2710 (96%)	106 (4%)	30	50

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	HIS
1	A	178	MET
1	A	204	LEU
1	A	205	GLU
1	A	226	ASN
1	A	236	PHE
1	A	271	VAL
1	A	334	ARG
1	A	422	VAL
1	A	425	THR
1	A	487	ASP
1	A	504	ASP
1	A	584	PHE
1	A	611	GLU
1	A	615	PHE
1	A	687	ASN

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Mol	Chain	Res	Type
1	A	693	LEU
1	A	707	HIS
1	A	708	HIS
1	A	711	ASN
1	A	738	PHE
2	B	55	LEU
2	B	74	PHE
2	B	82	GLU
2	B	88	ASP
2	B	107	GLU
2	B	175	MET
2	B	191	GLU
2	B	196	GLN
2	B	201	LEU
2	B	224	PHE
2	B	228	THR
2	B	285	GLN
2	B	301	ASP
2	B	376	PHE
2	B	421	VAL
2	B	424	LEU
2	B	441	ASP
2	B	608	ASP
2	B	656	LEU
2	B	711	THR
2	B	714	TRP
2	B	728	SER
2	B	740	ILE
2	B	749	GLN
2	B	754	LYS
2	B	776	ASN
2	B	814	THR
2	B	875	PHE
1	C	174	HIS
1	C	181	HIS
1	C	271	VAL
1	C	577	GLU
1	C	579	ILE
1	C	649	ASP
1	C	674	ASP
1	C	694	MET
1	C	695	THR

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Mol	Chain	Res	Type
2	D	8	GLN
2	D	16	PHE
2	D	88	ASP
2	D	99	ASP
2	D	100	ARG
2	D	173	ASP
2	D	175	MET
2	D	176	VAL
2	D	209	ILE
2	D	214	SER
2	D	240	GLU
2	D	268	SER
2	D	279	THR
2	D	285	GLN
2	D	301	ASP
2	D	358	THR
2	D	361	LYS
2	D	420	HIS
2	D	434	CYS
2	D	441	ASP
2	D	459	VAL
2	D	464	ASP
2	D	470	ASN
2	D	471	PHE
2	D	499	GLU
2	D	511	LYS
2	D	517	THR
2	D	528	TYR
2	D	552	VAL
2	D	608	ASP
2	D	617	GLU
2	D	643	ASN
2	D	655	PHE
2	D	714	TRP
2	D	720	ASN
2	D	723	HIS
2	D	740	ILE
2	D	749	GLN
2	D	750	LEU
2	D	762	LYS
2	D	764	GLU
2	D	765	VAL

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Mol	Chain	Res	Type
2	D	787	GLU
2	D	813	GLU
2	D	833	LEU
2	D	856	ASN
2	D	872	ASP
2	D	875	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	35	GLN
1	A	163	ASN
1	A	371	GLN
1	A	401	ASN
1	A	557	ASN
1	A	588	GLN
1	A	631	ASN
1	A	689	GLN
2	B	178	GLN
2	B	250	GLN
2	B	325	GLN
2	B	475	HIS
2	B	557	GLN
2	B	597	HIS
2	B	653	HIS
2	B	702	GLN
2	B	856	ASN
1	C	184	GLN
1	C	492	GLN
1	C	557	ASN
1	C	574	GLN
1	C	636	GLN
1	C	646	ASN
1	C	707	HIS
2	D	6	GLN
2	D	106	ASN
2	D	112	GLN
2	D	316	GLN
2	D	643	ASN
2	D	757	GLN
2	D	776	ASN

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Mol	Chain	Res	Type
2	D	777	ASN

### 5.3.3 RNA [i](#)

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](#)

There are no chain breaks in this entry.

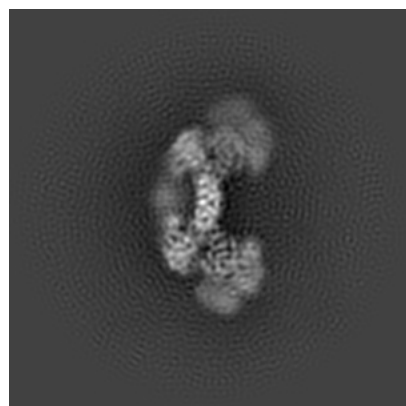
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-74435. These allow visual inspection of the internal detail of the map and identification of artifacts.

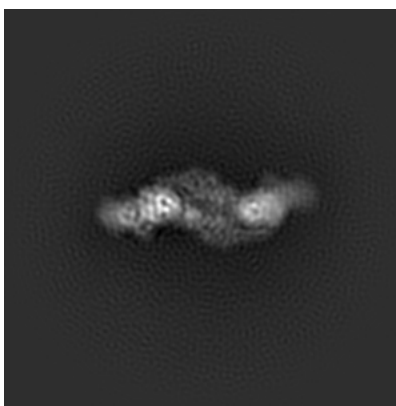
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

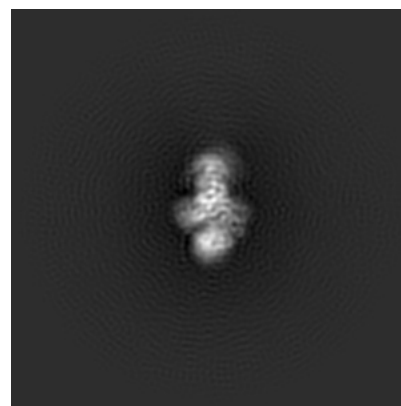
#### 6.1.1 Primary map



X

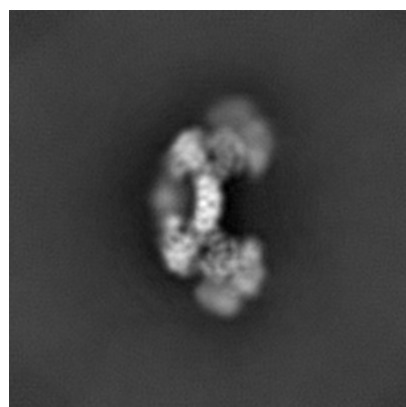


Y

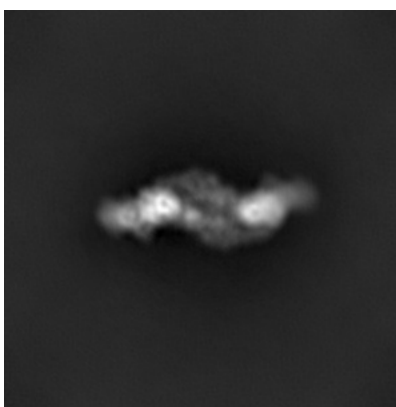


Z

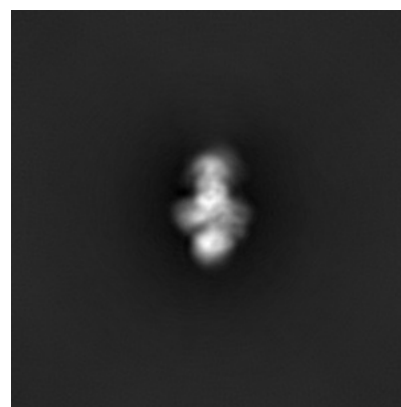
#### 6.1.2 Raw map



X



Y

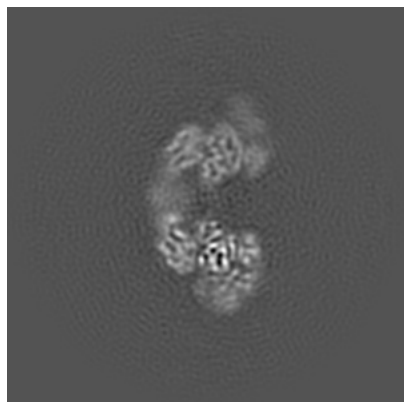


Z

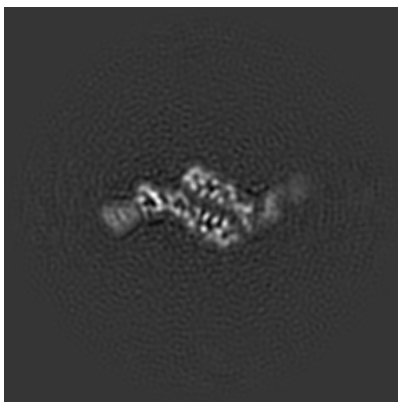
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

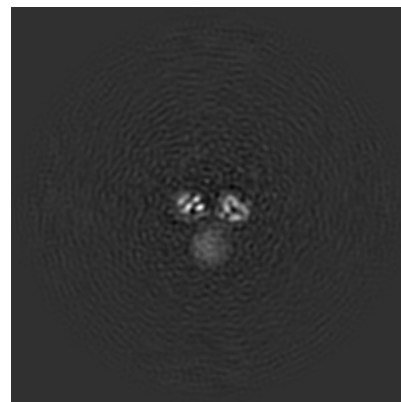
### 6.2.1 Primary map



X Index: 100

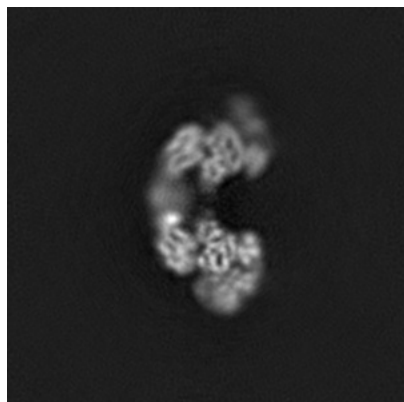


Y Index: 100

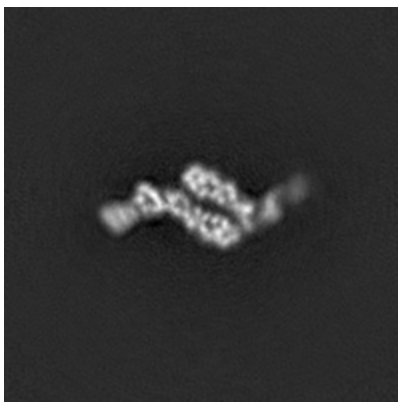


Z Index: 100

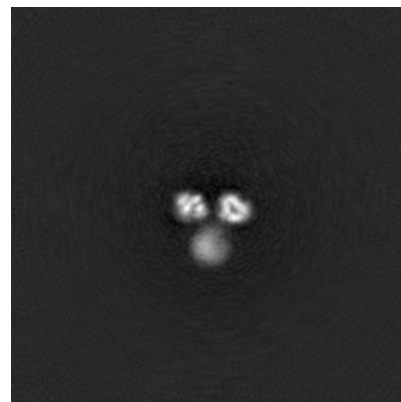
### 6.2.2 Raw map



X Index: 100



Y Index: 100

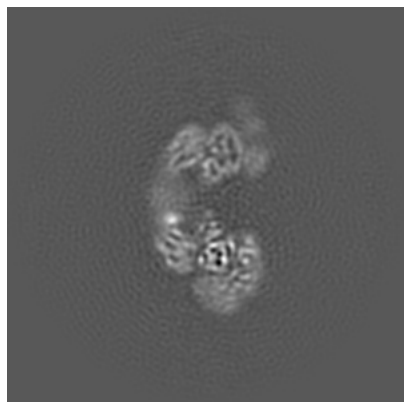


Z Index: 100

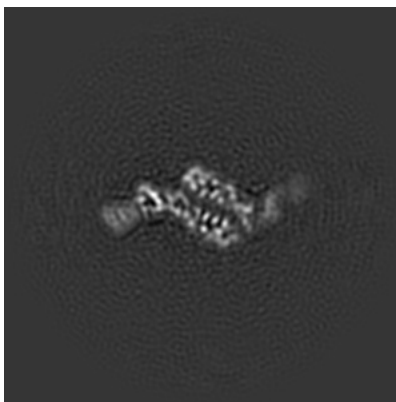
The images above show central slices of the map in three orthogonal directions.

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

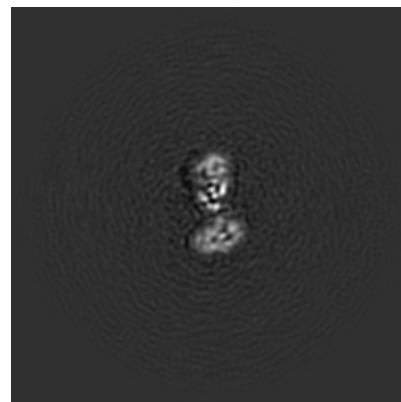
### 6.3.1 Primary map



X Index: 99

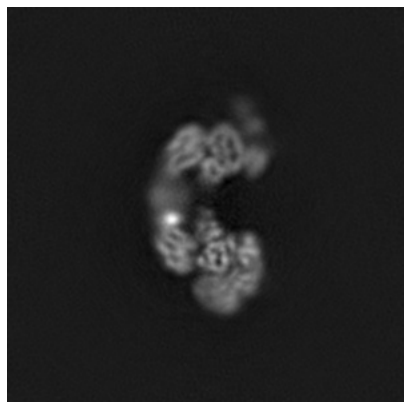


Y Index: 100

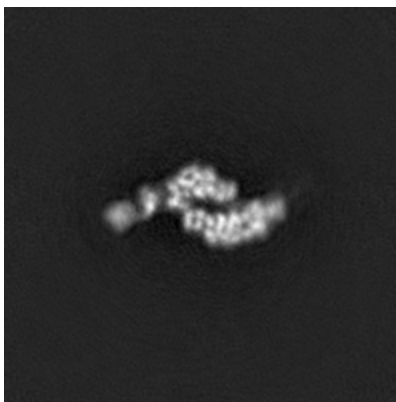


Z Index: 77

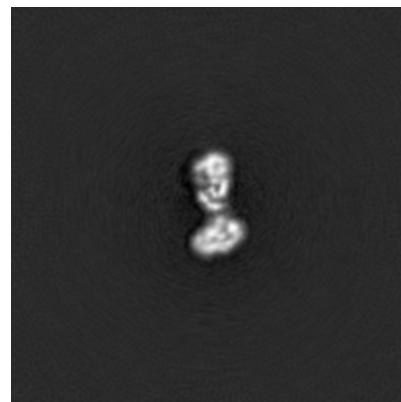
### 6.3.2 Raw map



X Index: 99



Y Index: 96



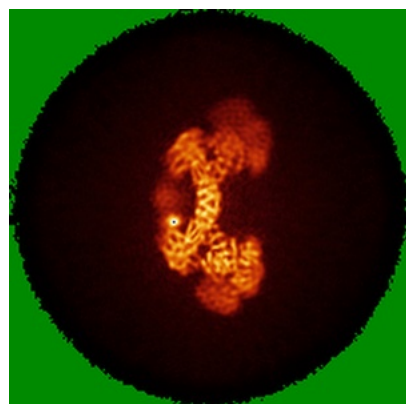
Z Index: 77

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

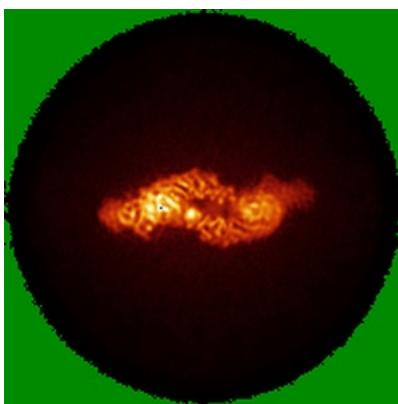


## 6.4 Orthogonal standard-deviation projections (False-color) [i](#)

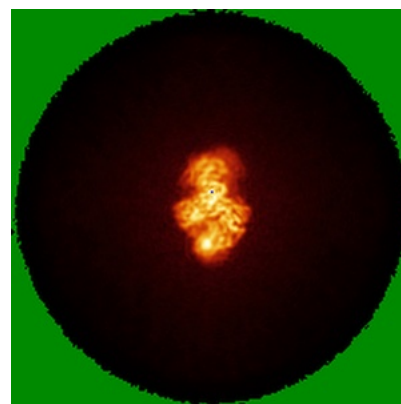
### 6.4.1 Primary map



X

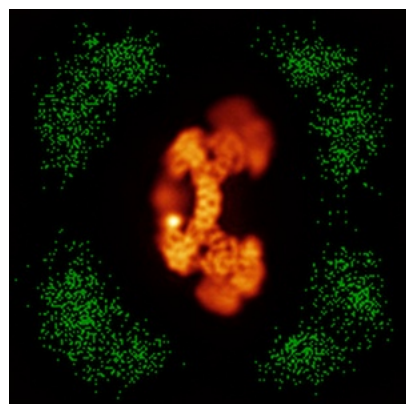


Y

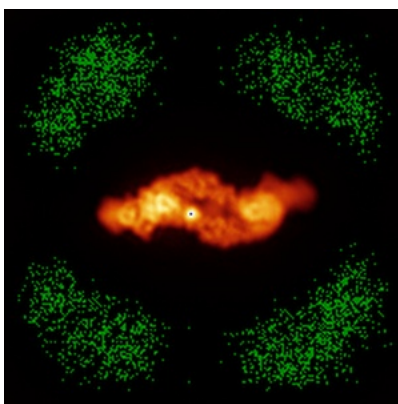


Z

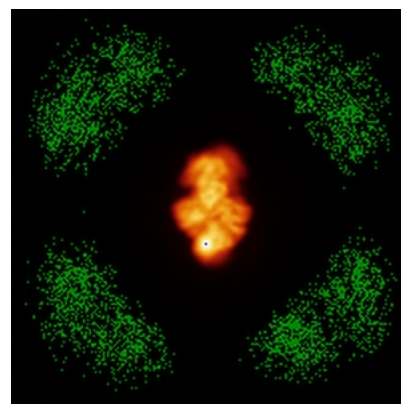
### 6.4.2 Raw map



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.

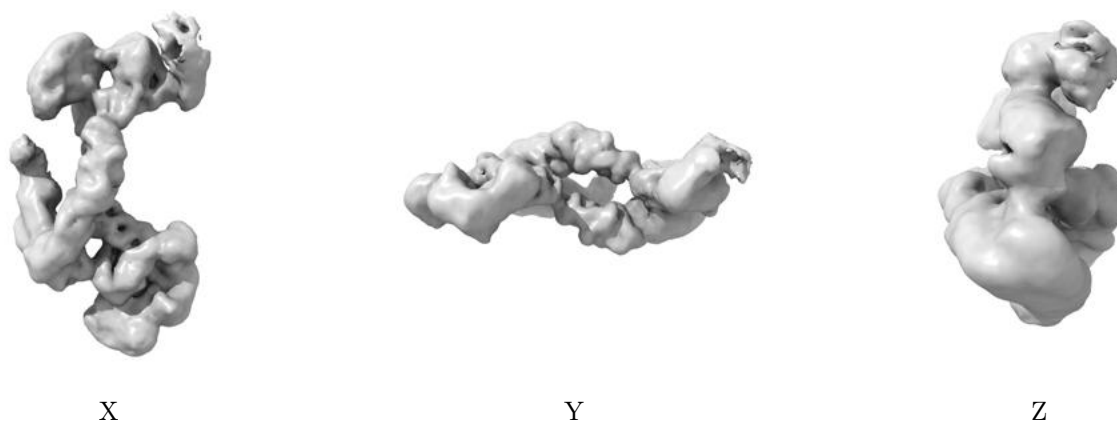
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.3. 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.

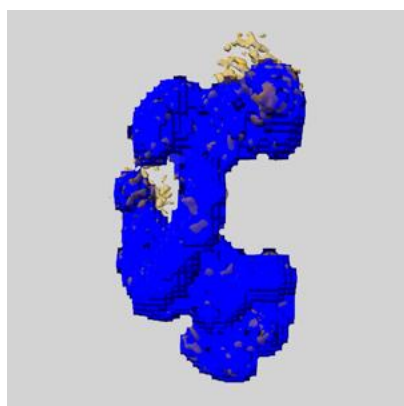
## 6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

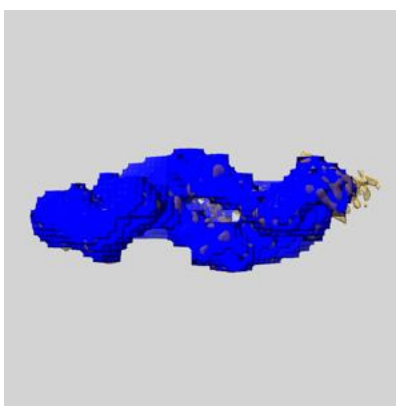
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

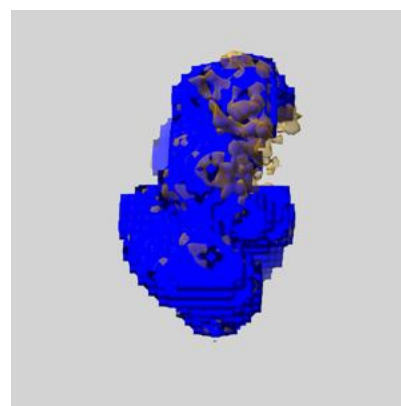
### 6.6.1 emd\_74435\_msk\_1.map [i](#)



X

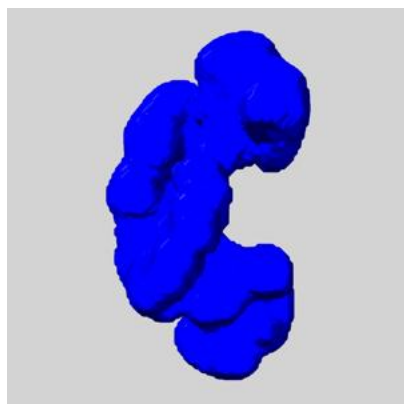


Y

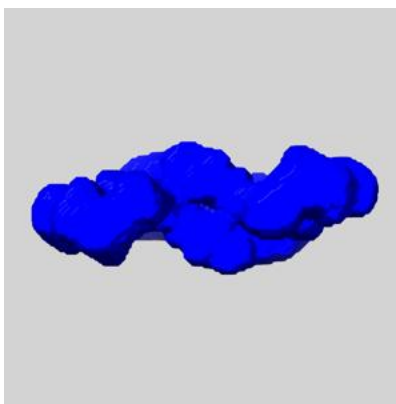


Z

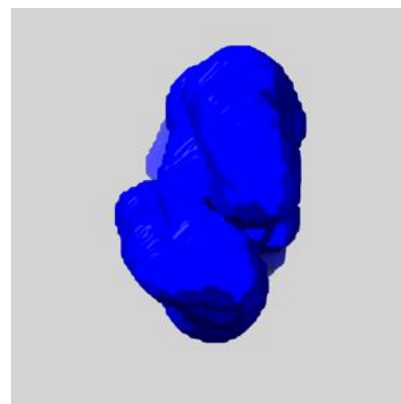
### 6.6.2 emd\_74435\_msk\_2.map [i](#)



X



Y

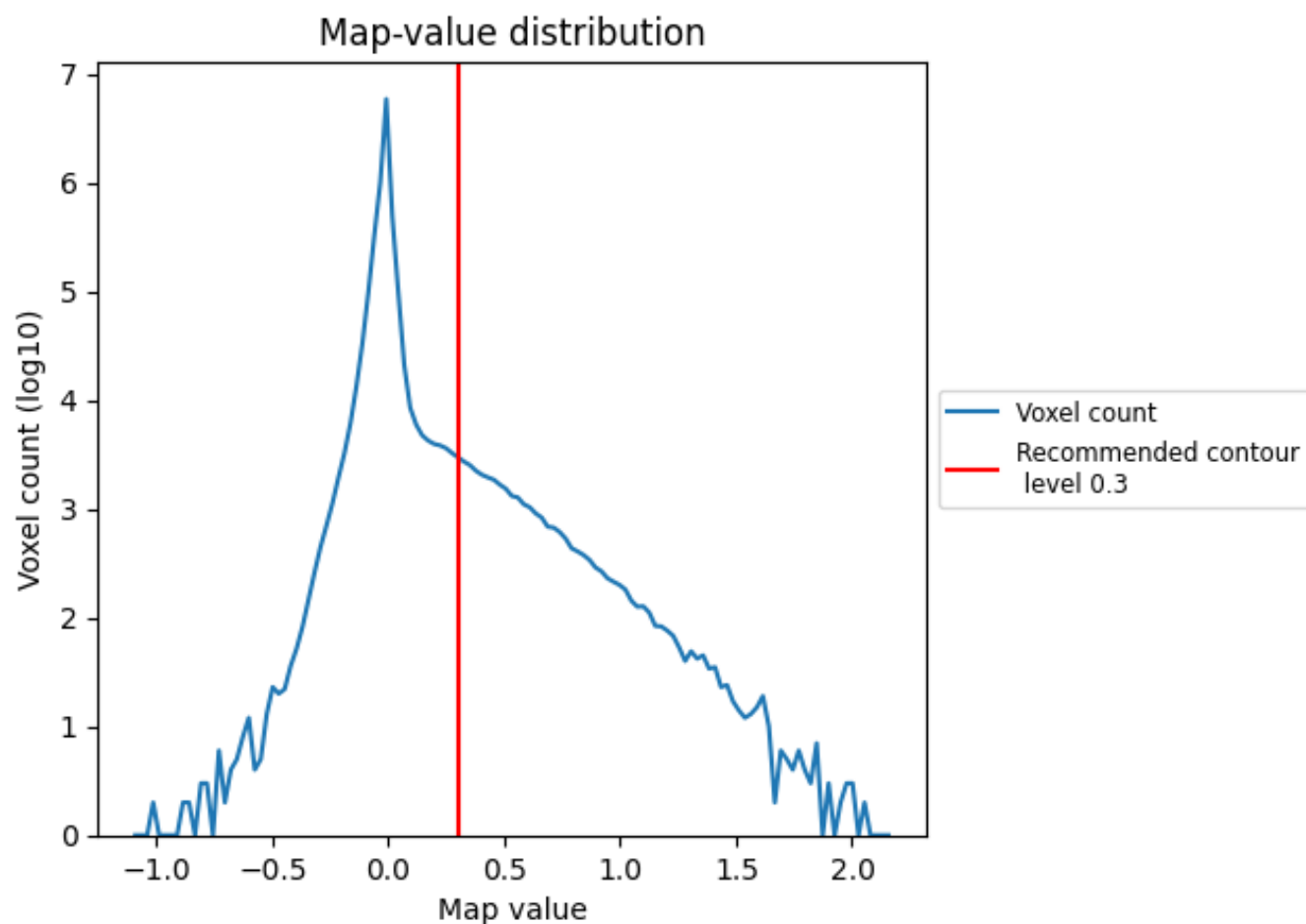


Z

## 7 Map analysis [i](#)

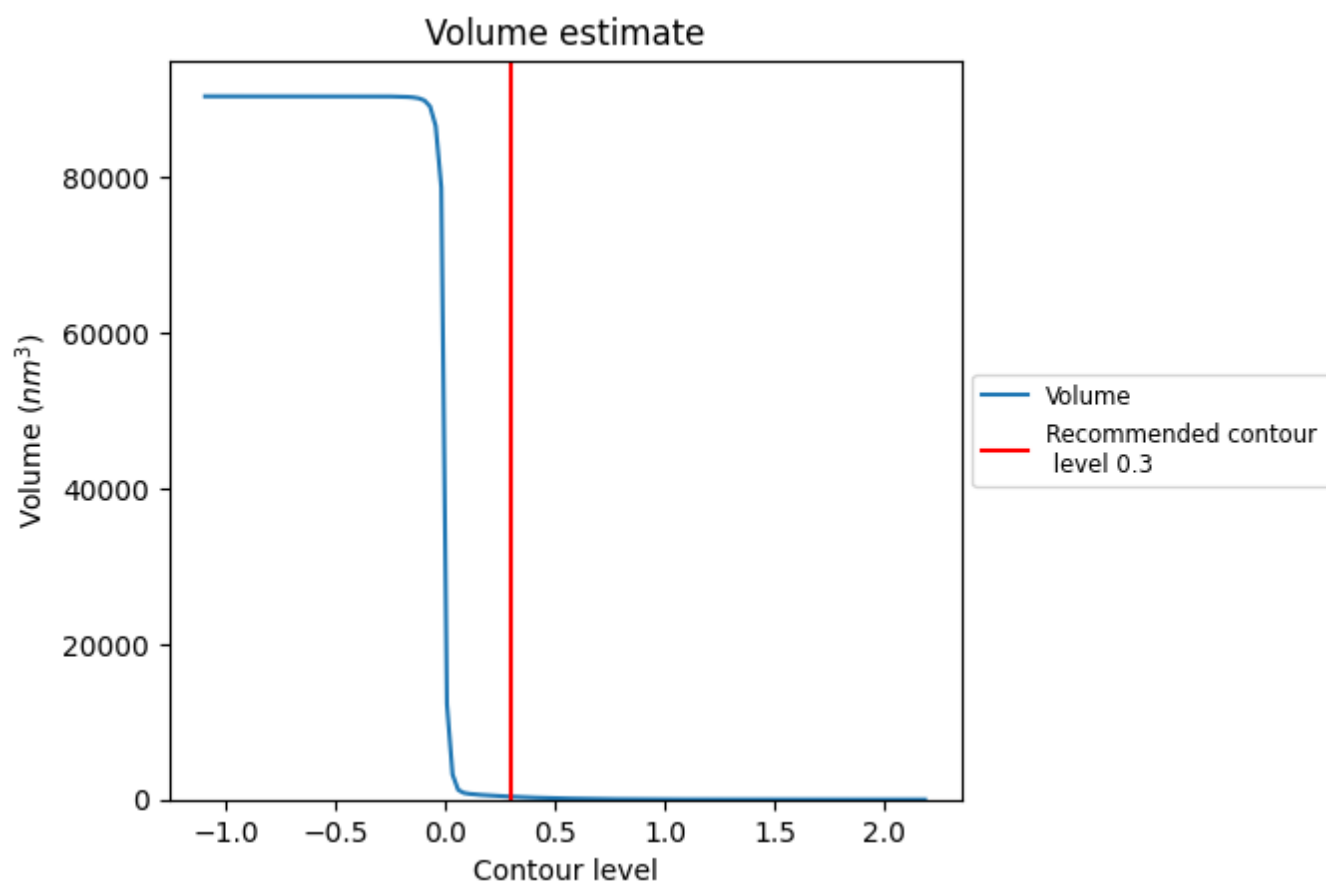
This section contains the results of statistical analysis of the map.

### 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.

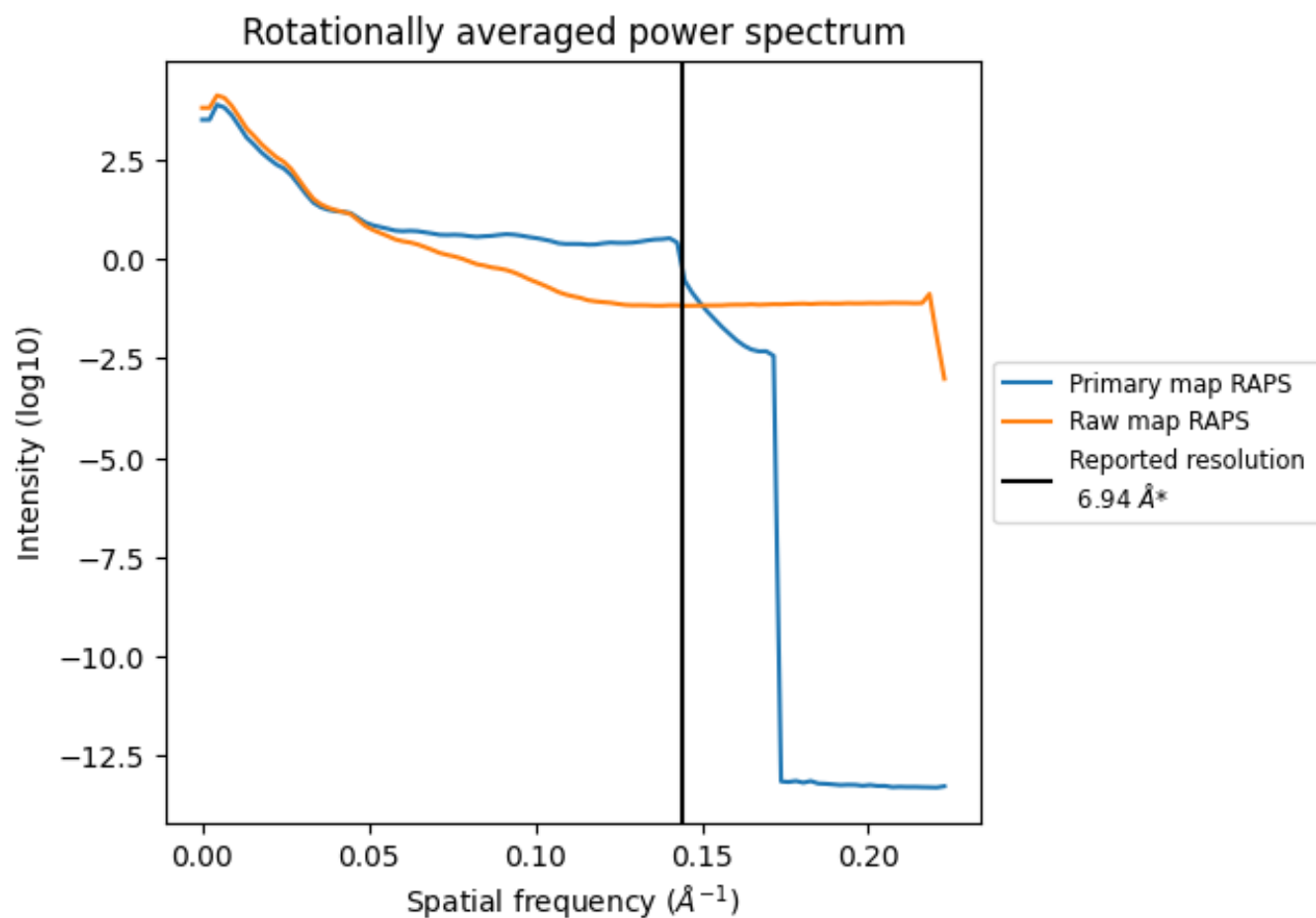
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 371 nm<sup>3</sup>; this corresponds to an approximate mass of 335 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 ⓘ

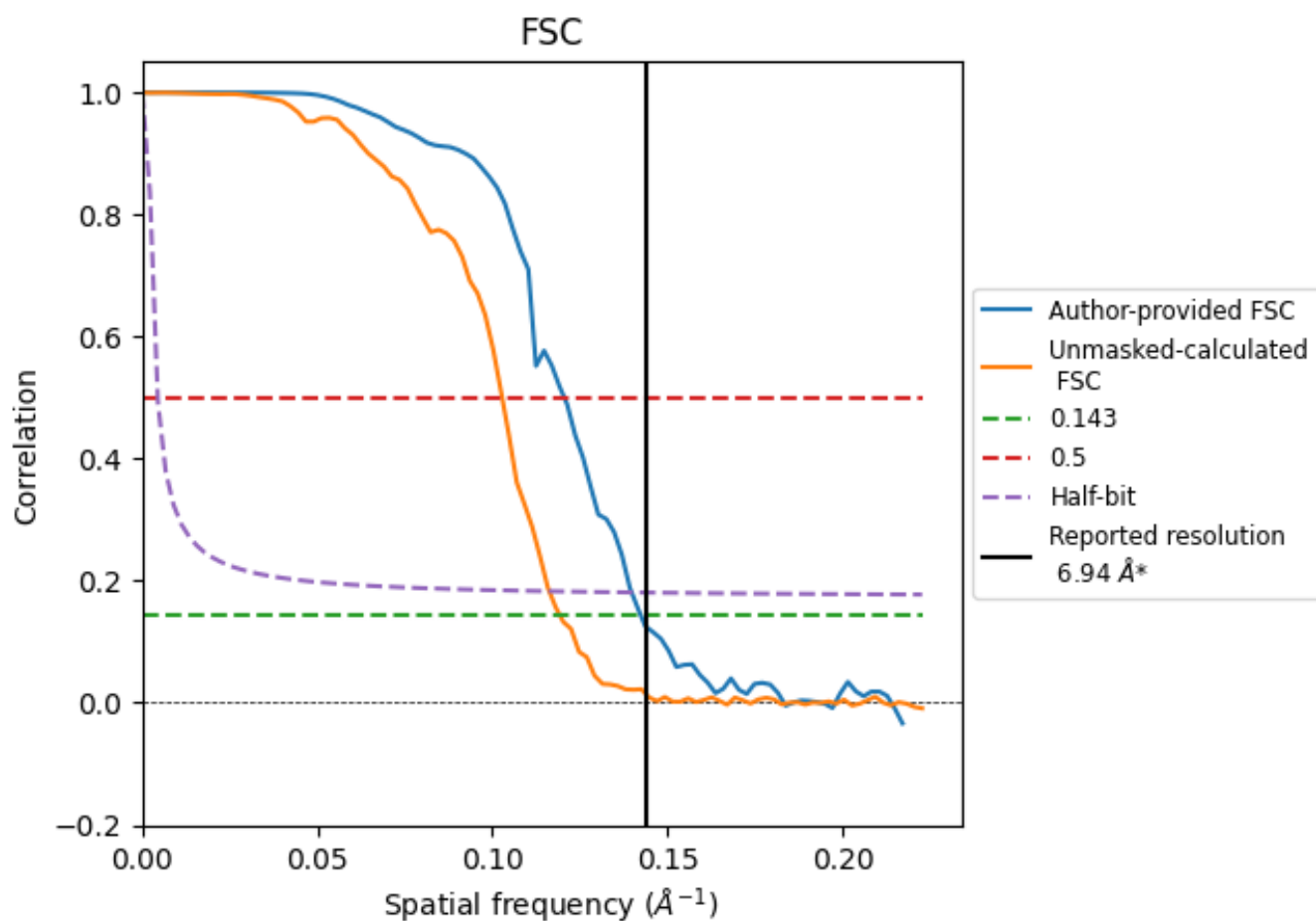


\*Reported resolution corresponds to spatial frequency of 0.144 Å<sup>-1</sup>

## 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 0.144  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	6.94	-	-
Author-provided FSC curve	7.02	8.29	7.15
Unmasked-calculated*	8.38	9.73	8.60

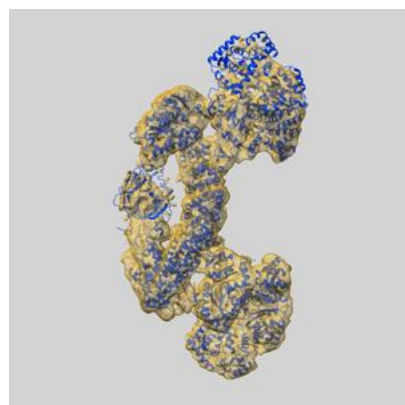
\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 8.38 differs from the reported value 6.94 by more than 10 %



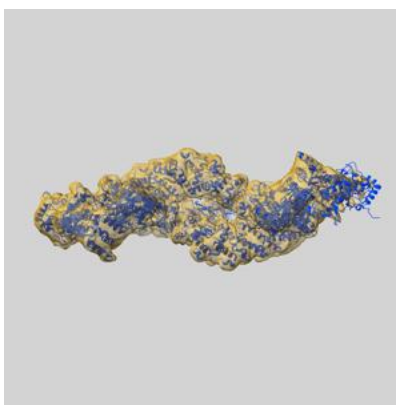
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-74435 and PDB model 9ZN5. Per-residue inclusion information can be found in section [3](#) on page [4](#).

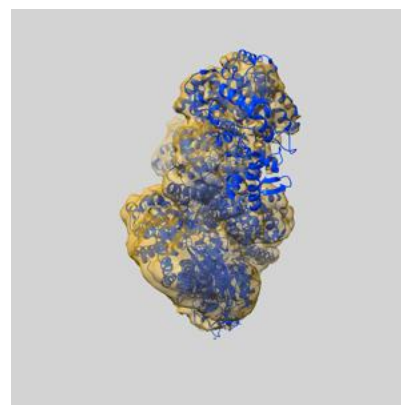
### 9.1 Map-model overlay [i](#)



X



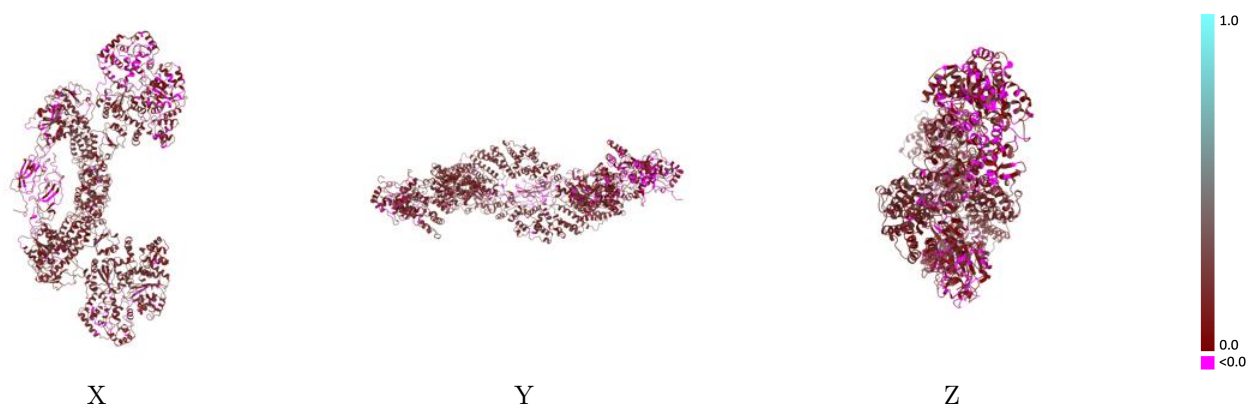
Y



Z

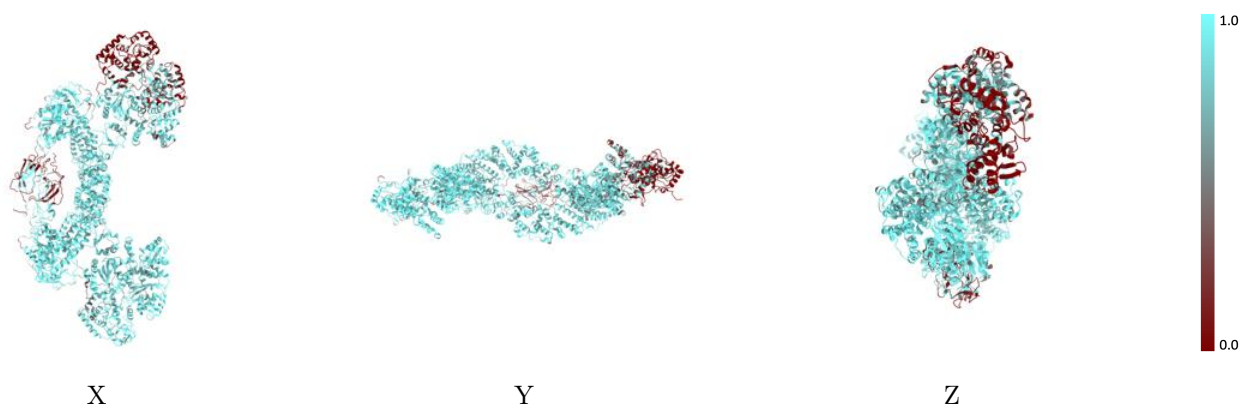
The images above show the 3D surface view of the map at the recommended contour level 0.3 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



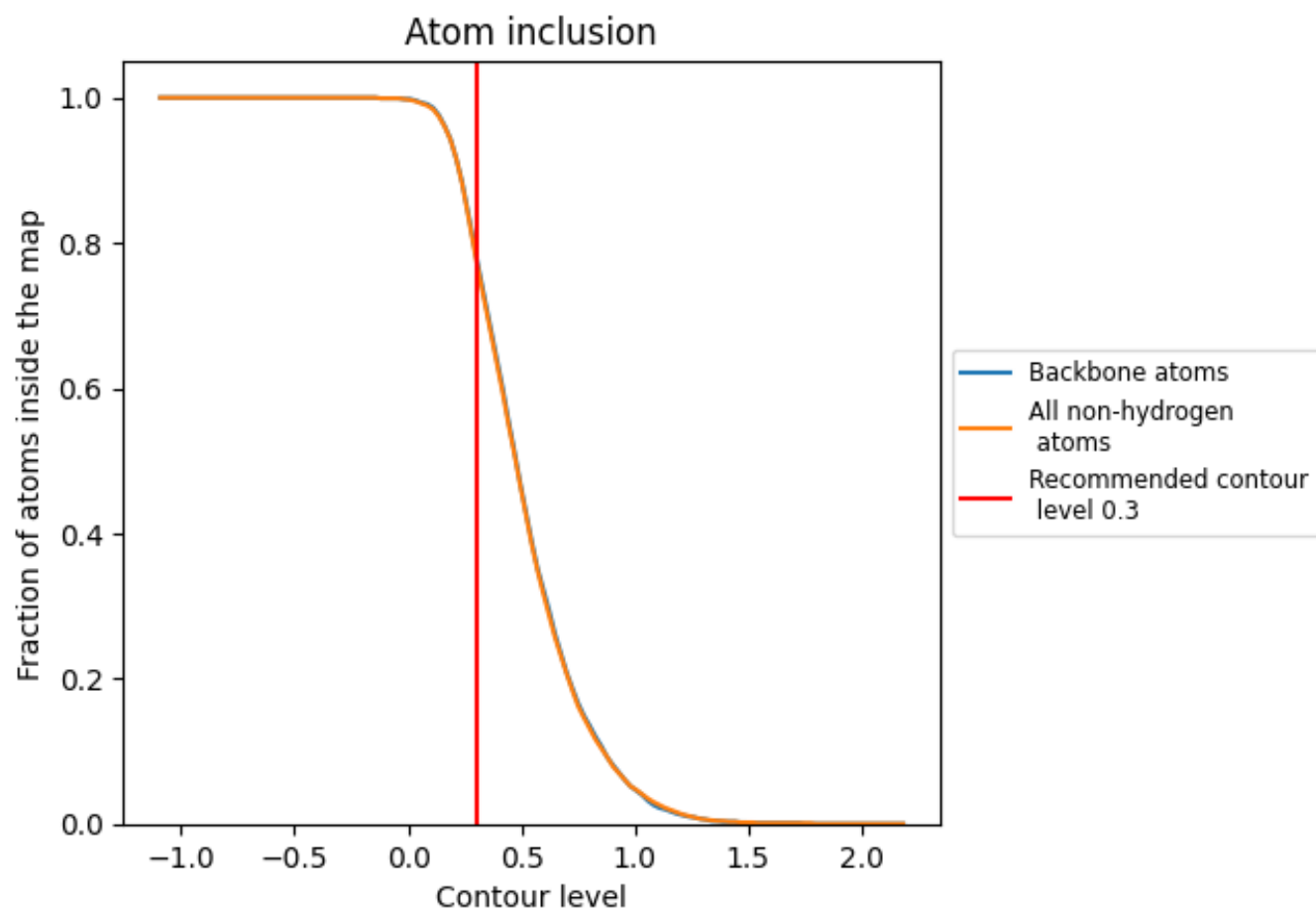
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.3).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 78% of all backbone atoms, 78% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7780	<div></div> 0.1480
A	<div></div> 0.9000	<div></div> 0.1790
B	<div></div> 0.8300	<div></div> 0.1500
C	<div></div> 0.5330	<div></div> 0.0990
D	<div></div> 0.8400	<div></div> 0.1610

