



Full wwPDB EM Validation Report ⓘ

Mar 17, 2026 – 02:53 PM UTC

PDB ID : 6ZYY / pdb_00006zyy
EMDB ID : EMD-11581
Title : Outer Dynein Arm-Shulin complex - Dyh3 motor region (Tetrahymena thermophila)
Authors : Mali, G.R.; Abid Ali, F.; Lau, C.K.; Begum, F.; Boulanger, J.; Howe, J.D.; Chen, Z.A.; Rappsilber, J.; Skehel, M.; Carter, A.P.
Deposited on : 2020-08-03
Resolution : 4.40 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:

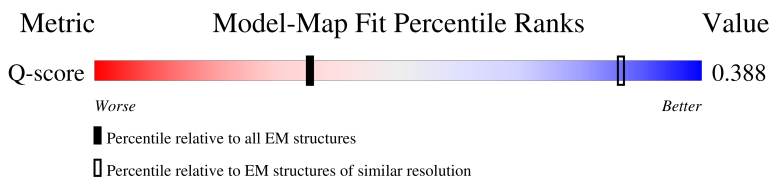
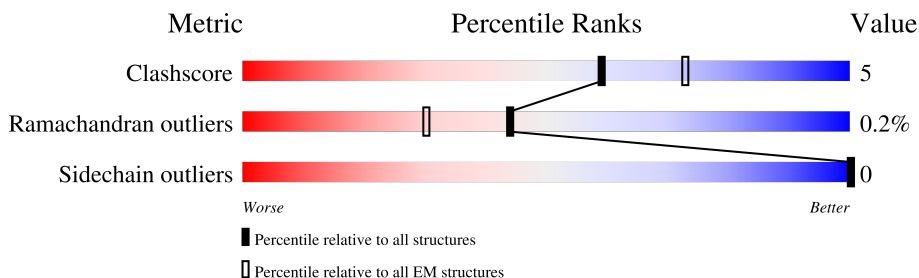
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

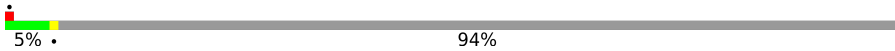
The reported resolution of this entry is 4.40 Å.

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	3132 (3.91 - 4.90)

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	C	4620	
2	Y	1200	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 22453 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 Dynein heavy chain, outer arm protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	2901	Total	C	N	O	S	0	0
			21875	13931	3736	4092	116		

- Molecule 2 is a protein called Shulin.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	Y	66	Total	C	N	O	0	0
			466	303	89	74		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

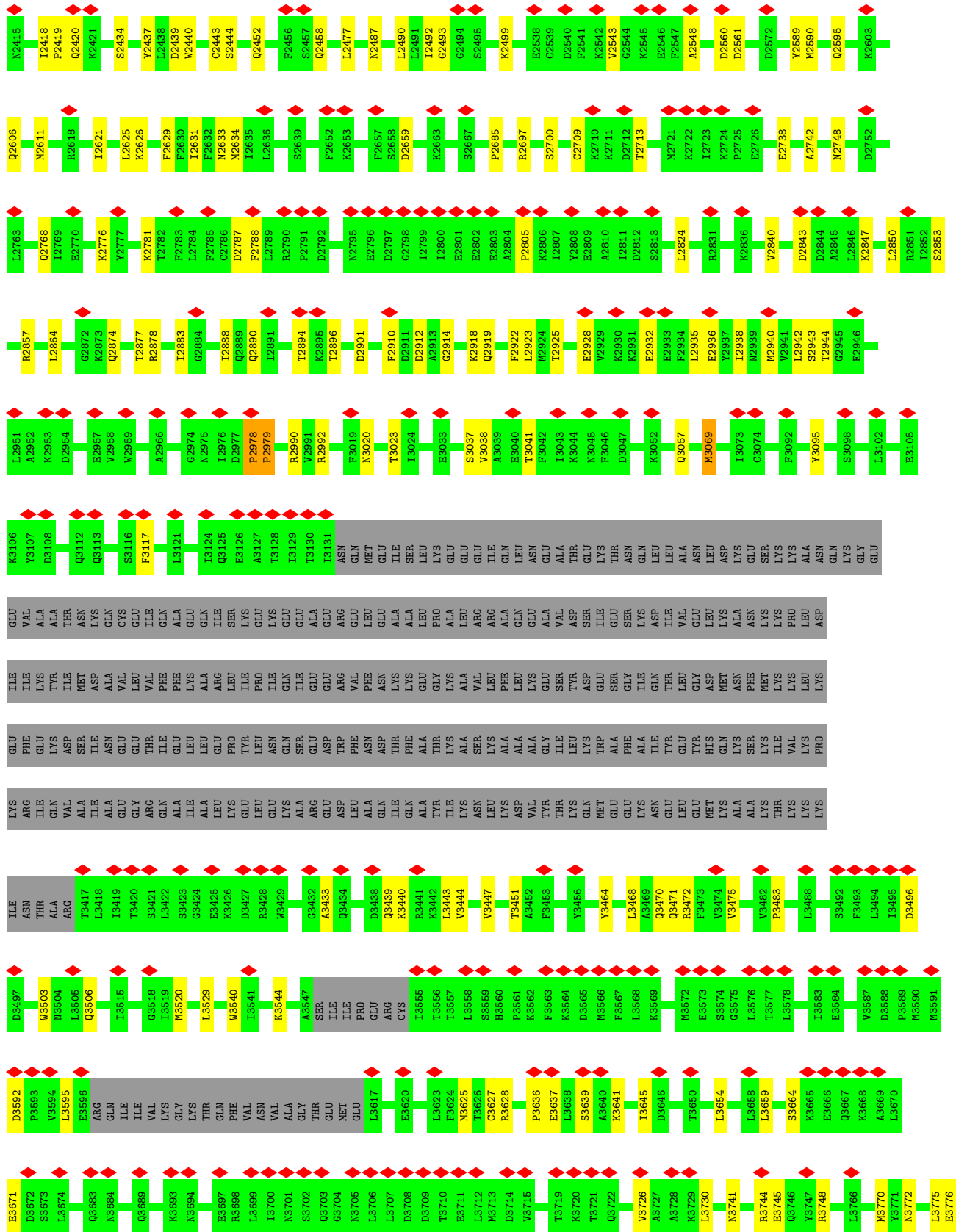


Mol	Chain	Residues	Atoms					AltConf
3	C	1	Total	C	N	O	P	0
			27	10	5	10	2	
3	C	1	Total	C	N	O	P	0
			27	10	5	10	2	
3	C	1	Total	C	N	O	P	0
			27	10	5	10	2	

-
- The image displays the chemical structure of Adenosine Triphosphate (ATP). It consists of an adenine base (a purine ring system with an amino group at N6) attached to a ribose sugar (a five-membered ring with hydroxyl groups at C2' and C3'). The ribose is linked to a chain of three phosphate groups (P1, P2, and P3) via phosphodiester bonds. The structure is labeled with atom names (N1-N9, O1-O9, C1-C5, C1P1, C2P1, C3P1, C4P1, C5P1, C1P2, C2P2, C3P2, C4P2, C5P2, C1P3, C2P3, C3P3, C4P3, C5P3) and shows the stereochemistry of the bonds.

Mol	Chain	Residues	Atoms					AltConf
4	C	1	Total	C	N	O	P	0
			31	10	5	13	3	







GLN	LEU	ASP	LYS	LEU	LEU	LEU	LYS	PHE	LEU	ILE	LYS	LYS	LYS	VAL	GLN	VAL	VAL
GLN	LEU	MET	GLY	GLN	GLN	GLN	LYS	ASN	ALA	THR	THR	GLN	GLN	THR	GLN	VAL	GLN
GLN	GLN	PHE	LEU	GLY	LEU	LEU	ASP	ASN	ASP	VAL	PRO	ASN	ASN	GLY	ASN	ASP	GLY
GLU	VAL	PHE	ARG	ASN	GLN	GLN	TYR	VAL	LEU	ILE	ALA	ASN	ASN	GLU	ASN	ASN	ARG
TRP	TRP	ASP	TYR	GLN	GLN	THR	LYS	THR	LEU	THR	THR	THR	ASN	ASN	GLN	ILE	LYS
GLU	GLU	ALA	ASP	THR	THR	THR	ASP	THR	THR	THR	ASN	ASN	ASN	GLU	GLN	THR	SER
THR	THR	GLY	SER	LEU	LEU	VAL	LYS	ILE	VAL	GLY	ILE	ILE	ILE	SER	ASN	THR	TYR
LYS	LYS	ASN	LYS	VAL	VAL	ASP	LYS	PHE	THR	THR	THR	THR	GLY	GLN	GLN	GLY	GLY
GLN	GLN	LEU	LYS	THR	THR	THR	LYS	THR	PHE	THR	THR	THR	GLN	GLN	GLN	ILE	PHE
VAL	VAL	ASN	GLU	ILE	GLU	ILE	GLU	ILE	ASN	ASN	ASN	ASN	ASN	ASN	ASN	ASN	ASN
ASP	ASP	GLU	GLY	GLU	GLU	GLU	GLU	GLY	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
LYS	LYS	LEU	ILE	GLU	GLU	GLU	GLU	GLY	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
HIS	HIS	LEU	ILE	LYS	LYS	LYS	LYS	THR	LEU	THR	THR	THR	THR	THR	THR	THR	THR
LEU	LEU	GLU	THR	ILE	ILE	ILE	ILE	SER	VAL	THR	THR	THR	THR	THR	THR	THR	THR
ASP	ASP	LEU	THR	ALA	ALA	ALA	ALA	GLN	ARG	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
LYS	LYS	LEU	THR	GLU	GLU	GLU	GLU	THR	LEU	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	ASN	ASN	ASN	ASN	ASN	ASN	ASN	ASN	ASN	ASN
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	SER	ALA	ALA	ALA	ALA	ALA	ALA	ALA	ALA	ALA
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
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TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
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GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
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GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU	GLU	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
GLN	GLN	LEU	THR	GLN	GLN	GLN	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
ALA	ALA	GLU	THR	ILE	ILE	ILE	ILE	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
TYR	TYR	LEU	THR	GLU	GLU	GLU											

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	49397	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 ($e^-/\text{\AA}^2$)	52	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.082	Depositor
Minimum map value	-0.069	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.015	Depositor
Map size (Å)	499.5, 499.5, 499.5	wwPDB
Map dimensions	450, 450, 450	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.11, 1.11, 1.11	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: ATP, ADP

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	C	0.32	2/22277 (0.0%)	0.59	15/30197 (0.0%)
2	Y	0.32	0/477	0.56	0/648
All	All	0.32	2/22754 (0.0%)	0.58	15/30845 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	2857	ARG	CD-NE	-5.51	1.38	1.46
1	C	3069	MET	CB-CG	-5.35	1.36	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	2857	ARG	NE-CZ-NH2	-12.99	107.51	119.20
1	C	2978	PRO	N-CA-CB	8.05	110.89	103.08
1	C	2805	PRO	N-CA-CB	7.71	110.06	103.35
1	C	4182	PRO	N-CA-CB	7.65	111.28	103.25
1	C	4188	PRO	N-CA-CB	7.64	111.28	103.25
1	C	3069	MET	CB-CG-SD	-7.42	90.44	112.70
1	C	2857	ARG	NH1-CZ-NH2	7.29	128.78	119.30
1	C	4180	PRO	N-CA-CB	7.24	110.85	103.25
1	C	3483	PRO	N-CA-CB	7.20	110.03	103.19
1	C	4546	PRO	N-CA-CB	7.14	110.75	103.25
1	C	2979	PRO	N-CA-CB	7.13	110.73	103.25
1	C	4433	PRO	CA-N-CD	-6.38	103.07	112.00
1	C	1323	LEU	N-CA-C	-6.26	102.78	110.61
1	C	2857	ARG	CG-CD-NE	-5.88	99.07	112.00
1	C	3069	MET	CG-SD-CE	-5.26	89.32	100.90

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	C	21875	0	20694	197	0
2	Y	466	0	390	5	0
3	C	81	0	36	0	0
4	C	31	0	12	0	0
All	All	22453	0	21132	202	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 5.

All (202) 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:C:2499:LYS:NZ	1:C:2611:MET:O	2.11	0.83
1:C:1314:SER:O	1:C:1318:VAL:N	2.14	0.79
1:C:4125:ASN:O	1:C:4128:SER:OG	2.00	0.77
1:C:2776:LYS:O	1:C:2781:LYS:NZ	2.18	0.76
1:C:3636:PRO:O	1:C:3639:SER:OG	2.03	0.74
1:C:2943:SER:O	1:C:2992:ARG:NH2	2.21	0.73
1:C:1986:LEU:O	1:C:1991:LYS:NZ	2.24	0.71
2:Y:1148:ASN:OD1	2:Y:1149:ILE:N	2.23	0.71
1:C:2843:ASP:OD2	1:C:2847:LYS:NZ	2.25	0.70
1:C:1862:MET:SD	1:C:1971:SER:OG	2.48	0.69
1:C:2322:LYS:NZ	1:C:2378:LEU:O	2.26	0.69
1:C:3664:SER:OG	1:C:3671:GLU:OE2	2.07	0.69
1:C:3789:SER:O	1:C:3798:ARG:NH2	2.26	0.68
1:C:2164:SER:OG	1:C:2273:GLU:OE2	2.12	0.67
1:C:3741:ASN:O	1:C:3745:GLU:N	2.27	0.67
1:C:4215:LEU:O	1:C:4218:THR:OG1	2.11	0.67
1:C:2434:SER:OG	1:C:2437:TYR:O	2.12	0.67
1:C:2595:GLN:N	1:C:2595:GLN:OE1	2.28	0.67
1:C:2027:PHE:CZ	1:C:2074:LEU:HD21	2.30	0.67
1:C:2340:GLN:NE2	1:C:2343:GLU:O	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2894:THR:OG1	1:C:2896:THR:O	2.12	0.66
1:C:2850:LEU:O	1:C:2853:SER:OG	2.12	0.66
1:C:2888:ILE:HG23	1:C:2923:LEU:HD23	1.78	0.65
1:C:2936:GLU:N	1:C:2936:GLU:OE1	2.30	0.65
1:C:4113:ILE:HG22	1:C:4117:PHE:CE2	2.32	0.65
1:C:2492:ILE:O	1:C:2634:MET:N	2.29	0.65
1:C:3037:SER:O	1:C:3041:THR:HG23	1.98	0.64
1:C:4461:LEU:CD2	1:C:4553:LEU:HD11	2.26	0.64
1:C:2890:GLN:NE2	1:C:2928:GLU:OE2	2.31	0.63
1:C:3776:GLU:OE1	1:C:3776:GLU:N	2.32	0.63
1:C:2768:GLN:N	1:C:2768:GLN:OE1	2.32	0.63
1:C:4022:GLY:O	1:C:4023:GLN:NE2	2.32	0.63
1:C:3870:GLU:O	1:C:3874:TRP:N	2.31	0.62
1:C:2910:PHE:O	1:C:2914:GLY:N	2.31	0.62
2:Y:1147:ARG:HD3	2:Y:1151:GLY:HA2	1.81	0.62
2:Y:1148:ASN:HB3	2:Y:1154:LEU:HD21	1.82	0.62
1:C:2543:VAL:HG12	1:C:2548:ALA:CB	2.29	0.62
1:C:2940:MET:SD	1:C:2944:THR:OG1	2.58	0.61
1:C:1589:GLU:OE1	1:C:1589:GLU:N	2.32	0.61
1:C:2543:VAL:HG12	1:C:2548:ALA:HB3	1.81	0.61
1:C:1932:GLU:O	1:C:1935:SER:OG	2.10	0.61
1:C:2211:VAL:O	1:C:2214:THR:OG1	2.13	0.61
1:C:2353:PRO:O	1:C:2356:SER:OG	2.16	0.61
1:C:1902:HIS:O	1:C:1903:ARG:NE	2.34	0.60
1:C:4209:ASP:OD1	1:C:4210:ASP:N	2.35	0.60
1:C:2621:ILE:O	1:C:2626:LYS:NZ	2.34	0.60
1:C:4537:SER:OG	1:C:4538:LYS:NZ	2.34	0.60
1:C:1801:LYS:N	1:C:1805:ASP:OD2	2.35	0.60
1:C:2443:CYS:SG	1:C:2444:SER:N	2.75	0.59
1:C:2560:ASP:OD1	1:C:2561:ASP:N	2.35	0.59
1:C:2008:ILE:HD11	1:C:2051:ILE:CG2	2.32	0.59
1:C:1487:VAL:O	1:C:1491:LYS:NZ	2.36	0.59
1:C:4143:ARG:NH1	1:C:4219:TYR:OH	2.35	0.59
1:C:3057:GLN:N	1:C:3057:GLN:OE1	2.35	0.59
1:C:1855:TYR:OH	1:C:1884:ASP:OD2	2.19	0.59
1:C:2685:PRO:O	1:C:4396:GLN:NE2	2.36	0.59
1:C:2748:ASN:C	1:C:2748:ASN:HD22	2.11	0.58
1:C:2912:ASP:O	1:C:2918:LYS:N	2.37	0.58
1:C:3069:MET:CE	1:C:3472:ARG:HB3	2.34	0.57
1:C:2175:THR:HG22	1:C:2180:PRO:HA	1.85	0.57
1:C:3095:TYR:HE1	1:C:3451:THR:HG23	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1827:ILE:HG22	1:C:1828:THR:H	1.70	0.56
1:C:4109:PHE:O	1:C:4113:ILE:HG21	2.05	0.56
1:C:3949:ASP:OD1	1:C:3950:ARG:N	2.37	0.56
1:C:2490:LEU:HD21	1:C:2611:MET:HG2	1.86	0.56
1:C:3117:PHE:CB	1:C:3433:ALA:HB2	2.36	0.56
1:C:3627:CYS:SG	1:C:3628:ARG:N	2.78	0.56
1:C:2477:LEU:HD12	1:C:2606:GLN:OE1	2.06	0.56
1:C:1700:SER:O	1:C:1704:GLN:NE2	2.39	0.55
1:C:2420:GLN:OE1	1:C:2420:GLN:N	2.38	0.55
1:C:4067:HIS:ND1	1:C:4068:GLU:O	2.40	0.55
1:C:4320:VAL:HG21	1:C:4411:TYR:CE2	2.42	0.55
1:C:4024:GLU:N	1:C:4024:GLU:OE1	2.40	0.55
1:C:1376:MET:O	1:C:1379:HIS:N	2.40	0.54
1:C:2452:GLN:O	1:C:2458:GLN:NE2	2.41	0.54
1:C:4522:GLU:OE1	1:C:4522:GLU:N	2.41	0.54
1:C:2697:ARG:O	1:C:2700:SER:OG	2.21	0.54
1:C:1925:GLU:OE1	1:C:1983:ARG:NH2	2.40	0.53
1:C:2439:ASP:OD1	1:C:2440:TRP:N	2.41	0.53
1:C:2625:LEU:HD21	1:C:2629:PHE:CE2	2.43	0.53
1:C:2008:ILE:HD11	1:C:2051:ILE:HG21	1.91	0.53
1:C:2324:GLY:O	1:C:2327:LEU:N	2.41	0.53
1:C:2490:LEU:HD22	1:C:2631:ILE:HG22	1.90	0.53
1:C:2925:THR:OG1	1:C:2928:GLU:OE1	2.12	0.53
1:C:1501:PHE:O	1:C:1504:VAL:HG12	2.09	0.52
1:C:1929:ILE:HG21	1:C:1934:LEU:HD21	1.91	0.52
1:C:3777:GLN:O	1:C:3781:LEU:HD23	2.09	0.52
1:C:1995:ARG:NH2	1:C:4161:ASP:OD1	2.42	0.52
1:C:2493:GLY:O	1:C:2499:LYS:NZ	2.42	0.52
1:C:2840:VAL:HG11	1:C:3038:VAL:HG22	1.91	0.52
1:C:1916:GLN:HG2	1:C:1964:ILE:HD11	1.92	0.52
1:C:2023:LEU:HD22	1:C:2074:LEU:HD23	1.92	0.51
1:C:3790:GLU:O	1:C:3798:ARG:NE	2.38	0.51
1:C:3801:ASN:O	1:C:3801:ASN:ND2	2.42	0.51
1:C:4163:GLU:O	1:C:4167:LEU:HD23	2.11	0.51
1:C:3520:MET:HG2	1:C:3645:ILE:HD11	1.93	0.51
1:C:4531:LEU:HD21	1:C:4534:CYS:SG	2.51	0.51
1:C:1970:VAL:O	1:C:1971:SER:OG	2.29	0.51
1:C:2787:ASP:OD1	1:C:2788:PHE:N	2.41	0.51
1:C:2397:GLY:O	1:C:2633:ASN:N	2.43	0.50
1:C:2919:GLN:N	1:C:2919:GLN:OE1	2.44	0.50
1:C:3069:MET:HE2	1:C:3472:ARG:HB3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Y:1156:HIS:ND1	2:Y:1157:HIS:O	2.44	0.50
1:C:2185:ILE:O	1:C:2589:TYR:OH	2.27	0.50
1:C:3529:LEU:HD13	1:C:3625:MET:HG3	1.93	0.50
1:C:2748:ASN:O	1:C:2748:ASN:ND2	2.39	0.49
1:C:2590:MET:SD	1:C:2590:MET:N	2.85	0.49
1:C:3095:TYR:CE1	1:C:3451:THR:HG23	2.47	0.49
1:C:4465:THR:HG1	1:C:4617:CYS:HG	1.60	0.49
1:C:2932:GLU:O	1:C:2935:LEU:N	2.42	0.49
1:C:2178:GLY:O	1:C:2179:SER:OG	2.30	0.48
1:C:2928:GLU:OE1	1:C:2928:GLU:N	2.47	0.48
1:C:4276:GLU:N	1:C:4276:GLU:OE1	2.46	0.48
1:C:2864:LEU:N	1:C:3023:THR:O	2.45	0.48
1:C:2180:PRO:O	1:C:2181:HIS:ND1	2.47	0.48
1:C:4304:ARG:NH2	1:C:4305:PRO:O	2.47	0.48
1:C:2357:VAL:O	1:C:2361:LEU:HD23	2.14	0.48
1:C:3592:ASP:O	1:C:3595:LEU:N	2.47	0.48
1:C:4385:THR:HG21	1:C:4419:TRP:HE1	1.78	0.47
1:C:2490:LEU:HD21	1:C:2611:MET:CG	2.43	0.47
1:C:4158:ASN:OD1	1:C:4160:SER:OG	2.32	0.47
1:C:4122:ASP:OD1	1:C:4124:SER:OG	2.32	0.47
1:C:1944:ILE:HG21	1:C:1972:TYR:OH	2.15	0.47
1:C:2709:CYS:CB	1:C:2713:THR:HG23	2.45	0.47
1:C:3744:ARG:O	1:C:3748:ARG:N	2.43	0.46
1:C:1857:THR:HG21	1:C:1997:VAL:HG22	1.97	0.46
1:C:4117:PHE:O	1:C:4120:LYS:NZ	2.40	0.46
1:C:4526:ILE:O	1:C:4553:LEU:HD13	2.16	0.46
1:C:2115:ASP:OD1	1:C:2116:VAL:N	2.48	0.46
1:C:3770:MET:N	1:C:3772:ASN:OD1	2.47	0.46
1:C:1819:ASP:OD1	1:C:1820:ILE:N	2.47	0.46
1:C:3520:MET:HG2	1:C:3645:ILE:CD1	2.46	0.46
1:C:2381:GLN:N	1:C:2381:GLN:OE1	2.49	0.46
1:C:2244:LEU:O	1:C:2247:VAL:HG12	2.16	0.45
1:C:2709:CYS:HB3	1:C:2713:THR:HG23	1.98	0.45
1:C:2738:GLU:O	1:C:2742:ALA:N	2.49	0.45
1:C:4190:SER:O	1:C:4190:SER:OG	2.34	0.45
1:C:4214:GLU:O	1:C:4218:THR:HG23	2.17	0.45
1:C:1765:SER:HA	1:C:1769:LEU:HD13	1.99	0.45
1:C:2324:GLY:HA2	1:C:2327:LEU:HB2	1.99	0.45
1:C:3464:TYR:CD1	1:C:3468:LEU:HD22	2.52	0.45
1:C:1735:VAL:HG11	1:C:1808:TRP:CZ3	2.52	0.45
1:C:1320:LYS:HA	1:C:1324:ALA:HB1	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2003:ASP:OD1	1:C:2004:ARG:N	2.50	0.44
1:C:2477:LEU:HD13	1:C:2487:ASN:OD1	2.17	0.44
1:C:2325:ASN:OD1	1:C:2328:ARG:NH2	2.48	0.44
1:C:3955:THR:HA	1:C:3958:PHE:CE1	2.52	0.44
1:C:4382:VAL:O	1:C:4386:LEU:HD23	2.18	0.44
1:C:2120:ILE:HG22	1:C:2124:ILE:HD12	2.00	0.44
1:C:2697:ARG:NH1	1:C:3020:ASN:OD1	2.51	0.44
1:C:2938:ILE:O	1:C:2942:LEU:HD12	2.18	0.44
1:C:3439:GLN:O	1:C:3443:LEU:N	2.42	0.44
2:Y:1147:ARG:HD3	2:Y:1151:GLY:CA	2.46	0.43
1:C:2418:ILE:HB	1:C:2419:PRO:HD2	2.00	0.43
1:C:2901:ASP:N	1:C:2901:ASP:OD1	2.52	0.43
1:C:4218:THR:HG21	1:C:4609:LYS:HG2	2.00	0.43
1:C:3811:TYR:CE1	1:C:3815:ASN:ND2	2.82	0.43
1:C:2206:ASP:N	1:C:2206:ASP:OD1	2.51	0.43
1:C:1983:ARG:NH2	1:C:2242:GLU:OE2	2.46	0.43
1:C:1400:TYR:O	1:C:1403:GLU:N	2.47	0.43
1:C:1877:GLY:O	1:C:1881:THR:OG1	2.28	0.43
1:C:2788:PHE:HE1	1:C:2853:SER:HB2	1.84	0.43
1:C:3503:TRP:O	1:C:3506:GLN:N	2.52	0.43
1:C:2177:LEU:HD12	1:C:2177:LEU:O	2.19	0.43
1:C:3790:GLU:OE1	1:C:3802:ILE:HG12	2.19	0.43
1:C:1521:LEU:HD12	1:C:1581:LEU:HD22	2.01	0.43
1:C:1917:SER:OG	1:C:1918:GLY:N	2.51	0.43
1:C:2874:GLN:O	1:C:2878:ARG:NH1	2.51	0.43
1:C:1772:LEU:HD23	1:C:1778:ARG:HA	2.01	0.42
1:C:1896:THR:OG1	1:C:1897:ASN:N	2.50	0.42
1:C:2748:ASN:C	1:C:2748:ASN:ND2	2.76	0.42
1:C:2877:THR:HG21	1:C:2923:LEU:CD1	2.49	0.42
1:C:3726:VAL:O	1:C:3730:LEU:N	2.44	0.42
1:C:3911:ILE:O	1:C:3947:ARG:NH2	2.53	0.42
1:C:4015:GLU:OE1	1:C:4015:GLU:N	2.51	0.42
1:C:3540:TRP:O	1:C:3544:LYS:N	2.50	0.42
1:C:4105:LEU:O	1:C:4108:THR:OG1	2.34	0.42
1:C:2210:GLY:O	1:C:2214:THR:HG23	2.20	0.42
1:C:2244:LEU:HD22	1:C:2247:VAL:HG11	2.01	0.42
1:C:3659:LEU:HD21	1:C:3744:ARG:HD2	2.02	0.42
1:C:2824:LEU:HD21	1:C:2883:ILE:CG2	2.50	0.42
1:C:2051:ILE:O	1:C:2055:LEU:HD23	2.20	0.42
1:C:3784:GLU:O	1:C:3787:ASP:N	2.53	0.42
1:C:2124:ILE:HG21	1:C:2132:ASN:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:4461:LEU:HD23	1:C:4553:LEU:HD11	2.01	0.41
1:C:1518:TRP:NE1	1:C:1522:GLU:OE2	2.53	0.41
1:C:3471:GLN:O	1:C:3475:VAL:HG13	2.20	0.41
1:C:2064:GLN:OE1	1:C:2066:ILE:HG23	2.20	0.41
1:C:2922:PHE:C	1:C:2923:LEU:HD22	2.45	0.41
1:C:4498:LEU:O	1:C:4501:VAL:HG22	2.20	0.41
1:C:2659:ASP:N	1:C:2659:ASP:OD1	2.54	0.41
1:C:4156:GLU:N	1:C:4156:GLU:OE1	2.53	0.41
1:C:2399:TYR:N	1:C:2633:ASN:OD1	2.54	0.41
1:C:3637:GLU:HB3	1:C:3641:LYS:HZ3	1.86	0.41
1:C:3440:LYS:O	1:C:3444:VAL:HG23	2.21	0.41
1:C:3447:VAL:O	1:C:3451:THR:HG22	2.21	0.41
1:C:3496:ASP:OD1	1:C:3496:ASP:N	2.54	0.41
1:C:3654:LEU:HD11	1:C:3775:LEU:HD11	2.03	0.41
1:C:2336:PHE:CZ	1:C:2337:ILE:HD11	2.56	0.40
1:C:2914:GLY:O	1:C:2990:ARG:NH1	2.54	0.40
1:C:3470:GLN:HG3	1:C:3471:GLN:N	2.34	0.40
1:C:3789:SER:O	1:C:3789:SER:OG	2.34	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	C	2877/4620 (62%)	2606 (91%)	265 (9%)	6 (0%)	43	77
2	Y	64/1200 (5%)	55 (86%)	9 (14%)	0	100	100
All	All	2941/5820 (50%)	2661 (90%)	274 (9%)	6 (0%)	44	77

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	2978	PRO
1	C	2979	PRO
1	C	4182	PRO
1	C	4188	PRO
1	C	4546	PRO
1	C	4121	VAL

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	C	2182/4196 (52%)	2182 (100%)	0	100	100
2	Y	33/1102 (3%)	33 (100%)	0	100	100
All	All	2215/5298 (42%)	2215 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	1567	ASN
1	C	1989	ASN
1	C	2245	ASN
1	C	2333	ASN
1	C	2483	HIS
1	C	2604	ASN
1	C	3065	ASN
1	C	3112	GLN
1	C	3517	ASN
1	C	3536	GLN
1	C	3543	ASN
1	C	3801	ASN
1	C	4125	ASN
1	C	4446	ASN
1	C	4562	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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
3	ADP	C	4704	-	28,29,29	1.34	4 (14%)	43,45,45	1.87	10 (23%)
4	ATP	C	4702	-	32,33,33	1.23	5 (15%)	48,52,52	1.80	9 (18%)
3	ADP	C	4703	-	28,29,29	1.32	4 (14%)	43,45,45	1.98	10 (23%)
3	ADP	C	4701	-	28,29,29	1.37	5 (17%)	43,45,45	1.85	10 (23%)

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
3	ADP	C	4704	-	-	3/16/32/32	0/3/3/3
4	ATP	C	4702	-	-	4/22/38/38	0/3/3/3
3	ADP	C	4703	-	-	3/16/32/32	0/3/3/3
3	ADP	C	4701	-	-	1/16/32/32	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	4704	ADP	C5-C4	4.37	1.46	1.39
3	C	4701	ADP	C5-C4	4.27	1.46	1.39
3	C	4703	ADP	C5-C4	4.19	1.46	1.39
4	C	4702	ATP	C5-C4	4.05	1.46	1.39
3	C	4703	ADP	C5-N7	-2.58	1.34	1.39
3	C	4704	ADP	C5-C6	2.57	1.48	1.41
4	C	4702	ATP	C5-N7	-2.52	1.34	1.39
3	C	4701	ADP	C5-C6	2.49	1.47	1.41
3	C	4701	ADP	C5-N7	-2.45	1.34	1.39
3	C	4704	ADP	C5-N7	-2.45	1.34	1.39
4	C	4702	ATP	C5-C6	2.34	1.47	1.41
3	C	4703	ADP	C5-C6	2.32	1.47	1.41
3	C	4701	ADP	C4-N9	-2.31	1.32	1.37
3	C	4701	ADP	C8-N7	2.28	1.36	1.31
3	C	4704	ADP	C8-N7	2.28	1.36	1.31
3	C	4703	ADP	C8-N7	2.11	1.35	1.31
4	C	4702	ATP	C4-N9	-2.09	1.33	1.37
4	C	4702	ATP	C8-N7	2.02	1.35	1.31

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	4703	ADP	C5-C4-N3	-6.20	118.17	126.72
3	C	4704	ADP	C5-C4-N3	-5.94	118.53	126.72
4	C	4702	ATP	C5-C4-N3	-5.93	118.55	126.72
3	C	4701	ADP	C5-C4-N3	-5.49	119.16	126.72
3	C	4703	ADP	N3-C4-N9	5.26	136.11	127.17
4	C	4702	ATP	N3-C4-N9	4.99	135.66	127.17
3	C	4704	ADP	N3-C4-N9	4.78	135.30	127.17
3	C	4701	ADP	N3-C4-N9	4.33	134.53	127.17
3	C	4703	ADP	C2-N3-C4	3.98	121.55	111.83
4	C	4702	ATP	C2-N3-C4	3.96	121.50	111.83
3	C	4704	ADP	C2-N3-C4	3.87	121.27	111.83
3	C	4701	ADP	C4-C5-N7	-3.71	106.34	110.58
3	C	4703	ADP	N3-C2-N1	-3.71	122.97	128.58
4	C	4702	ATP	N3-C2-N1	-3.71	122.97	128.58
3	C	4704	ADP	N3-C2-N1	-3.57	123.17	128.58
3	C	4701	ADP	C2-N3-C4	3.54	120.48	111.83
3	C	4704	ADP	C4-C5-N7	-3.35	106.75	110.58
3	C	4701	ADP	C4-N9-C8	3.31	109.21	105.74
3	C	4701	ADP	N3-C2-N1	-3.17	123.78	128.58
4	C	4702	ATP	C4-N9-C8	3.14	109.03	105.74

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	4703	ADP	C4-N9-C8	3.02	108.90	105.74
4	C	4702	ATP	C4-C5-N7	-2.99	107.16	110.58
3	C	4703	ADP	C4-C5-N7	-2.94	107.22	110.58
3	C	4704	ADP	C4-N9-C8	2.85	108.73	105.74
3	C	4701	ADP	C5-N7-C8	2.78	107.82	103.45
3	C	4704	ADP	C5-N7-C8	2.63	107.58	103.45
3	C	4703	ADP	C2'-C1'-N9	-2.57	106.91	113.30
3	C	4701	ADP	N9-C8-N7	-2.55	110.32	113.94
4	C	4702	ATP	C5-N7-C8	2.47	107.33	103.45
3	C	4703	ADP	C5-N7-C8	2.47	107.33	103.45
4	C	4702	ATP	N9-C8-N7	-2.30	110.67	113.94
3	C	4703	ADP	C3'-C2'-C1'	2.27	105.76	101.46
3	C	4704	ADP	N9-C8-N7	-2.23	110.77	113.94
3	C	4703	ADP	N9-C8-N7	-2.19	110.83	113.94
3	C	4701	ADP	C6-C5-N7	2.16	136.25	132.09
3	C	4701	ADP	O4'-C1'-N9	2.11	112.14	108.09
4	C	4702	ATP	N6-C6-N1	2.09	123.03	118.38
3	C	4704	ADP	O2A-PA-O1A	2.04	121.91	112.44
3	C	4704	ADP	C6-C5-N7	2.03	136.00	132.09

There are no chirality outliers.

All (11) torsion outliers are listed below:

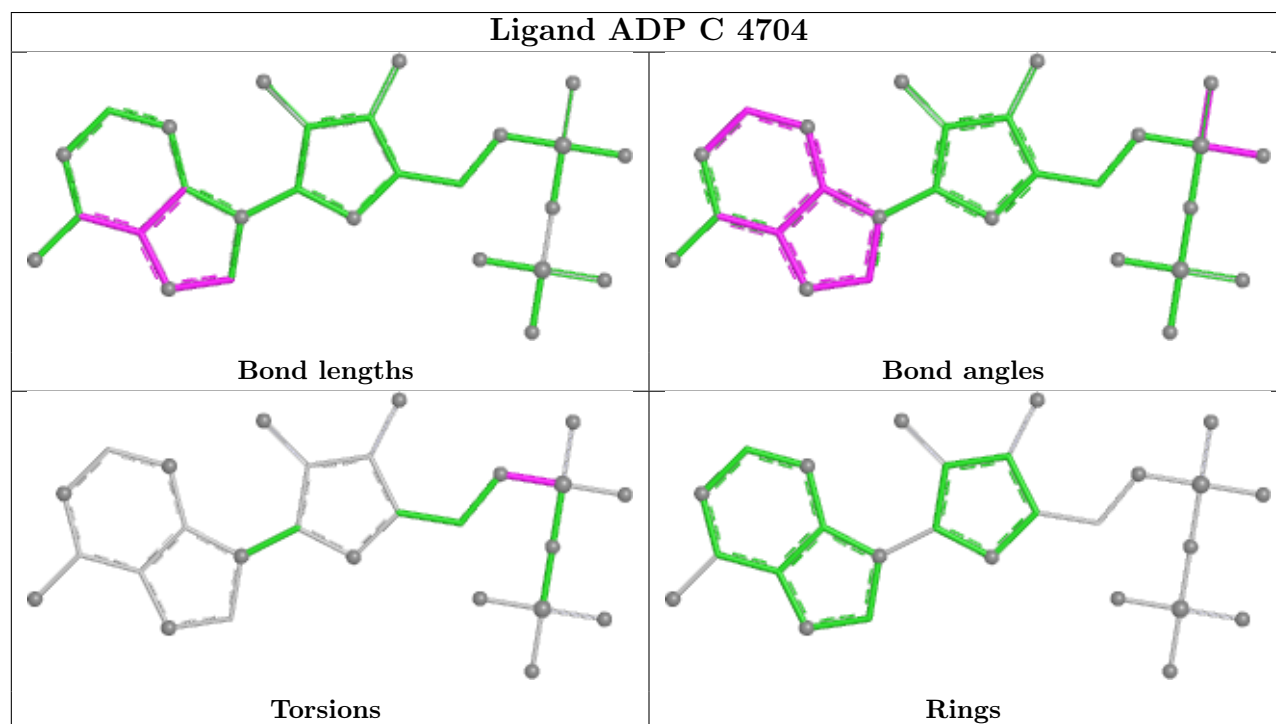
Mol	Chain	Res	Type	Atoms
3	C	4703	ADP	C5'-O5'-PA-O1A
3	C	4703	ADP	C5'-O5'-PA-O2A
3	C	4703	ADP	C5'-O5'-PA-O3A
3	C	4704	ADP	C5'-O5'-PA-O2A
3	C	4704	ADP	C5'-O5'-PA-O3A
4	C	4702	ATP	C5'-O5'-PA-O1A
4	C	4702	ATP	C5'-O5'-PA-O3A
4	C	4702	ATP	O4'-C4'-C5'-O5'
4	C	4702	ATP	C3'-C4'-C5'-O5'
3	C	4701	ADP	C5'-O5'-PA-O1A
3	C	4704	ADP	C5'-O5'-PA-O1A

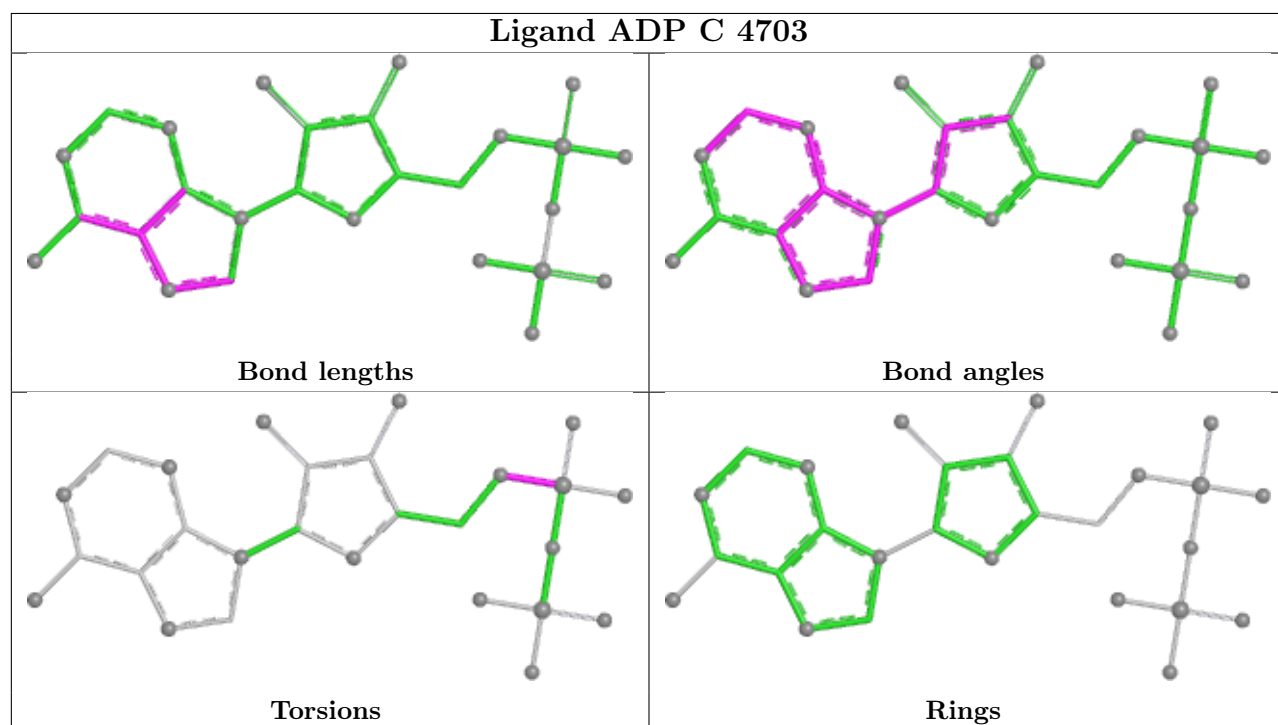
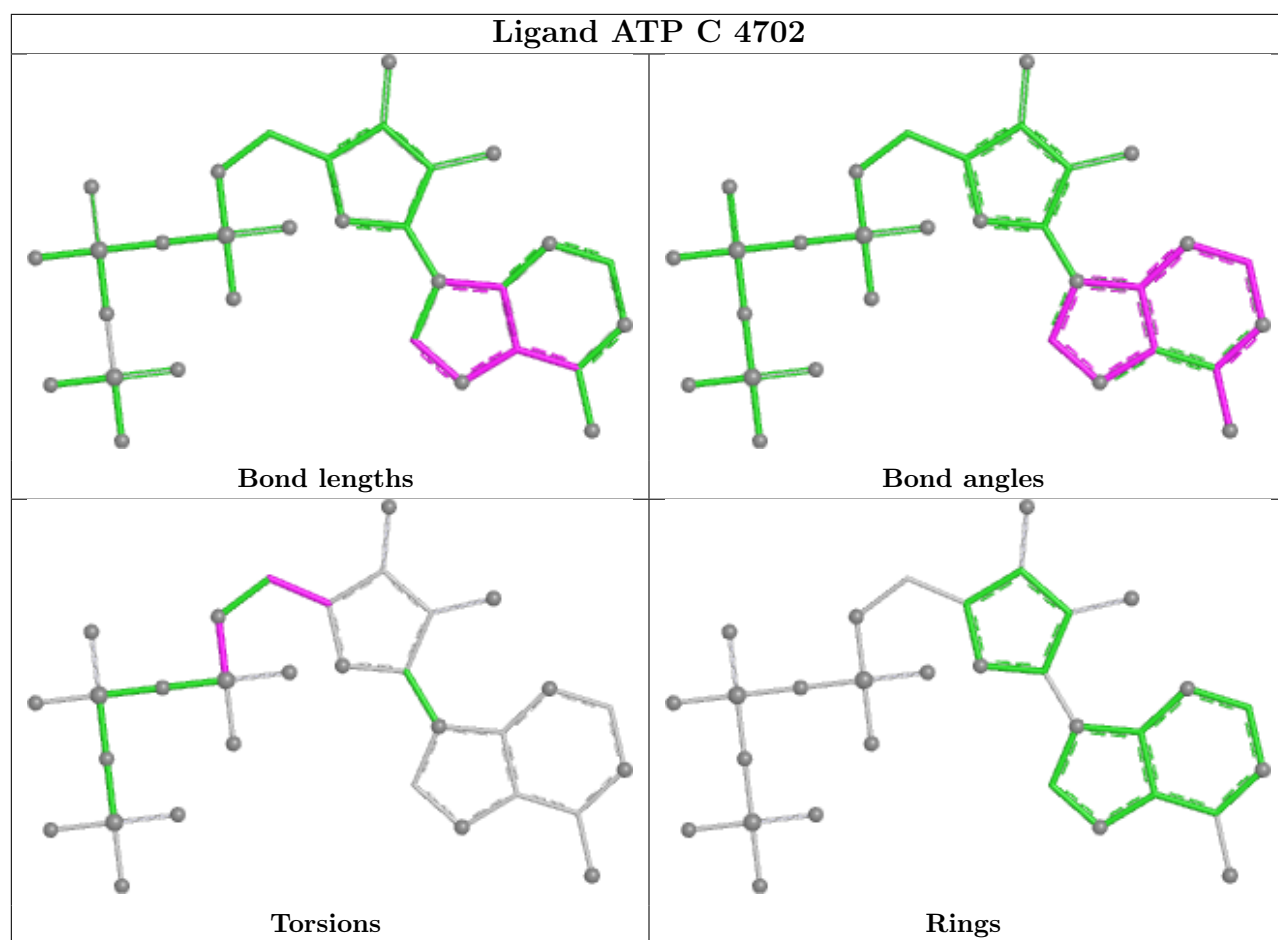
There are no ring outliers.

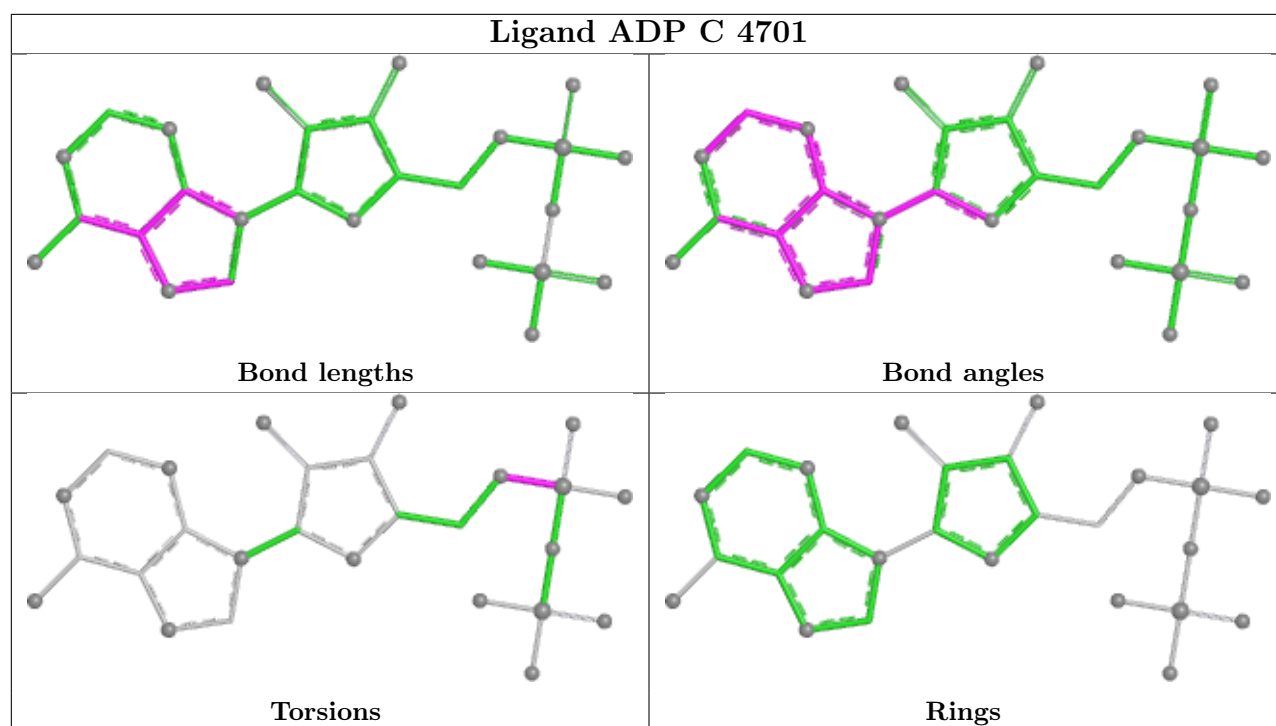
No monomer is involved in short contacts.

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.

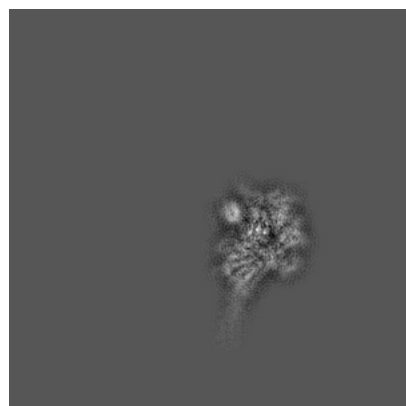
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11581. These allow visual inspection of the internal detail of the map and identification of artifacts.

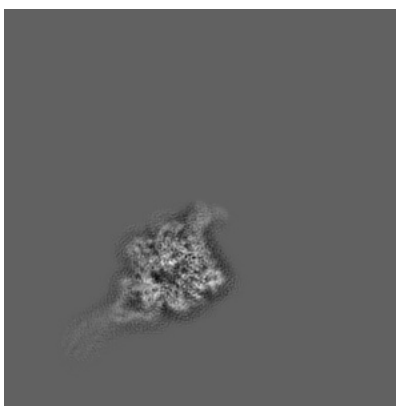
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

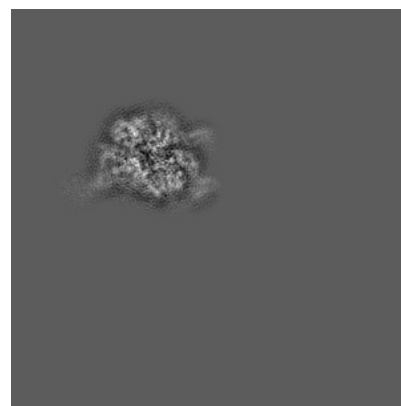
6.1.1 Primary map



X

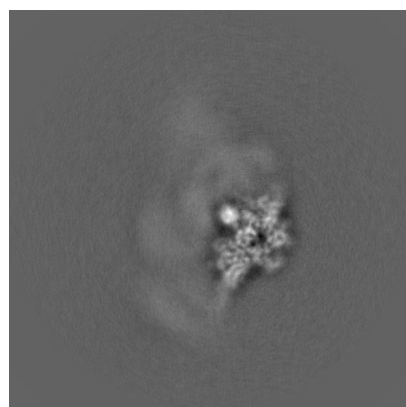


Y

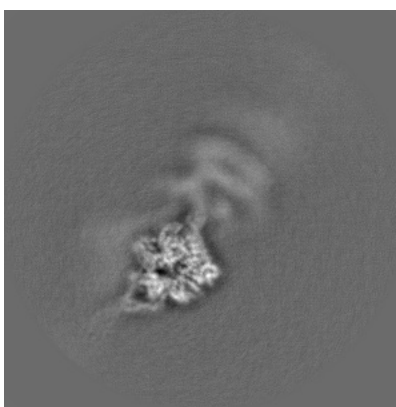


Z

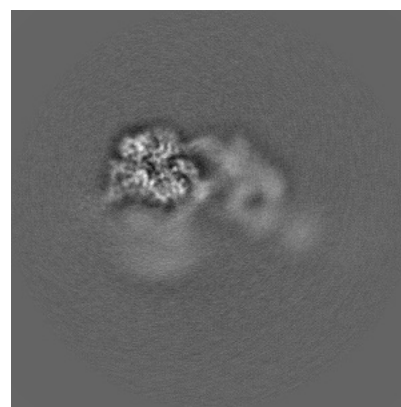
6.1.2 Raw map



X



Y

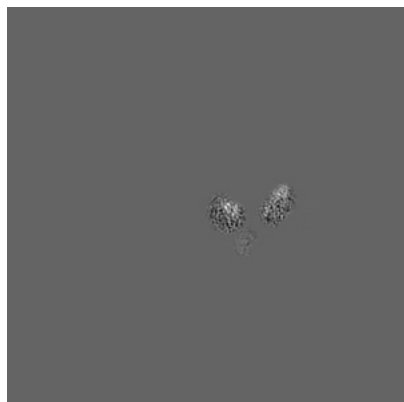


Z

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

6.2 Central slices [i](#)

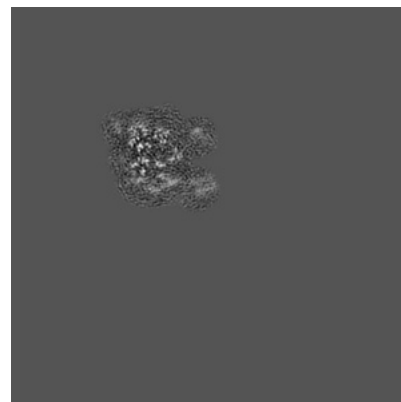
6.2.1 Primary map



X Index: 225

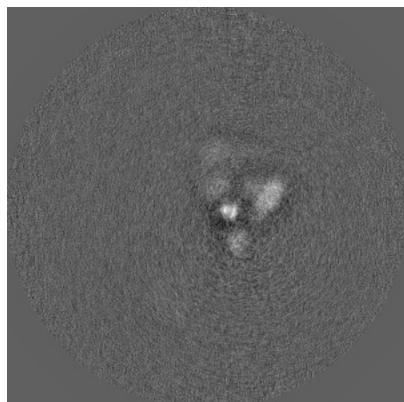


Y Index: 225

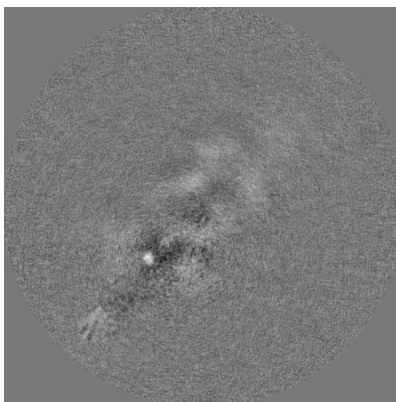


Z Index: 225

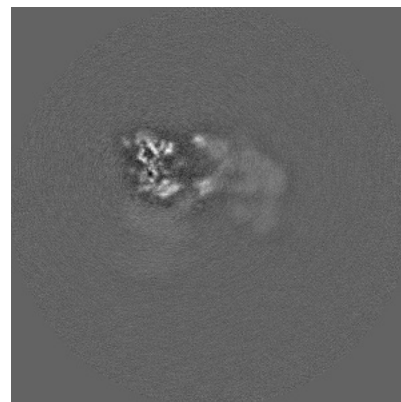
6.2.2 Raw map



X Index: 256



Y Index: 256

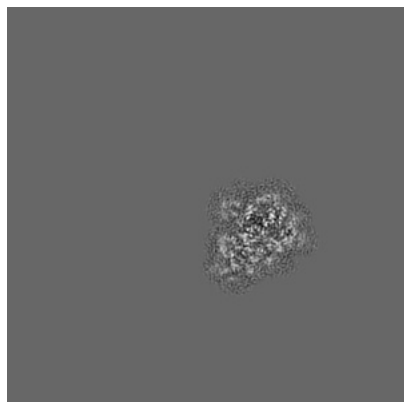


Z Index: 256

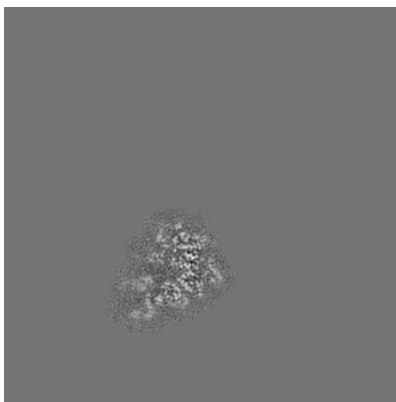
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

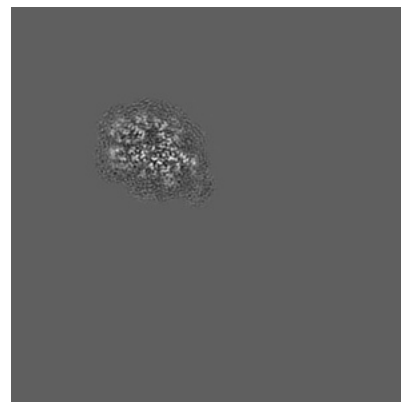
6.3.1 Primary map



X Index: 164

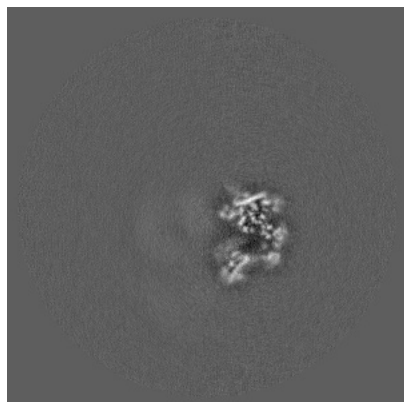


Y Index: 281

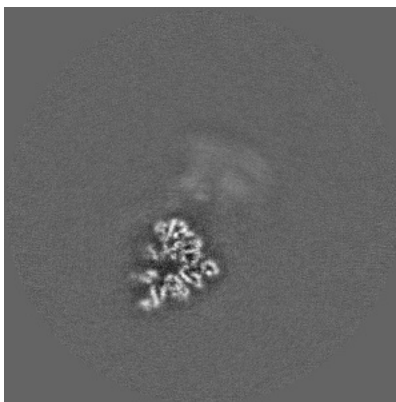


Z Index: 201

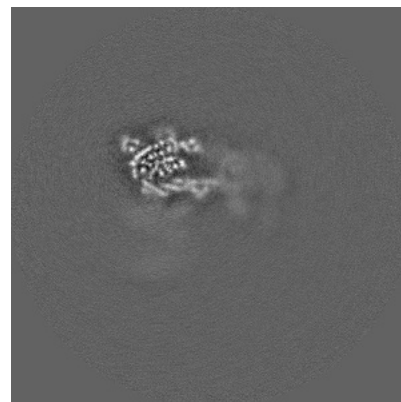
6.3.2 Raw map



X Index: 171



Y Index: 305

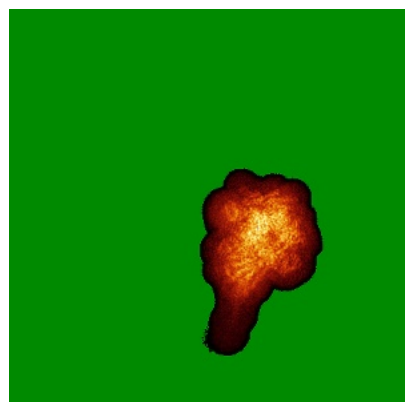


Z Index: 243

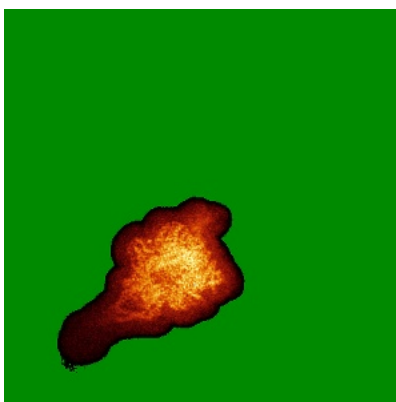
The images above show the largest variance slices of the map in three orthogonal directions.

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

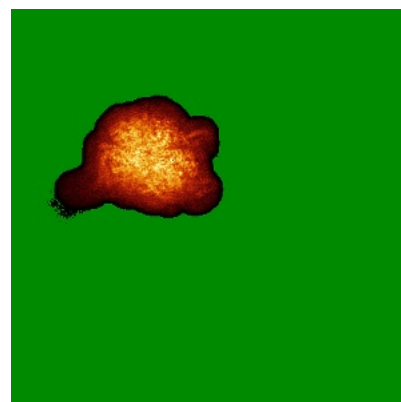
6.4.1 Primary map



X

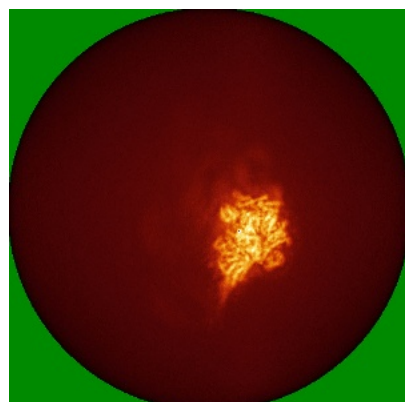


Y

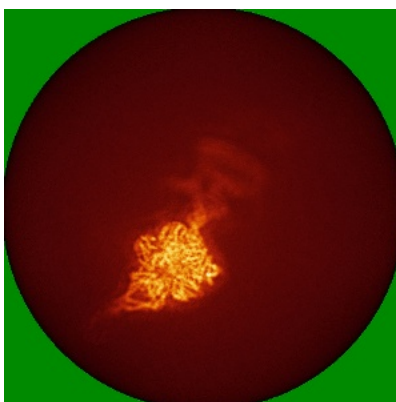


Z

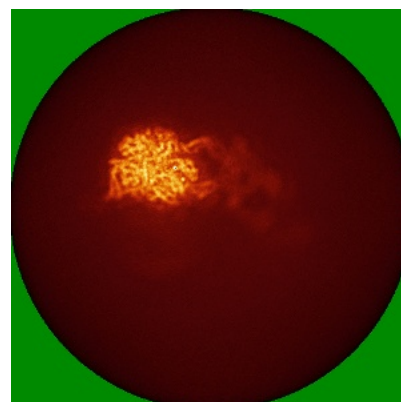
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.015. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

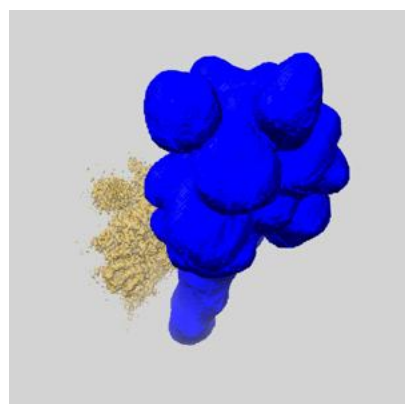
6.6 Mask visualisation [i](#)

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

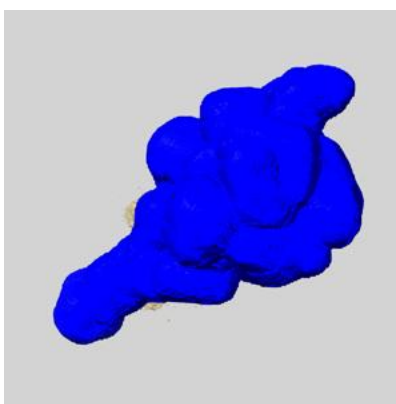
A mask typically either:

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

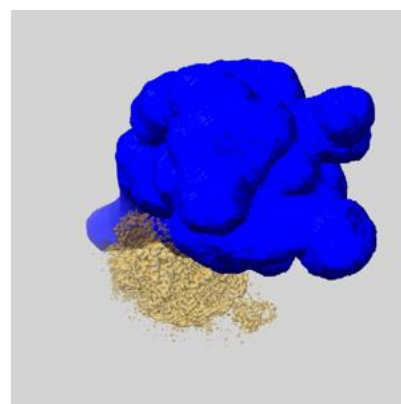
6.6.1 emd_11581_msk_2.map [i](#)



X

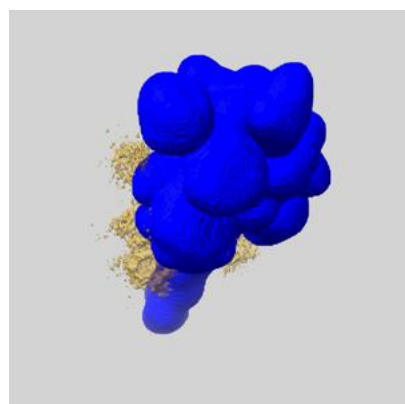


Y

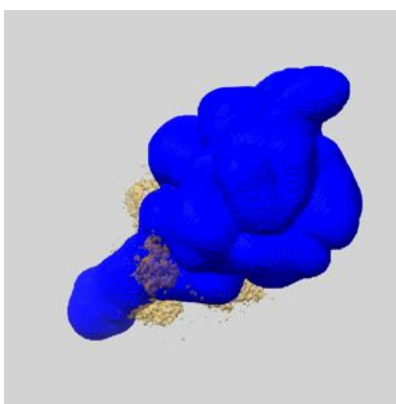


Z

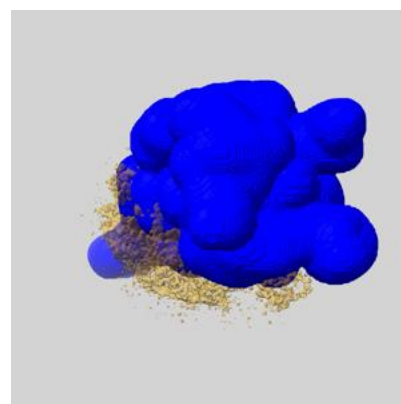
6.6.2 emd_11581_msk_1.map [i](#)



X



Y

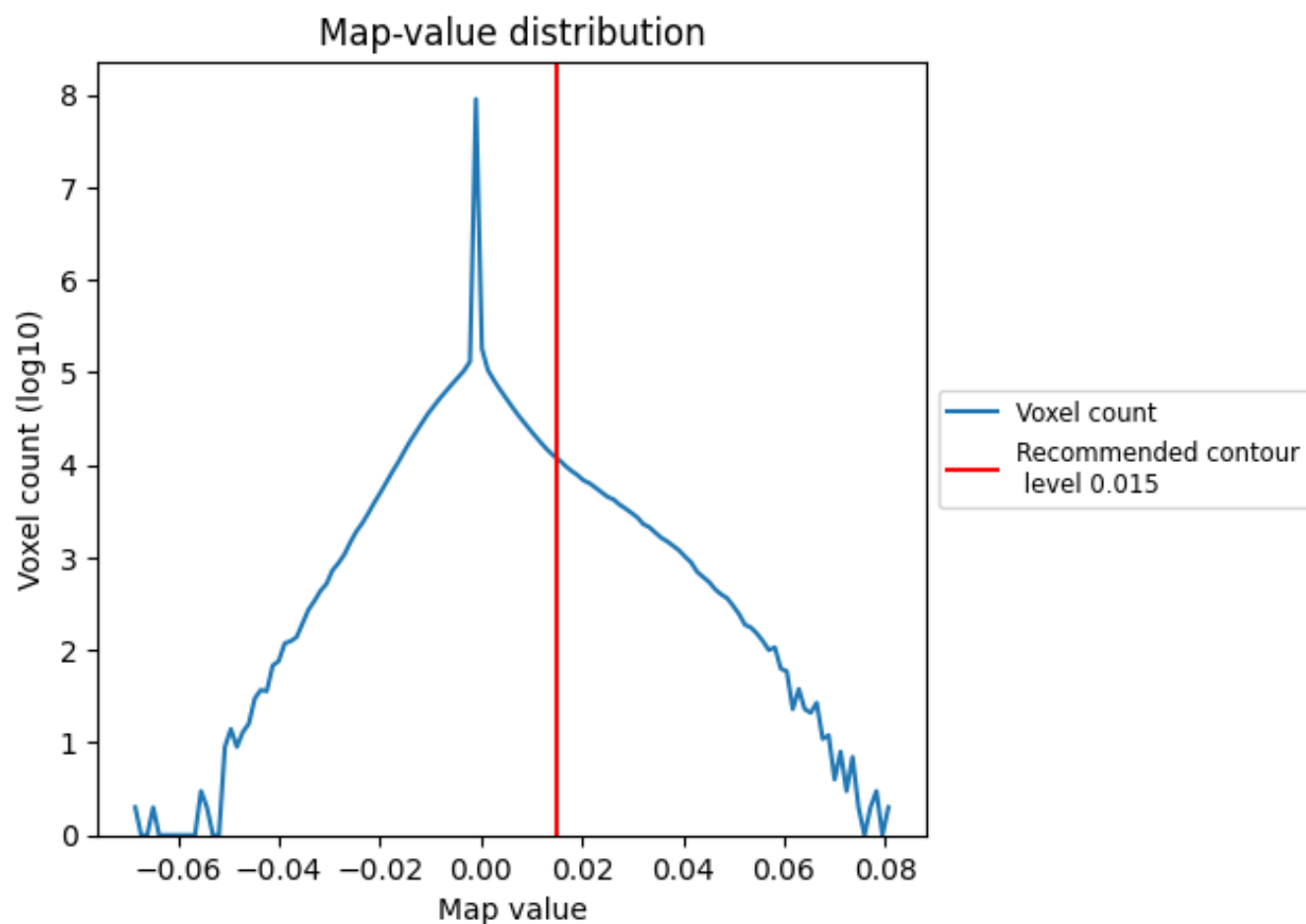


Z

7 Map analysis [i](#)

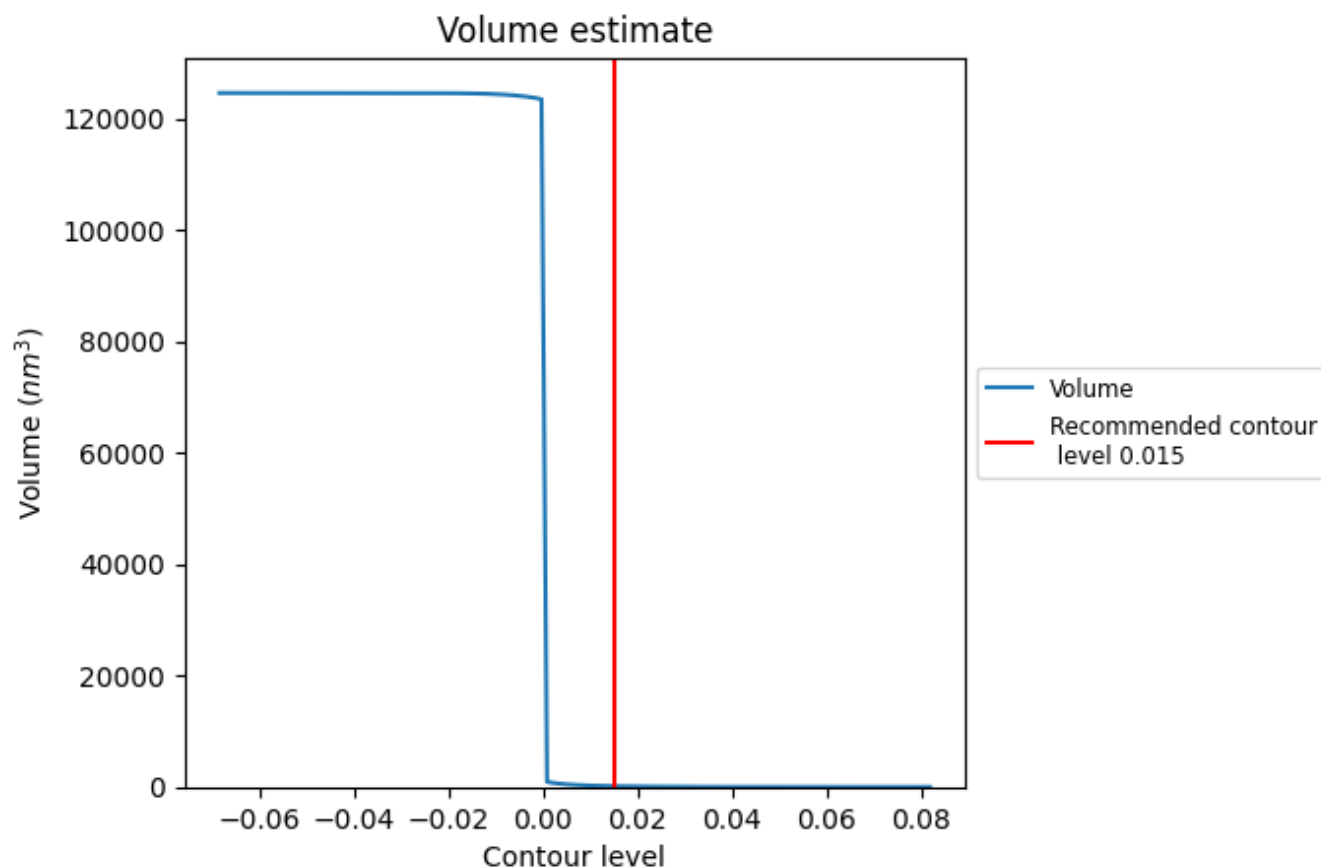
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

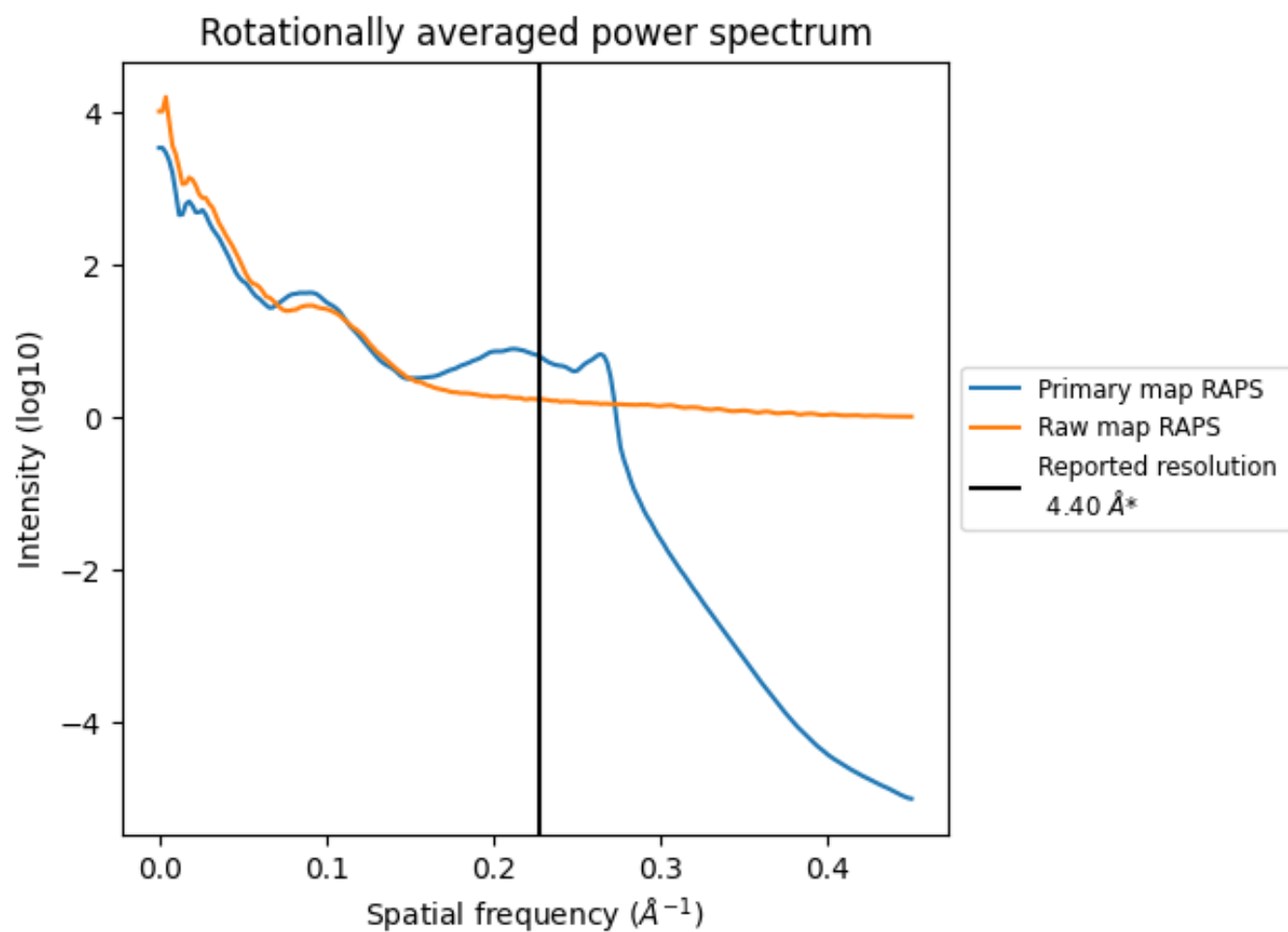
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 147 nm³; this corresponds to an approximate mass of 133 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 ⓘ

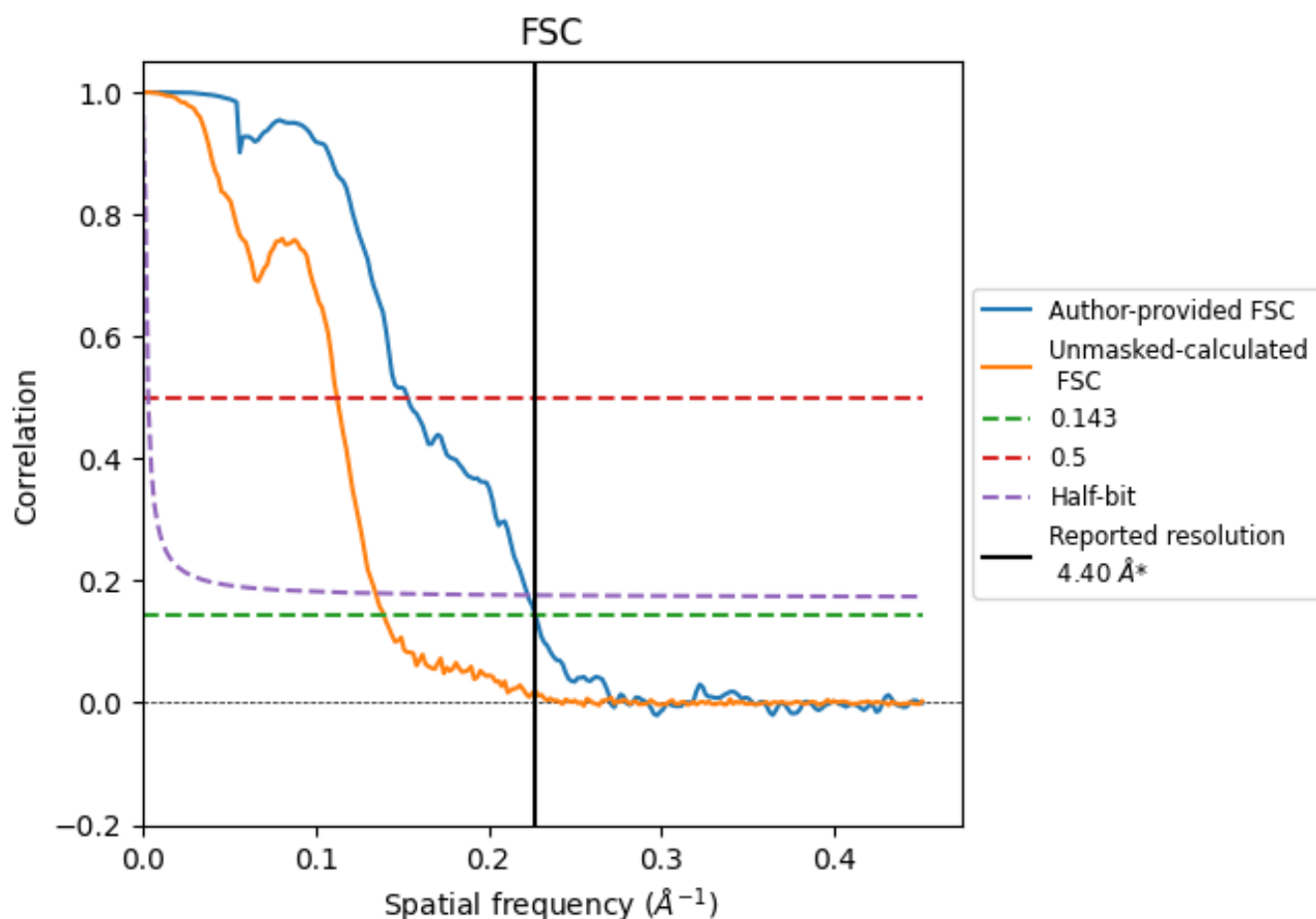


*Reported resolution corresponds to spatial frequency of 0.227 Å⁻¹

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 0.227 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

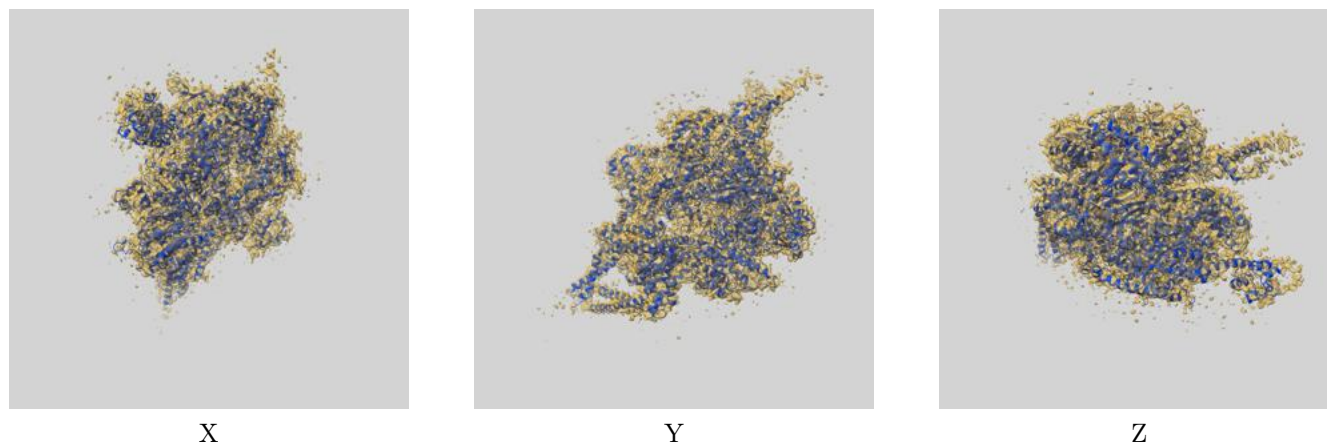
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.40	-	-
Author-provided FSC curve	4.40	6.53	4.49
Unmasked-calculated*	7.16	8.88	7.42

*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 7.16 differs from the reported value 4.4 by more than 10 %

9 Map-model fit [i](#)

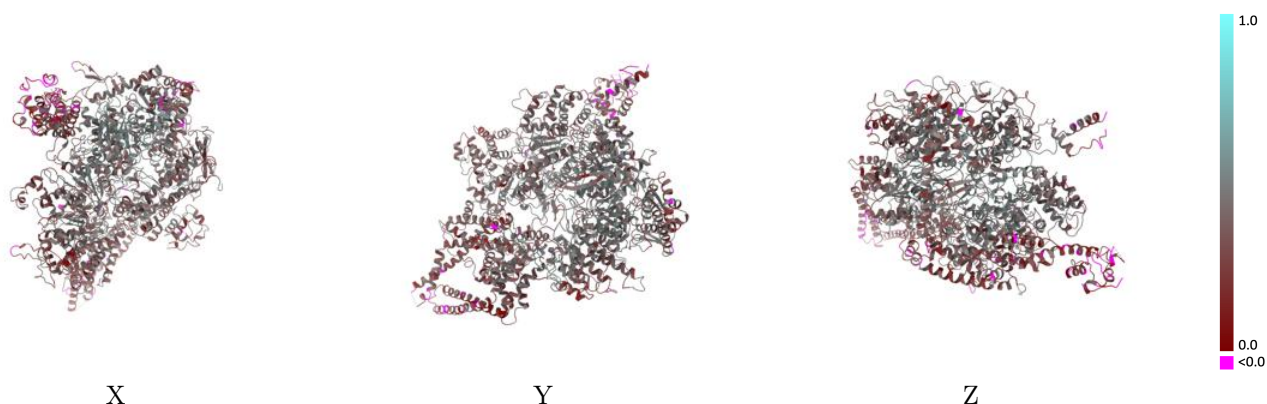
This section contains information regarding the fit between EMDB map EMD-11581 and PDB model 6ZYY. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)



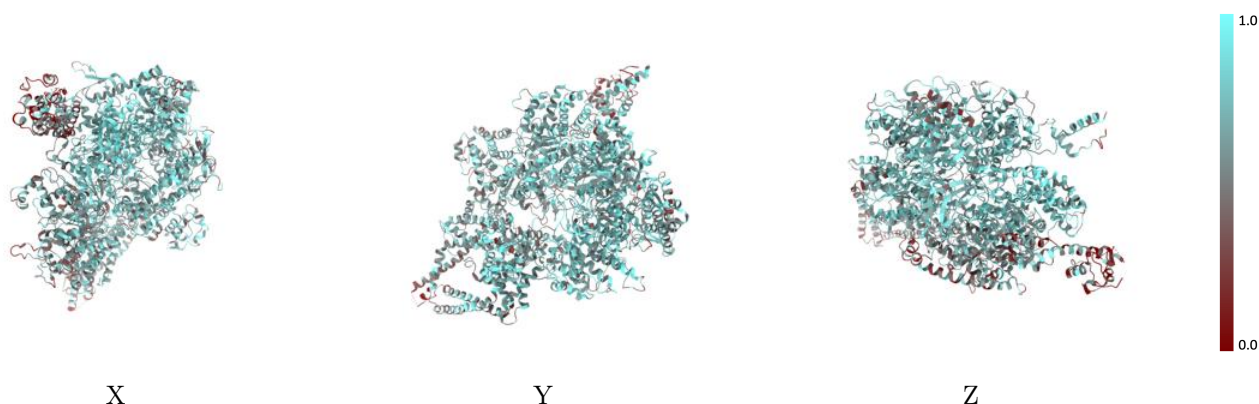
The images above show the 3D surface view of the map at the recommended contour level 0.015 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



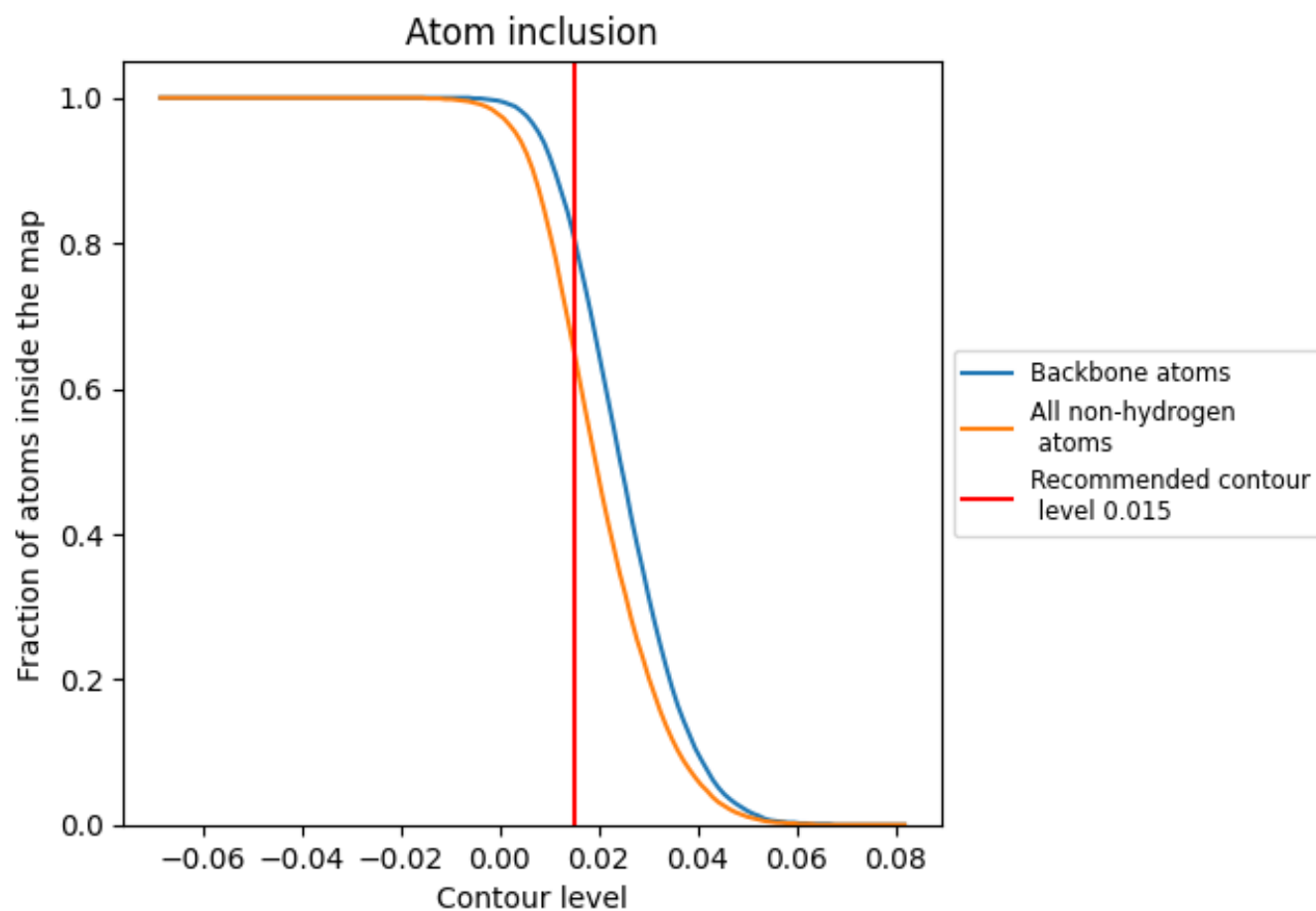
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.015).

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 65% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.015) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6540	0.3880
C	0.6550	0.3890
Y	0.6370	0.3590

