



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 06:02 AM UTC

PDB ID : 7TJY / pdb_00007tjy
EMDB ID : EMD-25946
Title : Yeast ATP synthase State 1catalytic(a) without exogenous ATP backbone model
Authors : Guo, H.; Rubinstein, J.L.
Deposited on : 2022-01-17
Resolution : 3.80 Å(reported)
Based on initial model : 2HLD

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:

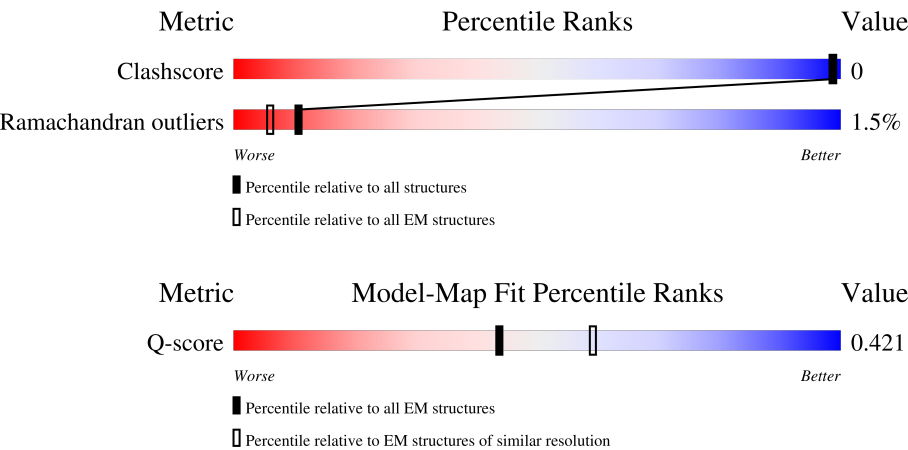
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.80 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	-
Q-score	-	25397	10198 (3.30 - 4.30)

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	0	76	7% 99% .
1	1	76	96% ..
1	2	76	97% ..
1	3	76	93% ...
1	4	76	8% 93% 5% .

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Mol	Chain	Length	Quality of chain
1	5	76	
1	6	76	
1	7	76	
1	8	76	
1	9	76	
2	A	510	
2	B	510	
2	C	510	
3	D	478	
3	E	478	
3	F	478	
4	G	278	
5	H	138	
6	I	61	
7	O	195	
8	T	249	
9	U	209	
10	V	173	
11	W	95	
12	X	92	
13	Y	59	
14	Z	48	

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 20228 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 ATP synthase subunit 9.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	0	75	Total	C	N	O	0	0
			300	150	75	75		
1	1	75	Total	C	N	O	0	0
			300	150	75	75		
1	2	75	Total	C	N	O	0	0
			300	150	75	75		
1	3	74	Total	C	N	O	0	0
			296	148	74	74		
1	4	75	Total	C	N	O	0	0
			300	150	75	75		
1	5	75	Total	C	N	O	0	0
			300	150	75	75		
1	6	74	Total	C	N	O	0	0
			296	148	74	74		
1	7	73	Total	C	N	O	0	0
			292	146	73	73		
1	8	75	Total	C	N	O	0	0
			300	150	75	75		
1	9	74	Total	C	N	O	0	0
			296	148	74	74		

- Molecule 2 is a protein called ATP synthase subunit alpha.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	A	499	Total	C	N	O	0	0
			1996	998	499	499		
2	B	505	Total	C	N	O	0	0
			2020	1010	505	505		
2	C	498	Total	C	N	O	0	0
			1992	996	498	498		

- Molecule 3 is a protein called ATP synthase subunit beta.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	D	470	Total	C	N	O	0	0
			1880	940	470	470		
3	E	468	Total	C	N	O	0	0
			1872	936	468	468		
3	F	469	Total	C	N	O	0	0
			1876	938	469	469		

- Molecule 4 is a protein called ATP synthase subunit gamma.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	G	265	Total	C	N	O	0	0
			1060	530	265	265		

- Molecule 5 is a protein called ATP synthase subunit delta.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	H	120	Total	C	N	O	0	0
			480	240	120	120		

- Molecule 6 is a protein called ATP synthase subunit epsilon.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	I	48	Total	C	N	O	0	0
			193	96	48	49		

- Molecule 7 is a protein called ATP synthase subunit 5.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	O	187	Total	C	N	O	0	0
			748	374	187	187		

- Molecule 8 is a protein called ATP synthase subunit a.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	T	224	Total	C	N	O	0	0
			897	448	224	225		

- Molecule 9 is a protein called ATP synthase subunit 4.

Mol	Chain	Residues	Atoms				AltConf	Trace
9	U	155	Total	C	N	O	0	0
			620	310	155	155		

- Molecule 10 is a protein called ATP synthase subunit d.

Mol	Chain	Residues	Atoms				AltConf	Trace
10	V	171	Total	C	N	O	0	0
			685	342	171	172		

- Molecule 11 is a protein called ATP synthase subunit f.

Mol	Chain	Residues	Atoms				AltConf	Trace
11	W	85	Total	C	N	O	0	0
			340	170	85	85		

- Molecule 12 is a protein called ATP synthase subunit H.

Mol	Chain	Residues	Atoms				AltConf	Trace
12	X	62	Total	C	N	O	0	0
			248	124	62	62		

- Molecule 13 is a protein called ATP synthase subunit J.

Mol	Chain	Residues	Atoms				AltConf	Trace
13	Y	37	Total	C	N	O	0	0
			148	74	37	37		

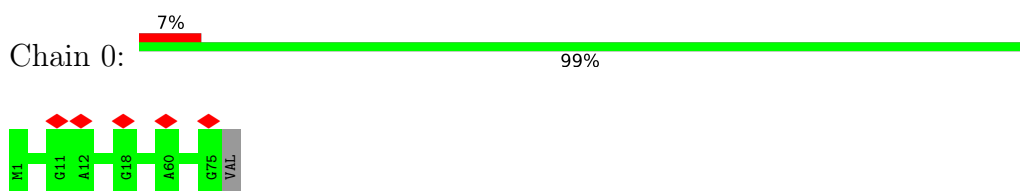
- Molecule 14 is a protein called ATP synthase protein 8.

Mol	Chain	Residues	Atoms				AltConf	Trace
14	Z	48	Total	C	N	O	0	0
			193	96	48	49		

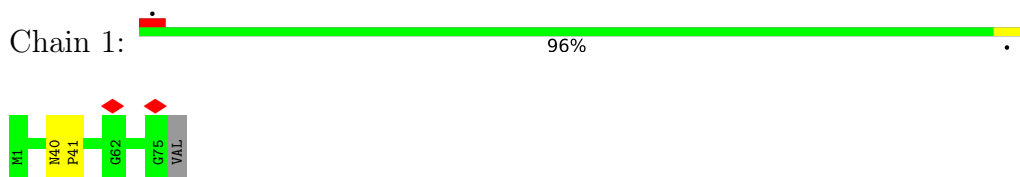
3 Residue-property plots [i](#)

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

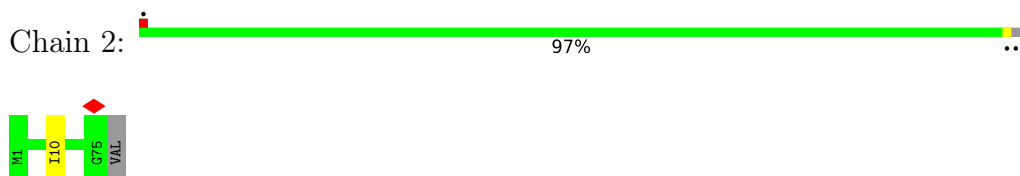
- Molecule 1: ATP synthase subunit 9



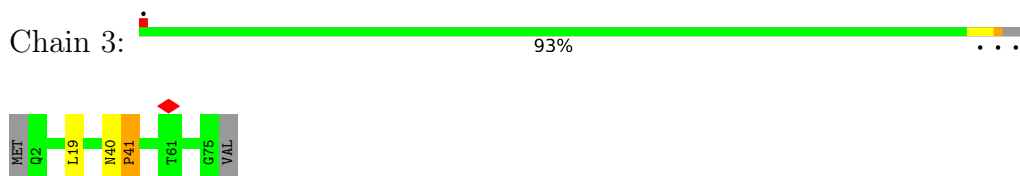
- Molecule 1: ATP synthase subunit 9



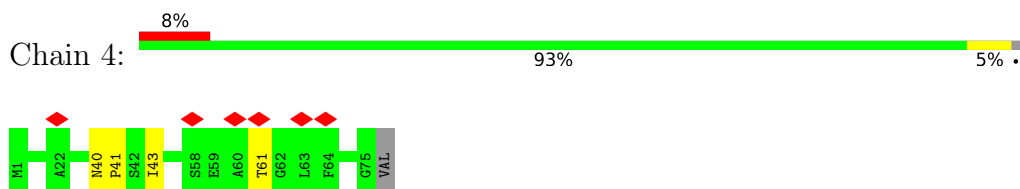
- Molecule 1: ATP synthase subunit 9



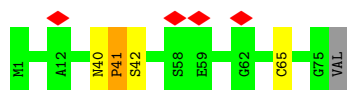
- Molecule 1: ATP synthase subunit 9



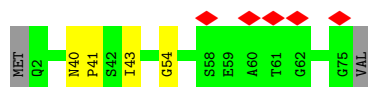
- Molecule 1: ATP synthase subunit 9



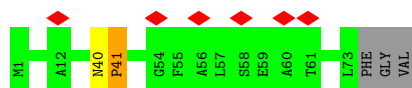
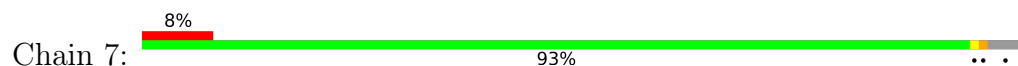
- Molecule 1: ATP synthase subunit 9



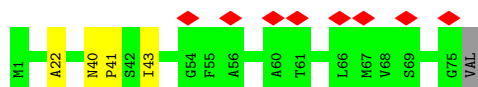
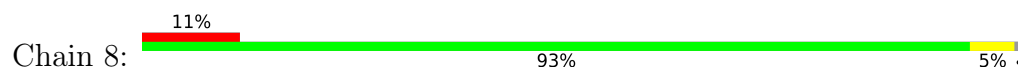
- Molecule 1: ATP synthase subunit 9



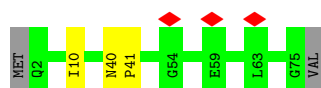
- Molecule 1: ATP synthase subunit 9



- Molecule 1: ATP synthase subunit 9



- Molecule 1: ATP synthase subunit 9

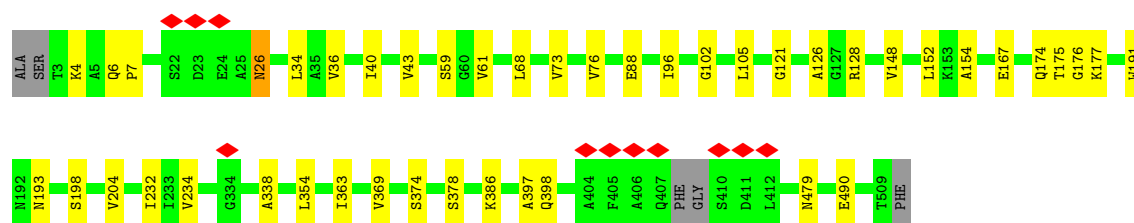


- Molecule 2: ATP synthase subunit alpha



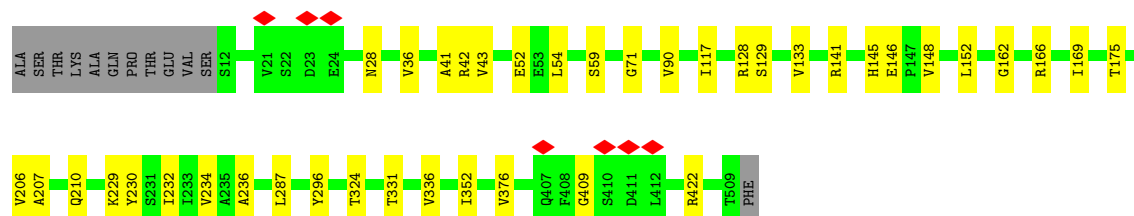
- Molecule 2: ATP synthase subunit alpha

Chain B:  90% 9% .



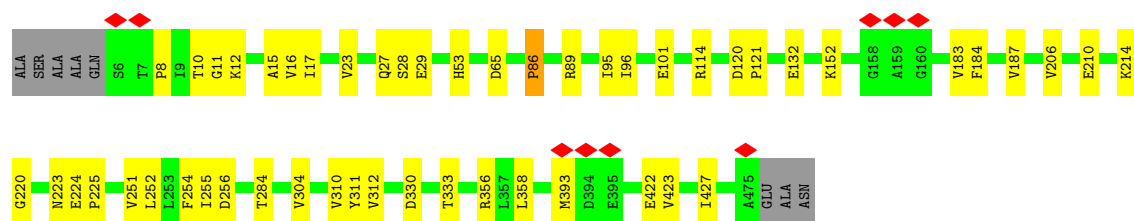
• Molecule 2: ATP synthase subunit alpha

Chain C:  90% 8% .



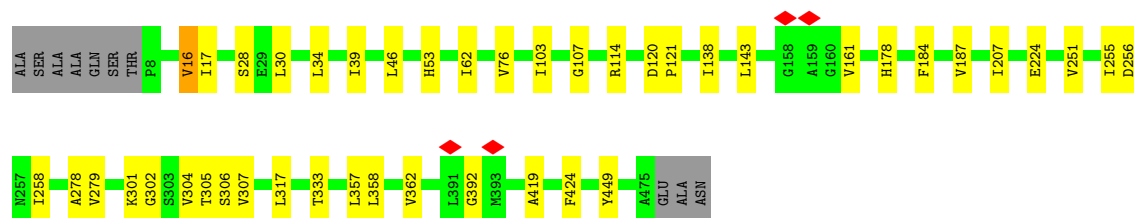
• Molecule 3: ATP synthase subunit beta

Chain D:  88% 10% .



• Molecule 3: ATP synthase subunit beta

Chain E:  89% 9% .



• Molecule 3: ATP synthase subunit beta

Chain F:  89% 9% .





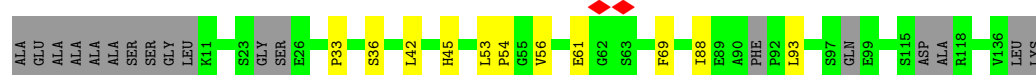
- Molecule 4: ATP synthase subunit gamma

Chain G: 90% 6% 5%



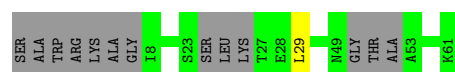
- Molecule 5: ATP synthase subunit delta

Chain H: 79% 8% 13%



- Molecule 6: ATP synthase subunit epsilon

Chain I: 77% 0% 21%



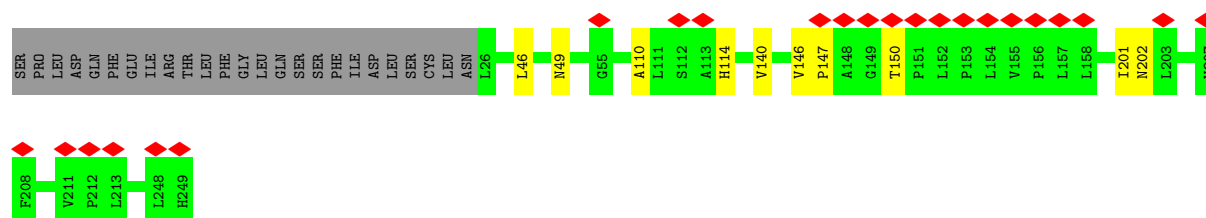
- Molecule 7: ATP synthase subunit 5

Chain O: 85% 11% 0%



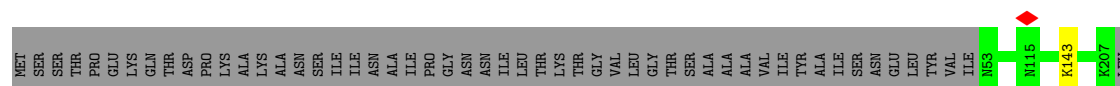
- Molecule 8: ATP synthase subunit a

Chain T: 9% 86% 0% 10%



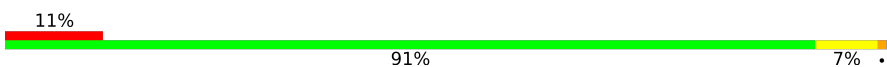
- Molecule 9: ATP synthase subunit 4

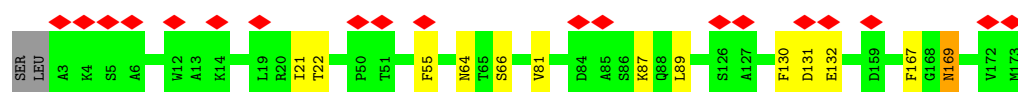
Chain U: 74% 0% 26%



LYS

- Molecule 10: ATP synthase subunit d

Chain V: 



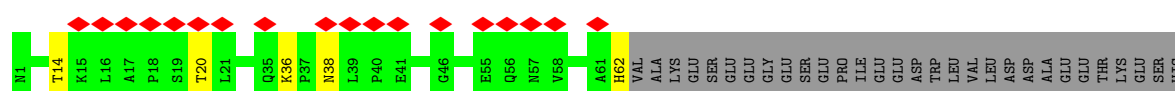
- Molecule 11: ATP synthase subunit f

Chain W: 



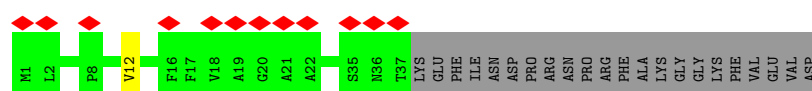
- Molecule 12: ATP synthase subunit H

Chain X: 

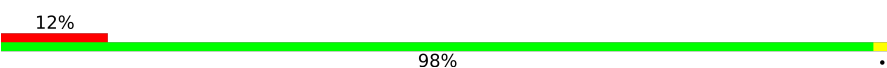


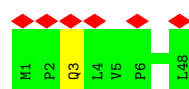
- Molecule 13: ATP synthase subunit J

Chain Y: 



- Molecule 14: ATP synthase protein 8

Chain Z: 



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	41506	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	39	Depositor
Minimum defocus (nm)	1100	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	133843	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.081	Depositor
Minimum map value	-0.558	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.107	Depositor
Recommended contour level	0.7	Depositor
Map size (Å)	344.96, 344.96, 344.96	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.3475, 1.3475, 1.3475	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	0	1.33	0/299	1.94	0/372
1	1	1.37	0/299	1.81	0/372
1	2	1.38	0/299	1.86	2/372 (0.5%)
1	3	1.37	0/295	1.74	3/367 (0.8%)
1	4	1.30	0/299	1.90	4/372 (1.1%)
1	5	1.30	0/299	1.97	3/372 (0.8%)
1	6	1.30	0/295	2.02	4/367 (1.1%)
1	7	1.28	0/291	2.01	1/362 (0.3%)
1	8	1.29	0/299	2.01	4/372 (1.1%)
1	9	1.33	0/295	1.94	2/367 (0.5%)
2	A	1.69	1/1994 (0.1%)	1.76	24/2489 (1.0%)
2	B	1.69	2/2018 (0.1%)	1.80	37/2519 (1.5%)
2	C	1.67	3/1991 (0.2%)	1.77	35/2487 (1.4%)
3	D	1.64	2/1879 (0.1%)	1.85	52/2347 (2.2%)
3	E	1.72	3/1871 (0.2%)	1.89	38/2337 (1.6%)
3	F	1.70	2/1875 (0.1%)	1.87	36/2342 (1.5%)
4	G	1.59	0/1058	1.70	10/1319 (0.8%)
5	H	1.55	0/475	1.73	9/585 (1.5%)
6	I	1.43	0/190	1.58	1/231 (0.4%)
7	O	1.64	0/747	1.84	17/932 (1.8%)
8	T	1.41	0/896	1.62	11/1117 (1.0%)
9	U	1.39	0/619	1.66	2/772 (0.3%)
10	V	1.42	0/684	1.70	8/852 (0.9%)
11	W	1.33	0/339	1.95	7/422 (1.7%)
12	X	1.40	0/247	2.06	3/307 (1.0%)
13	Y	1.25	0/147	1.79	1/182 (0.5%)
14	Z	1.37	0/192	1.60	1/237 (0.4%)
All	All	1.58	13/20192 (0.1%)	1.81	315/25172 (1.3%)

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	1	0	1
1	3	0	1
1	4	0	1
1	5	0	1
1	6	0	1
1	7	0	1
1	8	0	1
1	9	0	1
3	D	0	1
3	F	0	1
5	H	0	1
All	All	0	11

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	11	GLY	CA-C	-6.62	1.46	1.52
3	F	185	THR	CA-C	-5.64	1.45	1.52
2	C	207	ALA	CA-C	-5.64	1.46	1.52
2	B	374	SER	CA-C	-5.63	1.46	1.52
2	C	236	ALA	CA-C	-5.60	1.47	1.53
3	E	143	LEU	CA-C	-5.46	1.47	1.53
3	E	306	SER	CA-C	-5.44	1.46	1.52
2	C	162	GLY	CA-C	-5.23	1.47	1.52
3	F	253	LEU	CA-C	-5.19	1.46	1.52
3	D	15	ALA	CA-C	-5.14	1.46	1.52
2	B	73	VAL	CA-C	-5.12	1.46	1.52
2	A	36	VAL	CA-C	-5.04	1.46	1.52
3	E	251	VAL	CA-C	-5.03	1.47	1.52

All (315) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	362	VAL	N-CA-C	-16.26	98.46	112.12
3	E	121	PRO	CA-C-O	-9.70	113.92	120.90
3	F	255	ILE	N-CA-C	-9.60	94.15	108.17
7	O	55	HIS	N-CA-C	-9.60	103.67	114.62
4	G	194	PHE	N-CA-C	-9.46	96.58	109.54
7	O	176	VAL	N-CA-C	-9.40	94.59	108.12
3	E	304	VAL	N-CA-C	-9.37	99.84	110.05
2	B	96	ILE	N-CA-C	-9.14	99.88	110.21
3	E	161	VAL	N-CA-C	-8.96	103.66	112.17
11	W	6	PRO	N-CA-C	8.89	121.55	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	362	VAL	N-CA-C	-8.81	104.76	111.90
3	E	258	ILE	N-CA-C	-8.73	102.48	113.22
7	O	180	ILE	N-CA-C	-8.68	102.38	111.58
2	C	234	VAL	N-CA-C	-8.65	96.06	108.17
8	T	146	VAL	N-CA-C	-8.57	98.43	107.60
8	T	140	VAL	N-CA-C	-8.41	104.39	112.29
2	B	175	THR	N-CA-C	-8.39	104.08	112.97
2	A	232	ILE	N-CA-C	-8.34	96.11	108.45
2	B	234	VAL	N-CA-C	-8.20	96.64	108.53
3	E	305	THR	N-CA-C	-8.07	96.32	108.99
5	H	69	PHE	N-CA-C	-7.93	95.98	108.90
3	F	95	ILE	N-CA-C	-7.87	97.09	108.11
2	B	6	GLN	N-CA-C	-7.87	101.07	108.07
3	E	207	ILE	N-CA-C	-7.85	97.12	108.11
4	G	204	ASN	CA-C-N	7.84	125.18	120.24
4	G	204	ASN	C-N-CA	7.84	125.18	120.24
3	F	187	VAL	N-CA-C	-7.77	97.24	108.36
2	B	354	LEU	N-CA-C	-7.77	97.28	109.72
3	D	121	PRO	CA-C-O	-7.74	115.32	120.90
3	F	252	LEU	N-CA-C	-7.73	97.33	109.07
3	F	218	VAL	N-CA-C	-7.71	97.02	108.12
4	G	164	ILE	N-CA-C	-7.70	96.99	108.46
2	C	175	THR	N-CA-C	-7.68	103.22	112.59
7	O	14	GLY	CA-C-N	7.51	130.57	120.50
7	O	14	GLY	C-N-CA	7.51	130.57	120.50
3	D	89	ARG	N-CA-C	-7.49	103.27	113.30
3	E	114	ARG	N-CA-C	-7.45	97.97	109.52
3	D	254	PHE	N-CA-C	-7.40	96.48	109.06
2	A	352	ILE	N-CA-C	-7.35	97.53	108.12
3	D	16	VAL	N-CA-C	-7.26	96.63	107.37
2	B	369	VAL	N-CA-C	-7.24	105.95	112.90
3	F	39	ILE	N-CA-C	-7.21	98.05	108.36
3	D	53	HIS	N-CA-C	-7.20	97.66	109.40
2	C	28	ASN	N-CA-C	-7.19	96.45	108.75
2	C	336	VAL	N-CA-C	-7.17	105.73	112.83
7	O	33	ILE	N-CA-C	-7.13	105.68	111.81
2	B	73	VAL	N-CA-C	-7.10	97.88	108.46
7	O	84	TYR	N-CA-C	-7.10	102.96	111.69
2	C	169	ILE	N-CA-C	-7.05	97.70	107.99
3	D	10	THR	N-CA-C	-7.04	97.08	109.06
2	B	479	ASN	N-CA-C	-7.02	104.46	113.16
2	C	232	ILE	N-CA-C	-6.97	98.44	108.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	W	13	LYS	N-CA-C	-6.97	103.89	112.88
3	D	251	VAL	N-CA-C	-6.96	100.04	108.53
7	O	150	GLY	N-CA-C	-6.95	105.04	115.32
3	F	222	MET	N-CA-C	-6.92	104.71	113.02
2	A	479	ASN	N-CA-C	-6.89	104.88	113.28
11	W	46	THR	N-CA-C	-6.78	104.62	113.17
3	F	363	VAL	N-CA-C	-6.75	106.15	112.83
2	A	265	HIS	N-CA-C	-6.75	99.06	109.52
3	F	184	PHE	N-CA-C	-6.70	97.99	108.90
3	D	132	GLU	N-CA-C	-6.62	99.24	109.24
11	W	15	ILE	N-CA-C	-6.62	106.84	113.20
2	C	229	LYS	N-CA-C	-6.61	103.96	113.21
3	F	254	PHE	N-CA-C	-6.59	98.23	109.24
3	F	160	GLY	N-CA-C	-6.58	106.65	114.48
3	F	420	VAL	N-CA-C	-6.57	106.39	112.96
2	B	191	TRP	N-CA-C	-6.56	104.67	114.64
3	F	185	THR	N-CA-C	-6.55	97.69	108.76
2	B	378	SER	N-CA-C	-6.53	104.94	113.17
2	C	148	VAL	N-CA-C	-6.53	99.03	108.17
3	F	153	ILE	N-CA-C	-6.50	99.01	108.89
2	A	175	THR	N-CA-C	-6.47	105.26	112.57
3	E	449	TYR	N-CA-C	-6.45	102.63	110.88
10	V	167	PHE	N-CA-C	-6.43	99.55	109.14
2	A	296	TYR	N-CA-C	-6.43	101.20	110.40
2	B	128	ARG	N-CA-C	-6.43	99.34	109.50
3	D	304	VAL	N-CA-C	-6.42	98.80	108.17
11	W	36	GLN	N-CA-C	-6.40	99.61	109.14
14	Z	3	GLN	N-CA-C	-6.38	105.53	113.38
2	B	68	LEU	N-CA-C	-6.36	98.36	108.73
2	C	206	VAL	N-CA-C	-6.35	99.22	108.11
2	C	129	SER	N-CA-C	-6.34	99.93	108.86
7	O	56	LEU	N-CA-C	-6.31	104.73	112.88
2	A	117	ILE	N-CA-C	-6.30	106.85	112.90
3	D	358	LEU	N-CA-C	-6.29	104.86	113.30
3	E	333	THR	N-CA-C	-6.29	98.26	108.52
3	E	30	LEU	N-CA-C	-6.29	101.41	110.40
3	E	103	ILE	N-CA-C	-6.29	106.64	112.43
7	O	57	LEU	N-CA-C	-6.26	102.93	110.44
3	E	255	ILE	N-CA-C	-6.25	98.68	108.86
2	C	331	THR	N-CA-C	-6.25	100.39	110.32
3	E	120	ASP	CA-C-N	6.24	124.22	119.66
3	E	120	ASP	C-N-CA	6.24	124.22	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	133	VAL	N-CA-C	-6.22	99.08	108.17
2	A	61	VAL	N-CA-C	-6.20	103.29	110.05
2	B	26	ASN	N-CA-C	-6.19	97.62	110.80
2	A	99	VAL	N-CA-C	-6.18	104.12	109.19
3	D	393	MET	N-CA-C	-6.18	105.67	112.72
10	V	87	LYS	N-CA-C	-6.18	106.34	114.31
2	A	64	MET	N-CA-C	-6.16	100.04	109.41
3	D	423	VAL	N-CA-C	-6.16	106.71	113.43
7	O	58	LEU	N-CA-C	-6.15	105.09	112.59
3	D	252	LEU	N-CA-C	-6.13	97.75	110.80
3	D	312	VAL	N-CA-C	-6.12	101.06	107.60
3	D	427	ILE	N-CA-C	-6.11	101.59	107.76
4	G	2	THR	N-CA-C	-6.11	104.55	111.14
10	V	81	VAL	N-CA-C	-6.10	106.28	113.42
2	A	49	ILE	N-CA-C	-6.10	99.92	108.27
2	B	61	VAL	N-CA-C	-6.08	99.72	108.42
2	B	148	VAL	N-CA-C	-6.07	97.64	107.28
3	D	96	ILE	N-CA-C	-6.05	99.55	108.99
3	D	86	PRO	N-CA-C	-6.03	100.05	112.47
2	B	386	LYS	N-CA-C	-5.98	105.64	113.12
13	Y	12	VAL	N-CA-C	-5.96	105.89	113.22
8	T	150	THR	N-CA-C	-5.95	101.35	108.25
7	O	164	LYS	N-CA-C	-5.94	99.60	109.46
3	D	311	TYR	N-CA-C	-5.94	99.22	108.90
2	B	36	VAL	N-CA-C	-5.92	99.95	108.42
1	6	43	ILE	CA-C-N	5.91	128.47	120.38
1	6	43	ILE	C-N-CA	5.91	128.47	120.38
3	D	183	VAL	N-CA-C	-5.90	99.63	108.12
2	A	210	GLN	N-CA-C	-5.89	100.29	109.72
2	C	422	ARG	CA-C-N	5.89	126.52	119.98
2	C	422	ARG	C-N-CA	5.89	126.52	119.98
2	B	398	GLN	N-CA-C	-5.89	106.22	113.41
3	D	255	ILE	N-CA-C	-5.89	99.26	108.86
3	E	317	LEU	N-CA-C	-5.87	105.77	113.17
1	6	54	GLY	CA-C-N	5.86	128.06	120.44
1	6	54	GLY	C-N-CA	5.86	128.06	120.44
12	X	14	THR	N-CA-C	-5.84	99.52	108.76
2	C	324	THR	N-CA-C	-5.84	100.66	109.95
11	W	48	LEU	N-CA-C	-5.83	105.77	114.64
2	C	54	LEU	N-CA-C	-5.83	99.68	109.07
2	A	234	VAL	N-CA-C	-5.83	99.32	108.23
7	O	83	GLY	N-CA-C	-5.83	106.97	113.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	V	55	PHE	N-CA-C	-5.83	106.17	113.28
2	A	490	GLU	N-CA-C	-5.77	98.88	108.75
3	E	301	LYS	N-CA-C	-5.77	106.16	113.43
3	F	183	VAL	N-CA-C	-5.77	100.17	108.48
2	B	490	GLU	N-CA-C	-5.76	98.49	108.69
3	E	307	VAL	N-CA-C	-5.75	100.11	108.17
1	9	10	ILE	CA-C-N	5.75	126.36	119.98
1	9	10	ILE	C-N-CA	5.75	126.36	119.98
3	D	224	GLU	CA-C-N	5.74	126.29	120.38
3	D	224	GLU	C-N-CA	5.74	126.29	120.38
5	H	93	LEU	N-CA-C	5.72	117.52	111.28
2	B	76	VAL	N-CA-C	-5.72	100.10	108.11
3	F	257	ASN	N-CA-C	-5.71	95.03	111.00
3	E	46	LEU	N-CA-C	-5.70	99.12	108.76
2	B	34	LEU	N-CA-C	-5.70	98.67	110.80
2	C	236	ALA	N-CA-C	-5.68	97.74	107.23
2	A	32	ARG	N-CA-C	-5.68	99.17	108.76
3	E	107	GLY	CA-C-N	5.67	126.93	119.84
3	E	107	GLY	C-N-CA	5.67	126.93	119.84
3	D	120	ASP	CA-C-N	5.66	123.79	119.66
3	D	120	ASP	C-N-CA	5.66	123.79	119.66
2	B	177	LYS	CA-C-N	5.63	127.83	120.28
2	B	177	LYS	C-N-CA	5.63	127.83	120.28
2	C	296	TYR	CA-C-N	5.63	125.95	119.92
2	C	296	TYR	C-N-CA	5.63	125.95	119.92
3	D	95	ILE	N-CA-C	-5.63	97.79	107.24
3	F	452	ILE	CA-C-N	5.62	125.63	119.89
3	F	452	ILE	C-N-CA	5.62	125.63	119.89
2	C	376	VAL	N-CA-C	-5.60	105.39	113.07
3	E	39	ILE	N-CA-C	-5.58	99.71	107.80
1	5	41	PRO	N-CA-C	5.56	123.93	112.47
2	B	198	SER	N-CA-C	-5.55	105.82	112.59
3	E	302	GLY	N-CA-C	-5.54	103.15	111.19
3	E	184	PHE	N-CA-C	-5.54	99.70	108.73
3	F	46	LEU	N-CA-C	-5.54	99.87	108.90
1	8	43	ILE	CA-C-N	5.53	127.96	120.38
1	8	43	ILE	C-N-CA	5.53	127.96	120.38
3	D	220	GLY	N-CA-C	-5.53	100.07	113.18
7	O	170	GLU	N-CA-C	-5.53	99.66	109.06
3	E	424	PHE	N-CA-C	5.52	117.38	111.36
3	F	335	LEU	N-CA-C	-5.51	99.92	108.90
3	E	392	GLY	N-CA-C	-5.51	104.51	111.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	V	21	ILE	N-CA-C	-5.50	100.14	108.17
3	E	16	VAL	N-CA-C	-5.50	97.91	109.34
1	7	41	PRO	N-CA-C	5.49	123.78	112.47
3	D	225	PRO	CA-C-N	5.48	125.14	119.05
3	D	225	PRO	C-N-CA	5.48	125.14	119.05
3	F	249	GLN	N-CA-C	-5.48	101.02	109.52
3	F	53	HIS	N-CA-C	-5.48	100.25	109.07
3	F	207	ILE	N-CA-C	-5.47	100.70	108.58
1	4	43	ILE	CA-C-N	5.47	128.37	120.38
1	4	43	ILE	C-N-CA	5.47	128.37	120.38
3	D	214	LYS	N-CA-C	-5.46	105.05	112.26
2	B	232	ILE	N-CA-C	-5.46	100.44	108.85
3	D	11	GLY	N-CA-C	-5.46	102.77	111.18
2	A	422	ARG	CA-C-N	5.45	126.03	119.98
2	A	422	ARG	C-N-CA	5.45	126.03	119.98
3	D	206	VAL	N-CA-C	-5.45	105.22	110.72
3	E	62	ILE	N-CA-C	-5.45	100.75	109.20
5	H	56	VAL	N-CA-C	-5.45	100.48	108.11
2	C	128	ARG	N-CA-C	-5.45	101.08	109.52
2	C	409	GLY	N-CA-C	-5.45	107.75	115.43
4	G	168	ASP	N-CA-C	-5.44	102.79	109.65
3	F	114	ARG	N-CA-C	-5.43	100.32	109.07
3	D	330	ASP	N-CA-C	5.43	116.88	111.07
8	T	49	ASN	CA-C-N	5.42	130.05	122.08
8	T	49	ASN	C-N-CA	5.42	130.05	122.08
8	T	114	HIS	CA-C-N	5.41	128.54	120.31
8	T	114	HIS	C-N-CA	5.41	128.54	120.31
6	I	29	LEU	N-CA-C	-5.41	106.35	113.17
3	F	161	VAL	N-CA-C	-5.41	106.58	112.80
4	G	201	THR	N-CA-C	-5.40	106.06	113.30
3	F	284	THR	N-CA-C	-5.40	104.26	111.87
3	D	8	PRO	N-CA-C	-5.40	107.61	114.35
5	H	88	ILE	N-CA-C	5.39	115.60	110.53
2	C	210	GLN	N-CA-C	-5.39	101.78	110.14
3	F	259	PHE	N-CA-C	-5.39	105.57	111.82
2	A	198	SER	N-CA-C	-5.39	106.48	112.57
2	B	7	PRO	N-CA-C	-5.39	105.59	114.75
3	D	223	ASN	N-CA-C	-5.39	106.38	113.17
2	B	193	ASN	N-CA-C	-5.37	105.95	112.88
5	H	45	HIS	N-CA-C	-5.37	101.13	109.14
3	D	310	VAL	N-CA-C	-5.37	100.50	108.45
3	D	422	GLU	N-CA-C	-5.37	106.75	113.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	43	VAL	N-CA-C	-5.36	100.59	108.85
3	E	224	GLU	CA-C-N	5.36	125.90	120.38
3	E	224	GLU	C-N-CA	5.36	125.90	120.38
2	A	376	VAL	N-CA-C	-5.36	107.15	113.42
3	D	27	GLN	CA-C-N	5.36	131.78	121.54
3	D	27	GLN	C-N-CA	5.36	131.78	121.54
2	C	42	ARG	N-CA-C	-5.35	98.85	108.48
2	C	152	LEU	N-CA-C	-5.35	100.51	109.24
3	E	178	HIS	N-CA-C	-5.35	105.87	112.72
3	D	23	VAL	N-CA-C	-5.33	100.17	108.86
3	D	187	VAL	N-CA-C	-5.33	100.08	107.80
2	A	68	LEU	N-CA-C	-5.31	98.58	107.99
3	E	187	VAL	N-CA-C	-5.30	100.75	108.17
7	O	119	LEU	N-CA-C	-5.29	101.49	109.85
3	E	138	ILE	N-CA-C	-5.29	100.97	108.58
10	V	131	ASP	N-CA-C	-5.29	105.68	113.61
2	C	230	TYR	N-CA-C	-5.28	105.81	113.21
7	O	30	ASN	N-CA-C	-5.28	106.68	113.23
2	B	363	ILE	N-CA-C	-5.28	98.89	107.28
3	D	114	ARG	N-CA-C	-5.26	101.36	109.52
3	D	356	ARG	N-CA-C	-5.26	106.44	112.92
3	F	16	VAL	N-CA-C	-5.24	98.44	109.34
1	2	10	ILE	CA-C-N	5.23	125.79	119.98
1	2	10	ILE	C-N-CA	5.23	125.79	119.98
2	A	331	THR	N-CA-C	-5.23	100.00	108.52
2	C	117	ILE	N-CA-C	-5.23	108.18	113.20
2	C	166	ARG	N-CA-C	-5.23	100.65	108.96
3	E	419	ALA	N-CA-C	5.21	116.96	111.28
1	3	19	LEU	CA-C-N	5.21	127.51	120.38
1	3	19	LEU	C-N-CA	5.21	127.51	120.38
3	E	76	VAL	N-CA-C	-5.20	100.38	108.86
3	D	65	ASP	N-CA-C	-5.20	101.21	108.96
3	F	224	GLU	CA-C-N	5.20	125.73	120.38
3	F	224	GLU	C-N-CA	5.20	125.73	120.38
5	H	61	GLU	N-CA-C	-5.20	99.82	108.34
2	B	105	LEU	N-CA-C	-5.19	106.95	113.28
3	E	357	LEU	N-CA-C	-5.19	100.63	108.67
3	D	184	PHE	N-CA-C	-5.18	100.08	108.52
2	B	40	ILE	N-CA-C	-5.18	100.42	108.81
10	V	169	ASN	N-CA-C	-5.17	99.78	110.80
3	D	12	LYS	N-CA-C	-5.17	100.47	108.90
7	O	168	ILE	N-CA-C	-5.17	100.62	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	53	HIS	N-CA-C	-5.17	99.99	108.41
1	4	61	THR	CA-C-N	5.16	125.71	119.98
1	4	61	THR	C-N-CA	5.16	125.71	119.98
2	A	128	ARG	N-CA-C	-5.16	100.49	108.90
2	C	287	LEU	N-CA-C	-5.16	106.50	112.89
9	U	143	LYS	CA-C-N	5.15	127.05	120.56
9	U	143	LYS	C-N-CA	5.15	127.05	120.56
10	V	64	ASN	N-CA-C	-5.15	105.72	112.41
3	D	333	THR	N-CA-C	-5.14	99.73	108.26
2	C	352	ILE	N-CA-C	-5.14	98.66	109.34
5	H	53	LEU	CA-C-N	5.13	126.26	119.84
5	H	53	LEU	C-N-CA	5.13	126.26	119.84
3	D	101	GLU	N-CA-C	-5.13	102.39	110.14
3	D	210	GLU	N-CA-C	-5.13	104.66	112.04
3	D	152	LYS	CA-C-N	5.13	127.47	120.35
3	D	152	LYS	C-N-CA	5.13	127.47	120.35
4	G	127	LYS	N-CA-C	-5.12	106.97	113.43
1	8	22	ALA	CA-C-N	5.12	125.62	119.94
1	8	22	ALA	C-N-CA	5.12	125.62	119.94
3	E	17	ILE	N-CA-C	-5.12	100.52	107.99
3	F	213	SER	N-CA-C	-5.12	100.56	108.90
3	F	20	ILE	CA-C-N	5.11	128.22	120.75
3	F	20	ILE	C-N-CA	5.11	128.22	120.75
2	C	90	VAL	N-CA-C	-5.11	100.84	108.46
2	C	36	VAL	N-CA-C	-5.11	100.83	107.88
5	H	42	LEU	N-CA-C	-5.11	99.29	108.48
1	5	65	CYS	CA-C-N	5.10	127.11	120.28
1	5	65	CYS	C-N-CA	5.10	127.11	120.28
2	A	264	LYS	N-CA-C	-5.10	102.07	109.31
4	G	208	ASP	N-CA-C	-5.08	105.92	112.68
2	B	152	LEU	N-CA-C	-5.08	100.01	108.34
2	A	30	THR	N-CA-C	-5.07	101.03	108.79
4	G	74	ILE	N-CA-C	-5.07	100.91	108.46
2	C	41	ALA	N-CA-C	-5.06	102.40	109.69
3	F	62	ILE	N-CA-C	-5.06	102.33	109.21
1	3	41	PRO	N-CA-C	5.06	122.89	112.47
12	X	20	THR	CA-C-N	5.06	127.06	120.28
12	X	20	THR	C-N-CA	5.06	127.06	120.28
2	C	59	SER	N-CA-C	-5.05	107.15	113.72
3	D	427	ILE	CA-C-O	-5.05	116.17	119.15
3	D	284	THR	N-CA-C	-5.04	107.63	112.97
2	B	397	ALA	N-CA-C	-5.04	106.38	112.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	60	ARG	N-CA-C	-5.03	100.26	108.76
2	B	204	VAL	N-CA-C	-5.03	101.07	108.11
8	T	110	ALA	CA-C-N	5.03	127.02	120.28
8	T	110	ALA	C-N-CA	5.03	127.02	120.28
8	T	201	ILE	CA-C-N	5.03	131.14	121.54
8	T	201	ILE	C-N-CA	5.03	131.14	121.54
2	B	121	GLY	CA-C-N	5.02	126.11	119.84
2	B	121	GLY	C-N-CA	5.02	126.11	119.84
11	W	14	ASN	N-CA-C	-5.01	107.22	113.38
2	B	88	GLU	N-CA-C	-5.00	103.05	110.46
2	C	43	VAL	N-CA-C	-5.00	101.15	108.85

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	1	40	ASN	Peptide
1	3	40	ASN	Peptide
1	4	40	ASN	Peptide
1	5	40	ASN	Peptide
1	6	40	ASN	Peptide
1	7	40	ASN	Peptide
1	8	40	ASN	Peptide
1	9	40	ASN	Peptide
3	D	256	ASP	Peptide
3	F	345	TYR	Peptide
5	H	54	PRO	Peptide

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	0	300	0	95	0	0
1	1	300	0	95	0	0
1	2	300	0	95	0	0
1	3	296	0	91	0	0
1	4	300	0	95	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	5	300	0	95	0	0
1	6	296	0	91	0	0
1	7	292	0	91	0	0
1	8	300	0	95	0	0
1	9	296	0	91	0	0
2	A	1996	0	570	0	0
2	B	2020	0	575	2	0
2	C	1992	0	572	0	0
3	D	1880	0	538	0	0
3	E	1872	0	537	0	0
3	F	1876	0	537	0	0
4	G	1060	0	277	0	0
5	H	480	0	122	0	0
6	I	193	0	43	0	0
7	O	748	0	205	0	0
8	T	897	0	248	0	0
9	U	620	0	158	0	0
10	V	685	0	173	1	0
11	W	340	0	92	0	0
12	X	248	0	61	1	0
13	Y	148	0	40	0	0
14	Z	193	0	49	0	0
All	All	20228	0	5731	3	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 0.

All (3) 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)
10:V:22:THR:H	12:X:62:HIS:H	1.50	0.59
2:B:174:GLN:C	2:B:176:GLY:H	2.26	0.43
2:B:102:GLY:HA3	2:B:126:ALA:H	1.86	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	0	73/76 (96%)	73 (100%)	0	0	100	100
1	1	73/76 (96%)	70 (96%)	2 (3%)	1 (1%)	9	37
1	2	73/76 (96%)	72 (99%)	1 (1%)	0	100	100
1	3	72/76 (95%)	69 (96%)	2 (3%)	1 (1%)	9	37
1	4	73/76 (96%)	70 (96%)	2 (3%)	1 (1%)	9	37
1	5	73/76 (96%)	70 (96%)	1 (1%)	2 (3%)	4	27
1	6	72/76 (95%)	69 (96%)	2 (3%)	1 (1%)	9	37
1	7	71/76 (93%)	68 (96%)	2 (3%)	1 (1%)	9	37
1	8	73/76 (96%)	70 (96%)	2 (3%)	1 (1%)	9	37
1	9	72/76 (95%)	69 (96%)	2 (3%)	1 (1%)	9	37
2	A	495/510 (97%)	467 (94%)	25 (5%)	3 (1%)	21	54
2	B	501/510 (98%)	468 (93%)	27 (5%)	6 (1%)	10	40
2	C	496/510 (97%)	474 (96%)	17 (3%)	5 (1%)	12	43
3	D	468/478 (98%)	442 (94%)	22 (5%)	4 (1%)	14	45
3	E	466/478 (98%)	444 (95%)	15 (3%)	7 (2%)	8	36
3	F	467/478 (98%)	427 (91%)	33 (7%)	7 (2%)	8	36
4	G	261/278 (94%)	245 (94%)	9 (3%)	7 (3%)	4	27
5	H	110/138 (80%)	100 (91%)	8 (7%)	2 (2%)	6	33
6	I	42/61 (69%)	41 (98%)	1 (2%)	0	100	100
7	O	185/195 (95%)	171 (92%)	9 (5%)	5 (3%)	4	27
8	T	222/249 (89%)	212 (96%)	7 (3%)	3 (1%)	9	37
9	U	153/209 (73%)	152 (99%)	1 (1%)	0	100	100
10	V	169/173 (98%)	156 (92%)	8 (5%)	5 (3%)	3	25
11	W	83/95 (87%)	70 (84%)	5 (6%)	8 (10%)	0	7
12	X	60/92 (65%)	53 (88%)	5 (8%)	2 (3%)	3	24

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	Y	35/59 (59%)	34 (97%)	1 (3%)	0	100	100
14	Z	46/48 (96%)	43 (94%)	3 (6%)	0	100	100
All	All	4984/5321 (94%)	4699 (94%)	212 (4%)	73 (2%)	11	36

All (73) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	41	PRO
1	3	41	PRO
1	4	41	PRO
1	5	41	PRO
1	6	41	PRO
1	7	41	PRO
1	8	41	PRO
1	9	41	PRO
2	A	229	LYS
2	B	59	SER
2	B	338	ALA
3	D	28	SER
3	D	29	GLU
3	E	279	VAL
3	E	358	LEU
5	H	36	SER
2	B	4	LYS
2	B	154	ALA
2	B	167	GLU
2	C	141	ARG
3	E	28	SER
3	F	64	MET
3	F	451	ASN
4	G	58	LYS
4	G	72	GLU
7	O	61	ALA
10	V	130	PHE
10	V	169	ASN
11	W	11	SER
11	W	31	TYR
2	A	381	GLN
2	B	26	ASN
2	C	71	GLY
2	C	145	HIS

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Mol	Chain	Res	Type
3	E	34	LEU
3	F	131	ALA
5	H	33	PRO
7	O	51	PRO
10	V	89	LEU
11	W	39	ALA
11	W	42	ILE
11	W	44	ALA
12	X	38	ASN
1	5	42	SER
2	A	88	GLU
3	D	17	ILE
3	E	278	ALA
3	F	86	PRO
3	F	359	ASP
3	F	415	SER
4	G	131	ASN
7	O	10	VAL
7	O	16	GLU
8	T	46	LEU
8	T	202	ASN
12	X	36	LYS
3	E	16	VAL
3	E	256	ASP
4	G	56	GLU
10	V	66	SER
11	W	40	PRO
11	W	41	ALA
2	C	52	GLU
3	F	279	VAL
4	G	191	SER
10	V	132	GLU
4	G	181	PRO
8	T	147	PRO
3	D	86	PRO
11	W	5	ILE
2	C	146	GLU
7	O	165	GLY
4	G	123	PRO

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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.

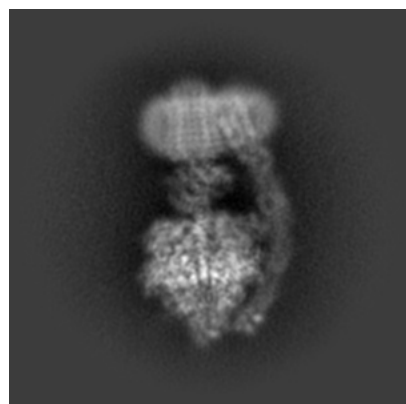
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-25946. These allow visual inspection of the internal detail of the map and identification of artifacts.

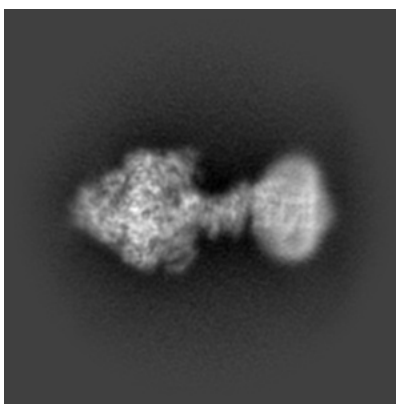
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

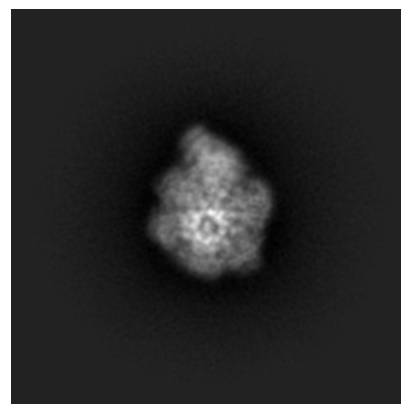
6.1.1 Primary map



X

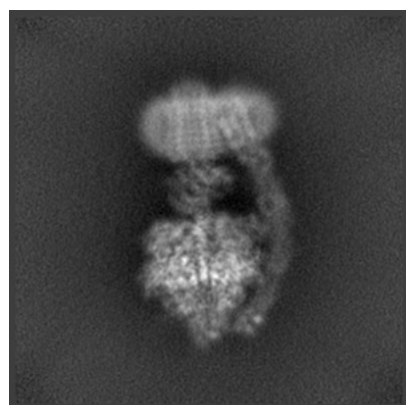


Y

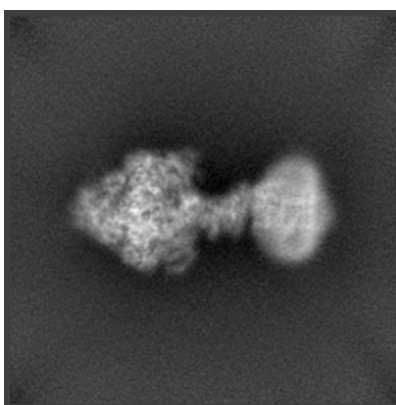


Z

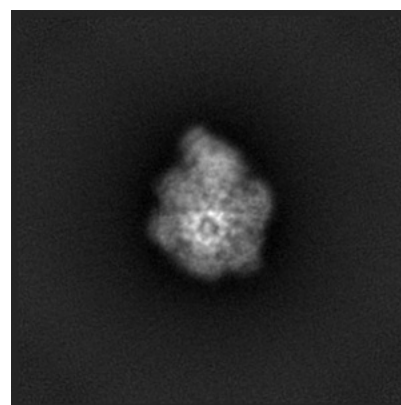
6.1.2 Raw map



X



Y

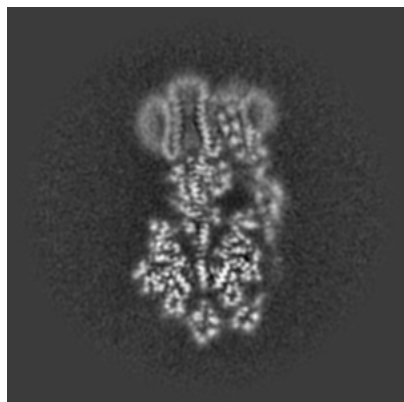


Z

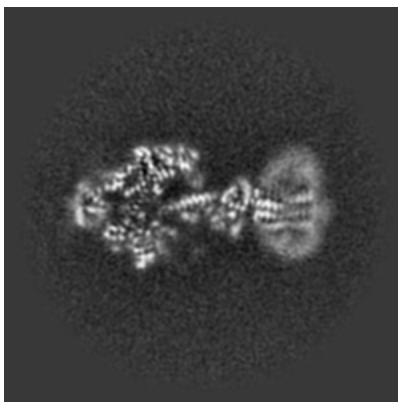
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

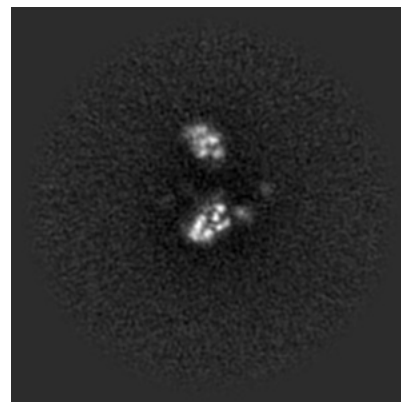
6.2.1 Primary map



X Index: 128

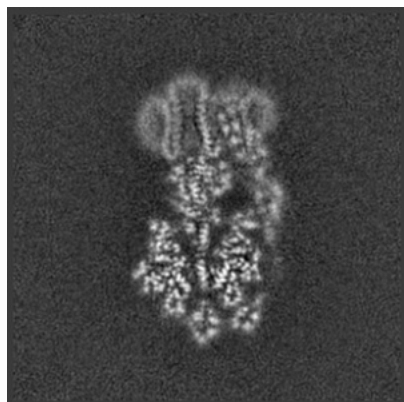


Y Index: 128

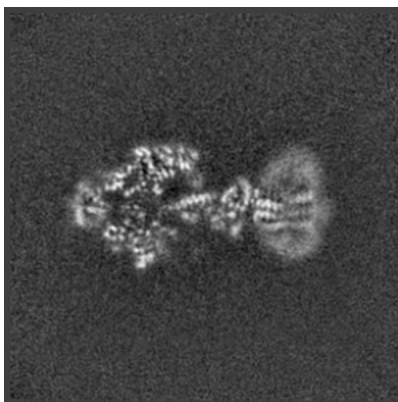


Z Index: 128

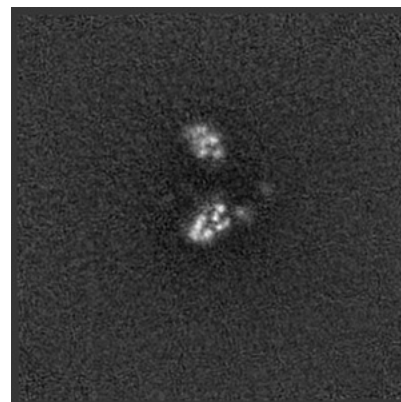
6.2.2 Raw map



X Index: 128



Y Index: 128

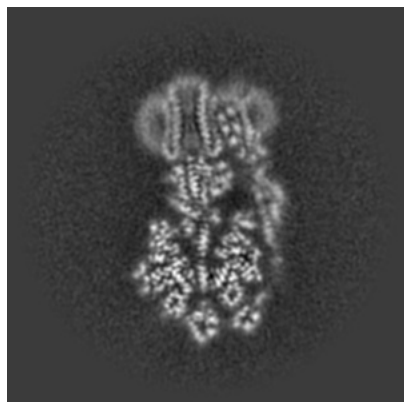


Z Index: 128

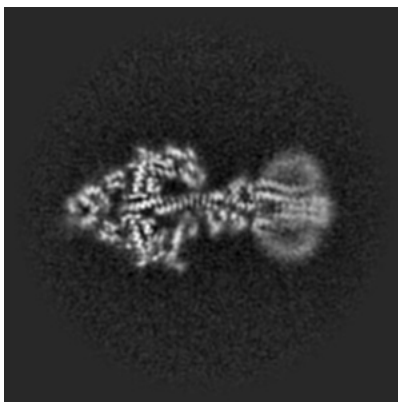
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

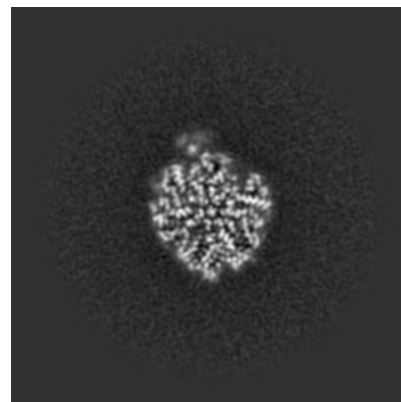
6.3.1 Primary map



X Index: 127

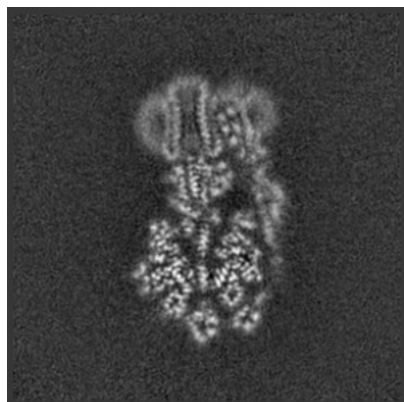


Y Index: 124

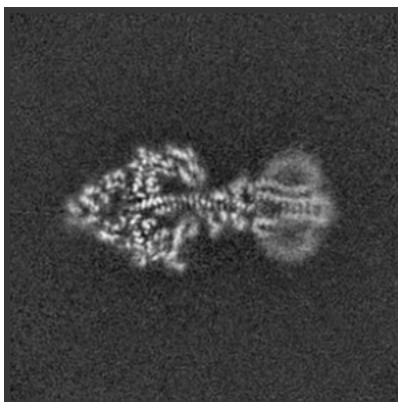


Z Index: 85

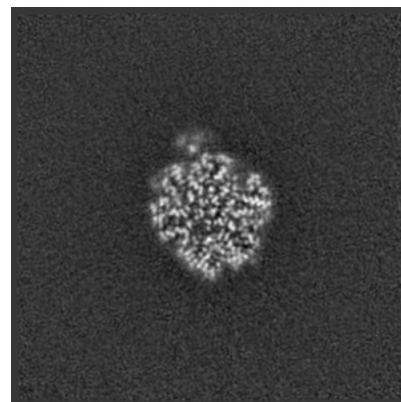
6.3.2 Raw map



X Index: 127



Y Index: 123

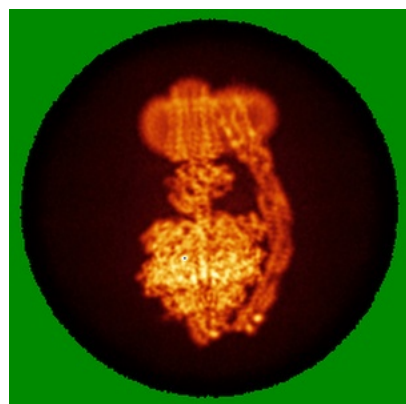


Z Index: 86

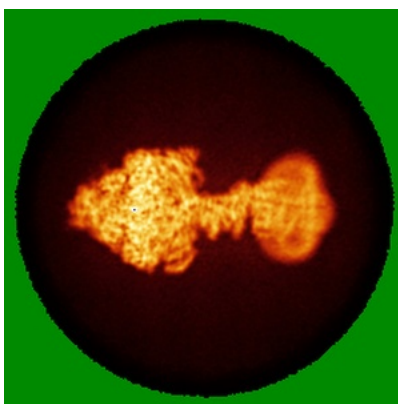
The images above show the largest variance slices of the map in three orthogonal directions.

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

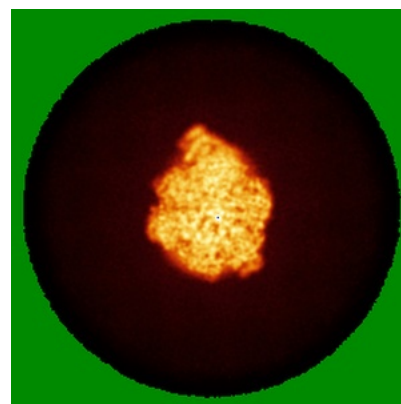
6.4.1 Primary map



X

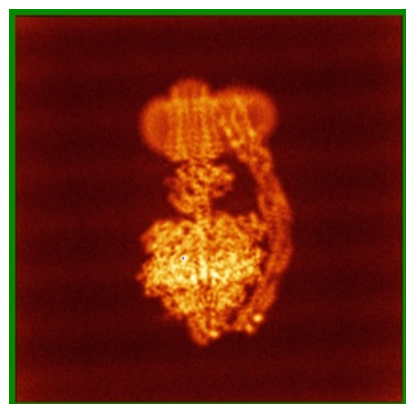


Y

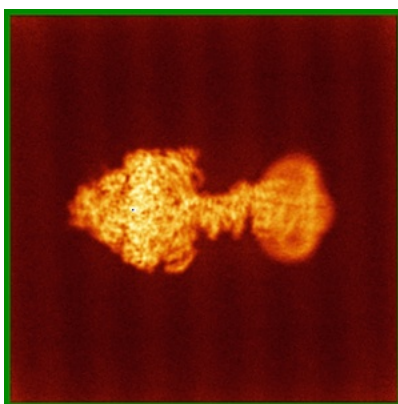


Z

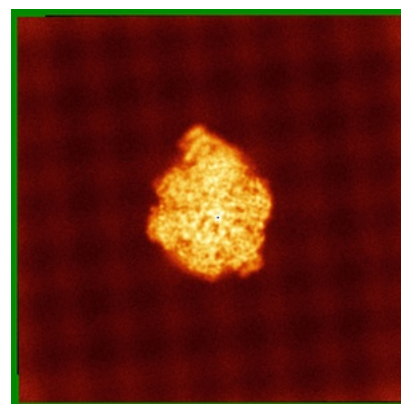
6.4.2 Raw map



X



Y

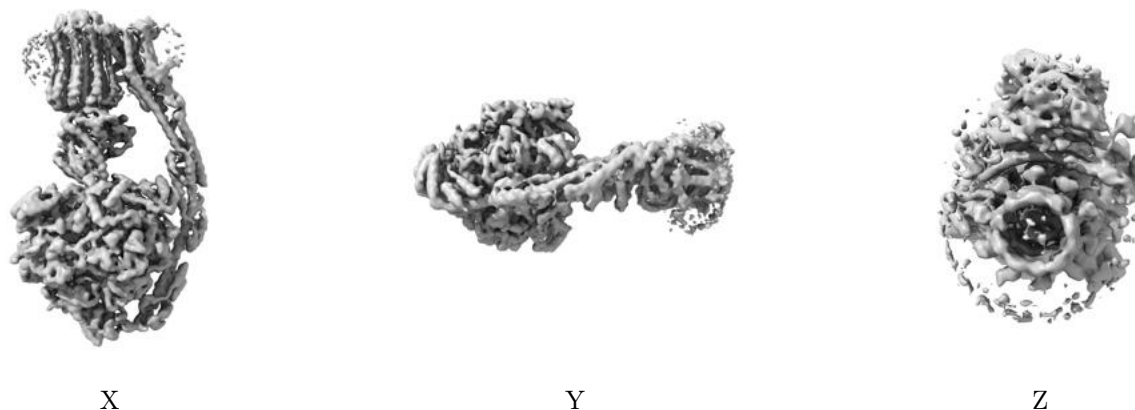


Z

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.

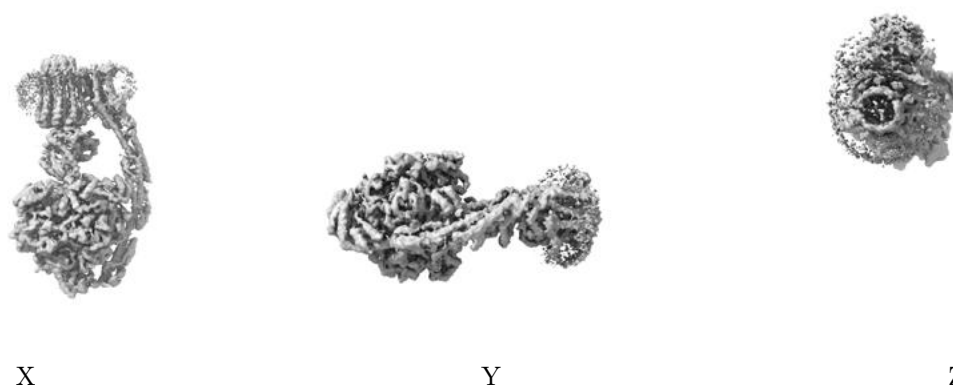
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.7. 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.

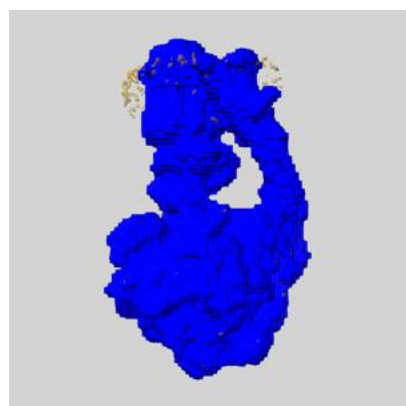
6.6 Mask visualisation [i](#)

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

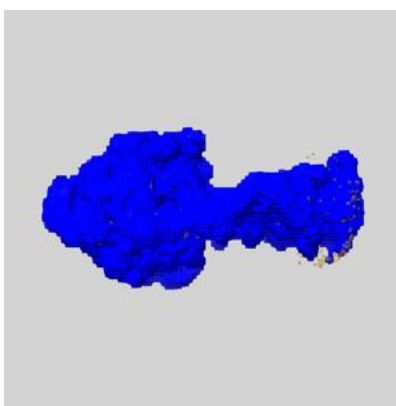
A mask typically either:

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

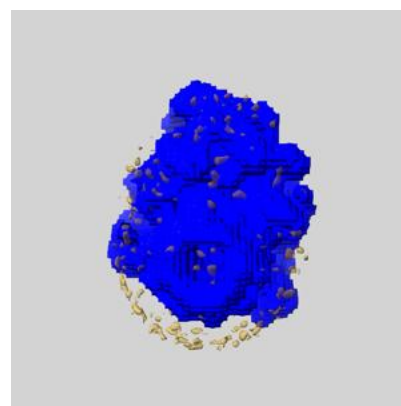
6.6.1 emd_25946_msk_1.map [i](#)



X



Y

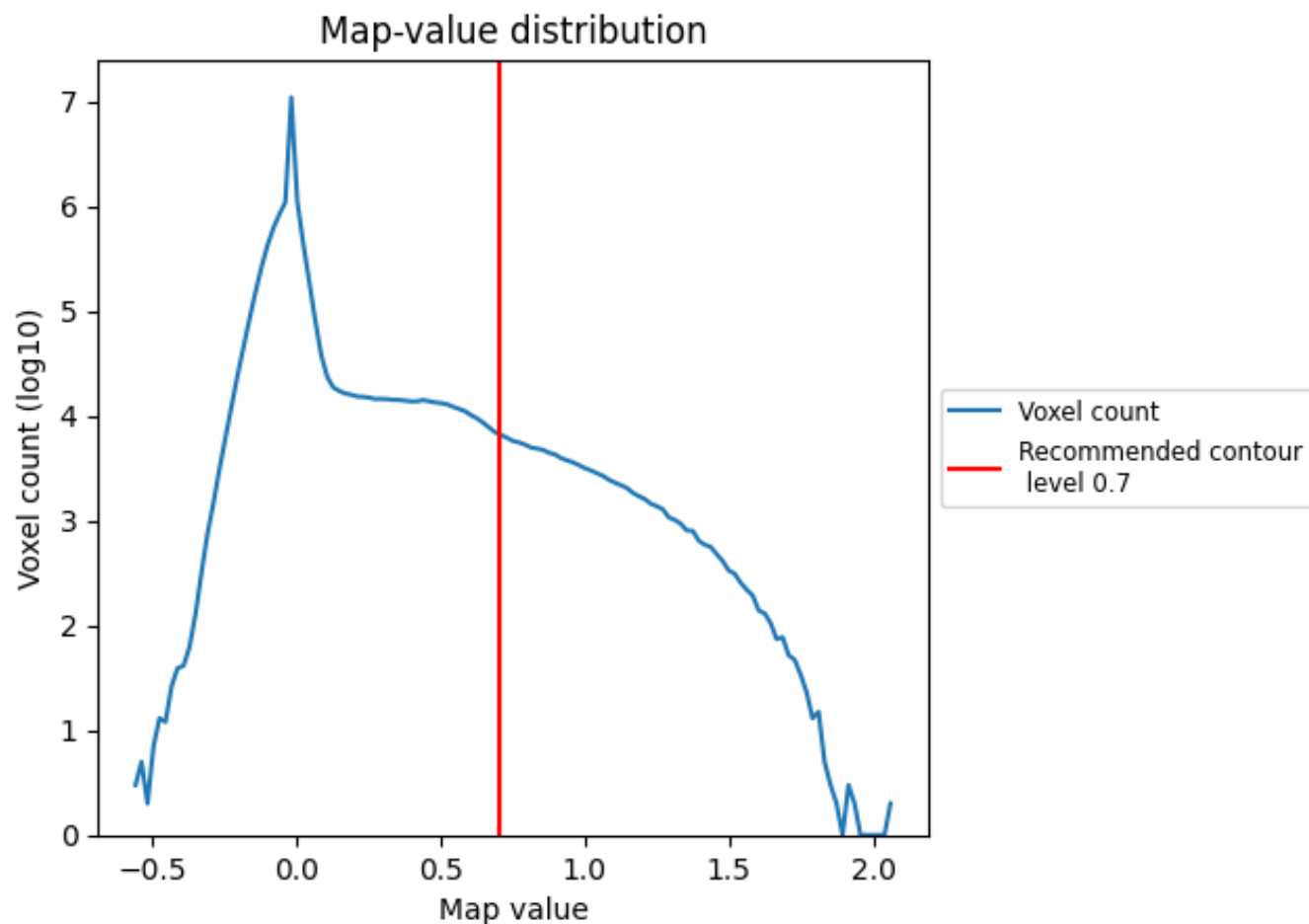


Z

7 Map analysis [i](#)

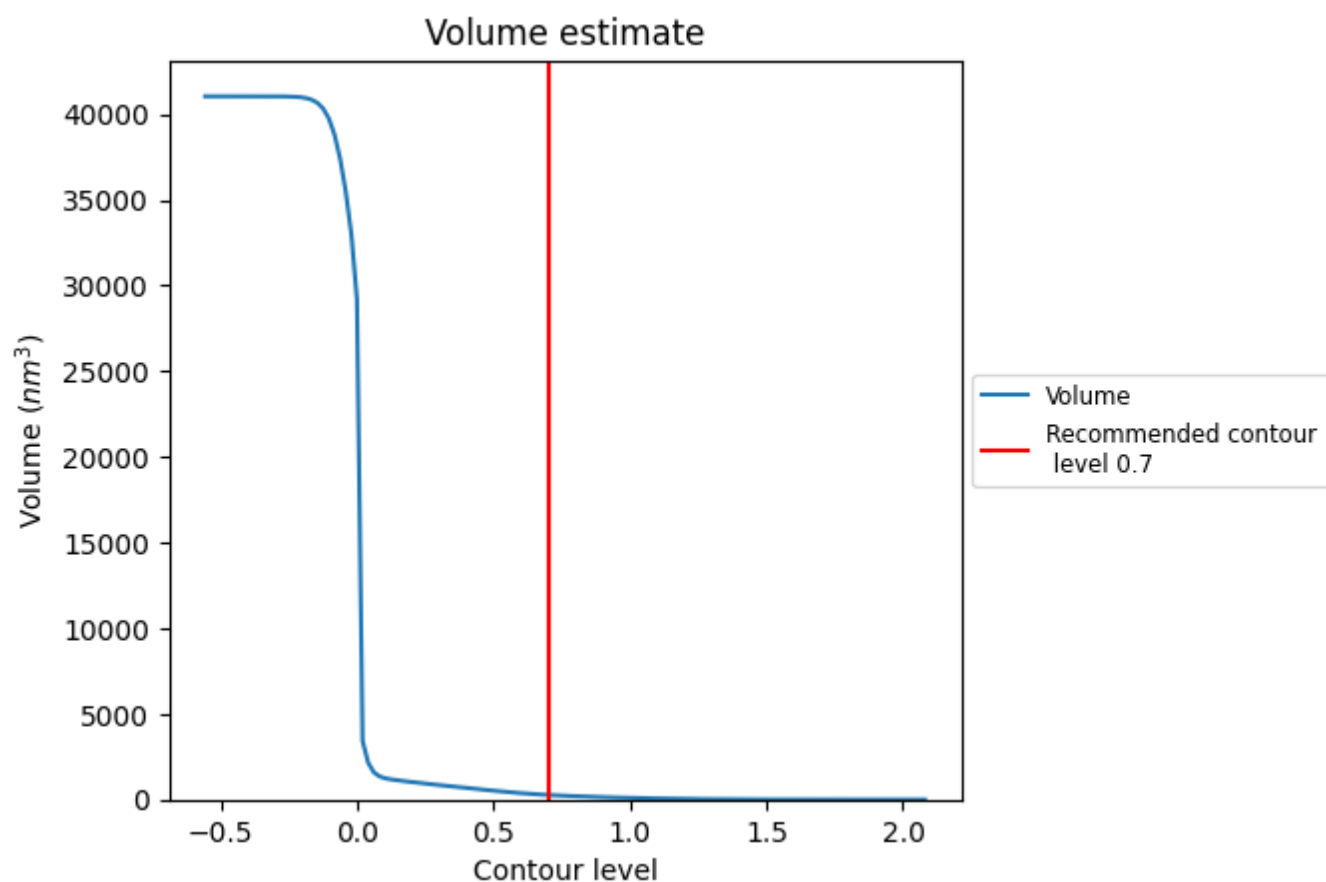
This section contains the results of statistical analysis of the map.

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.

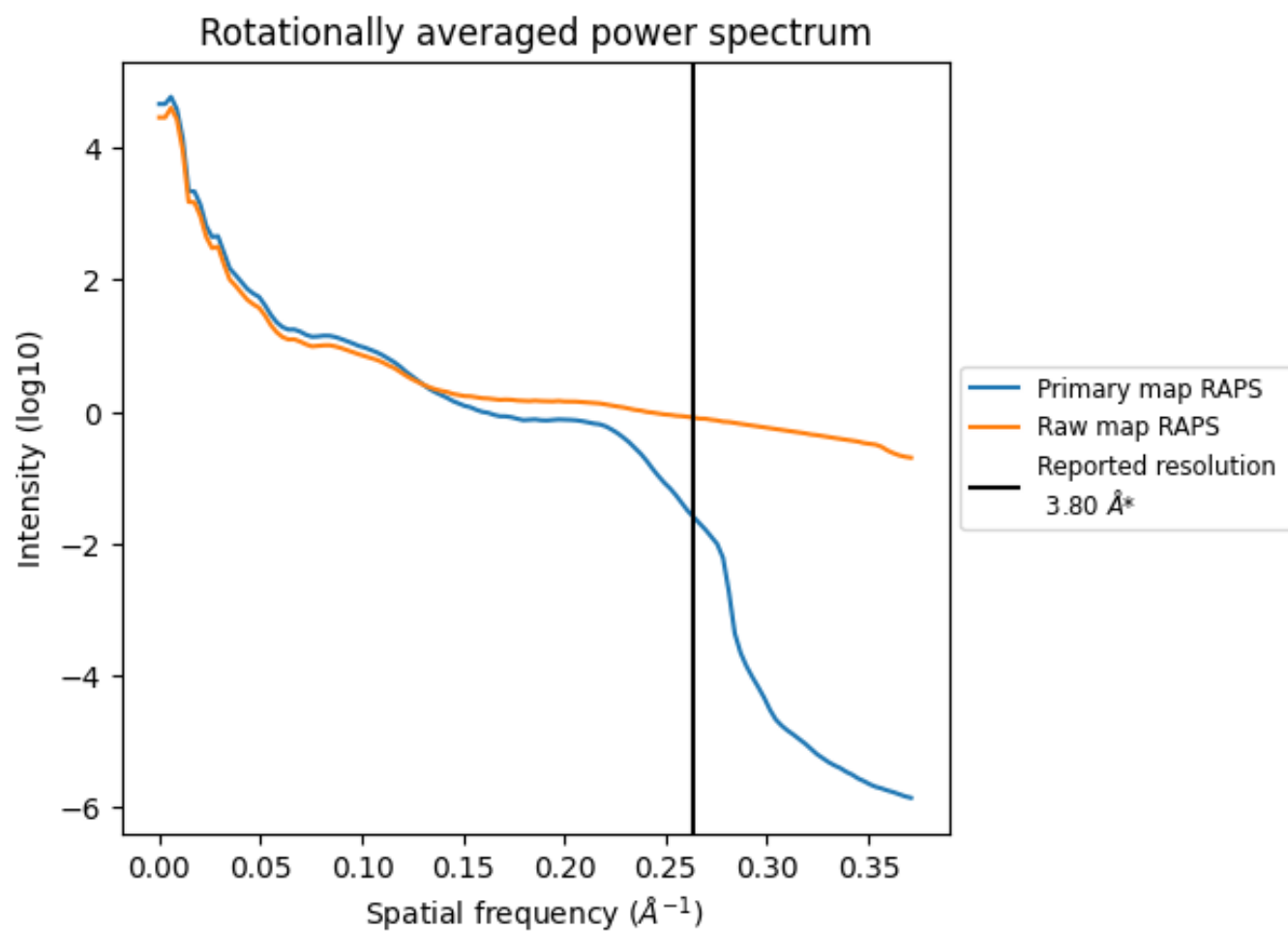
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 268 nm³; this corresponds to an approximate mass of 242 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 ⓘ

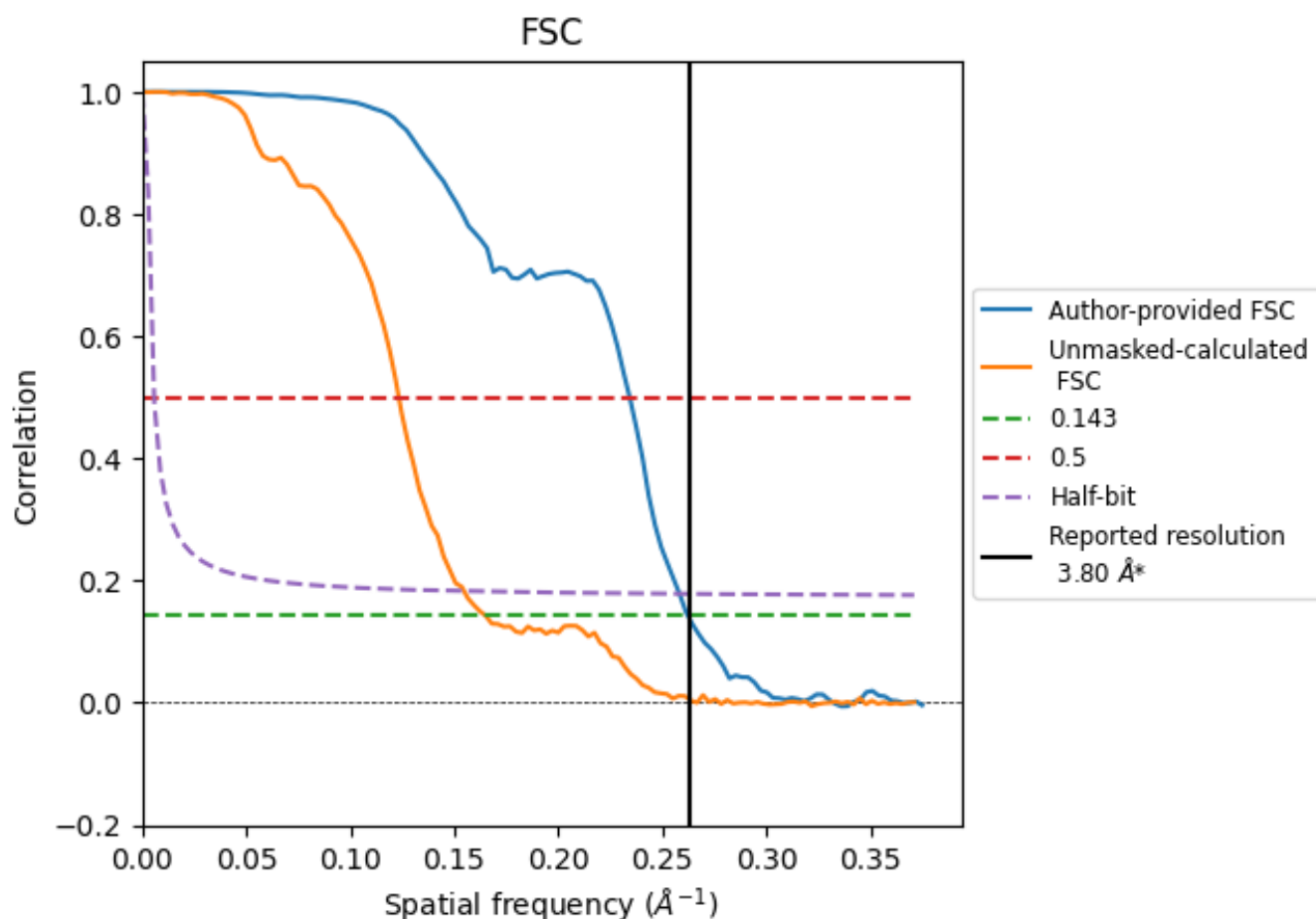


*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

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 $0.263 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

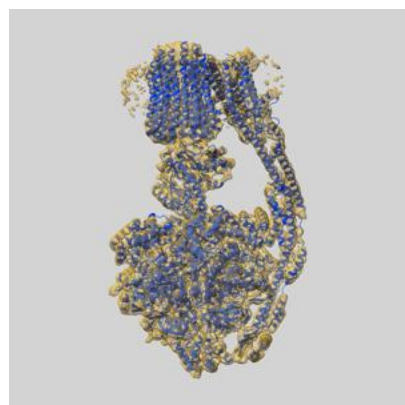
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.81	4.27	3.87
Unmasked-calculated*	6.07	8.11	6.46

*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 6.07 differs from the reported value 3.8 by more than 10 %

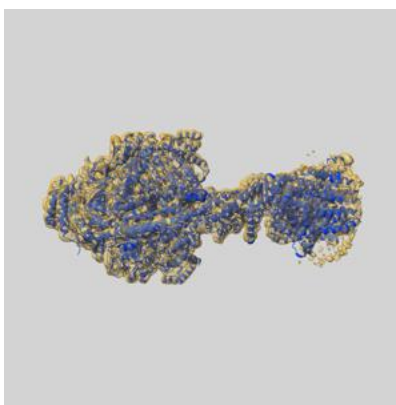
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-25946 and PDB model 7TJY. Per-residue inclusion information can be found in section [3](#) on page [7](#).

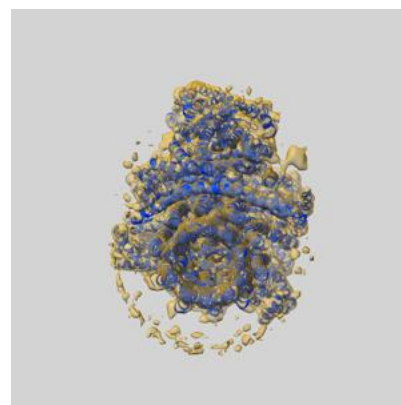
9.1 Map-model overlay [i](#)



X



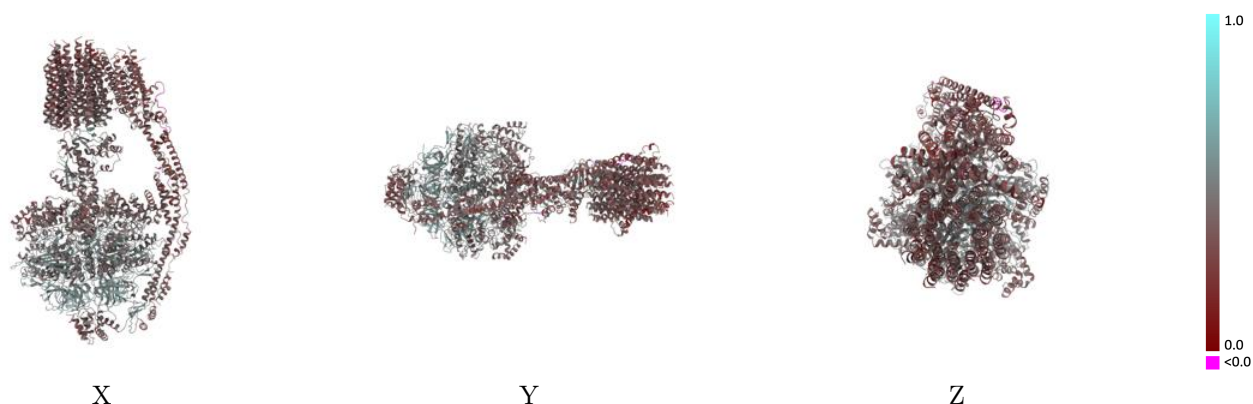
Y



Z

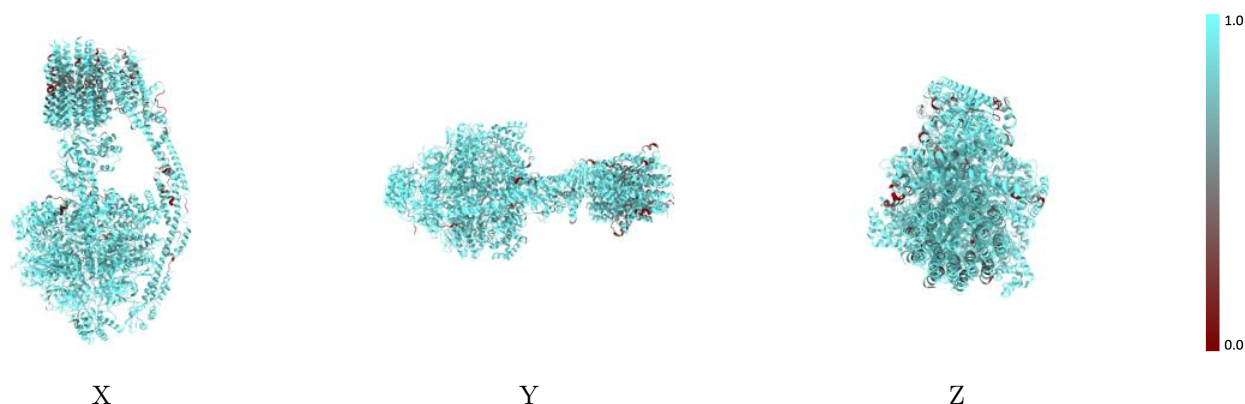
The images above show the 3D surface view of the map at the recommended contour level 0.7 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



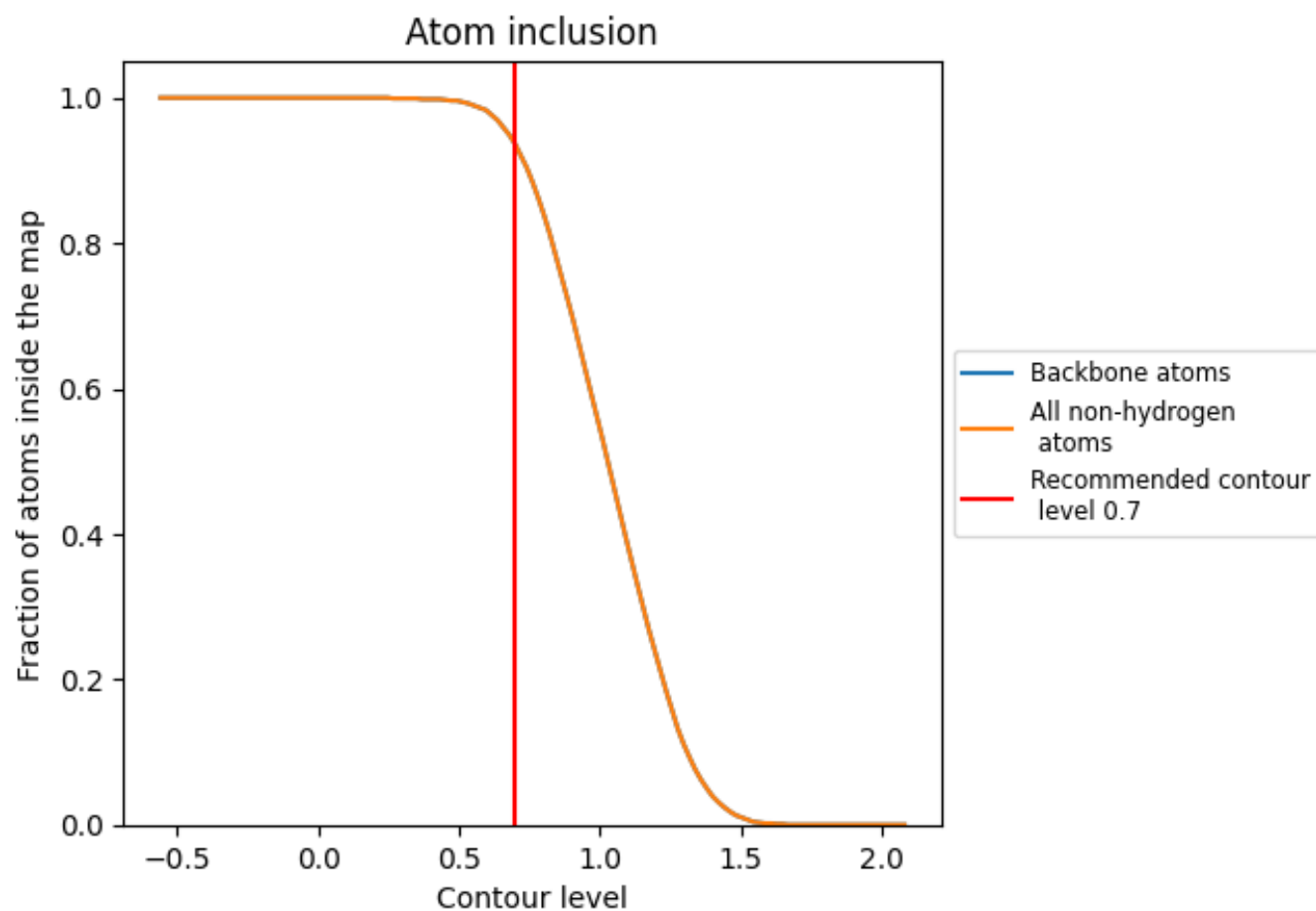
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 94% 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 (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9360	 0.4210
0	 0.8530	 0.3460
1	 0.9000	 0.3580
2	 0.9330	 0.3650
3	 0.9120	 0.3380
4	 0.8400	 0.3430
5	 0.8470	 0.3200
6	 0.8450	 0.3170
7	 0.8220	 0.3260
8	 0.8070	 0.3280
9	 0.8750	 0.3370
A	 0.9880	 0.4720
B	 0.9650	 0.4620
C	 0.9710	 0.4700
D	 0.9540	 0.4640
E	 0.9810	 0.4760
F	 0.9920	 0.4840
G	 0.9840	 0.4240
H	 0.9560	 0.4470
I	 1.0000	 0.4000
O	 0.9950	 0.4040
T	 0.8370	 0.3210
U	 0.9730	 0.3410
V	 0.8470	 0.3220
W	 0.6970	 0.2640
X	 0.6450	 0.3090
Y	 0.5680	 0.2820
Z	 0.8500	 0.3160

