



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 09:10 PM UTC

PDB ID : 7M2X / pdb_00007m2x
EMDB ID : EMD-23636
Title : Open conformation of the Yeast wild-type gamma-TuRC
Authors : Brilot, A.F.; Lyon, A.S.; Zelter, A.; Viswanath, S.; Maxwell, A.; MacCoss, M.J.; Muller, E.G.; Sali, A.; Davis, T.N.; Agard, D.A.
Deposited on : 2021-03-17
Resolution : 3.60 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:

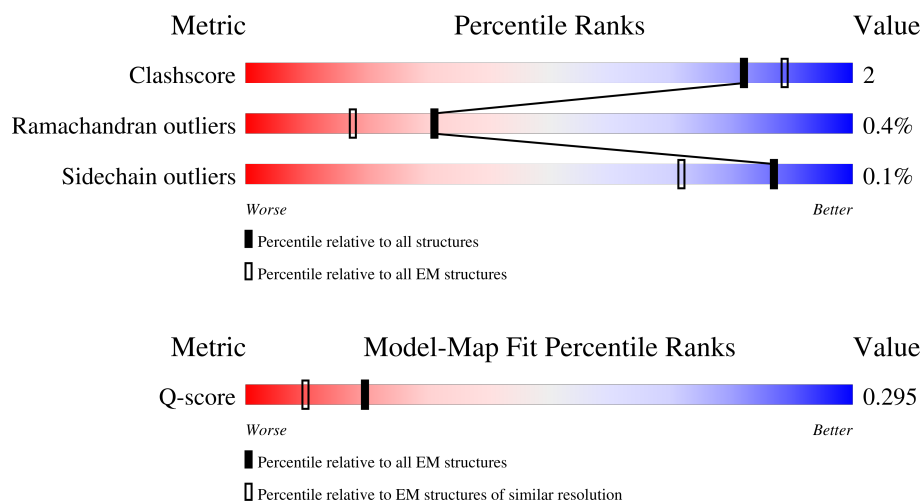
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 3.60 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	473	64% 82% 11% 6%
1	B	473	75% 80% 13% 6%
2	C	823	25% 76% 8% 15%
3	D	846	28% 74% 6% 20%

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Mol	Chain	Length	Quality of chain
4	U	220	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18704 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 Tubulin gamma chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	444	Total	C	N	O	S	0	0
			3474	2172	588	698	16		
1	B	446	Total	C	N	O	S	0	0
			3486	2179	590	700	17		

- Molecule 2 is a protein called Spindle pole body component SPC97.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	701	Total	C	N	O	S	0	0
			5853	3770	983	1071	29		

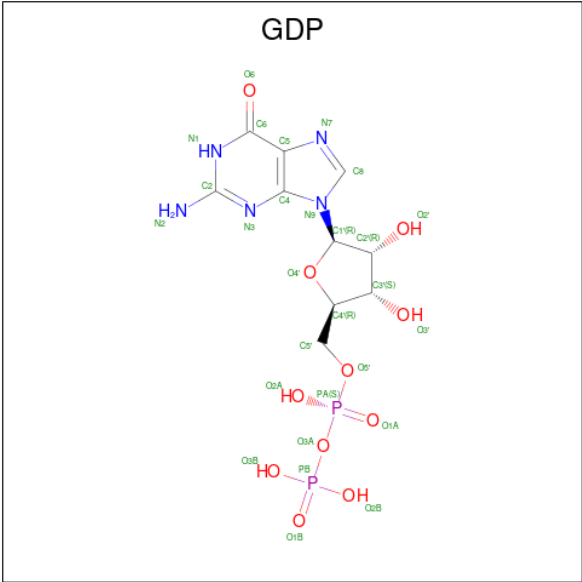
- Molecule 3 is a protein called Spindle pole body component SPC98.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	673	Total	C	N	O	S	0	0
			5560	3595	918	1031	16		

- Molecule 4 is a protein called Spindle pole body component 110.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	U	33	Total	C	N	O	S	0	0
			275	170	51	52	2		

- Molecule 5 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).

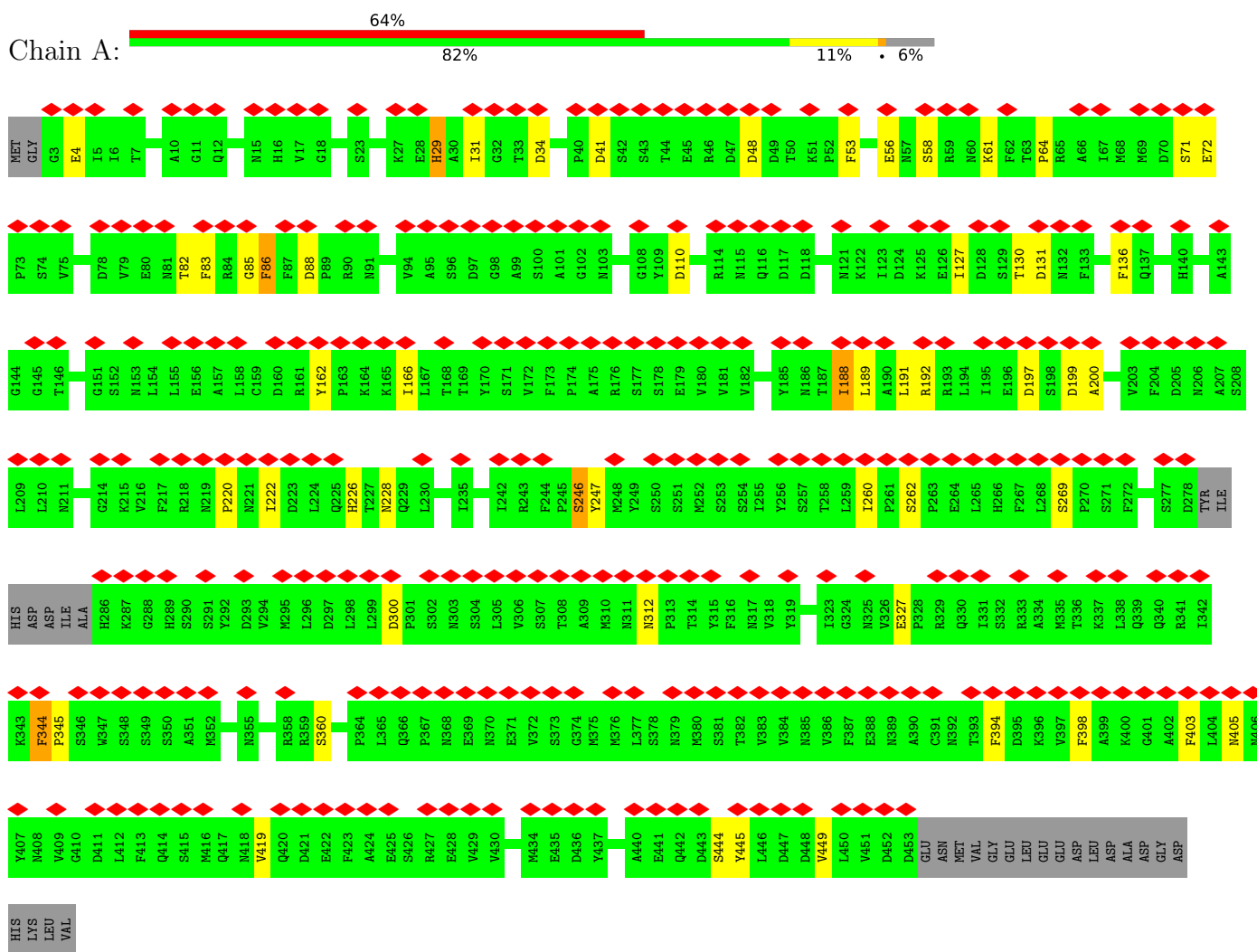


Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total	C	N	O	P	0
			28	10	5	11	2	
5	B	1	Total	C	N	O	P	0
			28	10	5	11	2	

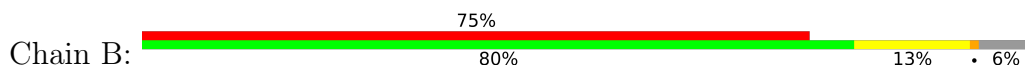
3 Residue-property plots

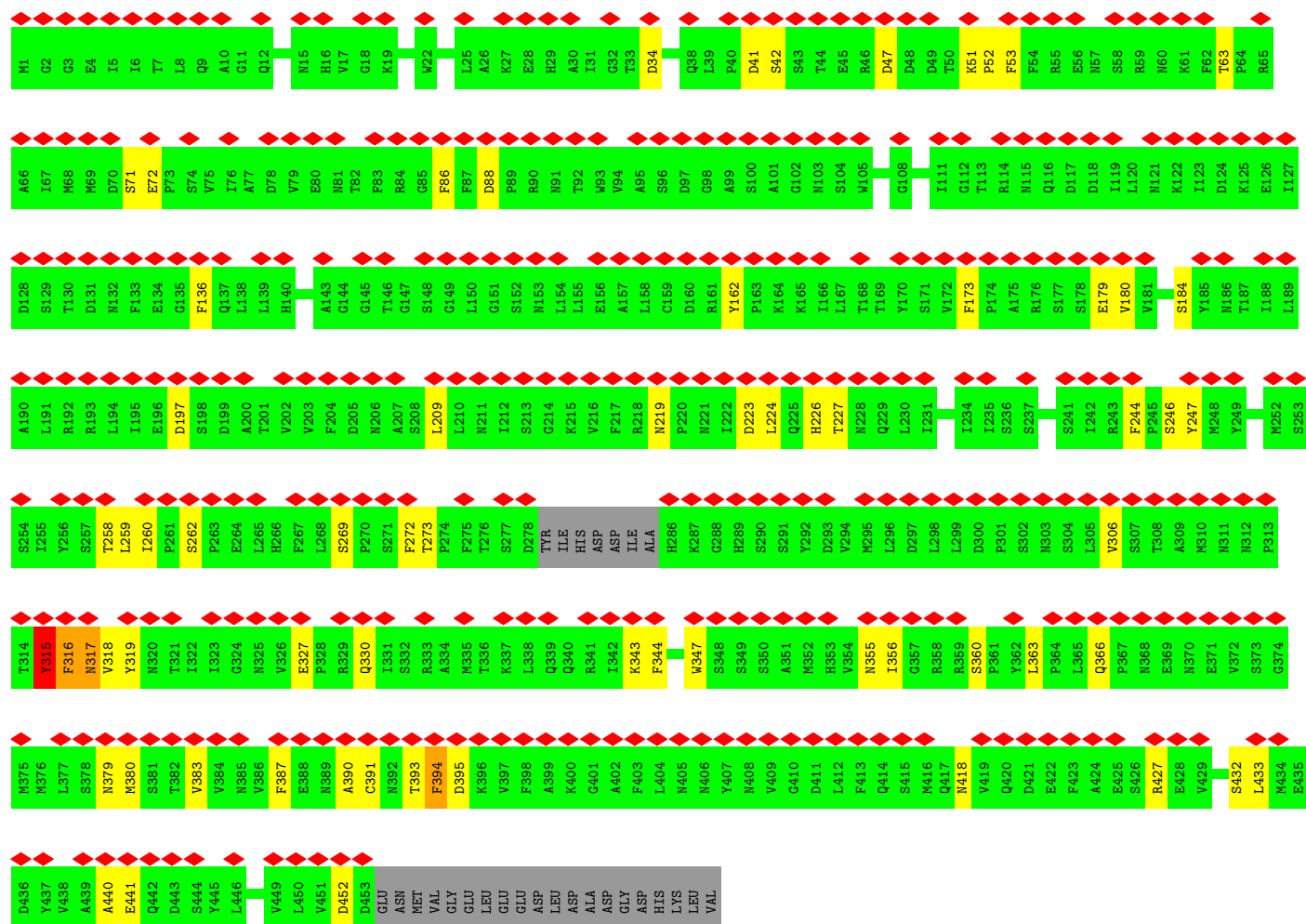
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

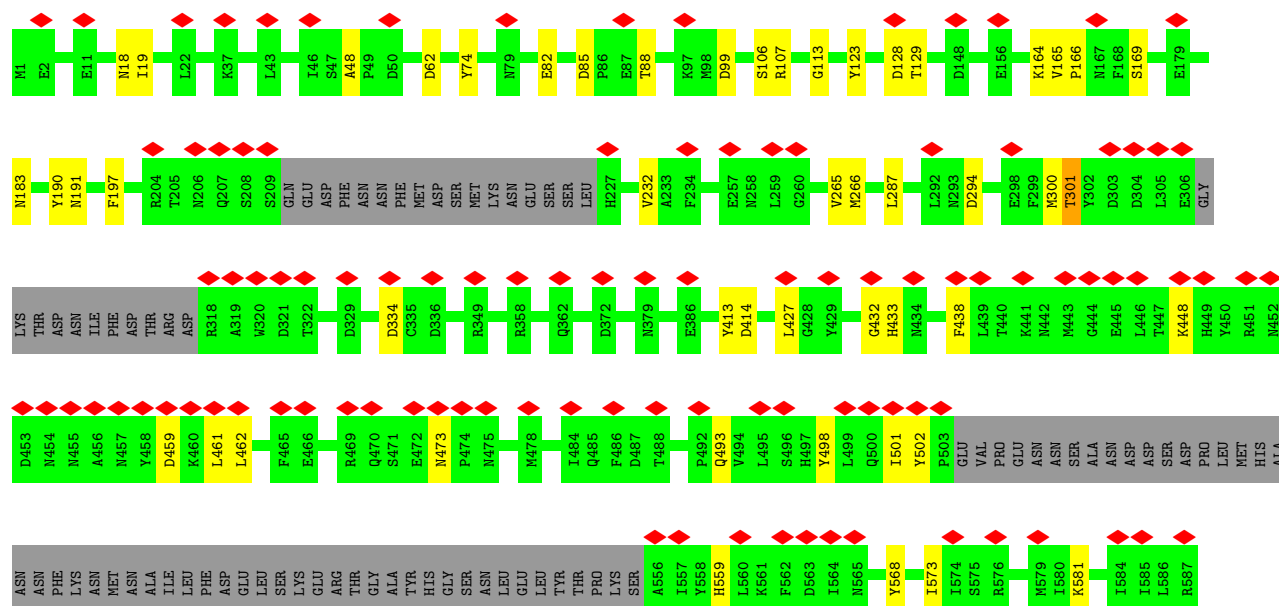
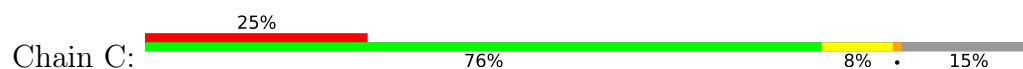


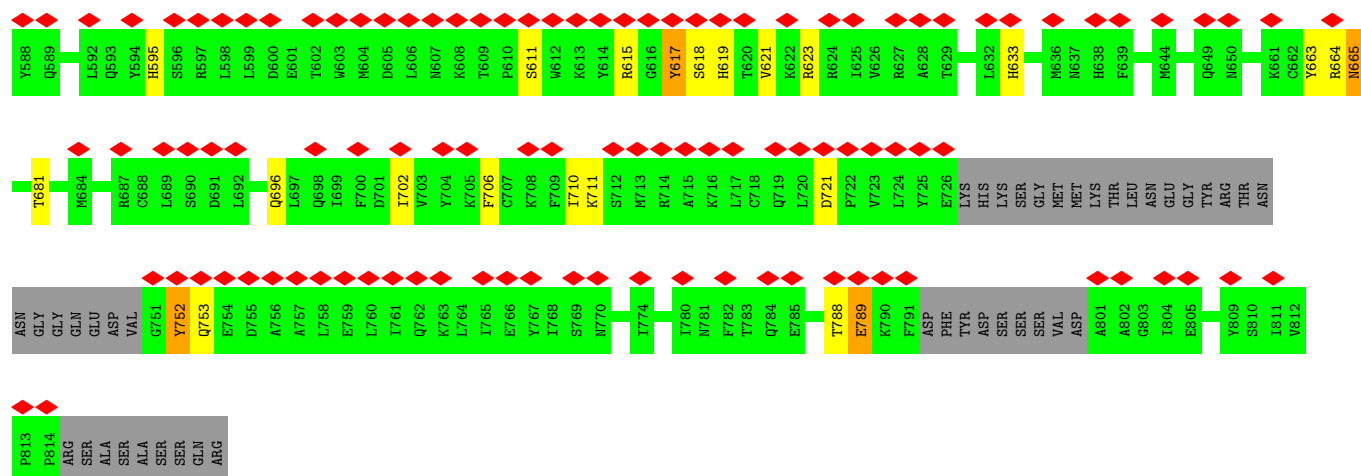
- Molecule 1: Tubulin gamma chain



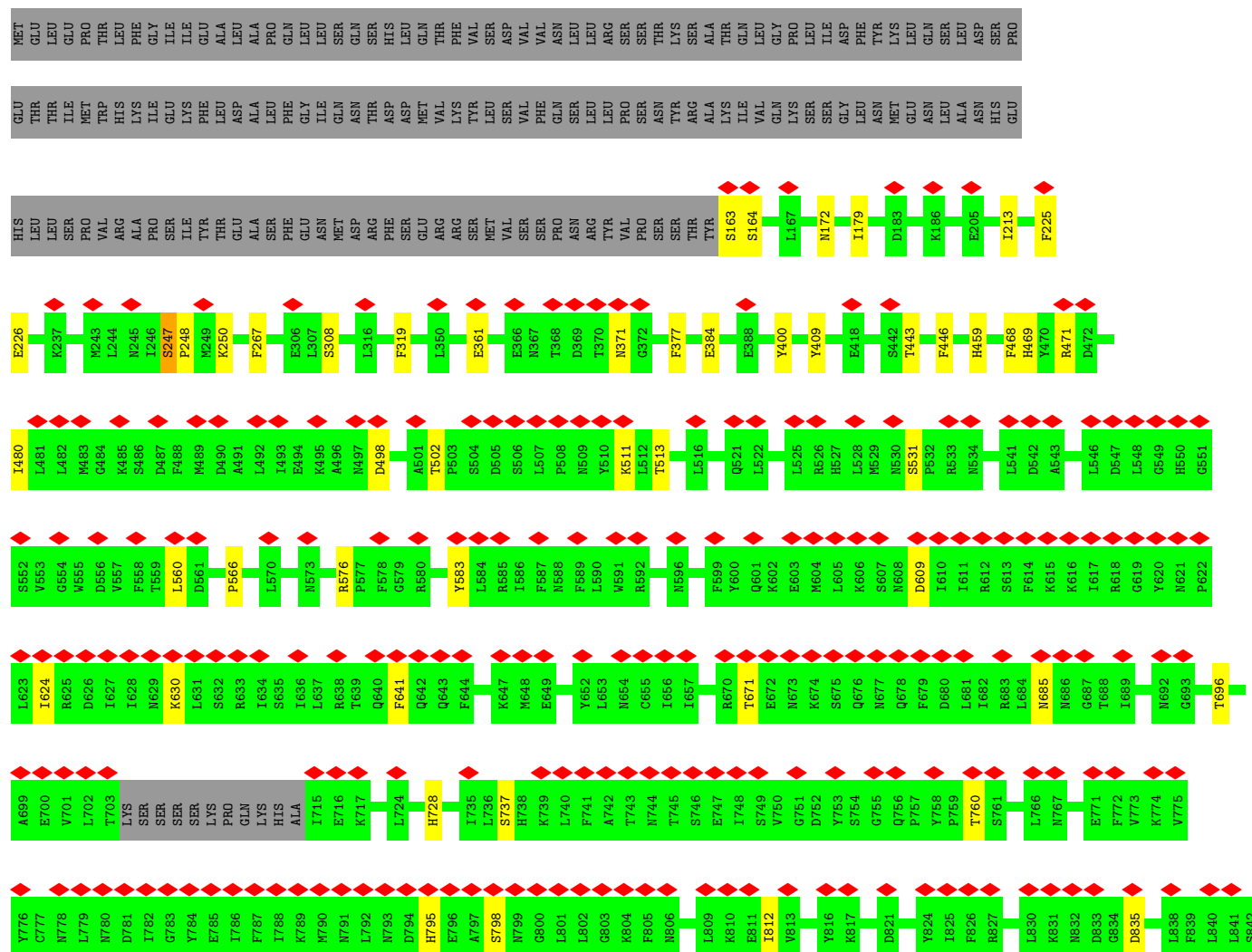


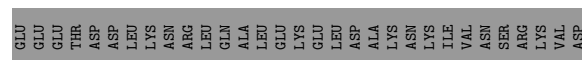
• Molecule 2: Spindle pole body component SPC97





• Molecule 3: Spindle pole body component SPC98





4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=0°, rise=0.01 Å, axial sym=C1	Depositor
Number of segments used	36251	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION; Final reconstruction in cis-TEM	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	80	Depositor
Minimum defocus (nm)	900	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	47214	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	23.834	Depositor
Minimum map value	-13.249	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.773	Depositor
Recommended contour level	7	Depositor
Map size (Å)	635.39844, 635.39844, 635.39844	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.4183, 1.4183, 1.4183	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GDP

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	1.19	8/3546 (0.2%)	1.36	46/4815 (1.0%)
1	B	1.12	1/3558 (0.0%)	1.38	56/4830 (1.2%)
2	C	1.29	7/5969 (0.1%)	1.27	52/8057 (0.6%)
3	D	1.30	6/5679 (0.1%)	1.22	34/7677 (0.4%)
4	U	1.17	0/278	1.21	3/368 (0.8%)
All	All	1.24	22/19030 (0.1%)	1.29	191/25747 (0.7%)

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	226	HIS	CB-CG	-7.07	1.40	1.50
2	C	559	HIS	CB-CG	-6.83	1.40	1.50
1	A	246	SER	CA-C	-6.77	1.50	1.53
1	A	29	HIS	CB-CG	-6.62	1.40	1.50
3	D	583	TYR	CB-CG	-6.18	1.38	1.51
1	A	29	HIS	CA-C	-6.11	1.45	1.52
2	C	498	TYR	CB-CG	-6.05	1.38	1.51
3	D	267	PHE	CB-CG	-5.89	1.37	1.50
2	C	633	HIS	CB-CG	-5.78	1.42	1.50
1	A	398	PHE	CG-CD1	-5.71	1.26	1.38
1	B	347	TRP	NE1-CE2	-5.67	1.31	1.37
2	C	197	PHE	CB-CG	-5.64	1.37	1.50
1	A	64	PRO	N-CD	-5.59	1.40	1.47
2	C	559	HIS	CG-CD2	-5.26	1.30	1.35
3	D	728	HIS	CB-CG	-5.21	1.42	1.50
1	A	4	GLU	CG-CD	-5.21	1.39	1.52
2	C	190	TYR	CB-CG	-5.19	1.40	1.51
3	D	400	TYR	CB-CG	-5.14	1.40	1.51
3	D	459	HIS	CB-CG	-5.12	1.43	1.50
3	D	795	HIS	CB-CG	-5.12	1.43	1.50
1	A	228	ASN	CB-CG	-5.10	1.39	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	433	HIS	CB-CG	-5.07	1.43	1.50

All (191) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	611	SER	N-CA-C	-11.87	98.90	113.50
1	A	86	PHE	CA-CB-CG	11.15	124.95	113.80
1	B	383	VAL	N-CA-C	-10.99	103.27	113.71
2	C	427	LEU	N-CA-C	-10.43	100.62	113.97
2	C	165	VAL	N-CA-C	9.74	118.84	107.73
3	D	566	PRO	N-CA-C	-8.96	100.99	110.58
2	C	615	ARG	N-CA-C	-8.73	104.67	114.62
1	B	42	SER	N-CA-C	-8.67	103.12	114.31
1	B	86	PHE	CA-CB-CG	8.41	122.21	113.80
1	B	71	SER	N-CA-C	-8.36	103.26	113.97
1	B	226	HIS	N-CA-C	-7.73	103.81	113.55
1	B	315	TYR	N-CA-C	7.71	127.22	110.80
1	A	71	SER	N-CA-C	-7.58	104.53	114.31
2	C	62	ASP	CA-CB-CG	7.56	120.16	112.60
3	D	377	PHE	CA-CB-CG	-7.35	106.45	113.80
2	C	473	ASN	CA-C-N	7.21	126.92	119.56
2	C	473	ASN	C-N-CA	7.21	126.92	119.56
1	B	269	SER	CA-C-N	7.19	127.36	120.31
1	B	269	SER	C-N-CA	7.19	127.36	120.31
1	B	88	ASP	CA-C-N	7.14	126.84	119.56
1	B	88	ASP	C-N-CA	7.14	126.84	119.56
3	D	247	SER	CA-C-N	7.09	126.79	119.56
3	D	247	SER	C-N-CA	7.09	126.79	119.56
1	B	197	ASP	N-CA-C	7.01	121.96	113.41
3	D	213	ILE	CA-C-N	6.98	127.38	119.92
3	D	213	ILE	C-N-CA	6.98	127.38	119.92
2	C	721	ASP	CA-C-N	6.96	126.66	119.56
2	C	721	ASP	C-N-CA	6.96	126.66	119.56
1	A	360	SER	CA-C-N	6.92	126.61	119.56
1	A	360	SER	C-N-CA	6.92	126.61	119.56
3	D	624	ILE	N-CA-C	6.87	117.47	111.90
1	B	272	PHE	N-CA-C	6.81	119.42	109.07
1	B	260	ILE	CA-C-N	6.79	126.49	119.56
1	B	260	ILE	C-N-CA	6.79	126.49	119.56
1	A	246	SER	CA-C-O	6.71	121.83	117.94
2	C	665	ASN	CA-C-N	6.70	126.61	119.85
2	C	665	ASN	C-N-CA	6.70	126.61	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	82	THR	N-CA-C	-6.68	103.80	112.68
1	A	312	ASN	CA-C-N	6.68	126.64	120.03
1	A	312	ASN	C-N-CA	6.68	126.64	120.03
1	A	162	TYR	CA-C-N	6.66	126.28	119.56
1	A	162	TYR	C-N-CA	6.66	126.28	119.56
2	C	85	ASP	CA-C-N	6.55	126.24	119.56
2	C	85	ASP	C-N-CA	6.55	126.24	119.56
2	C	232	VAL	N-CA-C	6.53	117.10	106.72
1	B	316	PHE	CA-C-N	6.49	133.93	121.54
1	B	316	PHE	C-N-CA	6.49	133.93	121.54
3	D	468	PHE	CA-CB-CG	-6.47	107.33	113.80
1	A	72	GLU	CA-C-N	6.46	126.15	119.56
1	A	72	GLU	C-N-CA	6.46	126.15	119.56
1	B	327	GLU	CA-C-N	6.46	126.97	119.47
1	B	327	GLU	C-N-CA	6.46	126.97	119.47
2	C	169	SER	N-CA-C	6.43	118.55	108.96
1	A	48	ASP	CA-CB-CG	6.42	119.02	112.60
2	C	448	LYS	N-CA-C	6.36	121.44	113.43
1	A	260	ILE	CA-C-N	6.35	126.90	120.04
1	A	260	ILE	C-N-CA	6.35	126.90	120.04
2	C	789	GLU	N-CA-C	-6.29	100.72	110.10
2	C	432	GLY	CA-C-N	6.25	128.57	120.44
2	C	432	GLY	C-N-CA	6.25	128.57	120.44
1	A	344	PHE	CA-C-N	6.21	126.18	120.03
1	A	344	PHE	C-N-CA	6.21	126.18	120.03
1	A	197	ASP	N-CA-C	6.18	121.88	113.97
1	A	88	ASP	CA-C-N	6.16	126.28	119.87
1	A	88	ASP	C-N-CA	6.16	126.28	119.87
3	D	576	ARG	CA-C-N	6.14	125.83	119.56
3	D	576	ARG	C-N-CA	6.14	125.83	119.56
1	B	63	THR	CA-C-N	6.12	126.14	119.90
1	B	63	THR	C-N-CA	6.12	126.14	119.90
2	C	621	VAL	N-CA-C	-6.10	107.00	112.12
1	B	360	SER	CA-C-N	6.04	126.15	119.87
1	B	360	SER	C-N-CA	6.04	126.15	119.87
3	D	469	HIS	CA-CB-CG	-6.04	107.76	113.80
1	A	226	HIS	N-CA-C	-6.02	105.60	113.12
1	A	110	ASP	CA-CB-CG	5.97	118.57	112.60
1	A	188	ILE	N-CA-C	5.97	116.72	110.62
3	D	179	ILE	CA-C-N	5.97	125.94	120.03
3	D	179	ILE	C-N-CA	5.97	125.94	120.03
1	B	418	ASN	N-CA-C	-5.94	105.62	112.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	294	ASP	CA-C-N	5.91	125.58	119.56
2	C	294	ASP	C-N-CA	5.91	125.58	119.56
1	A	222	ILE	N-CA-C	-5.88	105.65	111.88
1	A	269	SER	CA-C-N	5.82	126.01	120.31
1	A	269	SER	C-N-CA	5.82	126.01	120.31
1	B	363	LEU	CA-C-N	5.81	125.77	119.78
1	B	363	LEU	C-N-CA	5.81	125.77	119.78
1	A	56	GLU	N-CA-C	-5.80	101.46	110.10
3	D	319	PHE	CA-CB-CG	-5.79	108.01	113.80
1	B	306	VAL	CA-C-N	-5.78	113.38	121.99
1	B	306	VAL	C-N-CA	-5.78	113.38	121.99
1	A	58	SER	N-CA-C	-5.76	105.75	112.89
3	D	513	THR	N-CA-C	-5.75	104.91	111.07
1	B	318	VAL	N-CA-C	5.72	116.12	108.11
1	A	403	PHE	CA-CB-CG	-5.70	108.10	113.80
1	B	173	PHE	CA-C-N	5.69	125.62	119.76
1	B	173	PHE	C-N-CA	5.69	125.62	119.76
3	D	760	THR	N-CA-C	-5.68	105.00	111.14
2	C	19	ILE	CA-C-N	5.65	125.60	119.78
2	C	19	ILE	C-N-CA	5.65	125.60	119.78
2	C	752	TYR	N-CA-C	5.65	118.64	110.28
3	D	371	ASN	CA-CB-CG	-5.65	106.95	112.60
1	B	262	SER	CA-C-N	5.63	125.30	119.56
1	B	262	SER	C-N-CA	5.63	125.30	119.56
1	B	226	HIS	CA-C-N	5.57	128.30	120.28
1	B	226	HIS	C-N-CA	5.57	128.30	120.28
1	B	162	TYR	CA-C-N	5.57	125.19	119.56
1	B	162	TYR	C-N-CA	5.57	125.19	119.56
2	C	74	TYR	N-CA-C	-5.57	107.14	114.04
1	B	72	GLU	CA-C-N	5.56	125.23	119.56
1	B	72	GLU	C-N-CA	5.56	125.23	119.56
1	A	262	SER	CA-C-N	5.55	125.22	119.56
1	A	262	SER	C-N-CA	5.55	125.22	119.56
1	A	300	ASP	CA-C-N	5.54	125.21	119.56
1	A	300	ASP	C-N-CA	5.54	125.21	119.56
4	U	117	GLN	CA-C-N	5.54	125.30	119.76
4	U	117	GLN	C-N-CA	5.54	125.30	119.76
1	B	273	THR	CA-C-N	5.53	125.54	119.90
1	B	273	THR	C-N-CA	5.53	125.54	119.90
3	D	502	THR	CA-C-N	5.53	125.54	119.90
3	D	502	THR	C-N-CA	5.53	125.54	119.90
2	C	82	GLU	CA-C-N	5.50	125.17	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	82	GLU	C-N-CA	5.50	125.17	119.56
3	D	641	PHE	CA-CB-CG	-5.49	108.31	113.80
1	A	327	GLU	CA-C-N	5.48	125.15	119.56
1	A	327	GLU	C-N-CA	5.48	125.15	119.56
3	D	480	ILE	N-CA-C	5.48	116.27	111.56
2	C	573	ILE	N-CA-C	-5.47	106.47	111.67
3	D	609	ASP	CA-CB-CG	5.46	118.06	112.60
1	A	246	SER	CB-CA-C	-5.43	109.25	117.07
2	C	99	ASP	CA-C-N	5.43	125.09	119.56
2	C	99	ASP	C-N-CA	5.43	125.09	119.56
1	B	356	ILE	N-CA-C	5.41	116.90	111.00
2	C	113	GLY	N-CA-C	-5.41	105.84	112.77
3	D	696	THR	CA-C-N	5.40	125.38	120.03
3	D	696	THR	C-N-CA	5.40	125.38	120.03
1	B	366	GLN	CA-C-N	5.38	125.30	119.76
1	B	366	GLN	C-N-CA	5.38	125.30	119.76
1	A	419	VAL	N-CA-C	-5.38	106.56	111.67
3	D	384	GLU	CA-C-N	-5.37	113.51	122.21
3	D	384	GLU	C-N-CA	-5.37	113.51	122.21
3	D	531	SER	CA-C-N	5.36	125.30	119.78
3	D	531	SER	C-N-CA	5.36	125.30	119.78
2	C	462	LEU	N-CA-C	-5.36	105.33	111.07
3	D	308	SER	CA-C-N	5.32	125.89	119.98
3	D	308	SER	C-N-CA	5.32	125.89	119.98
2	C	48	ALA	CA-C-N	5.31	125.63	119.47
2	C	48	ALA	C-N-CA	5.31	125.63	119.47
2	C	166	PRO	N-CA-C	-5.31	102.77	111.21
2	C	617	TYR	N-CA-C	-5.28	106.34	114.16
1	B	390	ALA	O-C-N	5.28	127.79	122.09
1	A	53	PHE	CA-CB-CG	-5.24	108.56	113.80
1	B	344	PHE	CA-C-N	5.23	125.17	119.78
1	B	344	PHE	C-N-CA	5.23	125.17	119.78
2	C	498	TYR	N-CA-C	5.22	117.88	111.82
3	D	172	ASN	CA-C-N	5.21	124.93	119.82
3	D	172	ASN	C-N-CA	5.21	124.93	119.82
1	B	179	GLU	CA-C-N	5.20	129.55	120.86
1	B	179	GLU	C-N-CA	5.20	129.55	120.86
1	B	244	PHE	CA-C-N	5.20	125.17	120.03
1	B	244	PHE	C-N-CA	5.20	125.17	120.03
1	A	136	PHE	CA-CB-CG	5.18	118.98	113.80
2	C	573	ILE	CA-C-N	-5.17	116.27	122.22
2	C	573	ILE	C-N-CA	-5.17	116.27	122.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	219	ASN	CA-C-N	5.15	125.22	119.87
1	B	219	ASN	C-N-CA	5.15	125.22	119.87
1	B	395	ASP	N-CA-C	-5.14	106.85	113.02
2	C	88	THR	CA-C-N	5.11	125.01	119.85
2	C	88	THR	C-N-CA	5.11	125.01	119.85
2	C	568	TYR	N-CA-C	-5.11	103.34	109.83
1	B	136	PHE	CA-CB-CG	5.10	118.90	113.80
1	A	344	PHE	CB-CA-C	5.09	117.86	109.97
3	D	812	ILE	CB-CA-C	-5.08	105.46	111.97
1	A	166	ILE	N-CA-C	5.08	115.08	107.51
1	B	53	PHE	CA-CB-CG	-5.08	108.72	113.80
2	C	461	LEU	N-CA-C	-5.08	106.25	112.90
1	A	246	SER	N-CA-C	-5.07	104.62	108.78
2	C	164	LYS	N-CA-C	5.07	120.46	113.97
4	U	113	VAL	N-CA-CB	-5.06	107.37	111.67
1	A	398	PHE	CA-CB-CG	-5.06	108.74	113.80
2	C	287	LEU	N-CA-C	5.05	117.68	111.82
2	C	334	ASP	CA-CB-CG	-5.05	107.55	112.60
1	A	220	PRO	CA-C-N	-5.04	114.27	122.34
1	A	220	PRO	C-N-CA	-5.04	114.27	122.34
1	A	405	ASN	N-CA-C	5.04	120.08	113.88
2	C	438	PHE	CA-CB-CG	-5.03	108.78	113.80
2	C	493	GLN	N-CA-C	-5.02	105.51	111.69
3	D	737	SER	N-CA-C	-5.02	107.16	113.28
2	C	191	ASN	CA-CB-CG	-5.01	107.59	112.60
2	C	696	GLN	N-CA-C	-5.01	105.90	111.36
2	C	595	HIS	CA-CB-CG	-5.01	108.79	113.80
1	B	327	GLU	N-CA-C	5.00	116.19	109.24

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3474	0	3329	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3486	0	3344	43	0
2	C	5853	0	5914	19	0
3	D	5560	0	5575	11	0
4	U	275	0	280	0	0
5	A	28	0	12	0	0
5	B	28	0	12	0	0
All	All	18704	0	18466	93	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 (93) 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:B:316:PHE:HB2	1:B:379:ASN:O	1.46	1.14
1:B:315:TYR:HB3	1:B:343:LYS:O	1.52	1.08
1:B:316:PHE:CB	1:B:379:ASN:O	2.04	1.04
1:B:180:VAL:HG13	1:B:393:THR:HG21	1.40	1.04
1:A:188:ILE:HD12	1:A:394:PHE:CD1	2.02	0.93
1:A:188:ILE:CD1	1:A:394:PHE:CD1	2.56	0.87
1:B:184:SER:HB2	1:B:394:PHE:HB3	1.56	0.86
1:B:259:LEU:HA	1:B:317:ASN:ND2	1.90	0.85
1:B:180:VAL:HG13	1:B:393:THR:CG2	2.07	0.85
1:B:315:TYR:O	1:B:380:MET:HG2	1.86	0.76
1:B:180:VAL:CG1	1:B:393:THR:HG21	2.16	0.73
1:A:188:ILE:HD13	1:A:394:PHE:CG	2.24	0.72
1:A:188:ILE:CD1	1:A:394:PHE:CG	2.71	0.72
1:A:188:ILE:HD13	1:A:394:PHE:CD1	2.29	0.65
2:C:581:LYS:NZ	2:C:681:THR:OG1	2.30	0.63
1:B:387:PHE:O	1:B:391:CYS:SG	2.52	0.63
1:B:394:PHE:HE1	1:B:427:ARG:NE	1.97	0.61
1:B:180:VAL:HG22	1:B:393:THR:HG21	1.81	0.60
3:D:471:ARG:NH2	3:D:671:THR:OG1	2.35	0.60
1:B:258:THR:O	1:B:317:ASN:ND2	2.35	0.59
1:B:316:PHE:HB3	1:B:379:ASN:O	2.00	0.59
1:B:259:LEU:HA	1:B:317:ASN:HD21	1.68	0.59
1:B:246:SER:OG	1:B:247:TYR:N	2.35	0.58
1:A:246:SER:OG	1:A:247:TYR:N	2.39	0.56
2:C:459:ASP:N	2:C:459:ASP:OD1	2.38	0.55
1:B:316:PHE:HB2	1:B:380:MET:HA	1.88	0.55
1:B:394:PHE:C	1:B:394:PHE:CD1	2.85	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:163:SER:OG	3:D:164:SER:N	2.41	0.54
1:A:188:ILE:HD12	1:A:394:PHE:HD1	1.66	0.54
2:C:618:SER:OG	2:C:619:HIS:N	2.42	0.52
1:B:180:VAL:HG22	1:B:393:THR:CG2	2.40	0.52
1:B:394:PHE:CE1	1:B:427:ARG:CD	2.93	0.51
2:C:18:ASN:OD1	2:C:18:ASN:N	2.37	0.51
2:C:128:ASP:OD1	2:C:129:THR:N	2.43	0.51
2:C:663:TYR:O	2:C:665:ASN:N	2.44	0.51
1:B:394:PHE:HE1	1:B:427:ARG:CD	2.25	0.50
3:D:498:ASP:N	3:D:498:ASP:OD1	2.41	0.50
1:A:188:ILE:HG22	1:A:189:LEU:HD12	1.93	0.50
1:B:452:ASP:OD2	3:D:630:LYS:NZ	2.44	0.50
1:B:223:ASP:OD1	1:B:224:LEU:N	2.45	0.50
2:C:123:TYR:C	2:C:123:TYR:CD2	2.90	0.50
1:A:188:ILE:CD1	1:A:394:PHE:CB	2.89	0.50
1:B:394:PHE:C	1:B:394:PHE:HD1	2.19	0.49
1:B:180:VAL:C	1:B:393:THR:HG21	2.37	0.49
1:B:394:PHE:CE1	1:B:427:ARG:HD3	2.47	0.49
1:B:316:PHE:HB2	1:B:379:ASN:C	2.29	0.49
1:A:61:LYS:NZ	1:A:85:GLY:O	2.46	0.49
2:C:183:ASN:OD1	2:C:183:ASN:N	2.45	0.48
3:D:498:ASP:OD2	3:D:511:LYS:NZ	2.40	0.48
1:B:41:ASP:OD1	1:B:41:ASP:N	2.46	0.48
2:C:501:ILE:O	2:C:502:TYR:C	2.55	0.48
1:B:180:VAL:HG22	1:B:393:THR:CB	2.43	0.48
1:B:315:TYR:HE1	1:B:441:GLU:HA	1.79	0.47
2:C:788:THR:O	2:C:789:GLU:C	2.56	0.47
1:B:391:CYS:HA	1:B:394:PHE:CZ	2.49	0.46
3:D:409:TYR:C	3:D:409:TYR:CD2	2.94	0.46
3:D:685:ASN:OD1	3:D:685:ASN:N	2.46	0.46
3:D:250:LYS:NZ	3:D:361:GLU:OE2	2.49	0.46
1:A:29:HIS:O	1:A:31:ILE:N	2.49	0.46
1:A:188:ILE:CD1	1:A:394:PHE:HB2	2.46	0.46
1:B:180:VAL:CG2	1:B:393:THR:HG21	2.44	0.45
1:A:444:SER:O	1:A:445:TYR:C	2.60	0.45
1:B:319:TYR:OH	1:B:355:ASN:ND2	2.49	0.45
1:B:391:CYS:HA	1:B:394:PHE:CE2	2.51	0.45
1:B:209:LEU:HB3	1:B:227:THR:CG2	2.47	0.45
1:B:47:ASP:OD1	1:B:47:ASP:N	2.43	0.44
1:A:199:ASP:O	1:A:200:ALA:C	2.60	0.44
3:D:225:PHE:O	3:D:226:GLU:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:702:ILE:HG23	2:C:706:PHE:CZ	2.53	0.43
1:A:41:ASP:N	1:A:41:ASP:OD1	2.44	0.43
2:C:710:ILE:O	2:C:711:LYS:HB2	2.19	0.43
2:C:752:TYR:CD2	2:C:752:TYR:C	2.93	0.43
1:A:34:ASP:N	1:A:34:ASP:OD1	2.47	0.43
2:C:106:SER:O	2:C:107:ARG:C	2.60	0.42
1:A:449:VAL:HG22	2:C:623:ARG:HB3	2.00	0.42
1:B:51:LYS:N	1:B:52:PRO:HD2	2.34	0.42
1:A:191:LEU:O	1:A:192:ARG:C	2.63	0.42
1:B:184:SER:CB	1:B:394:PHE:HB3	2.39	0.41
1:A:131:ASP:OD1	1:A:131:ASP:N	2.53	0.41
1:A:344:PHE:HB2	1:A:345:PRO:HD2	2.02	0.41
1:A:127:ILE:O	1:A:130:THR:OG1	2.38	0.41
1:B:330:GLN:N	1:B:330:GLN:OE1	2.53	0.41
2:C:265:VAL:O	2:C:266:MET:C	2.63	0.41
2:C:300:MET:O	2:C:301:THR:CB	2.68	0.41
3:D:247:SER:HA	3:D:248:PRO:HD3	1.96	0.41
1:A:83:PHE:HB3	1:A:86:PHE:HB3	2.03	0.41
1:B:316:PHE:HZ	1:B:440:ALA:HB3	1.86	0.41
2:C:413:TYR:O	2:C:414:ASP:C	2.62	0.41
1:B:34:ASP:N	1:B:34:ASP:OD1	2.45	0.41
1:B:180:VAL:HG22	1:B:393:THR:HB	2.02	0.41
1:B:432:SER:O	1:B:433:LEU:C	2.64	0.40
3:D:443:THR:O	3:D:446:PHE:N	2.52	0.40
2:C:300:MET:O	2:C:301:THR:HB	2.21	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	440/473 (93%)	433 (98%)	7 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	442/473 (93%)	425 (96%)	15 (3%)	2 (0%)	24	57
2	C	689/823 (84%)	672 (98%)	14 (2%)	3 (0%)	30	61
3	D	669/846 (79%)	656 (98%)	10 (2%)	3 (0%)	30	61
4	U	31/220 (14%)	31 (100%)	0	0	100	100
All	All	2271/2835 (80%)	2217 (98%)	46 (2%)	8 (0%)	31	61

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	315	TYR
2	C	664	ARG
3	D	798	SER
2	C	301	THR
2	C	753	GLN
3	D	560	LEU
1	B	317	ASN
3	D	835	ASP

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	396/420 (94%)	396 (100%)	0	100	100
1	B	397/420 (94%)	395 (100%)	2 (0%)	81	80
2	C	657/766 (86%)	656 (100%)	1 (0%)	87	85
3	D	627/787 (80%)	627 (100%)	0	100	100
4	U	31/206 (15%)	31 (100%)	0	100	100
All	All	2108/2599 (81%)	2105 (100%)	3 (0%)	87	87

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	315	TYR
1	B	394	PHE
2	C	617	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	GLN
1	B	116	GLN
1	B	153	ASN
1	B	240	ASN
2	C	433	HIS
2	C	674	ASN
3	D	815	GLN

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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GDP	A	1400	-	29,30,30	1.18	3 (10%)	45,47,47	1.73	8 (17%)
5	GDP	B	501	-	29,30,30	1.10	2 (6%)	45,47,47	1.92	13 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GDP	A	1400	-	-	3/16/32/32	0/3/3/3
5	GDP	B	501	-	-	8/16/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	501	GDP	C4-N3	2.91	1.40	1.34
5	A	1400	GDP	C4-N3	2.86	1.40	1.34
5	A	1400	GDP	C5-C4	2.35	1.45	1.38
5	B	501	GDP	C5-C4	2.20	1.44	1.38
5	A	1400	GDP	C5'-C4'	2.10	1.57	1.51

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1400	GDP	C5-C4-N3	-5.52	119.61	128.39
5	B	501	GDP	C2-N3-C4	5.31	121.45	112.30
5	B	501	GDP	C5-C4-N3	-5.30	119.95	128.39
5	A	1400	GDP	C2-N3-C4	4.96	120.84	112.30
5	A	1400	GDP	N9-C4-N3	4.47	134.90	125.95
5	B	501	GDP	N9-C4-N3	4.20	134.35	125.95
5	B	501	GDP	C6-C5-N7	3.05	135.84	130.29
5	B	501	GDP	C8-N7-C5	2.89	109.41	104.26
5	A	1400	GDP	C6-C5-N7	2.82	135.43	130.29
5	B	501	GDP	O2B-PB-O3A	2.77	113.92	104.64
5	B	501	GDP	O2A-PA-O3A	2.67	114.50	107.27
5	A	1400	GDP	C8-N7-C5	2.51	108.72	104.26
5	A	1400	GDP	C5-C6-N1	2.47	119.55	113.25
5	B	501	GDP	O4'-C1'-C2'	-2.44	101.40	106.62
5	B	501	GDP	C2-N1-C6	-2.40	120.77	125.11
5	B	501	GDP	C4-C5-N7	-2.37	106.92	110.67
5	B	501	GDP	C5-C6-N1	2.34	119.22	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1400	GDP	C2-N1-C6	-2.20	121.11	125.11
5	B	501	GDP	O3B-PB-O3A	2.12	111.73	104.64
5	A	1400	GDP	C4-C5-N7	-2.11	107.33	110.67
5	B	501	GDP	O4'-C1'-N9	2.10	113.11	108.36

There are no chirality outliers.

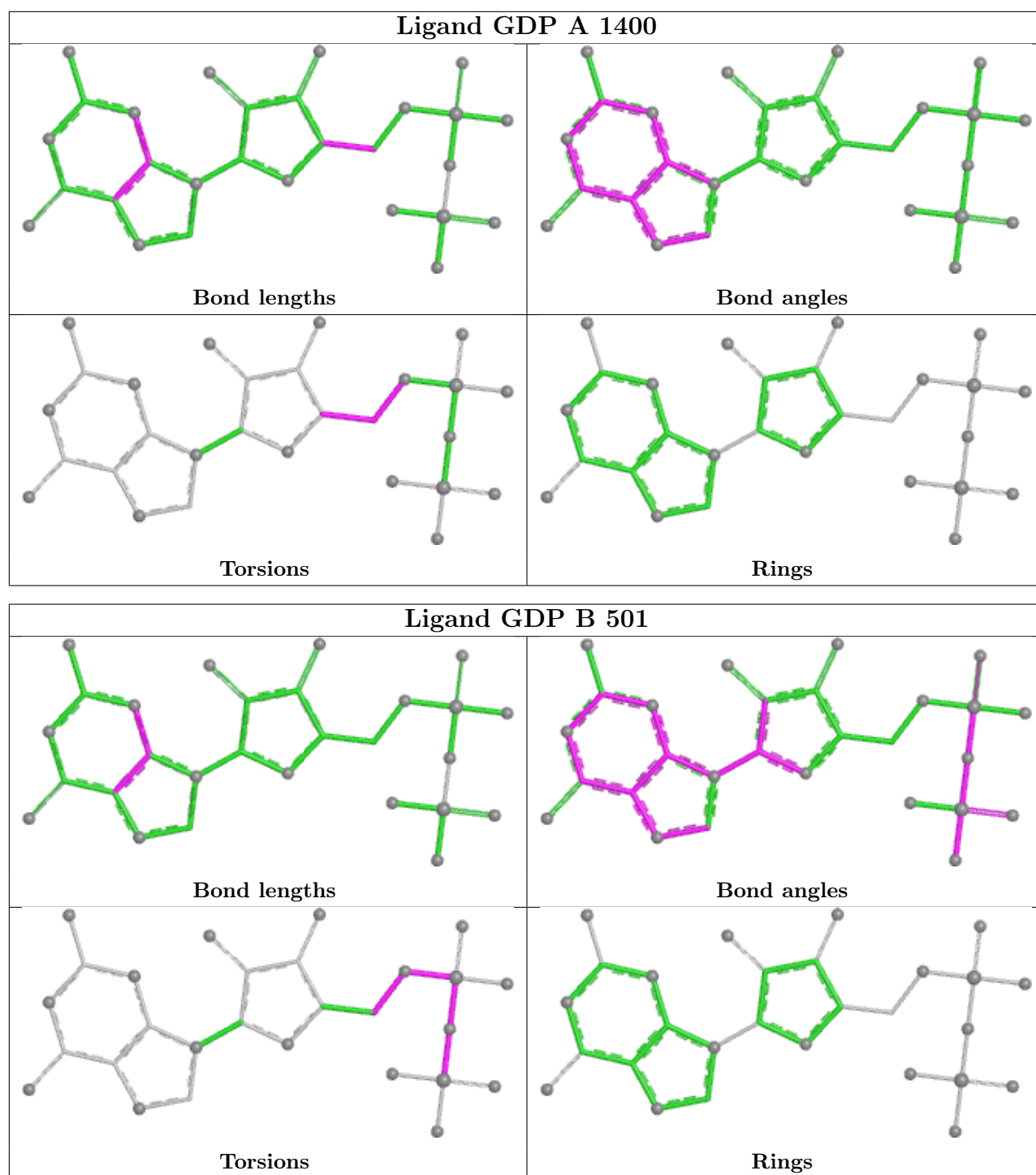
All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1400	GDP	O4'-C4'-C5'-O5'
5	B	501	GDP	C5'-O5'-PA-O3A
5	B	501	GDP	C5'-O5'-PA-O2A
5	B	501	GDP	C4'-C5'-O5'-PA
5	A	1400	GDP	C3'-C4'-C5'-O5'
5	A	1400	GDP	C4'-C5'-O5'-PA
5	B	501	GDP	C5'-O5'-PA-O1A
5	B	501	GDP	PA-O3A-PB-O1B
5	B	501	GDP	PA-O3A-PB-O2B
5	B	501	GDP	PA-O3A-PB-O3B
5	B	501	GDP	PB-O3A-PA-O2A

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.

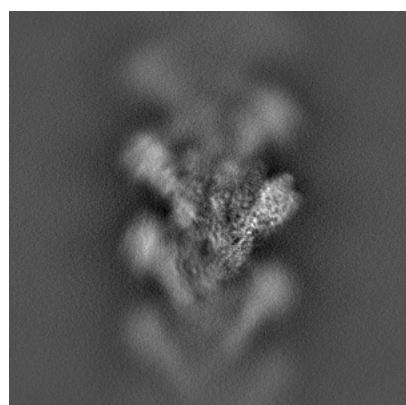
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23636. These allow visual inspection of the internal detail of the map and identification of artifacts.

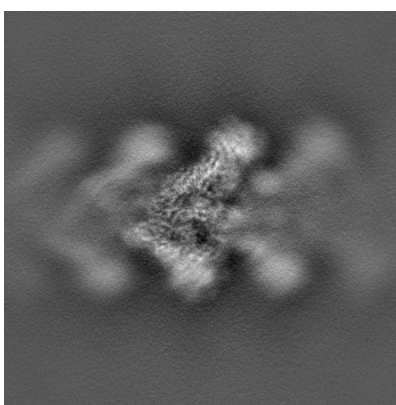
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

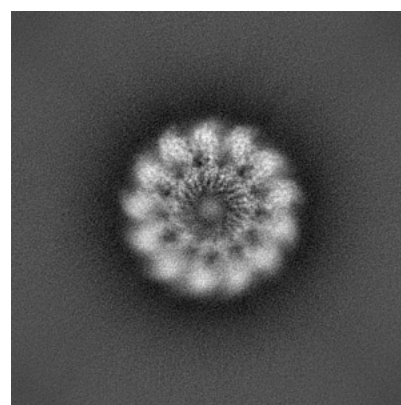
6.1.1 Primary map



X



Y

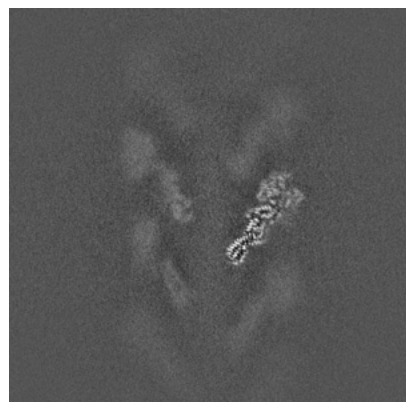


Z

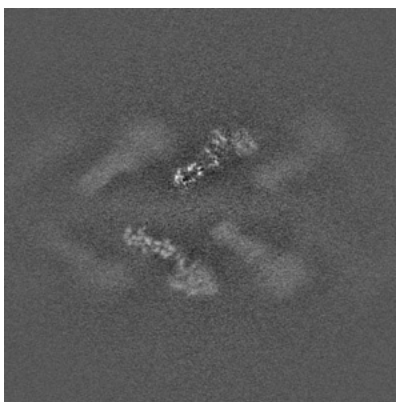
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

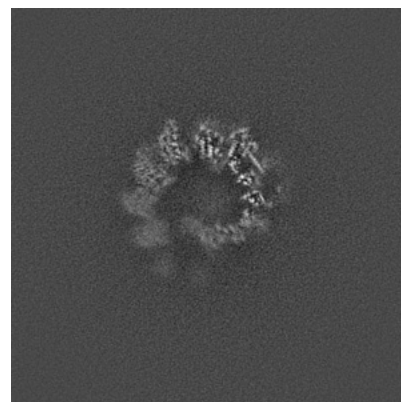
6.2.1 Primary map



X Index: 224



Y Index: 224

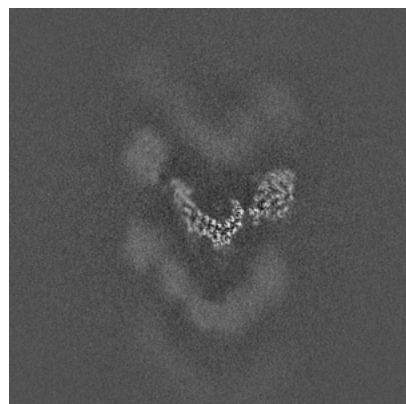


Z Index: 224

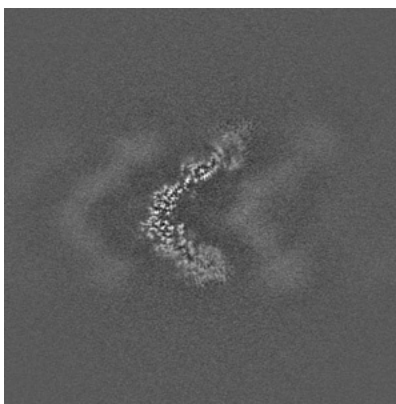
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

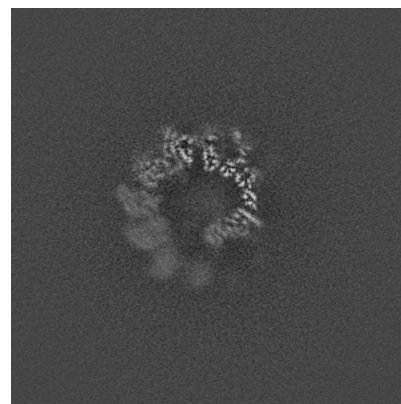
6.3.1 Primary map



X Index: 256



Y Index: 258

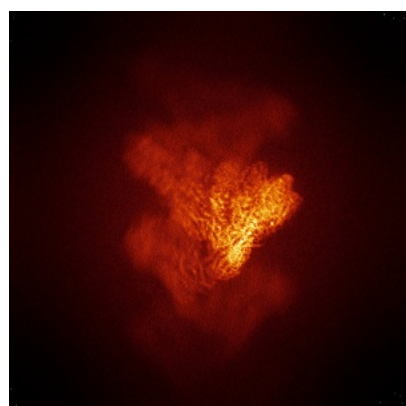


Z Index: 214

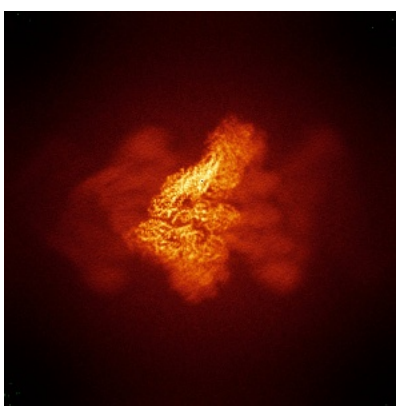
The images above show the largest variance slices of the map in three orthogonal directions.

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

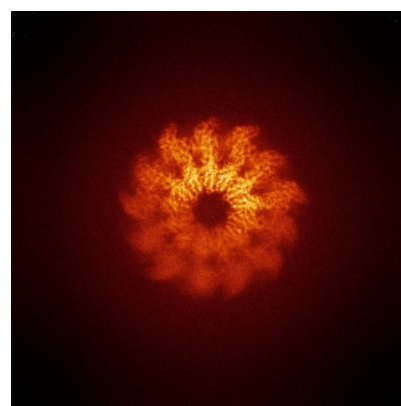
6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 7.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

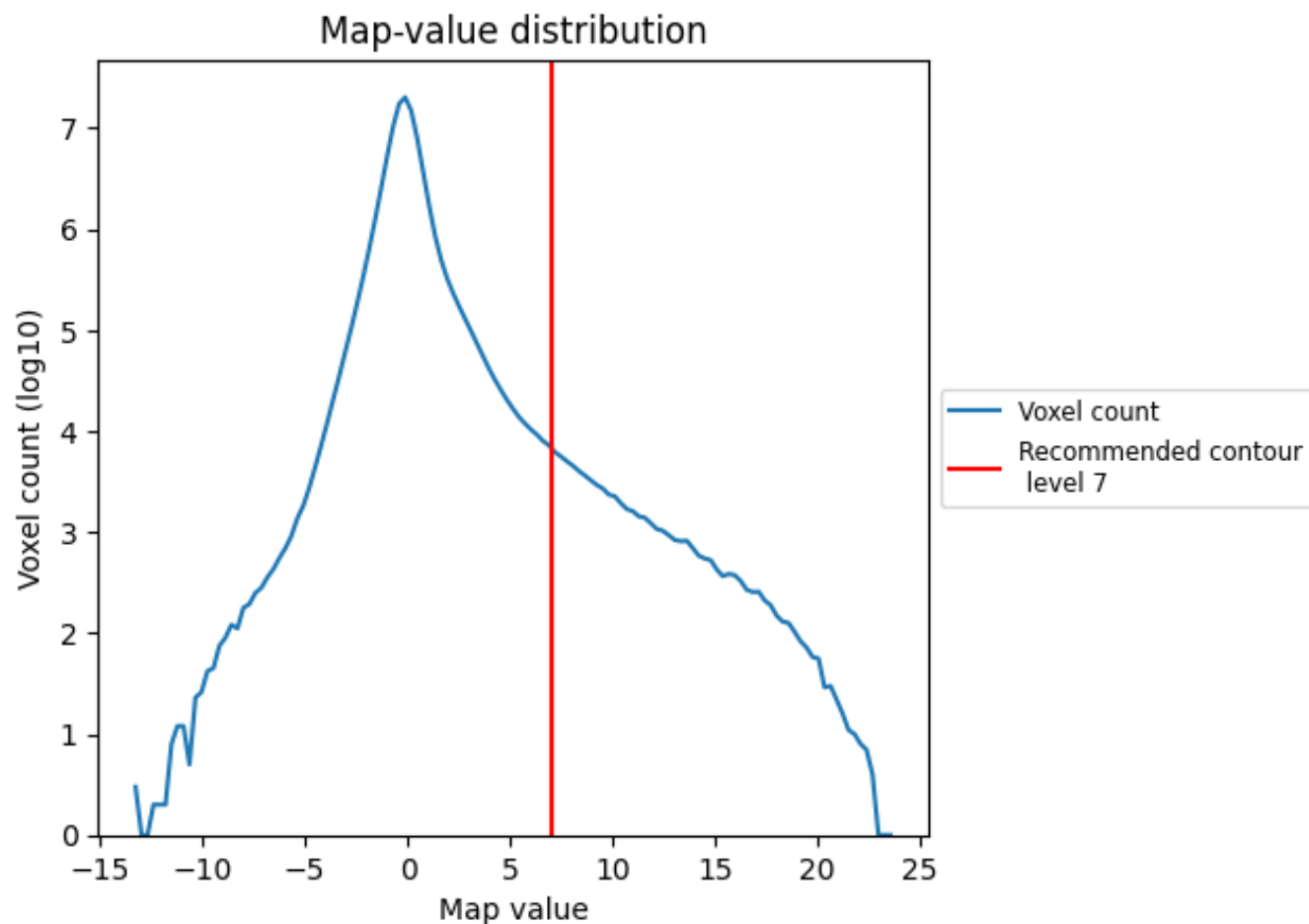
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

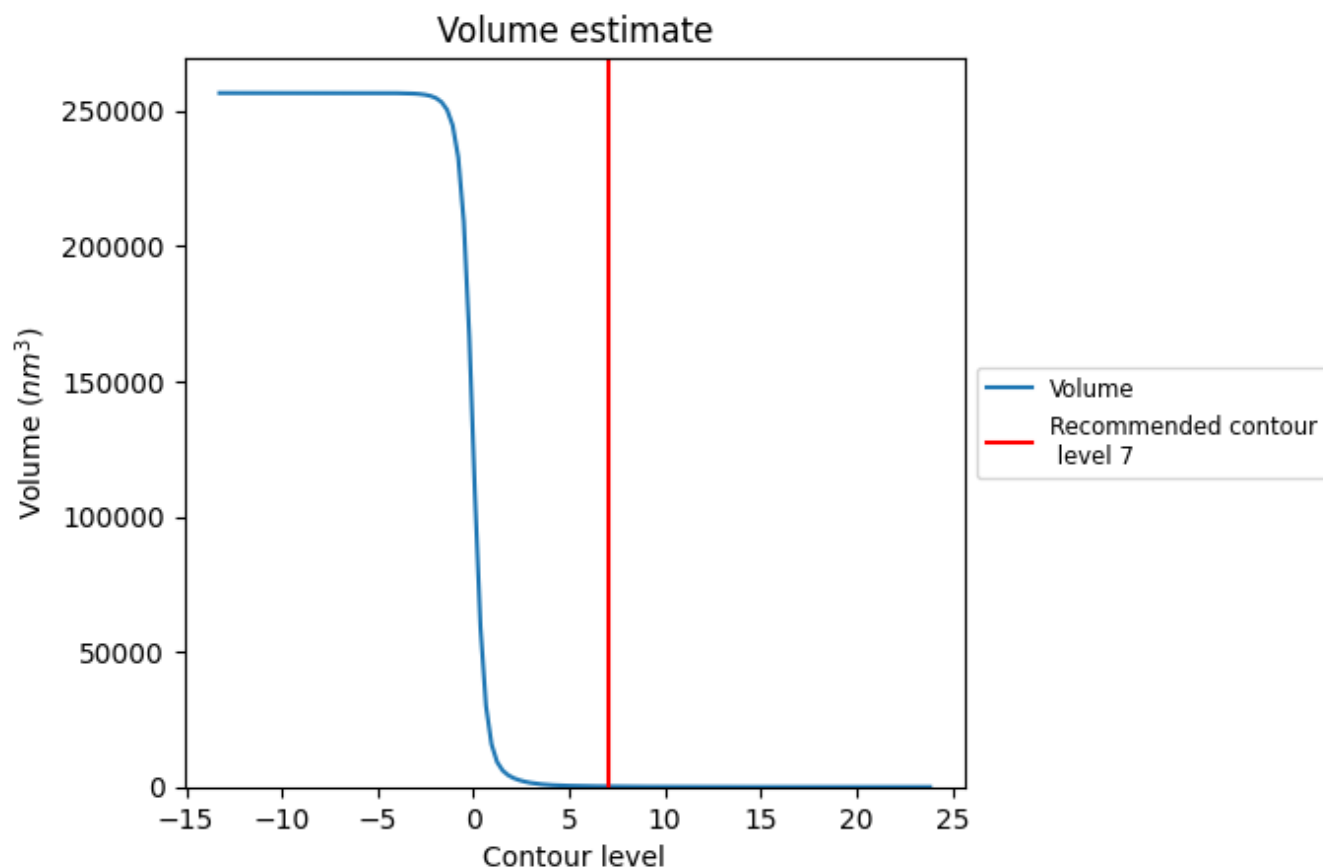
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

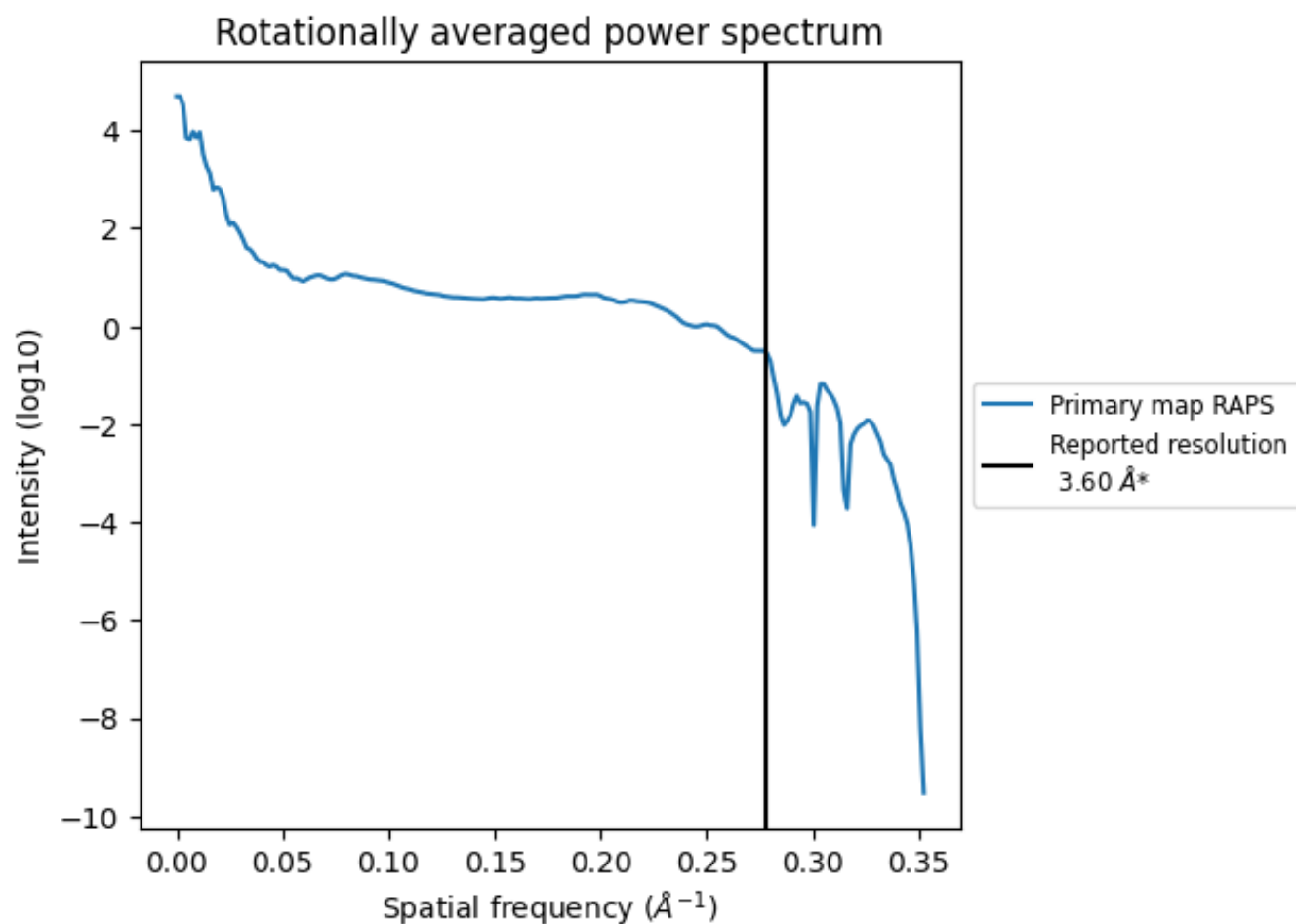
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 195 nm^3 ; this corresponds to an approximate mass of 176 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.278 Å⁻¹

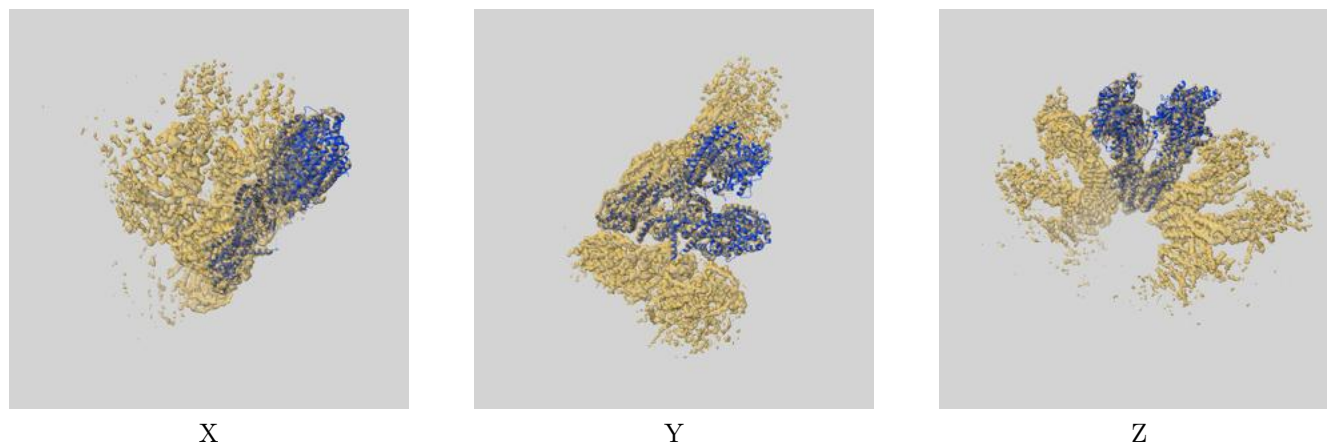
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23636 and PDB model 7M2X. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



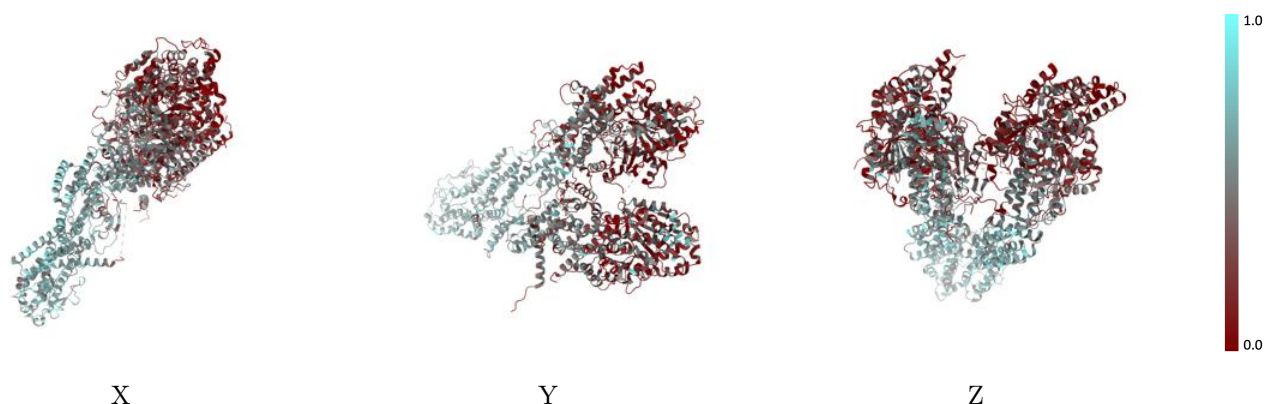
The images above show the 3D surface view of the map at the recommended contour level 7.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



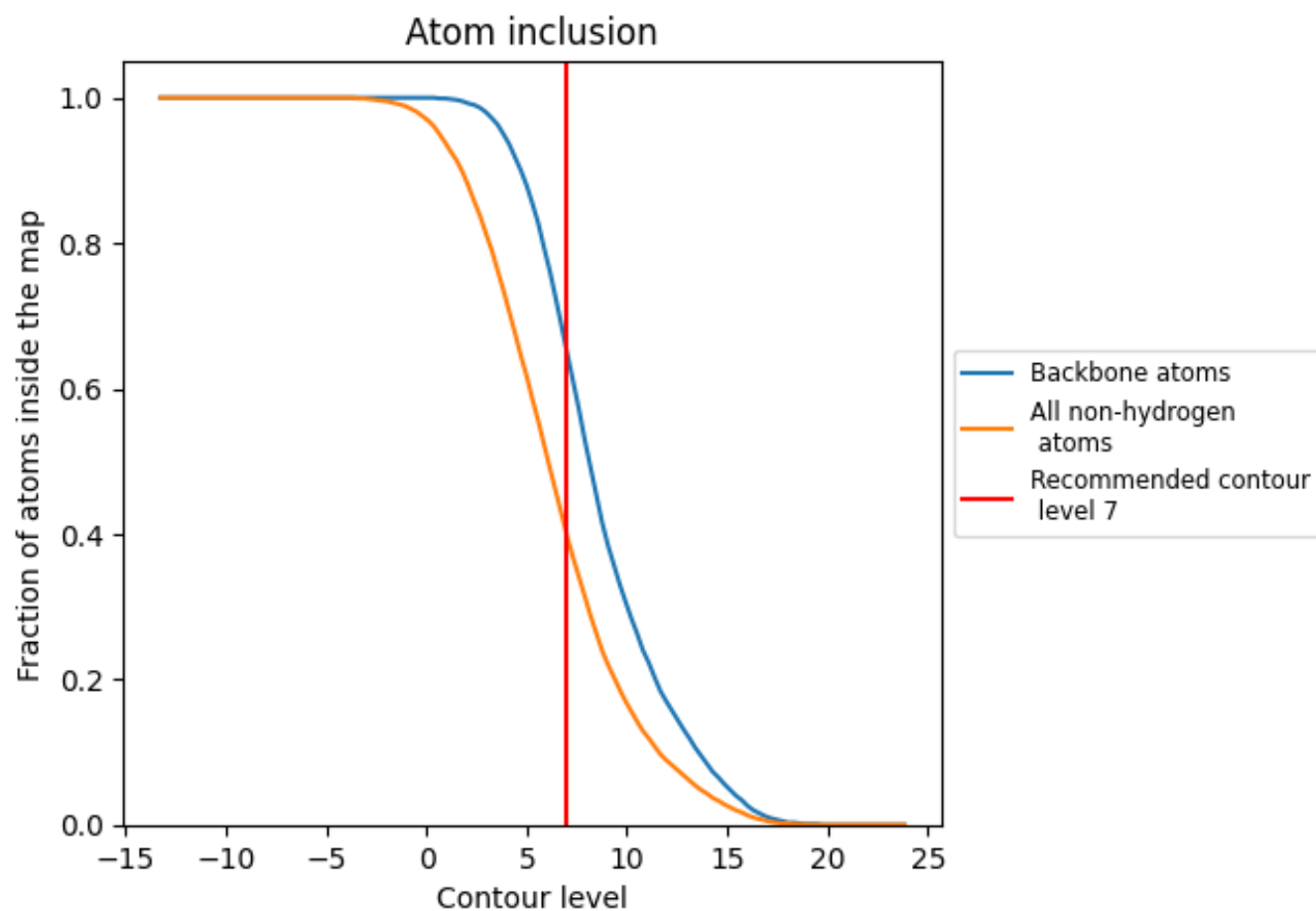
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 65% of all backbone atoms, 40% 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 (7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.3990	0.2950
A	0.3080	0.2550
B	0.2270	0.2490
C	0.4890	0.3210
D	0.4720	0.3210
U	0.3680	0.2640

