



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 04:49 AM UTC

PDB ID : 7M2Y / pdb_00007m2y
EMDB ID : EMD-23637
Title : Closed 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 : 4.03 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:

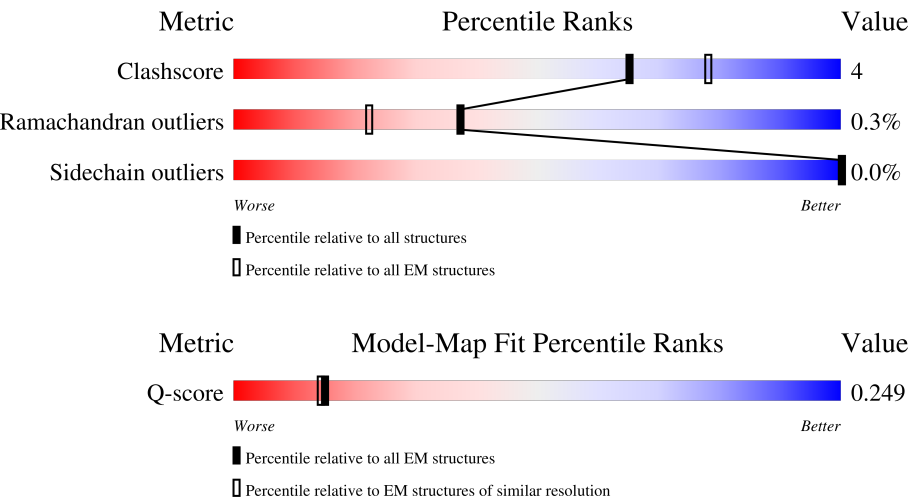
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 i

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

The reported resolution of this entry is 4.03 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	6618 (3.54 - 4.53)

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	
1	B	473	
2	C	846	
3	D	823	

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Mol	Chain	Length	Quality of chain
4	U	220	 13% 86%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GDP	A	501	-	-	X	-
5	GDP	B	501	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18560 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	441	Total	C	N	O	S	0	0
			3447	2156	582	693	16		
1	B	447	Total	C	N	O	S	0	0
			3491	2182	591	701	17		

- Molecule 2 is a protein called Spindle pole body component SPC98.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	656	Total	C	N	O	S	0	0
			5435	3518	899	1002	16		

- Molecule 3 is a protein called Spindle pole body component SPC97.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	703	Total	C	N	O	S	0	0
			5868	3781	984	1074	29		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
4	U	31	Total	C	N	O	S	0	0
			263	162	49	50	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	





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	20420	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)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	47214	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	21.845	Depositor
Minimum map value	-11.274	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.822	Depositor
Recommended contour level	3.4	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 [i](#)

5.1 Standard geometry [i](#)

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	1.13	1/3518 (0.0%)	1.41	49/4778 (1.0%)
1	B	1.09	1/3563 (0.0%)	1.44	58/4837 (1.2%)
2	C	1.30	9/5551 (0.2%)	1.25	34/7500 (0.5%)
3	D	1.23	5/5984 (0.1%)	1.23	31/8078 (0.4%)
4	U	1.17	0/266	1.22	3/352 (0.9%)
All	All	1.21	16/18882 (0.1%)	1.31	175/25545 (0.7%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	29	HIS	CB-CG	-8.94	1.37	1.50
3	D	190	TYR	CB-CG	-7.92	1.34	1.51
2	C	288	TYR	CB-CG	-7.46	1.35	1.51
3	D	78	PHE	CB-CG	-7.09	1.34	1.50
2	C	267	PHE	CB-CG	-6.51	1.35	1.50
2	C	469	HIS	CB-CG	-6.17	1.41	1.50
2	C	349	TRP	CB-CG	-5.78	1.32	1.50
2	C	757	PRO	N-CD	-5.49	1.40	1.47
2	C	550	HIS	CB-CG	-5.43	1.42	1.50
2	C	527	HIS	CB-CG	-5.38	1.42	1.50
3	D	433	HIS	CB-CG	-5.36	1.42	1.50
3	D	112	TYR	CB-CG	-5.29	1.40	1.51
3	D	197	PHE	CB-CG	-5.20	1.38	1.50
2	C	403	PHE	CB-CG	-5.05	1.39	1.50
2	C	596	ASN	CB-CG	-5.05	1.39	1.52
1	B	419	VAL	CB-CG2	-5.00	1.36	1.52

All (175) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	201	THR	N-CA-C	11.97	127.10	108.79
1	A	70	ASP	N-CA-C	10.96	124.32	111.71
3	D	51	GLN	N-CA-C	-10.21	101.33	113.88
2	C	792	LEU	N-CA-C	-9.91	101.32	113.41
3	D	804	ILE	N-CA-C	-8.90	101.30	111.00
3	D	87	GLU	N-CA-C	-8.87	103.29	114.56
1	A	88	ASP	CA-C-N	8.52	128.25	119.56
1	A	88	ASP	C-N-CA	8.52	128.25	119.56
1	B	88	ASP	CA-C-N	8.51	128.24	119.56
1	B	88	ASP	C-N-CA	8.51	128.24	119.56
3	D	488	THR	N-CA-C	-8.51	103.08	113.97
1	B	200	ALA	N-CA-C	8.50	120.54	111.28
1	A	236	SER	N-CA-C	-8.47	104.08	114.75
1	A	262	SER	CA-C-N	8.13	127.85	119.56
1	A	262	SER	C-N-CA	8.13	127.85	119.56
1	B	83	PHE	CA-CB-CG	-7.91	105.89	113.80
1	A	312	ASN	CA-C-N	7.83	127.99	120.31
1	A	312	ASN	C-N-CA	7.83	127.99	120.31
1	B	140	HIS	CA-CB-CG	7.68	121.48	113.80
1	A	63	THR	CA-C-N	7.67	127.59	119.85
1	A	63	THR	C-N-CA	7.67	127.59	119.85
1	B	96	SER	N-CA-C	7.64	120.98	108.99
3	D	502	TYR	CA-C-N	7.62	128.06	120.52
3	D	502	TYR	C-N-CA	7.62	128.06	120.52
2	C	621	ASN	CA-C-N	7.49	129.20	119.84
2	C	621	ASN	C-N-CA	7.49	129.20	119.84
1	B	219	ASN	CA-C-N	7.48	127.15	119.82
1	B	219	ASN	C-N-CA	7.48	127.15	119.82
2	C	342	TYR	N-CA-C	-7.34	103.36	111.36
1	A	173	PHE	CA-C-N	7.28	128.94	119.84
1	A	173	PHE	C-N-CA	7.28	128.94	119.84
1	B	293	ASP	CA-CB-CG	7.17	119.77	112.60
1	A	273	THR	CA-C-N	7.00	126.99	119.78
1	A	273	THR	C-N-CA	7.00	126.99	119.78
1	B	344	PHE	CA-C-N	6.94	126.87	120.21
1	B	344	PHE	C-N-CA	6.94	126.87	120.21
1	A	356	ILE	N-CA-C	6.89	120.34	108.90
1	B	394	PHE	CA-CB-CG	6.83	120.63	113.80
1	B	344	PHE	N-CA-C	6.81	114.26	108.13
1	B	162	TYR	CA-C-N	6.77	127.32	119.47
1	B	162	TYR	C-N-CA	6.77	127.32	119.47
1	A	338	LEU	N-CA-C	-6.73	103.41	111.69
1	A	269	SER	CA-C-N	6.72	126.69	120.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	SER	C-N-CA	6.72	126.69	120.03
3	D	473	ASN	CA-C-N	6.65	126.28	119.56
3	D	473	ASN	C-N-CA	6.65	126.28	119.56
1	B	244	PHE	CA-C-N	6.64	126.60	120.03
1	B	244	PHE	C-N-CA	6.64	126.60	120.03
1	B	71	SER	N-CA-C	-6.57	104.99	114.12
1	B	360	SER	CA-C-N	6.56	126.69	119.87
1	B	360	SER	C-N-CA	6.56	126.69	119.87
2	C	761	SER	N-CA-C	-6.55	104.22	111.36
2	C	213	ILE	CA-C-N	6.48	126.86	119.92
2	C	213	ILE	C-N-CA	6.48	126.86	119.92
1	B	269	SER	CA-C-N	6.43	126.61	120.31
1	B	269	SER	C-N-CA	6.43	126.61	120.31
1	B	200	ALA	CA-C-N	-6.43	109.86	121.62
1	B	200	ALA	C-N-CA	-6.43	109.86	121.62
3	D	294	ASP	CA-C-N	6.35	125.97	119.56
3	D	294	ASP	C-N-CA	6.35	125.97	119.56
1	A	197	ASP	N-CA-C	6.35	121.16	113.41
1	A	53	PHE	CA-CB-CG	-6.28	107.52	113.80
1	B	300	ASP	CA-C-N	6.26	125.94	119.56
1	B	300	ASP	C-N-CA	6.26	125.94	119.56
3	D	165	VAL	CA-C-N	6.25	125.93	119.56
3	D	165	VAL	C-N-CA	6.25	125.93	119.56
1	B	40	PRO	N-CA-C	-6.24	101.39	111.19
2	C	585	ARG	N-CA-C	-6.19	104.08	111.69
1	A	10	ALA	N-CA-C	6.17	118.35	108.41
2	C	480	ILE	N-CA-C	6.13	117.49	111.67
1	A	320	ASN	N-CA-C	6.09	119.44	108.48
1	B	173	PHE	CA-CB-CG	6.07	119.87	113.80
3	D	754	GLU	N-CA-C	6.05	118.67	111.71
3	D	99	ASP	CA-C-N	6.05	125.73	119.56
3	D	99	ASP	C-N-CA	6.05	125.73	119.56
1	A	363	LEU	CA-C-N	6.04	126.01	120.21
1	A	363	LEU	C-N-CA	6.04	126.01	120.21
2	C	465	ASN	CA-CB-CG	-5.97	106.63	112.60
1	A	69	MET	CA-C-N	5.96	128.87	120.28
1	A	69	MET	C-N-CA	5.96	128.87	120.28
2	C	531	SER	CA-C-N	5.95	125.57	119.56
2	C	531	SER	C-N-CA	5.95	125.57	119.56
1	A	344	PHE	CA-CB-CG	5.95	119.75	113.80
1	B	366	GLN	CA-C-N	5.93	125.55	119.56
1	B	366	GLN	C-N-CA	5.93	125.55	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	250	LYS	N-CA-C	-5.89	104.94	111.36
1	B	316	PHE	CA-CB-CG	-5.85	107.95	113.80
1	A	162	TYR	CA-C-N	5.84	125.46	119.56
1	A	162	TYR	C-N-CA	5.84	125.46	119.56
1	B	97	ASP	N-CA-C	-5.84	98.76	108.34
2	C	804	LYS	N-CA-C	-5.82	104.89	112.23
1	B	82	THR	N-CA-C	-5.79	105.23	112.93
2	C	390	VAL	CA-C-N	5.78	125.75	120.03
2	C	390	VAL	C-N-CA	5.78	125.75	120.03
1	B	230	LEU	N-CA-C	-5.73	105.39	112.90
1	B	327	GLU	CA-C-N	5.72	126.10	119.47
1	B	327	GLU	C-N-CA	5.72	126.10	119.47
4	U	119	LEU	N-CA-C	5.70	117.57	111.36
2	C	195	THR	N-CA-C	-5.69	101.77	110.14
1	A	150	LEU	N-CA-C	5.67	117.54	111.36
1	A	327	GLU	CA-C-N	5.63	126.01	119.47
1	A	327	GLU	C-N-CA	5.63	126.01	119.47
1	A	72	GLU	CA-C-N	5.63	125.30	119.56
1	A	72	GLU	C-N-CA	5.63	125.30	119.56
2	C	179	ILE	CA-C-N	5.61	125.59	120.03
2	C	179	ILE	C-N-CA	5.61	125.59	120.03
3	D	366	ILE	CA-C-N	5.59	125.57	120.03
3	D	366	ILE	C-N-CA	5.59	125.57	120.03
1	B	275	PHE	CA-CB-CG	-5.57	108.23	113.80
2	C	576	ARG	CA-C-N	5.54	126.77	119.84
2	C	576	ARG	C-N-CA	5.54	126.77	119.84
1	B	86	PHE	CA-C-N	-5.54	115.64	122.84
1	B	86	PHE	C-N-CA	-5.54	115.64	122.84
1	B	86	PHE	N-CA-C	5.54	116.99	111.07
1	A	360	SER	CA-C-N	5.53	125.62	119.87
1	A	360	SER	C-N-CA	5.53	125.62	119.87
3	D	142	ILE	CB-CA-C	-5.53	104.89	111.97
1	A	69	MET	N-CA-C	5.50	118.21	109.96
1	B	93	TRP	CA-C-N	-5.46	116.42	123.19
1	B	93	TRP	C-N-CA	-5.46	116.42	123.19
2	C	172	ASN	CA-C-N	5.46	125.13	119.56
2	C	172	ASN	C-N-CA	5.46	125.13	119.56
1	A	104	SER	N-CA-C	5.46	117.09	108.96
1	A	300	ASP	CA-C-N	5.45	126.65	119.84
1	A	300	ASP	C-N-CA	5.45	126.65	119.84
2	C	247	SER	CA-C-N	5.42	125.09	119.56
2	C	247	SER	C-N-CA	5.42	125.09	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	PHE	N-CA-C	5.42	118.29	109.24
4	U	117	GLN	CA-C-N	5.39	125.31	119.76
4	U	117	GLN	C-N-CA	5.39	125.31	119.76
3	D	82	GLU	CA-C-N	5.36	125.02	119.56
3	D	82	GLU	C-N-CA	5.36	125.02	119.56
1	A	99	ALA	N-CA-C	5.35	117.20	109.07
1	B	203	VAL	N-CA-C	5.35	116.18	108.48
2	C	573	ASN	N-CA-C	5.35	118.71	111.17
1	B	323	ILE	N-CA-C	5.34	115.82	108.12
2	C	637	LEU	CA-C-N	5.34	127.96	120.28
2	C	637	LEU	C-N-CA	5.34	127.96	120.28
1	B	262	SER	CA-C-N	5.31	125.63	119.47
1	B	262	SER	C-N-CA	5.31	125.63	119.47
2	C	499	ILE	CB-CA-C	-5.31	105.17	111.97
2	C	386	ASN	CA-CB-CG	-5.31	107.29	112.60
1	B	419	VAL	CB-CA-C	-5.29	105.09	112.02
1	B	231	ILE	N-CA-C	-5.29	105.38	110.72
3	D	19	ILE	CA-C-N	5.28	125.27	119.89
3	D	19	ILE	C-N-CA	5.28	125.27	119.89
3	D	164	LYS	N-CA-C	5.27	120.72	113.97
1	A	75	VAL	N-CA-C	-5.27	106.00	111.58
1	A	201	THR	N-CA-C	5.26	115.18	108.24
3	D	226	LEU	N-CA-C	5.24	116.99	111.28
1	B	199	ASP	CA-CB-CG	5.23	117.83	112.60
2	C	257	ILE	CB-CA-C	-5.20	105.32	111.97
1	A	203	VAL	N-CA-C	5.19	115.43	108.17
1	B	260	ILE	CA-C-N	5.19	124.80	119.56
1	B	260	ILE	C-N-CA	5.19	124.80	119.56
3	D	566	ILE	CA-C-N	5.18	125.08	119.85
3	D	566	ILE	C-N-CA	5.18	125.08	119.85
3	D	647	PHE	CA-C-N	-5.17	113.41	120.44
3	D	647	PHE	C-N-CA	-5.17	113.41	120.44
3	D	573	ILE	N-CA-C	-5.15	106.78	111.67
1	B	333	ARG	N-CA-C	-5.14	105.76	111.36
2	C	462	GLN	N-CA-C	-5.13	105.32	112.45
1	B	96	SER	CA-C-N	5.13	130.00	122.77
1	B	96	SER	C-N-CA	5.13	130.00	122.77
2	C	354	LEU	N-CA-C	5.12	117.59	109.24
1	A	39	LEU	CA-C-N	5.09	125.07	120.03
1	A	39	LEU	C-N-CA	5.09	125.07	120.03
1	A	354	VAL	N-CA-C	5.09	117.36	108.95
3	D	158	LEU	N-CA-C	5.08	117.60	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	352	MET	CA-C-N	-5.08	115.83	122.99
1	B	352	MET	C-N-CA	-5.08	115.83	122.99
3	D	42	SER	N-CA-C	5.07	119.31	113.18
1	A	244	PHE	N-CA-C	-5.03	98.15	108.21
1	B	11	GLY	N-CA-C	-5.03	107.17	111.95
2	C	558	PHE	N-CA-C	5.01	117.98	109.76

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	3447	0	3305	29	0
1	B	3491	0	3349	44	0
2	C	5435	0	5462	57	0
3	D	5868	0	5932	18	0
4	U	263	0	266	0	0
5	A	28	0	12	13	0
5	B	28	0	11	32	0
All	All	18560	0	18337	147	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 4.

All (147) 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:C:621:ASN:HB3	2:C:622:PRO:CD	1.17	1.46
1:B:174:PRO:CD	5:B:501:GDP:O3'	1.63	1.42
2:C:621:ASN:CB	2:C:622:PRO:CD	2.09	1.25
1:B:174:PRO:HD3	5:B:501:GDP:O3'	1.02	1.18
1:B:143:ALA:HB3	5:B:501:GDP:H4'	1.13	1.11
2:C:621:ASN:CB	2:C:622:PRO:HD2	1.77	1.08
1:B:143:ALA:HB3	5:B:501:GDP:C4'	1.89	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:613:SER:HA	2:C:616:LYS:CD	1.91	1.01
2:C:621:ASN:HB3	2:C:622:PRO:HD2	1.03	1.00
2:C:621:ASN:HB3	2:C:622:PRO:HD3	1.02	0.99
1:A:176:ARG:NH1	5:A:501:GDP:O2'	1.96	0.98
2:C:621:ASN:CB	2:C:622:PRO:HD3	1.85	0.98
1:A:141:SER:OG	5:A:501:GDP:H4'	1.63	0.98
1:B:174:PRO:CG	5:B:501:GDP:O3'	2.13	0.94
2:C:613:SER:HA	2:C:616:LYS:HG3	1.52	0.91
1:B:143:ALA:CB	5:B:501:GDP:H4'	2.00	0.90
2:C:613:SER:HA	2:C:616:LYS:CG	2.01	0.90
2:C:614:PHE:CE2	2:C:628:ILE:HG12	2.09	0.88
1:B:144:GLY:CA	5:B:501:GDP:O5'	2.25	0.85
1:A:141:SER:OG	5:A:501:GDP:C4'	2.24	0.84
1:A:146:THR:OG1	5:A:501:GDP:O1B	1.96	0.81
2:C:612:ARG:O	2:C:616:LYS:HG3	1.83	0.79
1:B:144:GLY:HA2	5:B:501:GDP:O5'	1.83	0.77
2:C:613:SER:HA	2:C:616:LYS:CE	2.15	0.76
2:C:614:PHE:CD2	2:C:628:ILE:HG12	2.22	0.75
1:B:174:PRO:HG3	5:B:501:GDP:O3'	1.88	0.74
1:A:176:ARG:CZ	5:A:501:GDP:O2'	2.36	0.73
1:B:16:HIS:HB2	5:B:501:GDP:O6	1.90	0.72
2:C:613:SER:HA	2:C:616:LYS:HE3	1.70	0.71
2:C:613:SER:CA	2:C:616:LYS:HE3	2.19	0.71
2:C:614:PHE:CE2	2:C:628:ILE:CG1	2.76	0.68
2:C:614:PHE:HE2	2:C:628:ILE:CG1	2.06	0.68
2:C:613:SER:CA	2:C:616:LYS:HG3	2.24	0.68
1:A:289:HIS:HB3	1:A:371:GLU:HG3	1.77	0.67
1:A:289:HIS:HB3	1:A:371:GLU:CD	2.19	0.67
2:C:612:ARG:O	2:C:616:LYS:CG	2.43	0.66
2:C:628:ILE:HG22	2:C:629:ASN:N	2.11	0.66
2:C:613:SER:HA	2:C:616:LYS:HD2	1.75	0.65
1:B:143:ALA:HB3	5:B:501:GDP:C3'	2.28	0.63
2:C:618:ARG:C	2:C:618:ARG:HD2	2.23	0.63
1:B:146:THR:HB	5:B:501:GDP:PB	2.38	0.62
1:B:143:ALA:CB	5:B:501:GDP:C4'	2.69	0.62
1:B:174:PRO:HB3	5:B:501:GDP:O2'	2.01	0.61
1:A:145:GLY:N	5:A:501:GDP:O2B	2.35	0.60
2:C:576:ARG:HB3	2:C:577:PRO:CD	2.31	0.60
1:B:174:PRO:HG3	5:B:501:GDP:C3'	2.31	0.59
2:C:611:ILE:O	2:C:615:LYS:HG2	2.02	0.59
3:D:52:ARG:NH2	3:D:101:SER:OG	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:VAL:HG21	5:B:501:GDP:HN22	1.68	0.58
1:B:306:VAL:O	1:B:307:SER:C	2.46	0.58
1:A:176:ARG:NH1	5:A:501:GDP:O3'	2.37	0.58
1:B:144:GLY:HA2	5:B:501:GDP:O3A	2.04	0.58
1:A:289:HIS:HB3	1:A:371:GLU:CG	2.34	0.57
1:A:224:LEU:HD13	5:A:501:GDP:C4	2.41	0.55
2:C:826:PHE:CE2	2:C:830:LEU:HD11	2.41	0.55
1:B:143:ALA:CB	5:B:501:GDP:C3'	2.84	0.55
2:C:614:PHE:HE2	2:C:628:ILE:HG13	1.71	0.54
2:C:576:ARG:HB3	2:C:577:PRO:HD2	1.90	0.54
2:C:288:TYR:OH	3:D:123:TYR:OH	2.26	0.53
1:A:224:LEU:HD22	5:A:501:GDP:C2	2.44	0.53
1:B:146:THR:N	5:B:501:GDP:O2B	2.34	0.53
1:B:143:ALA:HB3	5:B:501:GDP:O3'	2.09	0.52
1:A:443:ASP:OD1	1:A:443:ASP:N	2.37	0.52
1:B:144:GLY:HA2	5:B:501:GDP:PA	2.49	0.52
2:C:498:ASP:N	2:C:498:ASP:OD1	2.43	0.51
1:B:143:ALA:CB	5:B:501:GDP:O3'	2.59	0.50
3:D:167:ASN:OD1	3:D:167:ASN:N	2.44	0.50
2:C:618:ARG:C	2:C:618:ARG:CD	2.85	0.49
2:C:613:SER:OG	2:C:616:LYS:NZ	2.43	0.49
2:C:612:ARG:C	2:C:616:LYS:HE3	2.37	0.49
1:B:146:THR:CB	5:B:501:GDP:O2B	2.59	0.49
1:A:141:SER:HG	5:A:501:GDP:H4'	1.71	0.49
2:C:629:ASN:O	2:C:633:ARG:HG2	2.12	0.49
3:D:303:ASP:OD2	3:D:355:LYS:NZ	2.46	0.49
1:B:51:LYS:HB3	1:B:52:PRO:HD3	1.95	0.49
1:B:13:CYS:HA	5:B:501:GDP:C5	2.47	0.49
2:C:665:THR:O	2:C:669:GLN:NE2	2.47	0.48
2:C:487:ASP:N	2:C:487:ASP:OD1	2.46	0.48
1:B:146:THR:OG1	5:B:501:GDP:O2B	2.29	0.48
3:D:790:LYS:O	3:D:791:PHE:C	2.57	0.47
2:C:613:SER:N	2:C:616:LYS:HE3	2.29	0.47
1:A:191:LEU:O	1:A:192:ARG:C	2.57	0.47
2:C:685:ASN:OD1	2:C:685:ASN:N	2.47	0.47
1:B:201:THR:OG1	1:B:202:VAL:N	2.47	0.47
3:D:755:ASP:N	3:D:755:ASP:OD1	2.48	0.47
1:A:209:LEU:HD11	5:A:501:GDP:N2	2.30	0.47
2:C:554:GLY:O	2:C:556:ASP:N	2.48	0.46
1:A:103:ASN:OD1	5:A:501:GDP:O2B	2.34	0.46
2:C:793:ASN:O	2:C:795:HIS:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:94:ILE:O	3:D:95:ALA:C	2.59	0.45
1:B:205:ASP:OD1	1:B:205:ASP:C	2.58	0.45
1:B:224:LEU:HD22	5:B:501:GDP:C2	2.52	0.45
2:C:442:SER:OG	2:C:443:THR:N	2.50	0.45
1:A:340:GLN:O	1:A:341:ARG:CB	2.65	0.45
1:A:29:HIS:O	1:A:30:ALA:C	2.60	0.45
1:B:106:ALA:O	1:B:107:ASN:C	2.58	0.45
1:B:197:ASP:OD1	1:B:197:ASP:C	2.59	0.45
2:C:674:LYS:O	2:C:675:SER:C	2.59	0.45
1:B:229:GLN:OE1	1:B:229:GLN:N	2.47	0.44
2:C:163:SER:OG	2:C:164:SER:N	2.49	0.44
1:A:320:ASN:OD1	1:A:320:ASN:C	2.61	0.44
1:A:4:GLU:HG3	1:A:130:THR:HB	1.99	0.44
1:B:205:ASP:OD1	1:B:205:ASP:N	2.46	0.44
2:C:349:TRP:CE2	2:C:406:GLY:HA3	2.53	0.44
1:B:246:SER:OG	1:B:247:TYR:N	2.47	0.43
1:B:289:HIS:HB3	1:B:293:ASP:HB2	2.00	0.43
1:B:144:GLY:HA3	5:B:501:GDP:O5'	2.15	0.43
2:C:624:ILE:HG23	2:C:628:ILE:CD1	2.49	0.43
3:D:17:ASN:OD1	3:D:17:ASN:N	2.50	0.43
1:B:275:PHE:CD2	1:B:275:PHE:N	2.86	0.43
2:C:215:ASN:OD1	2:C:215:ASN:N	2.48	0.43
2:C:322:HIS:O	2:C:328:ARG:NE	2.52	0.43
2:C:613:SER:OG	2:C:616:LYS:CE	2.66	0.43
2:C:614:PHE:HA	2:C:787:PHE:CD1	2.53	0.43
3:D:249:LYS:NZ	3:D:334:ASP:OD1	2.39	0.43
2:C:691:LEU:HB2	2:C:694:ILE:HD11	2.01	0.42
1:A:88:ASP:OD1	1:A:88:ASP:C	2.59	0.42
1:B:141:SER:OG	5:B:501:GDP:C4'	2.68	0.42
1:B:144:GLY:CA	5:B:501:GDP:PA	3.08	0.42
2:C:624:ILE:HD13	2:C:624:ILE:HA	1.91	0.42
3:D:380:ASP:C	3:D:380:ASP:OD1	2.63	0.42
1:A:223:ASP:CG	1:A:224:LEU:H	2.27	0.42
1:A:394:PHE:CD2	1:A:394:PHE:C	2.93	0.42
2:C:624:ILE:HG23	2:C:628:ILE:HD11	2.01	0.42
3:D:813:PRO:HA	3:D:814:PRO:HD3	1.95	0.42
1:B:140:HIS:CG	1:B:141:SER:N	2.87	0.41
2:C:824:TYR:O	2:C:827:ARG:HG2	2.20	0.41
3:D:123:TYR:CD1	3:D:123:TYR:C	2.98	0.41
1:A:292:TYR:HB2	1:A:330:GLN:HB3	2.01	0.41
1:A:320:ASN:HD22	1:A:376:MET:HB2	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:HIS:CB	5:B:501:GDP:O6	2.64	0.41
1:A:12:GLN:OE1	5:A:501:GDP:O3B	2.37	0.41
1:A:201:THR:HG23	1:A:267:PHE:CD2	2.56	0.41
3:D:18:ASN:OD1	3:D:18:ASN:N	2.43	0.41
2:C:398:LEU:O	2:C:399:ALA:C	2.61	0.41
3:D:193:TYR:CD2	3:D:193:TYR:C	2.98	0.41
3:D:349:ARG:O	3:D:350:THR:C	2.63	0.41
3:D:145:PHE:CD2	3:D:145:PHE:C	2.98	0.41
3:D:647:PHE:O	3:D:651:VAL:HB	2.20	0.41
1:B:146:THR:HB	5:B:501:GDP:O1B	2.20	0.41
2:C:621:ASN:HB2	2:C:622:PRO:HD2	1.85	0.41
1:A:289:HIS:CB	1:A:371:GLU:HG3	2.47	0.41
1:B:141:SER:HG	5:B:501:GDP:C4'	2.33	0.41
2:C:572:LEU:O	2:C:573:ASN:C	2.64	0.41
2:C:310:ASP:OD1	2:C:311:THR:N	2.54	0.40
2:C:576:ARG:CB	2:C:577:PRO:CD	2.96	0.40
3:D:73:THR:O	3:D:76:ARG:NH2	2.55	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	437/473 (92%)	420 (96%)	15 (3%)	2 (0%)	24	60
1	B	443/473 (94%)	432 (98%)	11 (2%)	0	100	100
2	C	650/846 (77%)	634 (98%)	12 (2%)	4 (1%)	21	57
3	D	691/823 (84%)	681 (99%)	10 (1%)	0	100	100
4	U	29/220 (13%)	28 (97%)	1 (3%)	0	100	100
All	All	2250/2835 (79%)	2195 (98%)	49 (2%)	6 (0%)	37	69

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	555	TRP
2	C	628	ILE
1	A	135	GLY
1	A	341	ARG
2	C	621	ASN
2	C	627	ILE

5.3.2 Protein sidechains [i](#)

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	393/420 (94%)	393 (100%)	0	100	100
1	B	397/420 (94%)	397 (100%)	0	100	100
2	C	613/787 (78%)	612 (100%)	1 (0%)	87	87
3	D	659/766 (86%)	659 (100%)	0	100	100
4	U	30/206 (15%)	30 (100%)	0	100	100
All	All	2092/2599 (80%)	2091 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	624	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	GLN
1	B	9	GLN
1	B	366	GLN
2	C	208	GLN
2	C	423	ASN
2	C	444	ASN

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Mol	Chain	Res	Type
2	C	608	ASN
3	D	35	ASN
3	D	417	ASN
3	D	485	GLN
3	D	571	ASN
3	D	589	GLN
3	D	593	GLN
3	D	762	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	501	-	29,30,30	1.14	2 (6%)	45,47,47	1.80	8 (17%)
5	GDP	B	501	-	29,30,30	1.14	2 (6%)	45,47,47	1.80	8 (17%)

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	501	-	-	4/16/32/32	0/3/3/3
5	GDP	B	501	-	-	4/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	501	GDP	C5-C4	3.05	1.47	1.38
5	A	501	GDP	C5-C4	3.03	1.47	1.38
5	B	501	GDP	C6-N1	-2.03	1.35	1.38
5	A	501	GDP	C6-N1	-2.02	1.35	1.38

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GDP	C5-C4-N3	-5.74	119.26	128.39
5	B	501	GDP	C5-C4-N3	-5.73	119.27	128.39
5	B	501	GDP	C2-N3-C4	5.32	121.46	112.30
5	A	501	GDP	C2-N3-C4	5.27	121.38	112.30
5	A	501	GDP	N9-C4-N3	4.28	134.51	125.95
5	B	501	GDP	N9-C4-N3	4.27	134.49	125.95
5	B	501	GDP	C6-C5-N7	3.88	137.34	130.29
5	A	501	GDP	C6-C5-N7	3.86	137.32	130.29
5	A	501	GDP	C4-C5-N7	-2.65	106.47	110.67
5	B	501	GDP	C4-C5-N7	-2.64	106.49	110.67
5	B	501	GDP	O6-C6-C5	-2.29	120.50	126.53
5	A	501	GDP	O6-C6-C5	-2.27	120.54	126.53
5	B	501	GDP	C5-C6-N1	2.15	118.73	113.25
5	A	501	GDP	C5-C6-N1	2.14	118.70	113.25
5	B	501	GDP	O3'-C3'-C4'	-2.07	105.14	111.08
5	A	501	GDP	O3'-C3'-C4'	-2.07	105.14	111.08

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GDP	C5'-O5'-PA-O3A
5	A	501	GDP	C5'-O5'-PA-O2A
5	B	501	GDP	C5'-O5'-PA-O3A
5	B	501	GDP	C5'-O5'-PA-O2A

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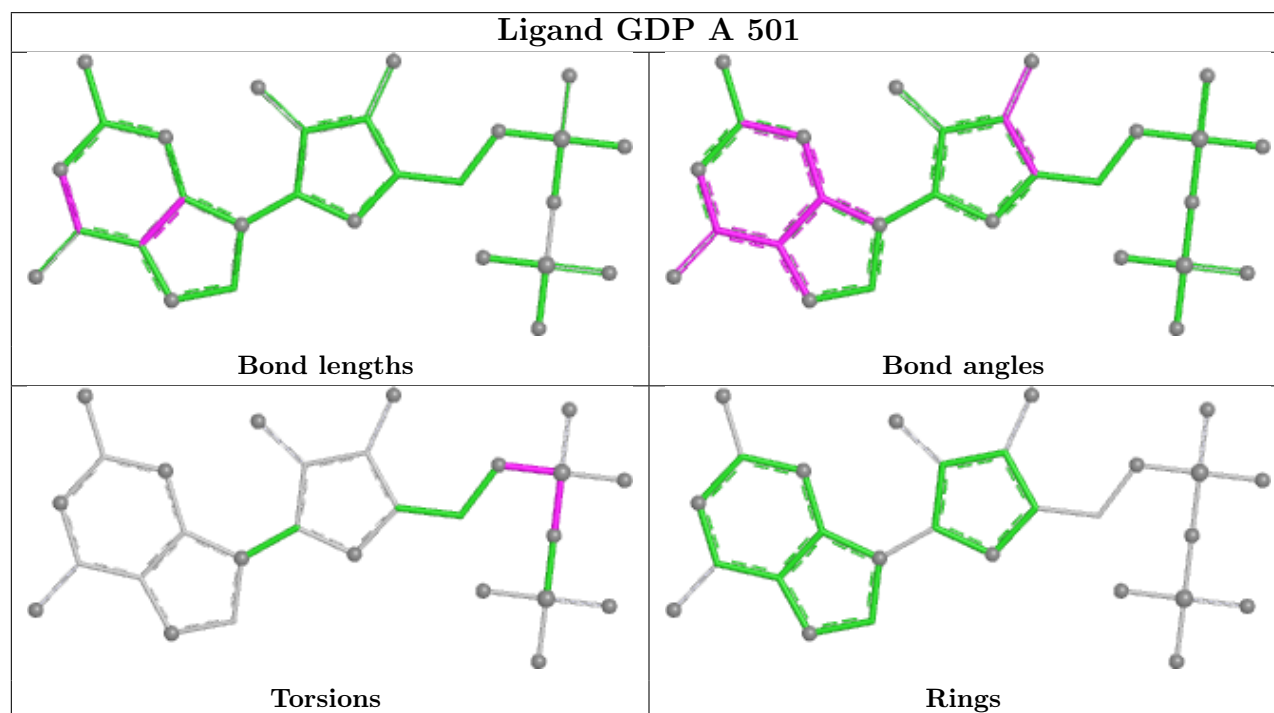
Mol	Chain	Res	Type	Atoms
5	A	501	GDP	C5'-O5'-PA-O1A
5	B	501	GDP	C5'-O5'-PA-O1A
5	A	501	GDP	PB-O3A-PA-O2A
5	B	501	GDP	PB-O3A-PA-O2A

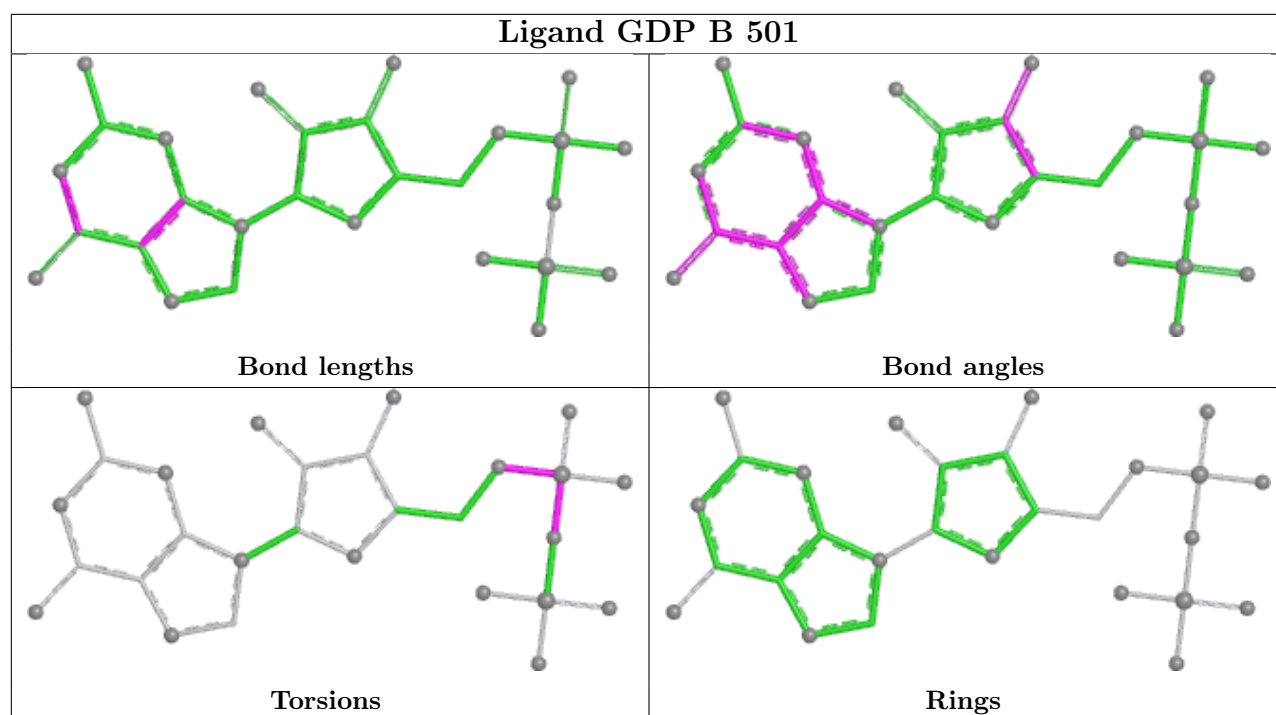
There are no ring outliers.

2 monomers are involved in 45 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	501	GDP	13	0
5	B	501	GDP	32	0

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.

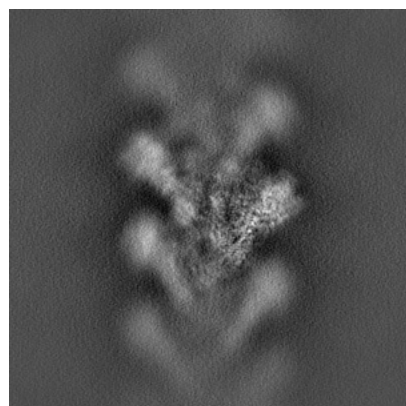
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23637. These allow visual inspection of the internal detail of the map and identification of artifacts.

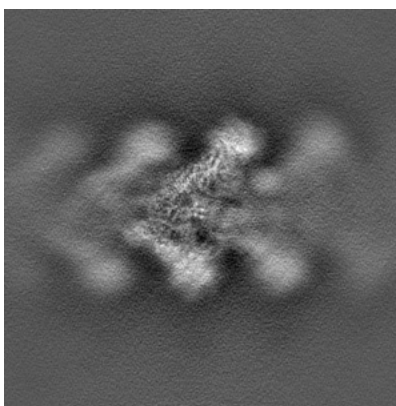
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

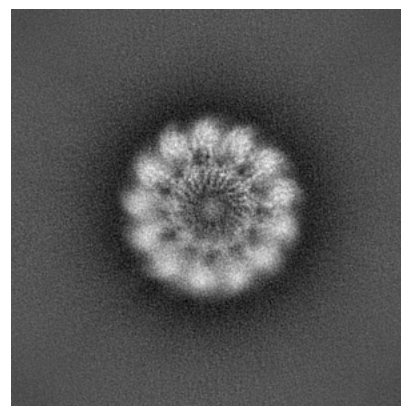
6.1.1 Primary map



X

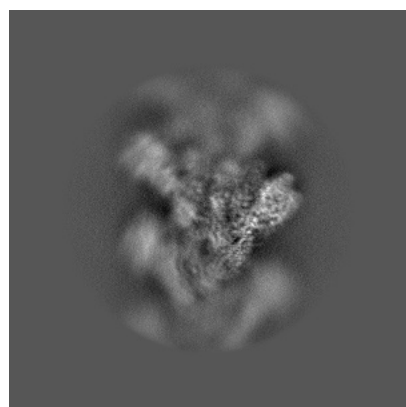


Y

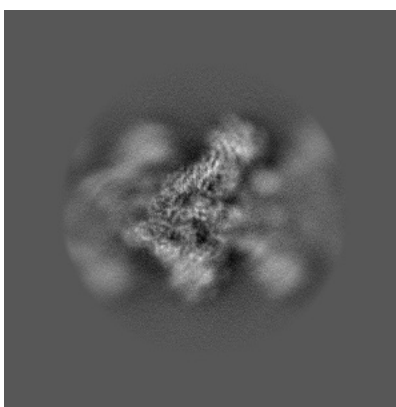


Z

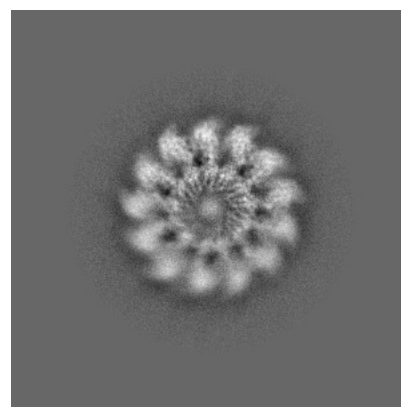
6.1.2 Raw map



X



Y

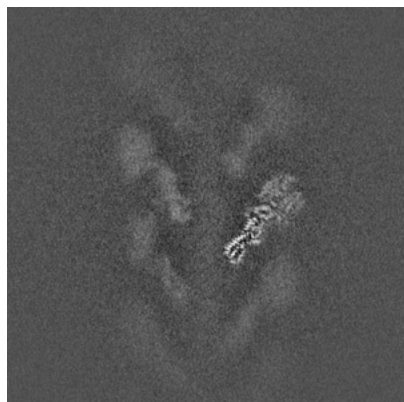


Z

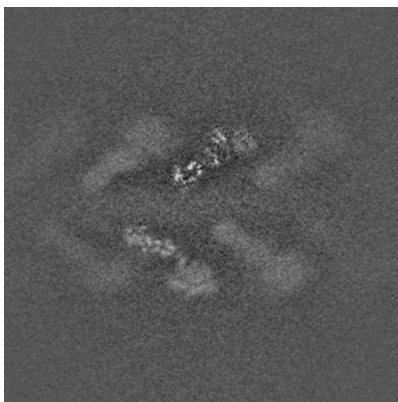
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

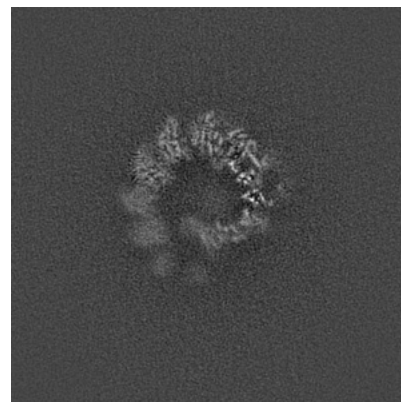
6.2.1 Primary map



X Index: 224

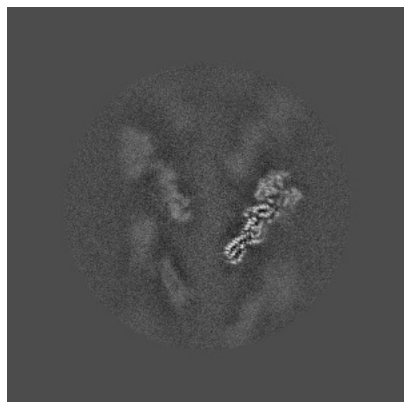


Y Index: 224

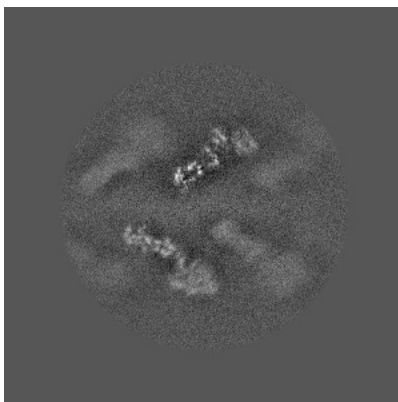


Z Index: 224

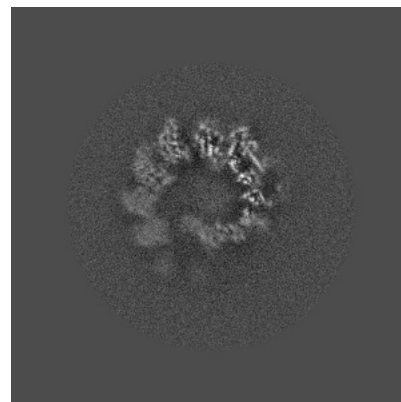
6.2.2 Raw map



X Index: 224



Y Index: 224

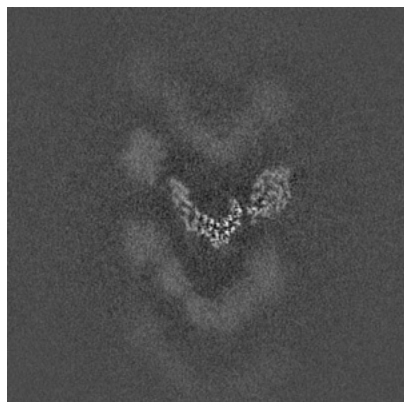


Z Index: 224

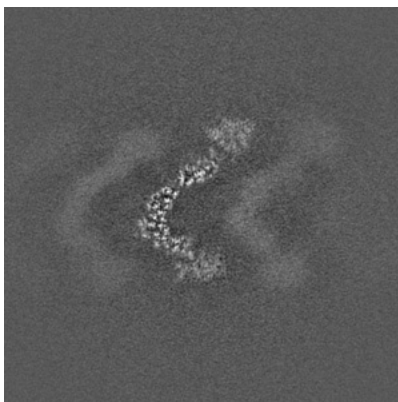
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

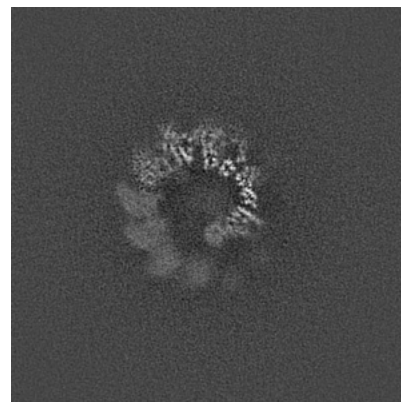
6.3.1 Primary map



X Index: 256

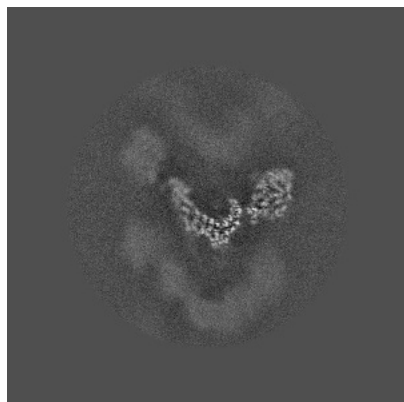


Y Index: 252

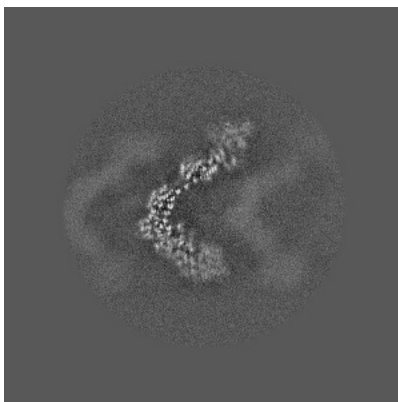


Z Index: 213

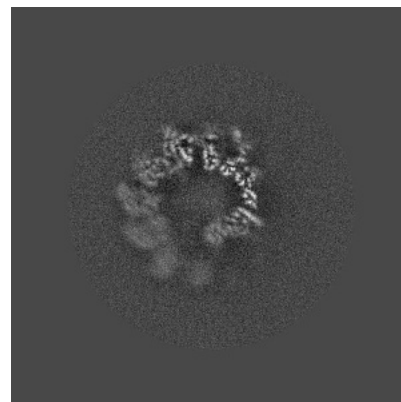
6.3.2 Raw map



X Index: 256



Y Index: 256

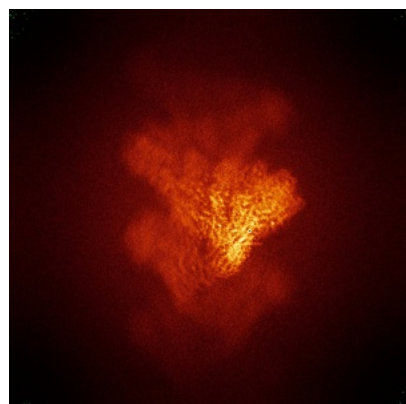


Z Index: 214

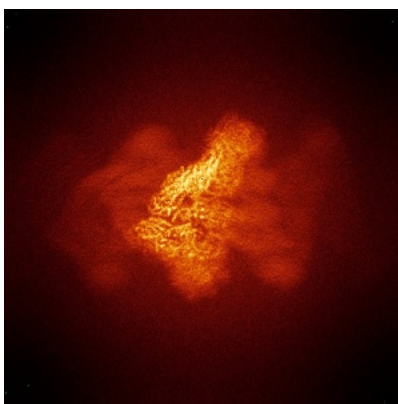
The images above show the largest variance slices of the map in three orthogonal directions.

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

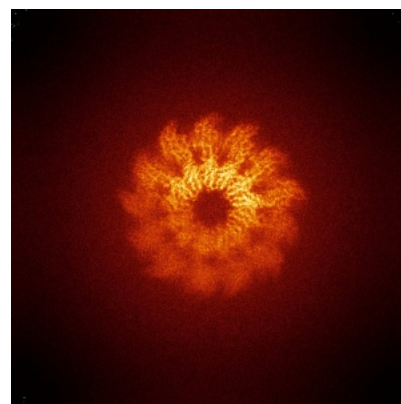
6.4.1 Primary map



X

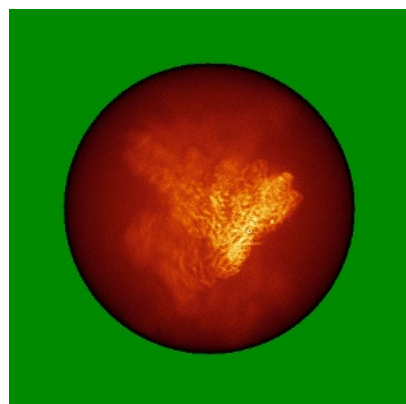


Y

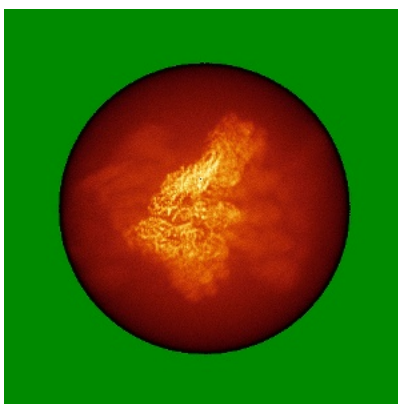


Z

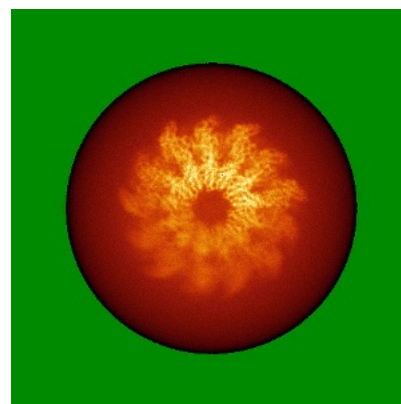
6.4.2 Raw map



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](#)

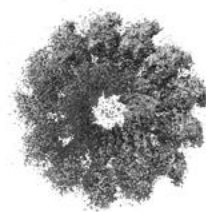
6.5.1 Primary map



X



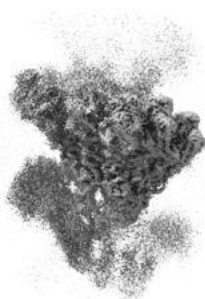
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 3.4. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

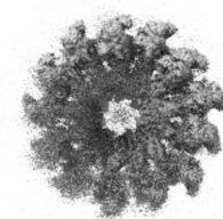
6.5.2 Raw map



X



Y



Z

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

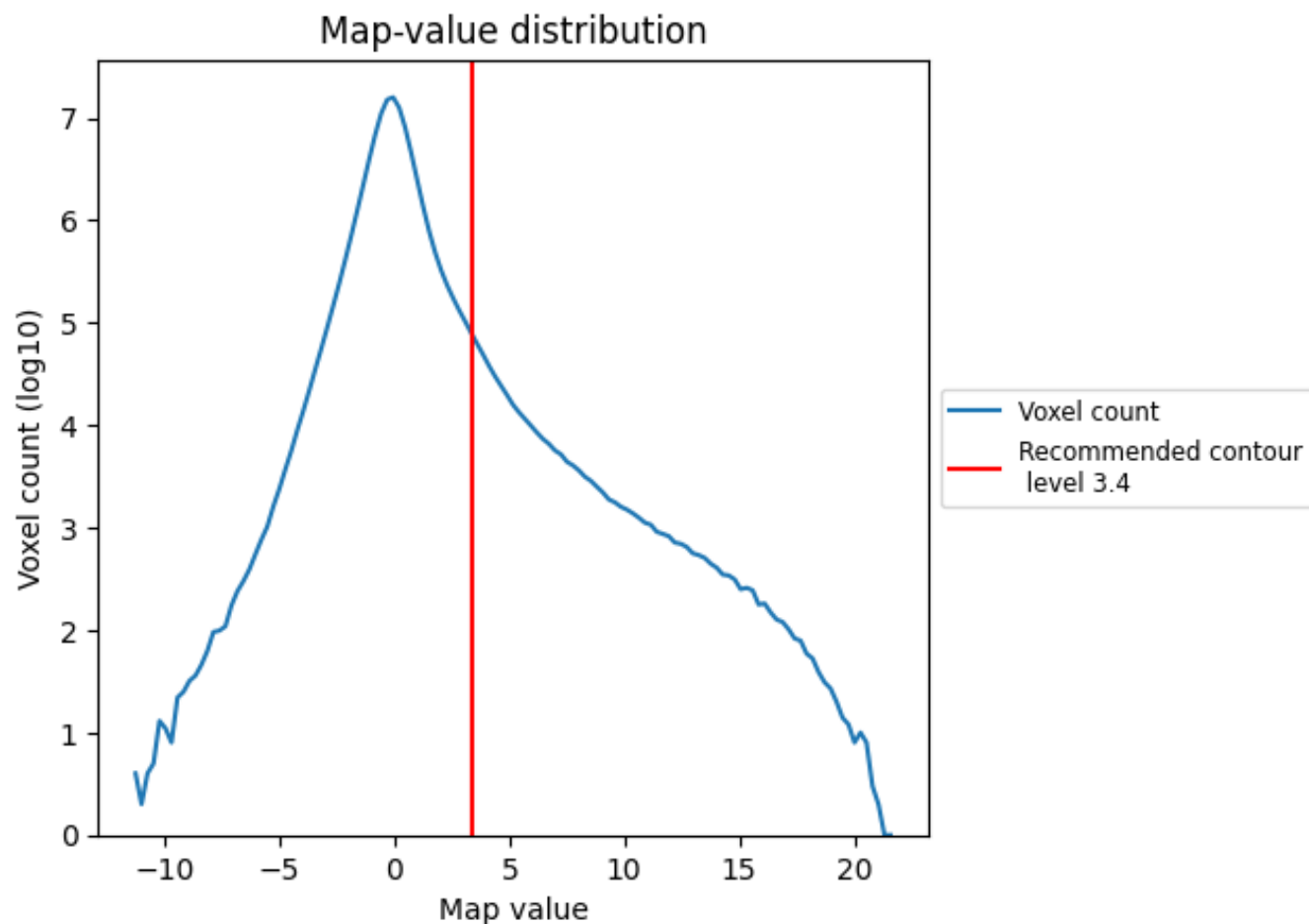
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

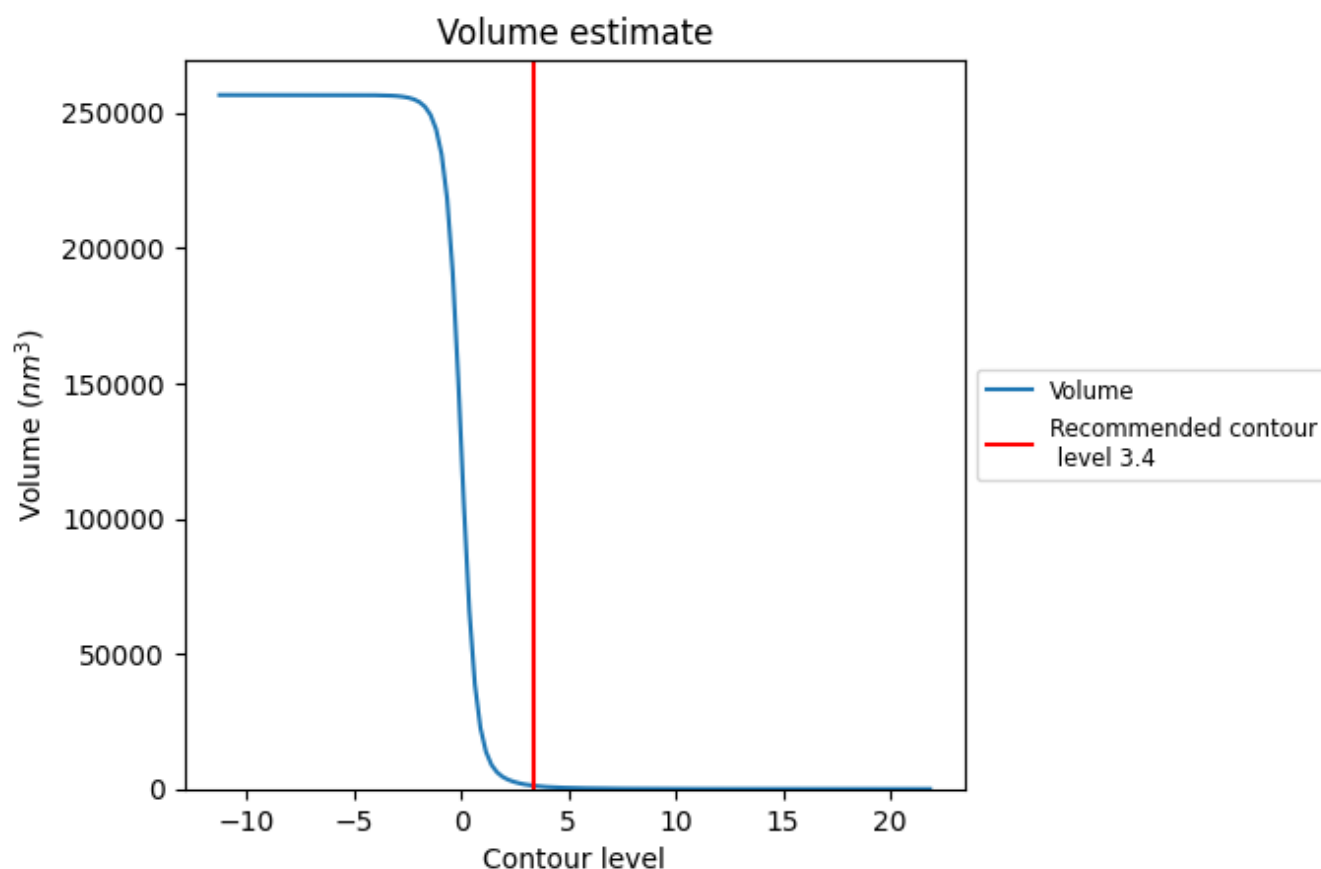
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

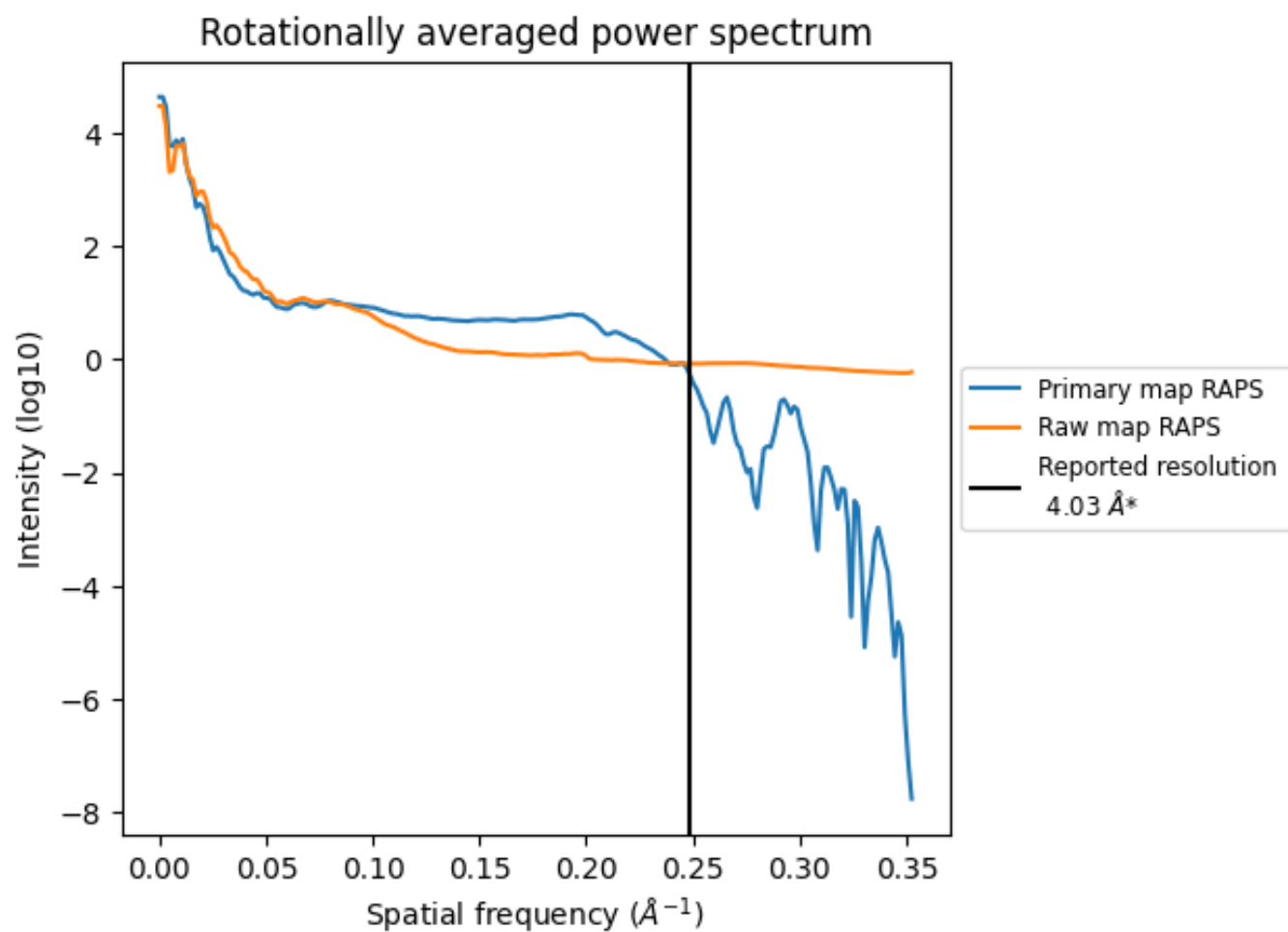
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1173 nm^3 ; this corresponds to an approximate mass of 1060 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 ⓘ

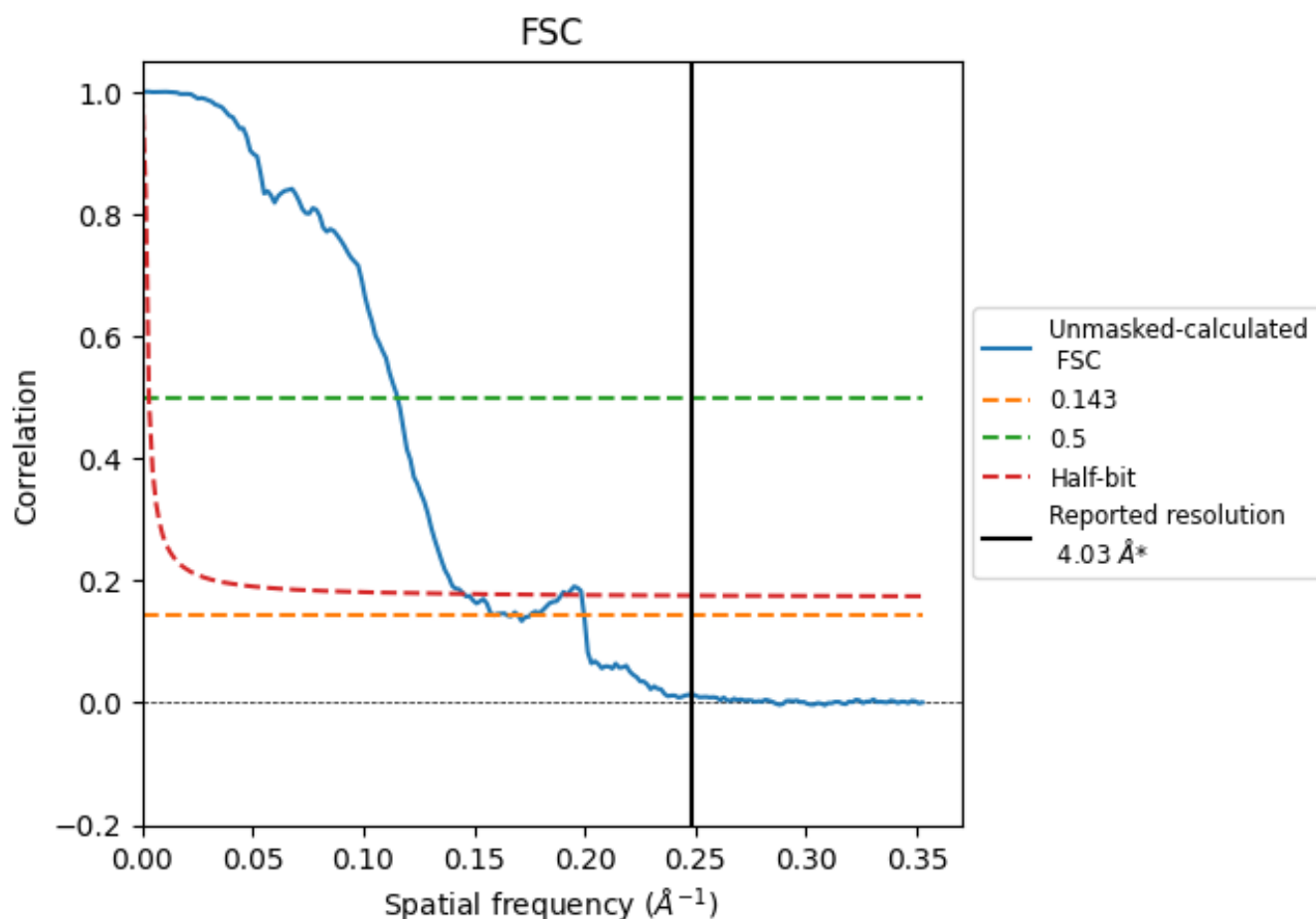


*Reported resolution corresponds to spatial frequency of 0.248 Å⁻¹

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 0.248 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

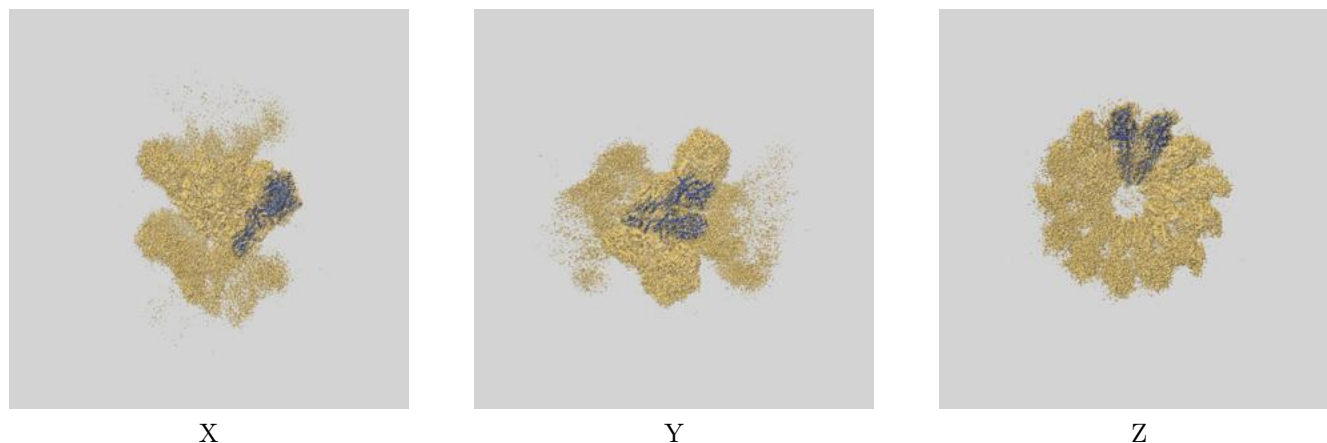
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.03	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	6.29	8.67	6.88

*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 6.29 differs from the reported value 4.03 by more than 10 %

9 Map-model fit [i](#)

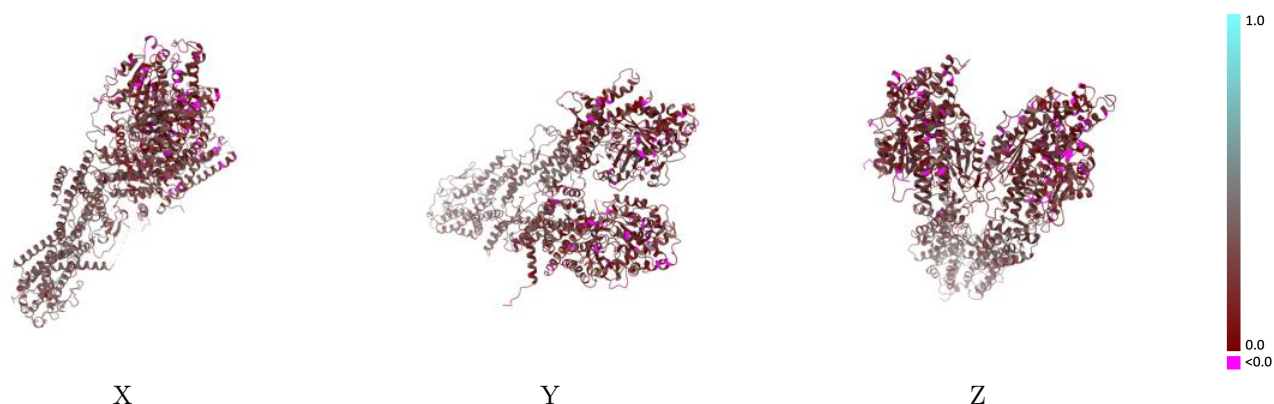
This section contains information regarding the fit between EMDB map EMD-23637 and PDB model 7M2Y. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 3.4 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](#)

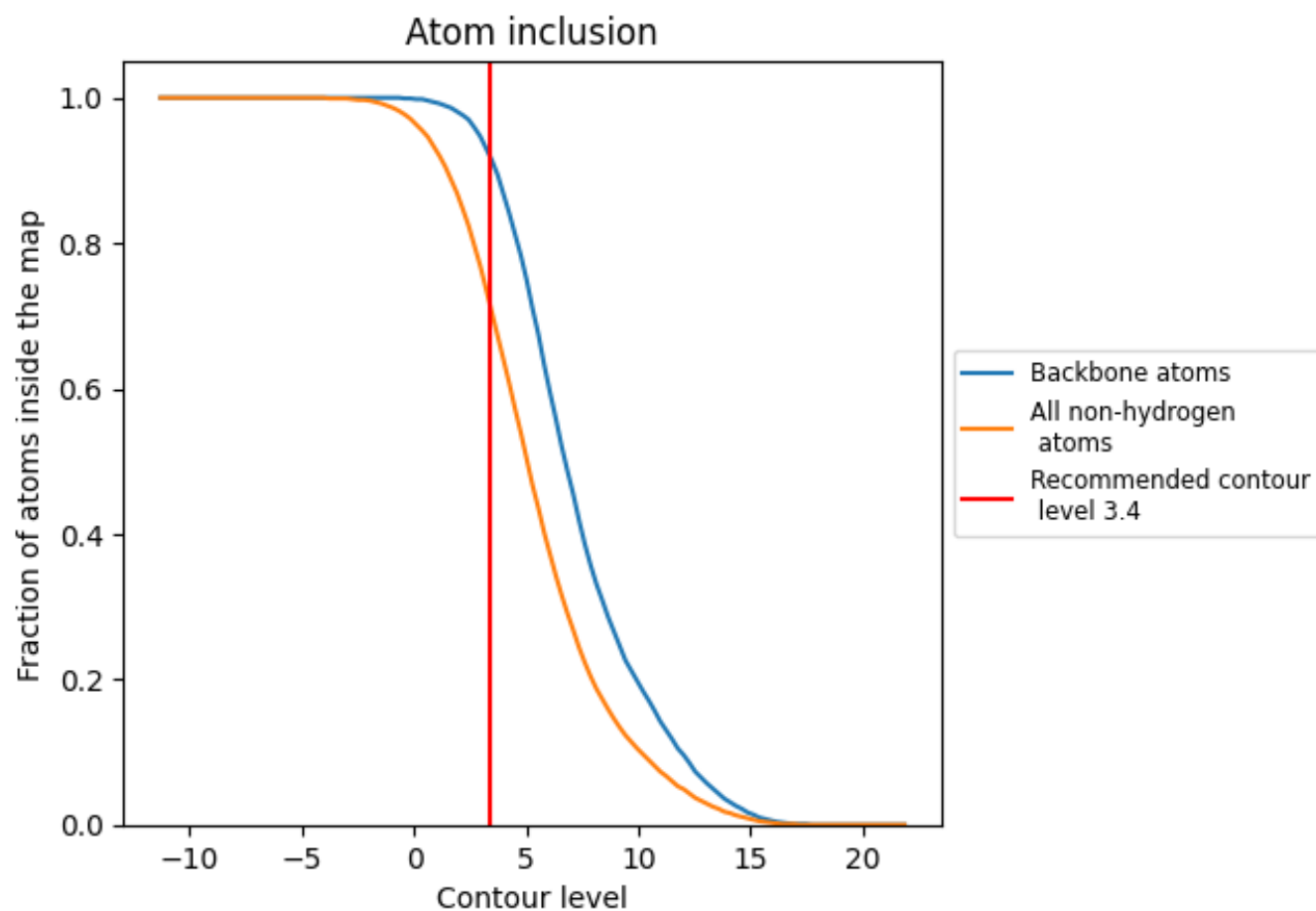


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 72% 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 (3.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7150	0.2490
A	0.6780	0.2060
B	0.6720	0.1950
C	0.7460	0.2780
D	0.7370	0.2830
U	0.6650	0.1940

