



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 10:41 PM UTC

PDB ID : 5SYE / pdb_00005sye
EMDB ID : EMD-8321
Title : Near-atomic resolution cryo-EM reconstruction of doubly bound Taxol- and p
eloruside-stabilized microtubule
Authors : Kellogg, E.H.; Nogales, E.
Deposited on : 2016-08-10
Resolution : 3.50 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

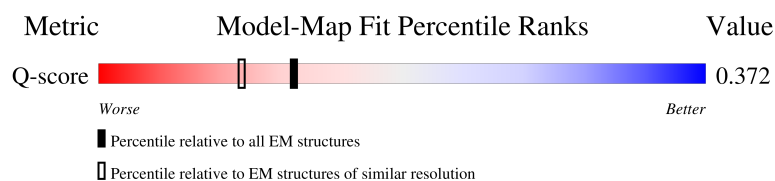
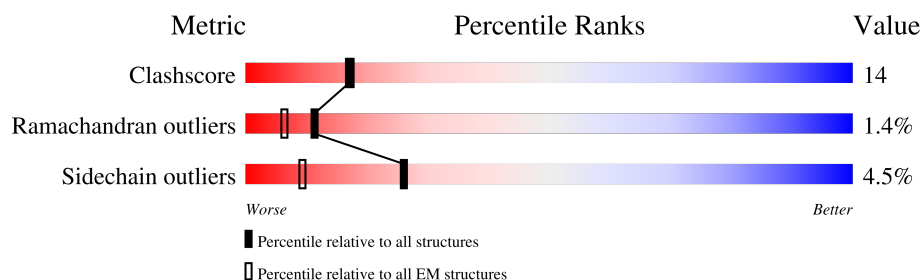
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	13950 (3.00 - 4.00)

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	437	 78% 18% . .
2	B	426	 83% 16% .

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
7	POU	B	503	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6864 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.

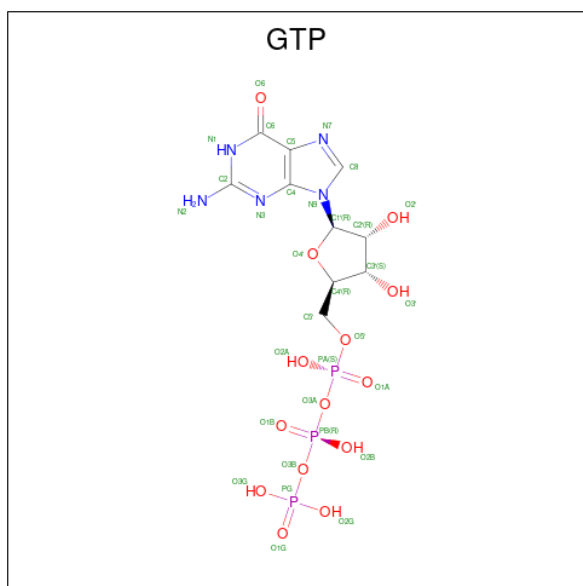
- Molecule 1 is a protein called Tubulin alpha chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	427	3351	2125	570	634	22	0	0

- Molecule 2 is a protein called Tubulin beta chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	426	3352	2106	575	645	26	0	0

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).

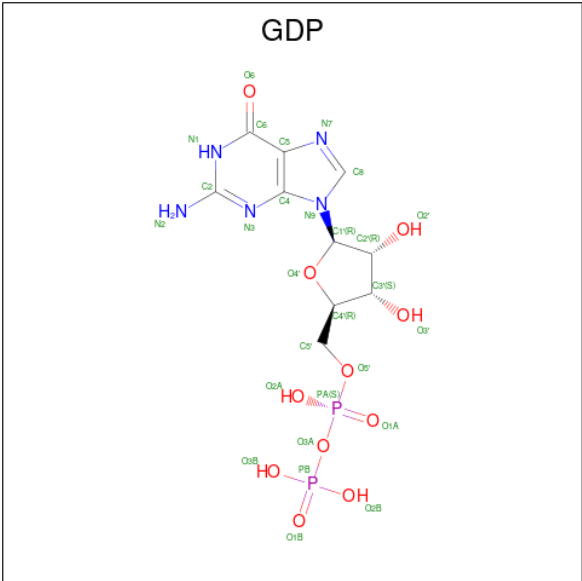


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	32	10	5	14	3	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

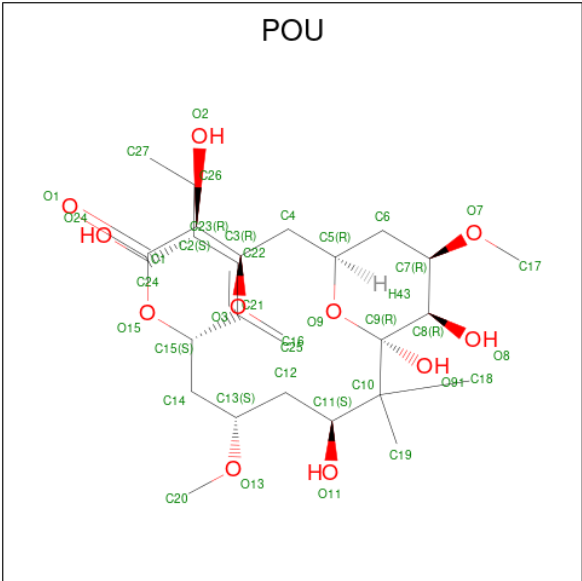
Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Mg	0
			1	1	

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



Mol	Chain	Residues	Atoms				AltConf
6	B	1	Total	C	N	O	0
			62	47	1	14	

- Molecule 7 is Peloruside A (CCD ID: POU) (formula: C₂₇H₄₈O₁₁).

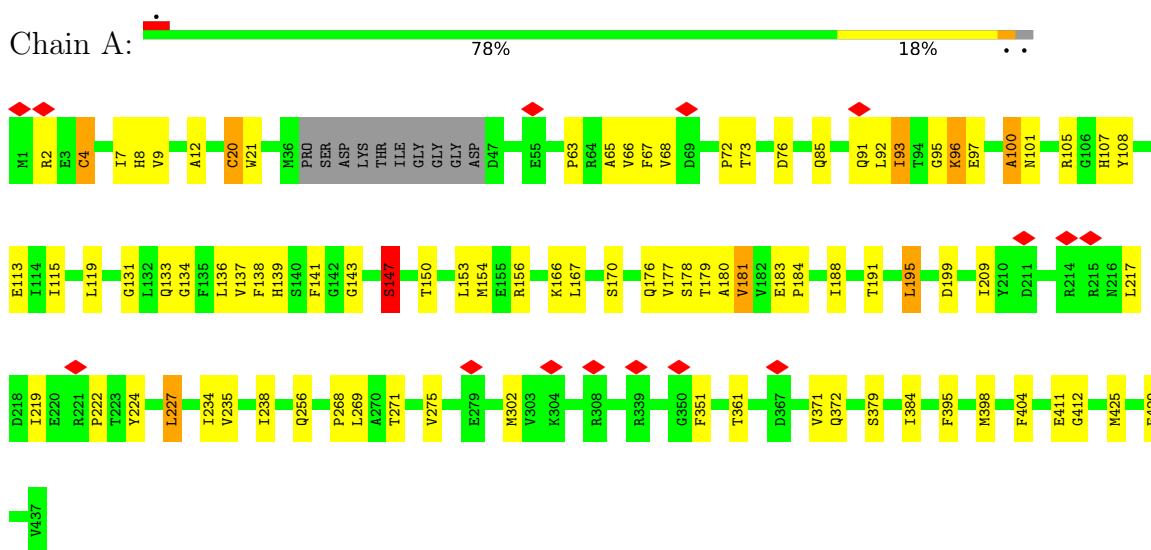


Mol	Chain	Residues	Atoms			AltConf
7	B	1	Total	C	O	0
			38	27	11	

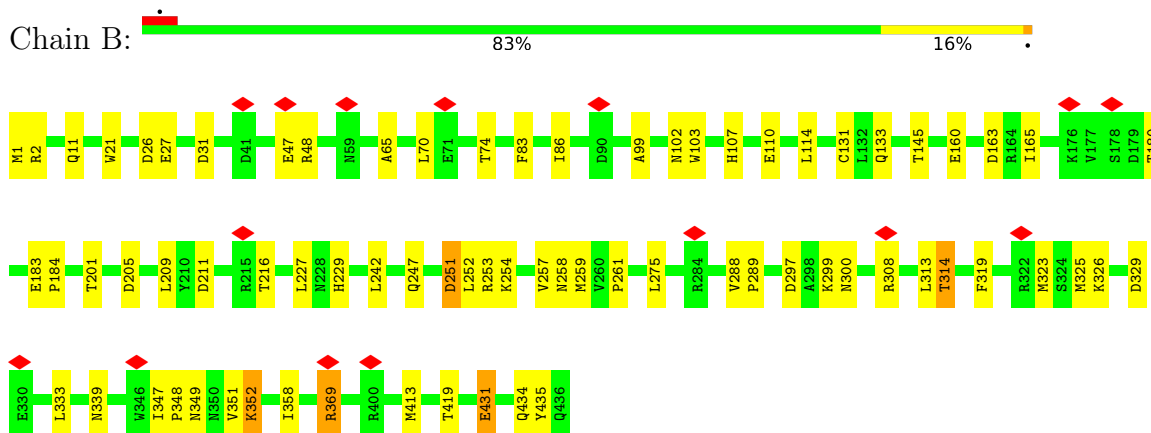
3 Residue-property plots

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 alpha chain



• Molecule 2: Tubulin beta chain



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-27.7°, rise=9.43 Å, axial sym=C1	Depositor
Number of segments used	17069	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{Å}^2$)	8	Depositor
Minimum defocus (nm)	1100	Depositor
Maximum defocus (nm)	3800	Depositor
Magnification	27500	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	23.832	Depositor
Minimum map value	-10.062	Depositor
Average map value	0.065	Depositor
Map value standard deviation	1.162	Depositor
Recommended contour level	5.59	Depositor
Map size (Å)	675.84, 675.84, 675.84	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.32, 1.32, 1.32	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: TA1, GTP, GDP, MG, POU

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.51	0/3427	0.94	1/4650 (0.0%)
2	B	0.55	0/3427	0.87	0/4641
All	All	0.53	0/6854	0.91	1/9291 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4	CYS	N-CA-C	5.16	116.82	109.14

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3351	0	3265	114	0
2	B	3352	0	3237	146	0
3	A	32	0	12	1	0
4	A	1	0	0	0	0
5	B	28	0	12	0	0
6	B	62	0	51	12	0
7	B	38	0	48	23	0
All	All	6864	0	6625	186	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 14.

All (186) 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:A:398:MET:HE2	2:B:348:PRO:CD	1.16	1.55
1:A:398:MET:CE	2:B:348:PRO:HD2	1.07	1.53
2:B:165:ILE:CD1	2:B:253:ARG:HG3	1.50	1.37
1:A:73:THR:HA	2:B:48:ARG:NH2	1.42	1.33
1:A:404:PHE:CE2	2:B:261:PRO:HB3	1.65	1.31
2:B:165:ILE:HD11	2:B:253:ARG:CG	1.60	1.29
2:B:163:ASP:O	2:B:253:ARG:NH2	1.66	1.25
1:A:404:PHE:CD2	2:B:261:PRO:HA	1.72	1.25
2:B:297:ASP:OD1	7:B:503:POU:C2	1.90	1.18
2:B:299:LYS:NZ	7:B:503:POU:H34	1.54	1.18
1:A:404:PHE:CE2	2:B:261:PRO:CB	2.28	1.16
2:B:297:ASP:OD1	7:B:503:POU:O2	1.61	1.15
2:B:229:HIS:NE2	6:B:502:TA1:O14	1.81	1.13
1:A:404:PHE:CE2	2:B:261:PRO:CA	2.32	1.12
2:B:26:ASP:OD2	6:B:502:TA1:H331	1.54	1.07
1:A:224:TYR:CE1	2:B:325:MET:HG3	1.88	1.06
1:A:398:MET:HE3	2:B:348:PRO:HD2	1.39	1.04
1:A:404:PHE:CE2	2:B:261:PRO:HA	1.92	1.03
1:A:178:SER:HB3	2:B:349:ASN:ND2	1.76	1.01
1:A:398:MET:HE1	2:B:347:ILE:HG23	1.38	1.01
1:A:100:ALA:HB1	2:B:254:LYS:HA	1.43	0.97
1:A:95:GLY:O	2:B:1:MET:HE2	1.65	0.95
1:A:404:PHE:HE2	2:B:261:PRO:CB	1.80	0.94
1:A:404:PHE:CZ	2:B:261:PRO:HB3	2.03	0.94
1:A:181:VAL:CG2	2:B:349:ASN:O	2.16	0.94
2:B:308:ARG:HH11	7:B:503:POU:C16	1.82	0.92
1:A:105:ARG:HH22	2:B:253:ARG:NH1	1.68	0.92
1:A:398:MET:HE2	2:B:348:PRO:HD3	1.52	0.92
1:A:7:ILE:HD12	1:A:137:VAL:HG22	1.52	0.91
2:B:297:ASP:OD1	7:B:503:POU:H44	1.68	0.91
1:A:73:THR:HA	2:B:48:ARG:HH21	1.07	0.89
2:B:299:LYS:HZ3	7:B:503:POU:H34	1.38	0.89
2:B:308:ARG:NH1	7:B:503:POU:O3	2.06	0.89
2:B:299:LYS:NZ	7:B:503:POU:C6	2.36	0.88
1:A:404:PHE:HE2	2:B:261:PRO:CA	1.79	0.87
2:B:297:ASP:CG	7:B:503:POU:H44	2.01	0.86
1:A:398:MET:HE1	2:B:347:ILE:CG2	2.05	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:THR:HG21	1:A:425:MET:HE1	1.58	0.85
1:A:404:PHE:HD2	2:B:261:PRO:HA	1.35	0.85
1:A:73:THR:CA	2:B:48:ARG:NH2	2.35	0.84
1:A:105:ARG:HH22	2:B:253:ARG:HH11	1.24	0.82
2:B:299:LYS:HZ1	7:B:503:POU:H34	1.41	0.81
1:A:73:THR:HA	2:B:48:ARG:HH22	1.42	0.80
1:A:73:THR:CA	2:B:48:ARG:HH21	1.93	0.80
1:A:181:VAL:N	2:B:258:ASN:OD1	2.14	0.80
1:A:180:ALA:HB1	2:B:258:ASN:HD21	1.45	0.80
2:B:229:HIS:HE1	6:B:502:TA1:H132	1.47	0.80
1:A:100:ALA:HB1	2:B:254:LYS:CA	2.11	0.80
1:A:95:GLY:O	2:B:1:MET:CE	2.26	0.78
2:B:165:ILE:CD1	2:B:253:ARG:CG	2.38	0.78
1:A:181:VAL:CG2	2:B:351:VAL:O	2.31	0.78
1:A:224:TYR:CD1	2:B:325:MET:HG3	2.19	0.77
2:B:70:LEU:HD23	2:B:114:LEU:HD22	1.66	0.76
1:A:181:VAL:HG21	2:B:349:ASN:O	1.84	0.76
2:B:165:ILE:HD11	2:B:253:ARG:HG3	0.76	0.75
1:A:181:VAL:CG1	2:B:349:ASN:O	2.34	0.75
1:A:72:PRO:HD2	2:B:2:ARG:NH1	2.03	0.74
1:A:72:PRO:HG2	2:B:2:ARG:NH1	2.03	0.74
1:A:76:ASP:OD2	2:B:48:ARG:NH1	2.21	0.73
1:A:139:HIS:HB3	1:A:150:THR:HG21	1.69	0.73
2:B:165:ILE:HG13	2:B:253:ARG:HE	1.52	0.73
2:B:247:GLN:OE1	2:B:325:MET:HE3	1.91	0.71
1:A:181:VAL:HG23	2:B:351:VAL:O	1.89	0.71
1:A:178:SER:OG	2:B:349:ASN:HB3	1.91	0.70
1:A:398:MET:CE	2:B:348:PRO:CD	2.02	0.70
1:A:181:VAL:HG22	2:B:349:ASN:O	1.91	0.69
2:B:275:LEU:HD12	2:B:300:ASN:OD1	1.92	0.69
1:A:176:GLN:HB3	2:B:333:LEU:HD11	1.73	0.68
2:B:229:HIS:HE2	6:B:502:TA1:C30	2.05	0.67
1:A:224:TYR:CE1	2:B:325:MET:CG	2.72	0.66
2:B:2:ARG:HH22	2:B:48:ARG:CZ	2.08	0.66
1:A:72:PRO:CG	2:B:2:ARG:NH1	2.58	0.66
1:A:72:PRO:CG	2:B:2:ARG:HH12	2.08	0.66
1:A:177:VAL:CG2	2:B:329:ASP:OD2	2.42	0.66
1:A:12:ALA:HB2	3:A:501:GTP:C8	2.32	0.65
2:B:165:ILE:HD11	2:B:253:ARG:CD	2.26	0.65
2:B:299:LYS:HZ2	7:B:503:POU:C5	2.10	0.65
1:A:100:ALA:HA	2:B:254:LYS:HE3	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ASP:OD2	2:B:48:ARG:NH2	2.30	0.64
1:A:178:SER:CB	2:B:349:ASN:ND2	2.59	0.64
2:B:308:ARG:NH1	7:B:503:POU:C16	2.58	0.64
1:A:96:LYS:HG2	2:B:131:CYS:CB	2.28	0.62
2:B:299:LYS:HZ2	7:B:503:POU:H43	1.66	0.61
2:B:369:ARG:HB2	6:B:502:TA1:O13	2.00	0.61
2:B:369:ARG:CB	6:B:502:TA1:O13	2.48	0.61
6:B:502:TA1:H463	6:B:502:TA1:H261	1.82	0.61
2:B:165:ILE:CG1	2:B:253:ARG:HE	2.13	0.60
2:B:319:PHE:CD2	2:B:323:MET:HE1	2.36	0.60
1:A:177:VAL:HG21	2:B:329:ASP:OD2	2.02	0.60
1:A:72:PRO:HG2	2:B:2:ARG:HH12	1.67	0.59
1:A:72:PRO:CD	2:B:2:ARG:NH1	2.65	0.59
1:A:7:ILE:HG23	1:A:66:VAL:HB	1.85	0.58
2:B:297:ASP:OD1	7:B:503:POU:C1	2.51	0.58
1:A:180:ALA:HB1	2:B:258:ASN:ND2	2.16	0.57
1:A:404:PHE:HE2	2:B:261:PRO:N	2.02	0.57
1:A:178:SER:HB3	2:B:349:ASN:CG	2.28	0.57
2:B:247:GLN:OE1	2:B:325:MET:CE	2.54	0.55
1:A:398:MET:CE	2:B:347:ILE:HG23	2.24	0.55
1:A:181:VAL:HG11	2:B:349:ASN:O	2.04	0.55
1:A:217:LEU:HG	1:A:275:VAL:HG12	1.88	0.55
2:B:11:GLN:HA	2:B:74:THR:HG21	1.88	0.55
2:B:133:GLN:HG3	2:B:252:LEU:HD12	1.88	0.54
1:A:76:ASP:OD2	2:B:48:ARG:CZ	2.56	0.54
1:A:234:ILE:CD1	1:A:302:MET:HE2	2.37	0.54
2:B:133:GLN:NE2	2:B:251:ASP:OD2	2.40	0.54
1:A:209:ILE:HG21	1:A:227:LEU:HD13	1.90	0.54
1:A:404:PHE:CD2	2:B:261:PRO:CA	2.58	0.54
1:A:139:HIS:CB	1:A:150:THR:HG21	2.38	0.53
2:B:319:PHE:CG	2:B:323:MET:HE1	2.44	0.53
2:B:308:ARG:HD3	7:B:503:POU:H15	1.90	0.53
1:A:269:LEU:HD21	1:A:384:ILE:HG12	1.91	0.53
1:A:234:ILE:HD12	1:A:302:MET:HE2	1.90	0.52
2:B:299:LYS:NZ	7:B:503:POU:C5	2.70	0.52
6:B:502:TA1:H171	6:B:502:TA1:H442	1.91	0.52
1:A:199:ASP:OD1	1:A:256:GLN:NE2	2.43	0.51
2:B:308:ARG:HH11	7:B:503:POU:H7	1.72	0.51
1:A:178:SER:CB	2:B:349:ASN:CG	2.84	0.51
2:B:247:GLN:OE1	2:B:325:MET:SD	2.68	0.51
1:A:181:VAL:HG13	2:B:349:ASN:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:ARG:HG2	1:A:133:GLN:HE21	1.76	0.50
1:A:4:CYS:SG	1:A:134:GLY:N	2.85	0.50
1:A:72:PRO:HD2	2:B:2:ARG:HH11	1.77	0.49
1:A:178:SER:OG	2:B:349:ASN:CB	2.60	0.49
1:A:398:MET:CE	2:B:347:ILE:HA	2.43	0.49
2:B:297:ASP:OD2	7:B:503:POU:H43	2.12	0.49
1:A:72:PRO:HB2	2:B:2:ARG:HH12	1.77	0.49
1:A:72:PRO:CB	2:B:2:ARG:HH12	2.26	0.48
1:A:398:MET:HE2	2:B:348:PRO:N	2.12	0.48
2:B:165:ILE:CG1	2:B:253:ARG:CG	2.92	0.48
6:B:502:TA1:C25	6:B:502:TA1:H101	2.44	0.48
1:A:21:TRP:HZ2	1:A:65:ALA:HB2	1.79	0.48
1:A:63:PRO:O	1:A:91:GLN:NE2	2.31	0.47
2:B:21:TRP:CZ2	2:B:65:ALA:HB2	2.50	0.47
2:B:165:ILE:HD13	2:B:253:ARG:HG3	1.74	0.47
6:B:502:TA1:H131	6:B:502:TA1:H381	1.96	0.47
1:A:136:LEU:HD23	1:A:167:LEU:HB2	1.97	0.47
2:B:83:PHE:O	2:B:86:ILE:HG12	2.14	0.47
1:A:68:VAL:HG13	1:A:93:ILE:HG21	1.98	0.46
1:A:143:GLY:O	1:A:147:SER:HB2	2.16	0.46
2:B:103:TRP:CH2	2:B:107:HIS:CG	3.03	0.46
2:B:70:LEU:HD12	2:B:145:THR:HG22	1.98	0.46
2:B:183:GLU:N	2:B:184:PRO:HD2	2.31	0.46
1:A:154:MET:HE1	1:A:166:LYS:HB3	1.97	0.46
1:A:7:ILE:HG22	1:A:9:VAL:HG23	1.97	0.46
1:A:115:ILE:HD11	1:A:153:LEU:HA	1.96	0.46
1:A:7:ILE:N	1:A:136:LEU:O	2.44	0.46
1:A:67:PHE:CB	1:A:92:LEU:HD23	2.46	0.45
1:A:179:THR:O	2:B:352:LYS:HA	2.17	0.45
2:B:308:ARG:HG2	7:B:503:POU:H15	1.98	0.45
6:B:502:TA1:H463	6:B:502:TA1:C26	2.47	0.45
1:A:67:PHE:HB3	1:A:92:LEU:HD23	1.99	0.45
1:A:68:VAL:HG13	1:A:93:ILE:CG2	2.47	0.45
1:A:97:GLU:OE1	2:B:131:CYS:SG	2.75	0.44
2:B:253:ARG:O	2:B:257:VAL:HG23	2.18	0.44
2:B:229:HIS:CE1	6:B:502:TA1:O14	2.60	0.44
7:B:503:POU:H27	7:B:503:POU:O15	2.18	0.44
1:A:224:TYR:HE1	2:B:325:MET:SD	2.41	0.44
1:A:178:SER:OG	2:B:349:ASN:CG	2.61	0.44
2:B:259:MET:HE1	2:B:314:THR:O	2.18	0.43
1:A:224:TYR:CD1	2:B:325:MET:CG	2.97	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2:ARG:HD2	2:B:242:LEU:HD22	2.00	0.43
1:A:271:THR:HA	1:A:302:MET:HG3	2.00	0.43
2:B:99:ALA:O	2:B:102:ASN:HB3	2.19	0.43
2:B:103:TRP:CE3	2:B:413:MET:HE1	2.54	0.43
2:B:209:LEU:HB3	2:B:227:LEU:HD22	2.00	0.43
1:A:115:ILE:HD11	1:A:153:LEU:CA	2.48	0.42
1:A:101:ASN:OD1	2:B:254:LYS:HE2	2.19	0.42
1:A:181:VAL:CA	2:B:258:ASN:OD1	2.67	0.42
1:A:188:ILE:HD12	1:A:395:PHE:HB2	2.02	0.42
2:B:313:LEU:HD21	2:B:435:TYR:HB3	2.02	0.42
1:A:119:LEU:HD21	1:A:156:ARG:HB3	2.02	0.42
1:A:224:TYR:CE1	2:B:325:MET:SD	3.13	0.41
1:A:398:MET:HE2	2:B:347:ILE:HA	1.99	0.41
1:A:398:MET:HE2	2:B:348:PRO:HD2	0.43	0.41
1:A:96:LYS:HG2	2:B:131:CYS:HB3	2.01	0.41
2:B:299:LYS:NZ	7:B:503:POU:H31	2.36	0.41
2:B:299:LYS:HZ1	7:B:503:POU:C6	2.20	0.41
2:B:431:GLU:O	2:B:434:GLN:HB2	2.21	0.41
1:A:8:HIS:CE1	1:A:138:PHE:CD1	3.09	0.41
2:B:70:LEU:HD13	2:B:110:GLU:HG3	2.02	0.41
7:B:503:POU:H45	7:B:503:POU:H12	1.85	0.41
2:B:165:ILE:HD13	2:B:252:LEU:HB2	2.02	0.41
1:A:222:PRO:HG2	2:B:326:LYS:HB2	2.03	0.41
2:B:288:VAL:N	2:B:289:PRO:HD2	2.36	0.40
1:A:183:GLU:N	1:A:184:PRO:CD	2.84	0.40
1:A:20:CYS:SG	1:A:235:VAL:HG11	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	423/437 (97%)	378 (89%)	35 (8%)	10 (2%)	4	29
2	B	424/426 (100%)	409 (96%)	13 (3%)	2 (0%)	24	57
All	All	847/863 (98%)	787 (93%)	48 (6%)	12 (1%)	11	38

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	LEU
2	B	369	ARG
1	A	113	GLU
1	A	141	PHE
1	A	147	SER
1	A	100	ALA
1	A	108	TYR
1	A	412	GLY
1	A	131	GLY
2	B	358	ILE
1	A	181	VAL
1	A	268	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	360/368 (98%)	342 (95%)	18 (5%)	22	48
2	B	367/367 (100%)	352 (96%)	15 (4%)	27	53
All	All	727/735 (99%)	694 (96%)	33 (4%)	26	50

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	CYS
1	A	85	GLN
1	A	93	ILE
1	A	96	LYS

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Mol	Chain	Res	Type
1	A	107	HIS
1	A	147	SER
1	A	170	SER
1	A	195	LEU
1	A	219	ILE
1	A	227	LEU
1	A	238	ILE
1	A	351	PHE
1	A	361	THR
1	A	371	VAL
1	A	372	GLN
1	A	379	SER
1	A	411	GLU
1	A	429	GLU
2	B	27	GLU
2	B	31	ASP
2	B	47	GLU
2	B	160	GLU
2	B	180	THR
2	B	201	THR
2	B	205	ASP
2	B	211	ASP
2	B	216	THR
2	B	251	ASP
2	B	314	THR
2	B	339	ASN
2	B	352	LYS
2	B	419	THR
2	B	431	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	18	ASN
1	A	186	ASN
1	A	192	HIS
1	A	301	GLN
2	B	14	ASN
2	B	309	HIS
2	B	339	ASN
2	B	349	ASN

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Mol	Chain	Res	Type
2	B	350	ASN
2	B	380	ASN
2	B	385	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](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	$\# Z > 2$
6	TA1	B	502	-	68,68,68	2.04	17 (25%)	105,105,105	1.50	16 (15%)
3	GTP	A	501	4	33,34,34	1.44	5 (15%)	50,54,54	1.61	6 (12%)
5	GDP	B	501	-	29,30,30	1.29	4 (13%)	45,47,47	1.68	11 (24%)
7	POU	B	503	-	38,39,39	2.20	9 (23%)	44,57,57	2.12	10 (22%)

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
6	TA1	B	502	-	-	9/41/127/127	0/7/7/7
3	GTP	A	501	4	-	5/22/38/38	0/3/3/3
5	GDP	B	501	-	-	5/16/32/32	0/3/3/3
7	POU	B	503	-	-	7/54/76/76	0/1/2/2

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	503	POU	C22-C21	9.88	1.55	1.33
6	B	502	TA1	C06-C05	6.33	1.49	1.38
6	B	502	TA1	C08-C07	-5.88	1.25	1.38
7	B	503	POU	C2-C1	-5.30	1.38	1.52
6	B	502	TA1	C18-C10	5.05	1.68	1.57
6	B	502	TA1	C05-C04	4.44	1.46	1.39
6	B	502	TA1	C36-C31	3.69	1.45	1.39
6	B	502	TA1	C45-C24	3.28	1.59	1.54
6	B	502	TA1	O02-C03	3.26	1.41	1.34
3	A	501	GTP	C6-N1	-3.24	1.32	1.38
6	B	502	TA1	C46-C45	3.22	1.60	1.53
3	A	501	GTP	C4-N9	-3.18	1.30	1.38
5	B	501	GDP	C6-N1	-3.12	1.33	1.38
6	B	502	TA1	C43-C01	3.00	1.60	1.54
3	A	501	GTP	C5-N7	-2.99	1.33	1.39
6	B	502	TA1	C11-C10	2.92	1.61	1.54
6	B	502	TA1	C04-C03	-2.81	1.43	1.50
6	B	502	TA1	C25-C24	2.79	1.38	1.34
6	B	502	TA1	C43-C26	2.76	1.58	1.52
3	A	501	GTP	PG-O3G	-2.67	1.44	1.54
7	B	503	POU	O91-C9	2.65	1.44	1.39
5	B	501	GDP	C4-N9	-2.58	1.31	1.38
7	B	503	POU	O11-C11	2.38	1.47	1.43
6	B	502	TA1	C41-C42	2.38	1.43	1.38
7	B	503	POU	C23-C22	2.32	1.53	1.49
7	B	503	POU	C14-C13	2.30	1.57	1.52
5	B	501	GDP	C5-C4	2.27	1.44	1.38
6	B	502	TA1	C01-C45	2.21	1.65	1.56
7	B	503	POU	C9-C8	-2.19	1.51	1.53
3	A	501	GTP	C5-C4	2.17	1.44	1.38
5	B	501	GDP	O4'-C1'	2.17	1.47	1.42
6	B	502	TA1	C26-C25	2.11	1.55	1.51
7	B	503	POU	C3-C2	-2.10	1.48	1.52
7	B	503	POU	C14-C15	2.09	1.56	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	502	TA1	C07-C06	2.09	1.42	1.38

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	503	POU	C13-C12-C11	-6.94	103.99	114.42
6	B	502	TA1	C07-C08-C09	5.89	127.50	120.24
6	B	502	TA1	C06-C05-C04	-5.82	114.64	120.36
7	B	503	POU	C15-O15-C1	5.50	126.48	116.58
3	A	501	GTP	C5-C4-N3	-5.37	119.85	128.39
7	B	503	POU	O9-C9-C8	5.19	112.88	105.25
3	A	501	GTP	C2-N3-C4	4.56	120.16	112.30
5	B	501	GDP	C5-C4-N3	-4.47	121.27	128.39
3	A	501	GTP	N9-C4-N3	4.11	134.16	125.95
5	B	501	GDP	C2-N3-C4	3.84	118.91	112.30
6	B	502	TA1	C05-C04-C03	-3.84	111.94	120.40
5	B	501	GDP	N9-C4-N3	3.77	133.49	125.95
7	B	503	POU	O15-C15-C14	3.72	114.44	106.65
6	B	502	TA1	C09-C04-C03	3.37	127.83	120.40
3	A	501	GTP	C6-C5-N7	3.33	136.36	130.29
5	B	501	GDP	C6-C5-N7	3.22	136.15	130.29
6	B	502	TA1	C17-C18-C20	3.08	109.49	102.58
7	B	503	POU	C3-C4-C5	-3.07	107.68	114.70
7	B	503	POU	O15-C15-C21	3.04	117.80	109.59
5	B	501	GDP	O6-C6-C5	-2.85	119.01	126.53
6	B	502	TA1	O04-C11-C14	-2.78	102.24	108.13
6	B	502	TA1	C45-C01-C02	2.63	114.81	111.93
6	B	502	TA1	O01-C01-C43	2.54	113.53	107.04
5	B	501	GDP	C8-N9-C4	2.52	110.76	106.03
6	B	502	TA1	C08-C09-C04	-2.47	117.94	120.36
3	A	501	GTP	C4-C5-N7	-2.38	106.89	110.67
5	B	501	GDP	O2B-PB-O3A	2.29	112.32	104.64
6	B	502	TA1	C10-C18-C17	-2.29	102.20	106.55
7	B	503	POU	C25-C21-C22	-2.27	117.73	123.56
7	B	503	POU	O9-C5-C6	2.25	112.98	109.02
7	B	503	POU	C4-C3-C2	-2.25	108.18	113.26
3	A	501	GTP	O6-C6-C5	-2.25	120.59	126.53
5	B	501	GDP	N9-C8-N7	-2.22	109.28	113.40
5	B	501	GDP	C1'-N9-C4	-2.21	119.95	126.49
6	B	502	TA1	C38-C37-C29	-2.19	117.29	120.78
6	B	502	TA1	O11-C27-O12	2.19	127.91	123.95
6	B	502	TA1	O06-C15-C11	2.17	92.95	90.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	502	TA1	C14-C11-C15	-2.16	83.15	85.38
6	B	502	TA1	C43-C01-C02	-2.11	107.92	111.67
7	B	503	POU	O91-C9-O9	-2.11	104.24	109.47
5	B	501	GDP	C6-C5-C4	-2.08	115.69	118.83
5	B	501	GDP	C5-C6-N1	2.01	118.38	113.25
6	B	502	TA1	C31-C30-N01	-2.01	113.31	117.04

There are no chirality outliers.

All (26) torsion outliers are listed below:

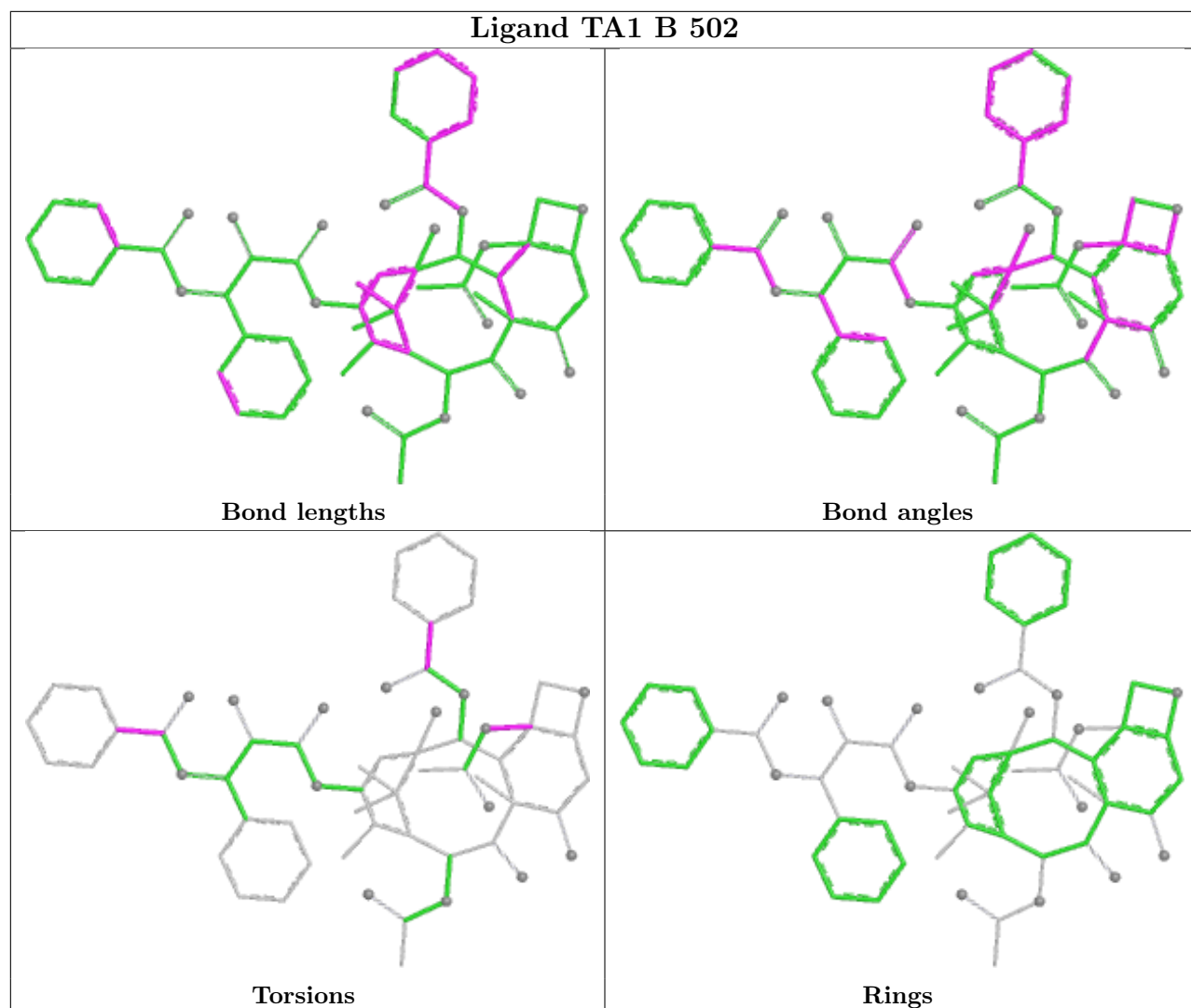
Mol	Chain	Res	Type	Atoms
3	A	501	GTP	C5'-O5'-PA-O3A
3	A	501	GTP	C5'-O5'-PA-O1A
3	A	501	GTP	C5'-O5'-PA-O2A
5	B	501	GDP	C5'-O5'-PA-O3A
5	B	501	GDP	C5'-O5'-PA-O1A
5	B	501	GDP	C5'-O5'-PA-O2A
7	B	503	POU	C25-C21-C22-C23
7	B	503	POU	C15-C21-C22-C23
6	B	502	TA1	O02-C03-C04-C05
6	B	502	TA1	O02-C03-C04-C09
6	B	502	TA1	O03-C03-C04-C09
6	B	502	TA1	N01-C30-C31-C36
6	B	502	TA1	O03-C03-C04-C05
6	B	502	TA1	O14-C30-C31-C36
6	B	502	TA1	N01-C30-C31-C32
6	B	502	TA1	O14-C30-C31-C32
7	B	503	POU	C8-C7-O7-C17
7	B	503	POU	C13-C14-C15-O15
3	A	501	GTP	PB-O3B-PG-O3G
7	B	503	POU	C14-C13-O13-C20
7	B	503	POU	C2-C3-O3-C16
5	B	501	GDP	PB-O3A-PA-O2A
6	B	502	TA1	C15-C11-O04-C12
5	B	501	GDP	C4'-C5'-O5'-PA
7	B	503	POU	C6-C7-O7-C17
3	A	501	GTP	PB-O3A-PA-O1A

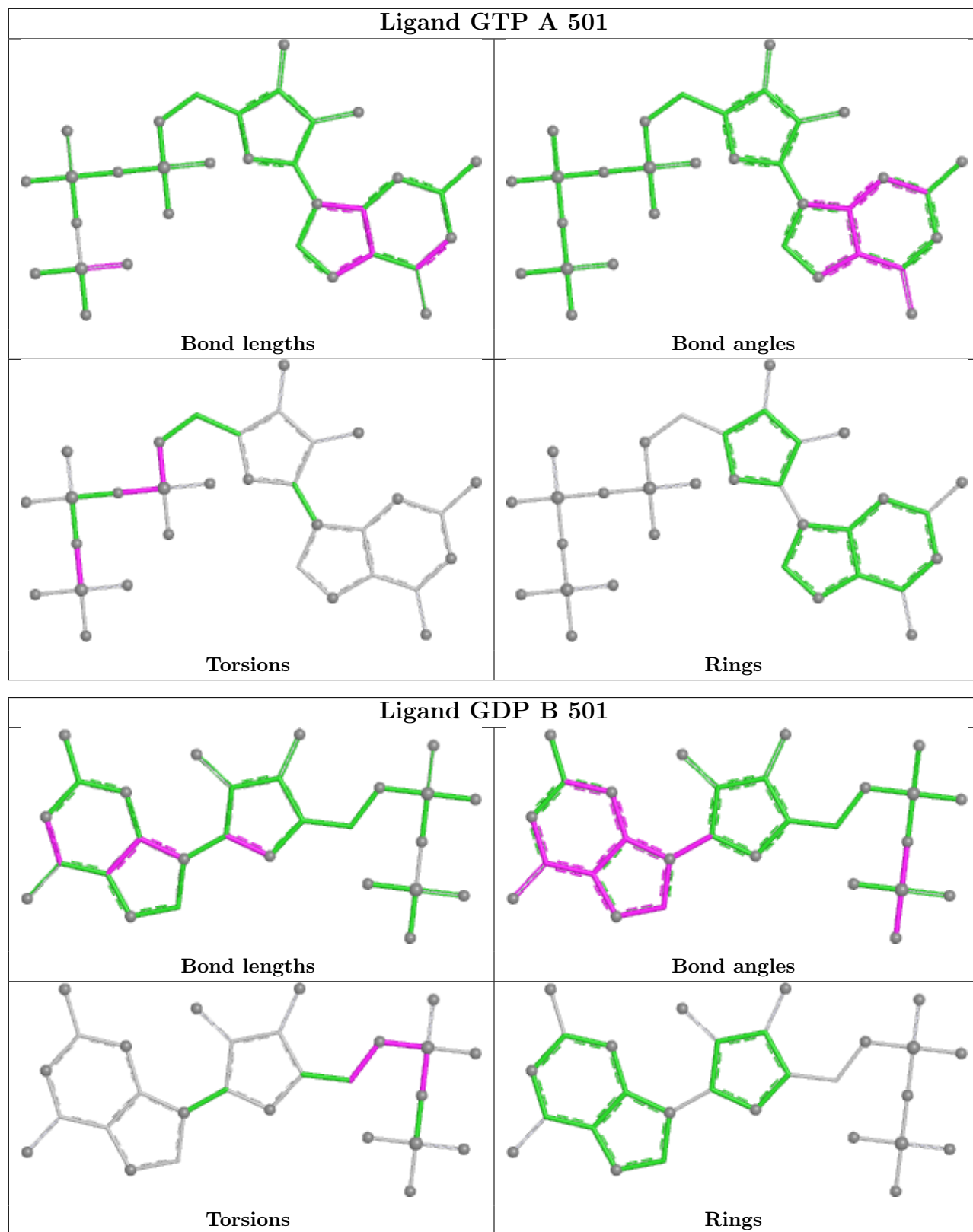
There are no ring outliers.

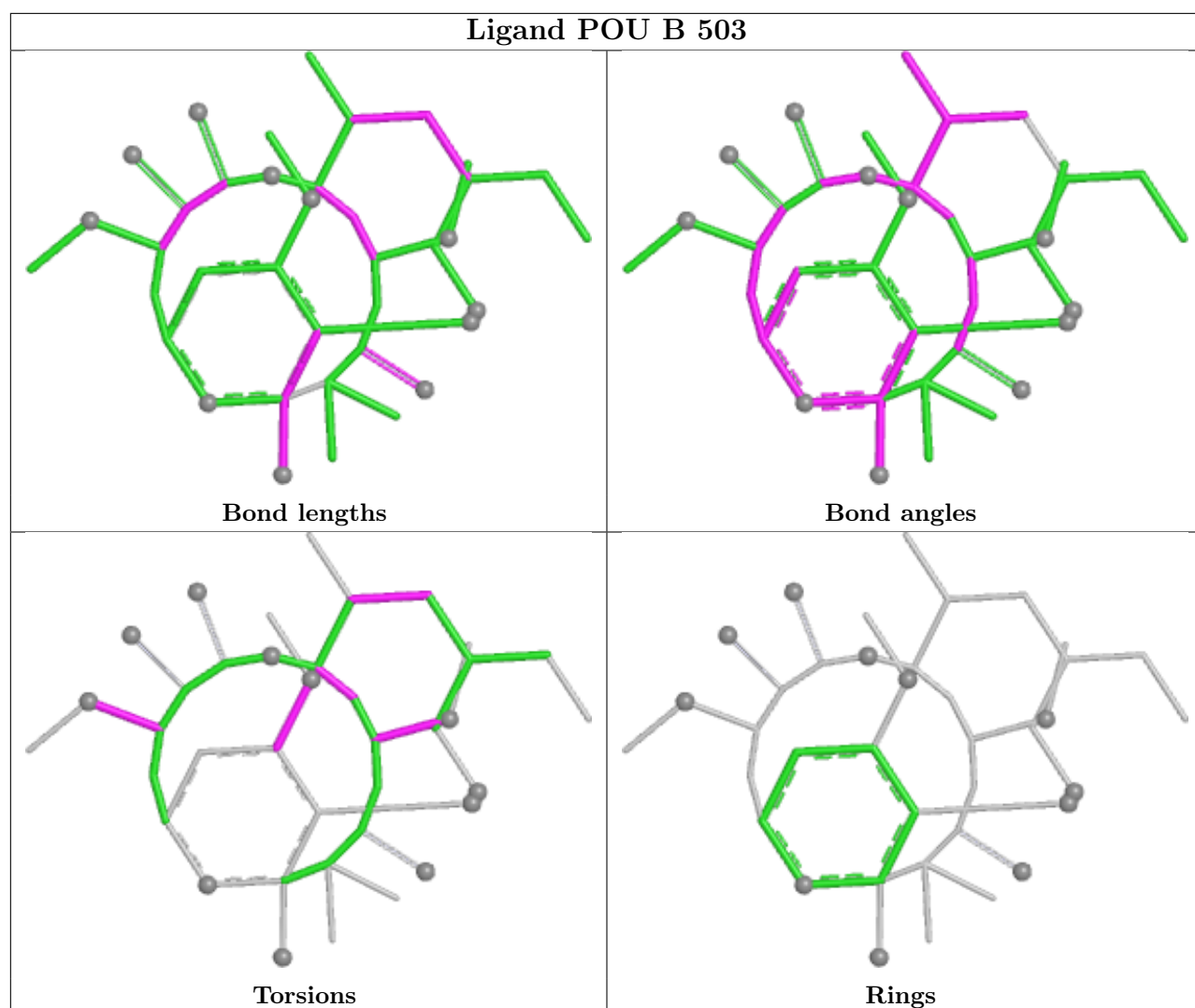
3 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	502	TA1	12	0
3	A	501	GTP	1	0
7	B	503	POU	23	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.

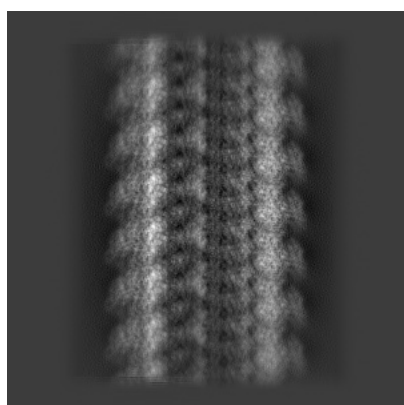
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8321. These allow visual inspection of the internal detail of the map and identification of artifacts.

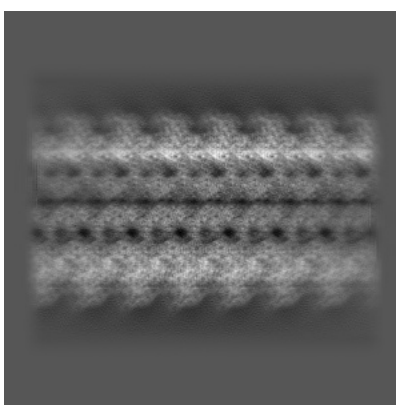
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

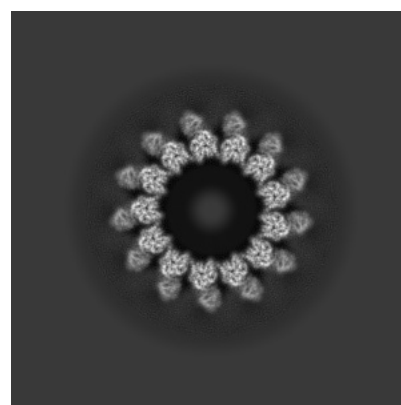
6.1.1 Primary map



X



Y

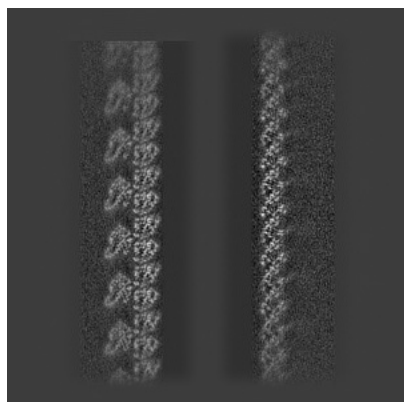


Z

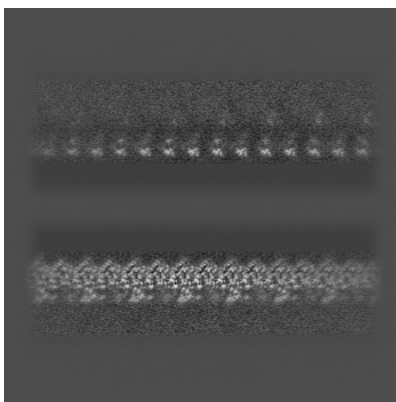
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

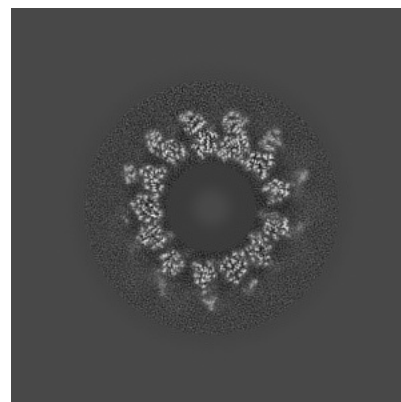
6.2.1 Primary map



X Index: 256



Y Index: 256

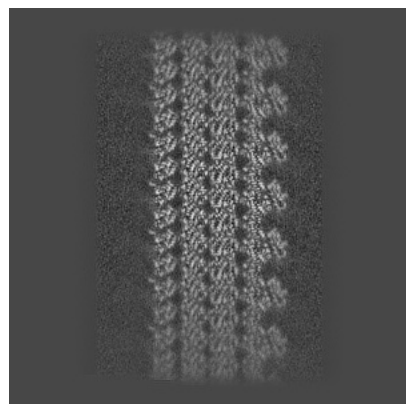


Z Index: 256

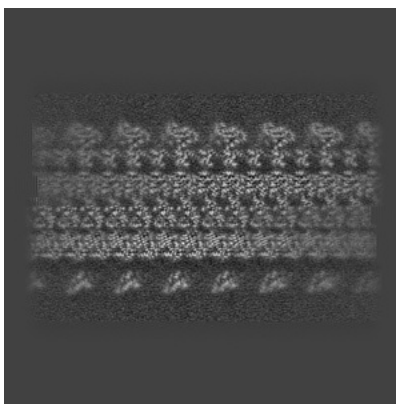
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

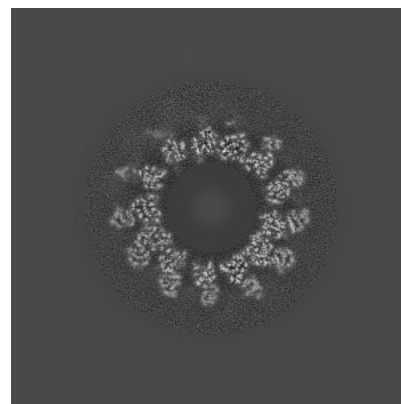
6.3.1 Primary map



X Index: 331



Y Index: 184

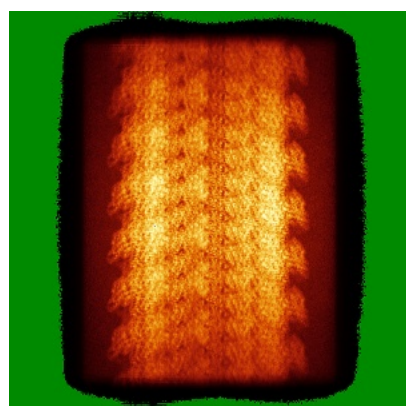


Z Index: 284

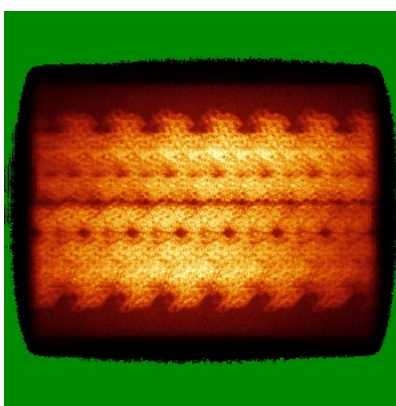
The images above show the largest variance slices of the map in three orthogonal directions.

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

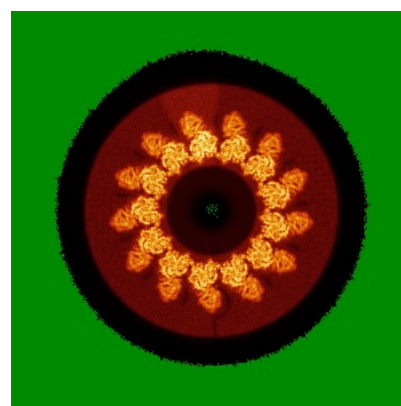
6.4.1 Primary map



X



Y

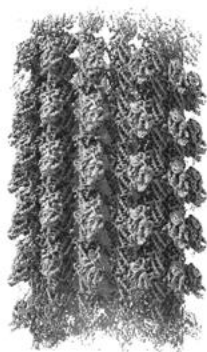


Z

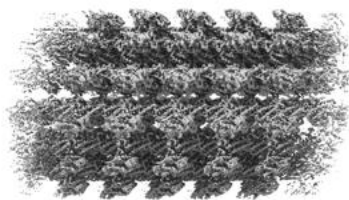
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

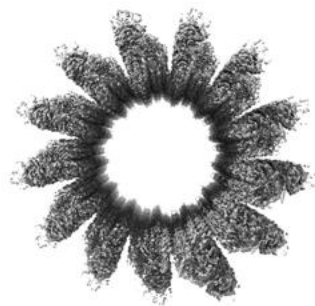
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 5.59. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

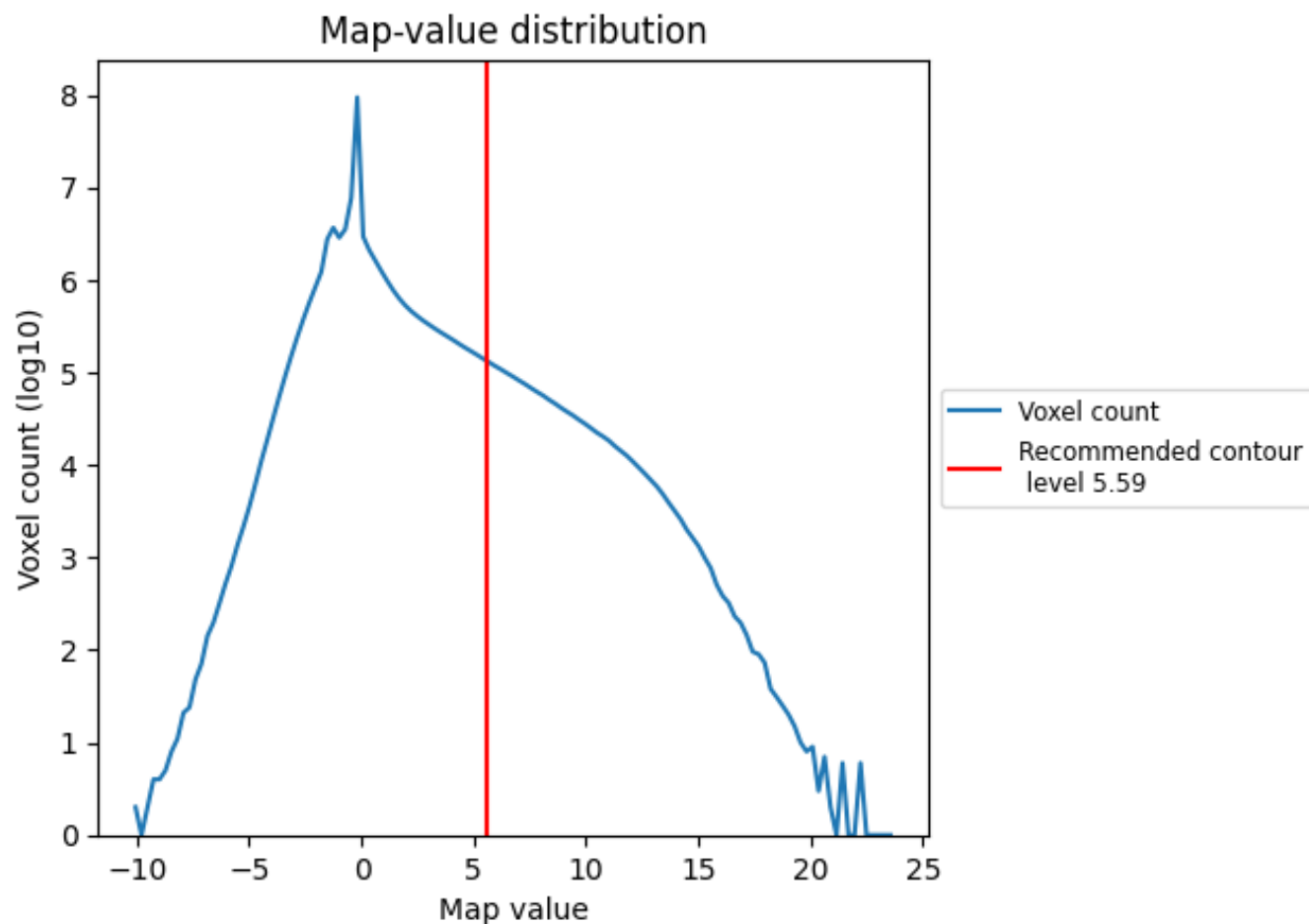
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

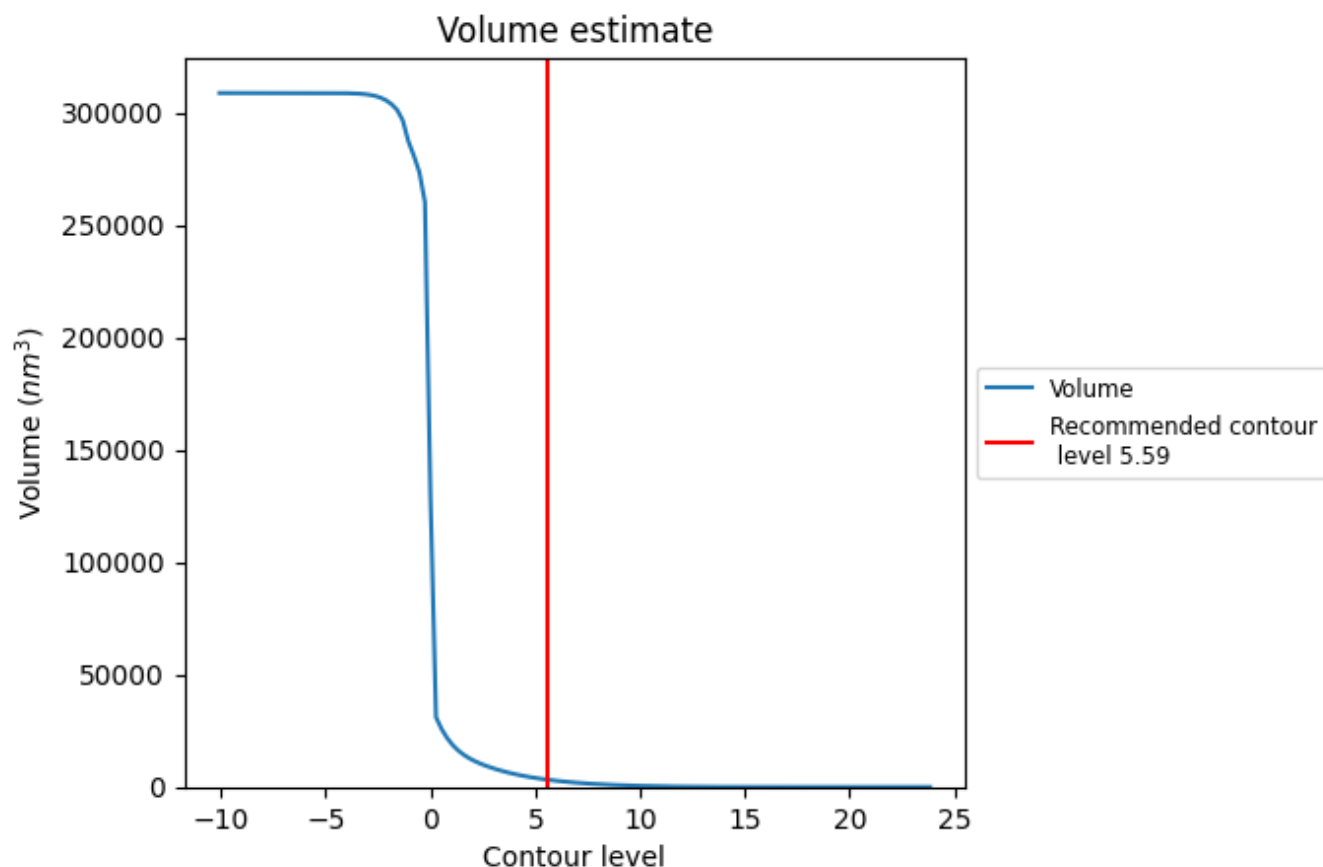
This section contains the results of statistical analysis of the map.

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.

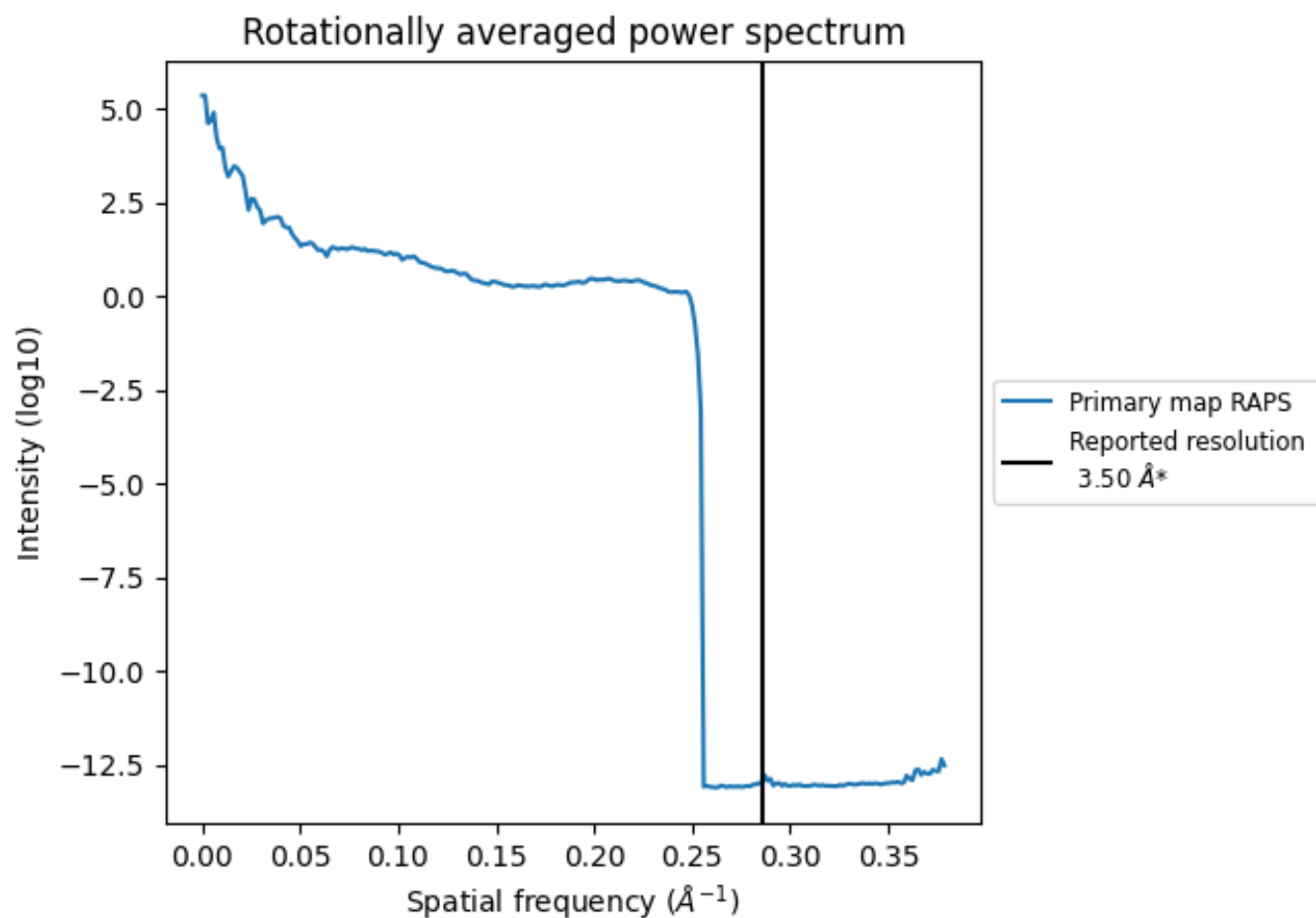
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 3216 nm³; this corresponds to an approximate mass of 2905 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 0.286 Å⁻¹

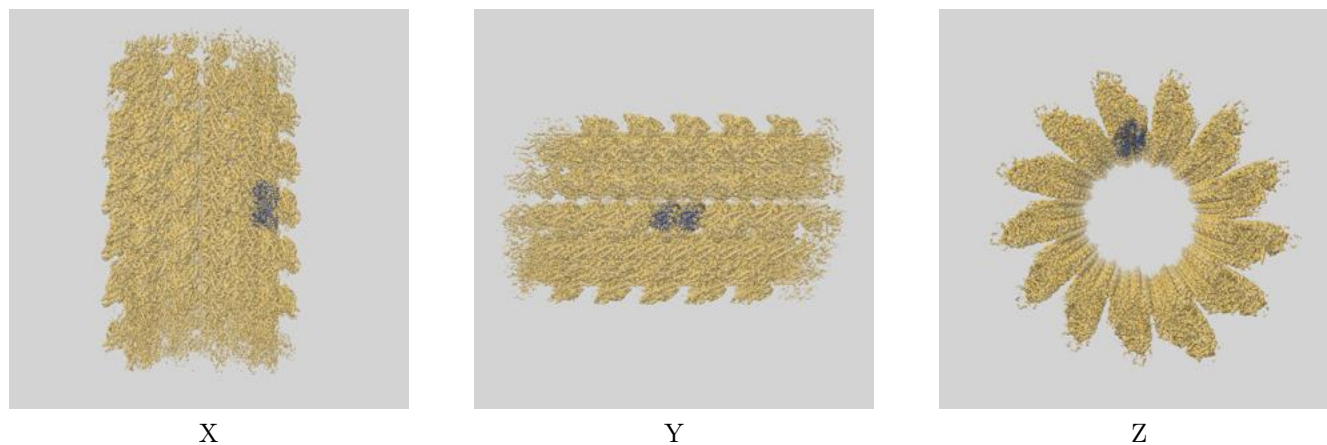
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

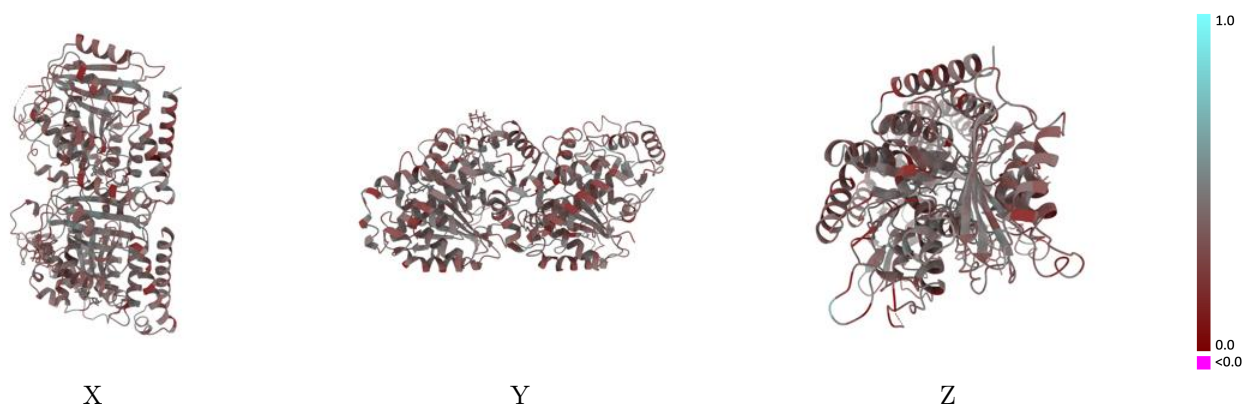
This section contains information regarding the fit between EMDB map EMD-8321 and PDB model 5SYE. Per-residue inclusion information can be found in [section 3](#) on [page 7](#).

9.1 Map-model overlay [i](#)



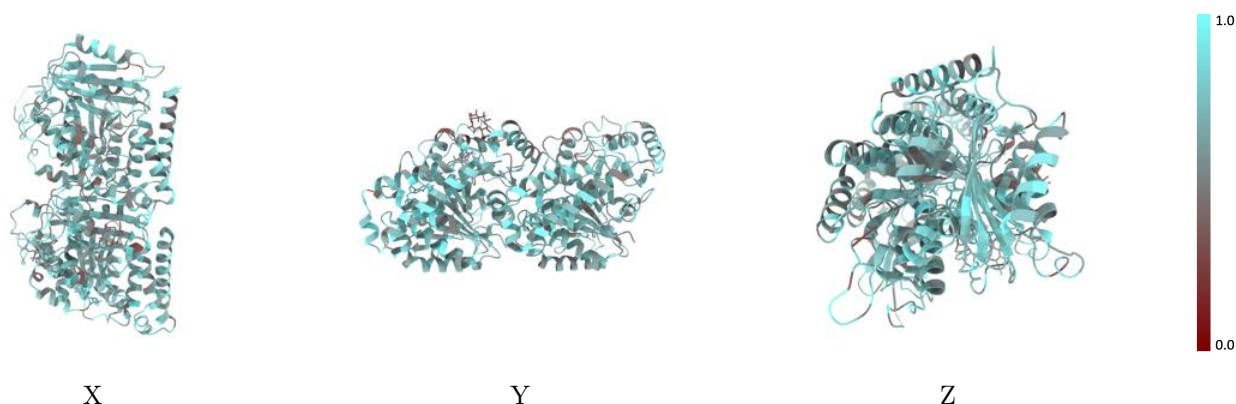
The images above show the 3D surface view of the map at the recommended contour level 5.59 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



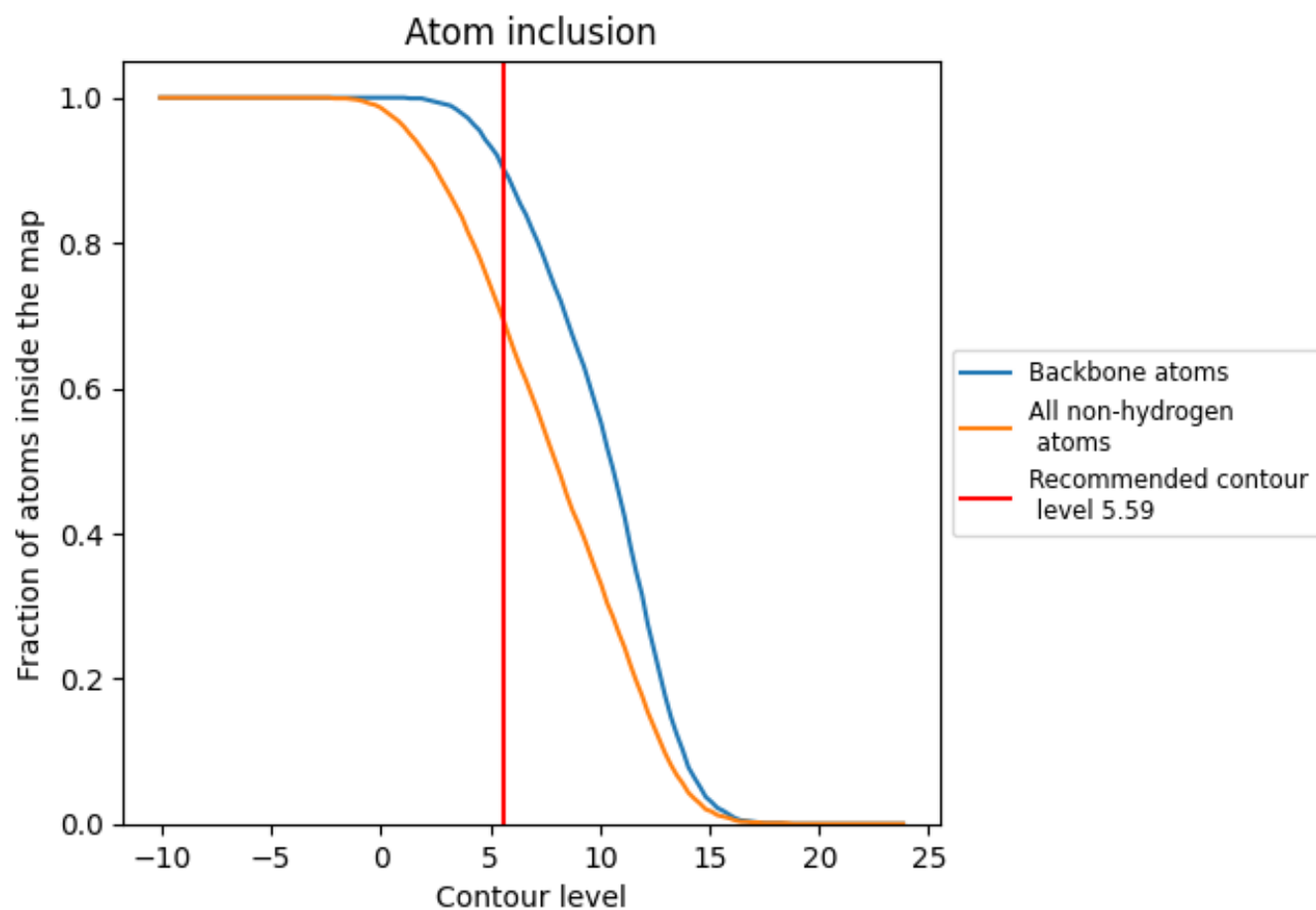
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (5.59).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 69% 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 (5.59) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6940	0.3720
A	0.6900	0.3640
B	0.6980	0.3790

