



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 06:18 AM UTC

PDB ID : 7UQK / pdb_00007uqk
EMDB ID : EMD-26697
Title : Cryo-EM structure of the *S. cerevisiae* chromatin remodeler Yta7 hexamer bound to ADP
Authors : Wang, F.; Feng, X.; Li, H.
Deposited on : 2022-04-19
Resolution : 3.10 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:

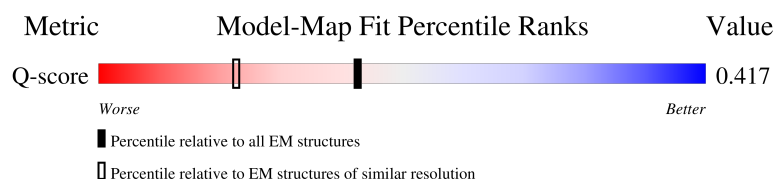
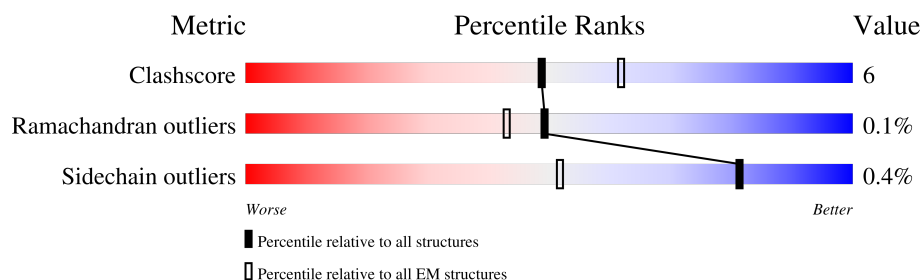
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.10 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	14724 (2.60 - 3.60)

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	1416	
1	B	1416	
1	C	1416	
1	D	1416	

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Mol	Chain	Length	Quality of chain
1	E	1416	15% 33% 6% 61%
1	F	1416	9% 35% • 60%
1	G	1416	11% 9% • 86%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 29142 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 ATPase histone chaperone YTA7.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	576	Total	C	N	O	S	0	0
			4591	2945	780	850	16		
1	B	579	Total	C	N	O	S	0	0
			4614	2959	783	855	17		
1	C	578	Total	C	N	O	S	0	0
			4607	2954	782	854	17		
1	D	576	Total	C	N	O	S	0	0
			4586	2940	780	849	17		
1	E	558	Total	C	N	O	S	0	0
			4448	2855	752	824	17		
1	F	567	Total	C	N	O	S	0	0
			4519	2903	763	836	17		
1	G	197	Total	C	N	O	S	0	0
			1642	1046	288	301	7		

There are 259 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	HIS	-	expression tag	UNP P40340
A	-34	HIS	-	expression tag	UNP P40340
A	-33	HIS	-	expression tag	UNP P40340
A	-32	HIS	-	expression tag	UNP P40340
A	-31	HIS	-	expression tag	UNP P40340
A	-30	HIS	-	expression tag	UNP P40340
A	-29	HIS	-	expression tag	UNP P40340
A	-28	HIS	-	expression tag	UNP P40340
A	-27	HIS	-	expression tag	UNP P40340
A	-26	HIS	-	expression tag	UNP P40340
A	-25	THR	-	expression tag	UNP P40340
A	-24	SER	-	expression tag	UNP P40340
A	-23	GLY	-	expression tag	UNP P40340
A	-22	SER	-	expression tag	UNP P40340
A	-21	MET	-	expression tag	UNP P40340
A	-20	ASP	-	expression tag	UNP P40340

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	TYR	-	expression tag	UNP P40340
A	-18	LYS	-	expression tag	UNP P40340
A	-17	ASP	-	expression tag	UNP P40340
A	-16	HIS	-	expression tag	UNP P40340
A	-15	ASP	-	expression tag	UNP P40340
A	-14	GLY	-	expression tag	UNP P40340
A	-13	ASP	-	expression tag	UNP P40340
A	-12	TYR	-	expression tag	UNP P40340
A	-11	LYS	-	expression tag	UNP P40340
A	-10	ASP	-	expression tag	UNP P40340
A	-9	HIS	-	expression tag	UNP P40340
A	-8	ASP	-	expression tag	UNP P40340
A	-7	ILE	-	expression tag	UNP P40340
A	-6	ASP	-	expression tag	UNP P40340
A	-5	TYR	-	expression tag	UNP P40340
A	-4	LYS	-	expression tag	UNP P40340
A	-3	ASP	-	expression tag	UNP P40340
A	-2	ASP	-	expression tag	UNP P40340
A	-1	ASP	-	expression tag	UNP P40340
A	0	ASP	-	expression tag	UNP P40340
A	1	LYS	-	expression tag	UNP P40340
B	-36	HIS	-	expression tag	UNP P40340
B	-35	HIS	-	expression tag	UNP P40340
B	-34	HIS	-	expression tag	UNP P40340
B	-33	HIS	-	expression tag	UNP P40340
B	-32	HIS	-	expression tag	UNP P40340
B	-31	HIS	-	expression tag	UNP P40340
B	-30	HIS	-	expression tag	UNP P40340
B	-29	HIS	-	expression tag	UNP P40340
B	-28	HIS	-	expression tag	UNP P40340
B	-27	HIS	-	expression tag	UNP P40340
B	-26	THR	-	expression tag	UNP P40340
B	-25	SER	-	expression tag	UNP P40340
B	-24	GLY	-	expression tag	UNP P40340
B	-23	SER	-	expression tag	UNP P40340
B	-22	MET	-	expression tag	UNP P40340
B	-21	ASP	-	expression tag	UNP P40340
B	-20	TYR	-	expression tag	UNP P40340
B	-19	LYS	-	expression tag	UNP P40340
B	-18	ASP	-	expression tag	UNP P40340
B	-17	HIS	-	expression tag	UNP P40340
B	-16	ASP	-	expression tag	UNP P40340

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	GLY	-	expression tag	UNP P40340
B	-14	ASP	-	expression tag	UNP P40340
B	-13	TYR	-	expression tag	UNP P40340
B	-12	LYS	-	expression tag	UNP P40340
B	-11	ASP	-	expression tag	UNP P40340
B	-10	HIS	-	expression tag	UNP P40340
B	-9	ASP	-	expression tag	UNP P40340
B	-8	ILE	-	expression tag	UNP P40340
B	-7	ASP	-	expression tag	UNP P40340
B	-6	TYR	-	expression tag	UNP P40340
B	-5	LYS	-	expression tag	UNP P40340
B	-4	ASP	-	expression tag	UNP P40340
B	-3	ASP	-	expression tag	UNP P40340
B	-2	ASP	-	expression tag	UNP P40340
B	-1	ASP	-	expression tag	UNP P40340
B	0	LYS	-	expression tag	UNP P40340
C	-36	HIS	-	expression tag	UNP P40340
C	-35	HIS	-	expression tag	UNP P40340
C	-34	HIS	-	expression tag	UNP P40340
C	-33	HIS	-	expression tag	UNP P40340
C	-32	HIS	-	expression tag	UNP P40340
C	-31	HIS	-	expression tag	UNP P40340
C	-30	HIS	-	expression tag	UNP P40340
C	-29	HIS	-	expression tag	UNP P40340
C	-28	HIS	-	expression tag	UNP P40340
C	-27	HIS	-	expression tag	UNP P40340
C	-26	THR	-	expression tag	UNP P40340
C	-25	SER	-	expression tag	UNP P40340
C	-24	GLY	-	expression tag	UNP P40340
C	-23	SER	-	expression tag	UNP P40340
C	-22	MET	-	expression tag	UNP P40340
C	-21	ASP	-	expression tag	UNP P40340
C	-20	TYR	-	expression tag	UNP P40340
C	-19	LYS	-	expression tag	UNP P40340
C	-18	ASP	-	expression tag	UNP P40340
C	-17	HIS	-	expression tag	UNP P40340
C	-16	ASP	-	expression tag	UNP P40340
C	-15	GLY	-	expression tag	UNP P40340
C	-14	ASP	-	expression tag	UNP P40340
C	-13	TYR	-	expression tag	UNP P40340
C	-12	LYS	-	expression tag	UNP P40340
C	-11	ASP	-	expression tag	UNP P40340

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	HIS	-	expression tag	UNP P40340
C	-9	ASP	-	expression tag	UNP P40340
C	-8	ILE	-	expression tag	UNP P40340
C	-7	ASP	-	expression tag	UNP P40340
C	-6	TYR	-	expression tag	UNP P40340
C	-5	LYS	-	expression tag	UNP P40340
C	-4	ASP	-	expression tag	UNP P40340
C	-3	ASP	-	expression tag	UNP P40340
C	-2	ASP	-	expression tag	UNP P40340
C	-1	ASP	-	expression tag	UNP P40340
C	0	LYS	-	expression tag	UNP P40340
D	-36	HIS	-	expression tag	UNP P40340
D	-35	HIS	-	expression tag	UNP P40340
D	-34	HIS	-	expression tag	UNP P40340
D	-33	HIS	-	expression tag	UNP P40340
D	-32	HIS	-	expression tag	UNP P40340
D	-31	HIS	-	expression tag	UNP P40340
D	-30	HIS	-	expression tag	UNP P40340
D	-29	HIS	-	expression tag	UNP P40340
D	-28	HIS	-	expression tag	UNP P40340
D	-27	HIS	-	expression tag	UNP P40340
D	-26	THR	-	expression tag	UNP P40340
D	-25	SER	-	expression tag	UNP P40340
D	-24	GLY	-	expression tag	UNP P40340
D	-23	SER	-	expression tag	UNP P40340
D	-22	MET	-	expression tag	UNP P40340
D	-21	ASP	-	expression tag	UNP P40340
D	-20	TYR	-	expression tag	UNP P40340
D	-19	LYS	-	expression tag	UNP P40340
D	-18	ASP	-	expression tag	UNP P40340
D	-17	HIS	-	expression tag	UNP P40340
D	-16	ASP	-	expression tag	UNP P40340
D	-15	GLY	-	expression tag	UNP P40340
D	-14	ASP	-	expression tag	UNP P40340
D	-13	TYR	-	expression tag	UNP P40340
D	-12	LYS	-	expression tag	UNP P40340
D	-11	ASP	-	expression tag	UNP P40340
D	-10	HIS	-	expression tag	UNP P40340
D	-9	ASP	-	expression tag	UNP P40340
D	-8	ILE	-	expression tag	UNP P40340
D	-7	ASP	-	expression tag	UNP P40340
D	-6	TYR	-	expression tag	UNP P40340

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	LYS	-	expression tag	UNP P40340
D	-4	ASP	-	expression tag	UNP P40340
D	-3	ASP	-	expression tag	UNP P40340
D	-2	ASP	-	expression tag	UNP P40340
D	-1	ASP	-	expression tag	UNP P40340
D	0	LYS	-	expression tag	UNP P40340
E	-36	HIS	-	expression tag	UNP P40340
E	-35	HIS	-	expression tag	UNP P40340
E	-34	HIS	-	expression tag	UNP P40340
E	-33	HIS	-	expression tag	UNP P40340
E	-32	HIS	-	expression tag	UNP P40340
E	-31	HIS	-	expression tag	UNP P40340
E	-30	HIS	-	expression tag	UNP P40340
E	-29	HIS	-	expression tag	UNP P40340
E	-28	HIS	-	expression tag	UNP P40340
E	-27	HIS	-	expression tag	UNP P40340
E	-26	THR	-	expression tag	UNP P40340
E	-25	SER	-	expression tag	UNP P40340
E	-24	GLY	-	expression tag	UNP P40340
E	-23	SER	-	expression tag	UNP P40340
E	-22	MET	-	expression tag	UNP P40340
E	-21	ASP	-	expression tag	UNP P40340
E	-20	TYR	-	expression tag	UNP P40340
E	-19	LYS	-	expression tag	UNP P40340
E	-18	ASP	-	expression tag	UNP P40340
E	-17	HIS	-	expression tag	UNP P40340
E	-16	ASP	-	expression tag	UNP P40340
E	-15	GLY	-	expression tag	UNP P40340
E	-14	ASP	-	expression tag	UNP P40340
E	-13	TYR	-	expression tag	UNP P40340
E	-12	LYS	-	expression tag	UNP P40340
E	-11	ASP	-	expression tag	UNP P40340
E	-10	HIS	-	expression tag	UNP P40340
E	-9	ASP	-	expression tag	UNP P40340
E	-8	ILE	-	expression tag	UNP P40340
E	-7	ASP	-	expression tag	UNP P40340
E	-6	TYR	-	expression tag	UNP P40340
E	-5	LYS	-	expression tag	UNP P40340
E	-4	ASP	-	expression tag	UNP P40340
E	-3	ASP	-	expression tag	UNP P40340
E	-2	ASP	-	expression tag	UNP P40340
E	-1	ASP	-	expression tag	UNP P40340

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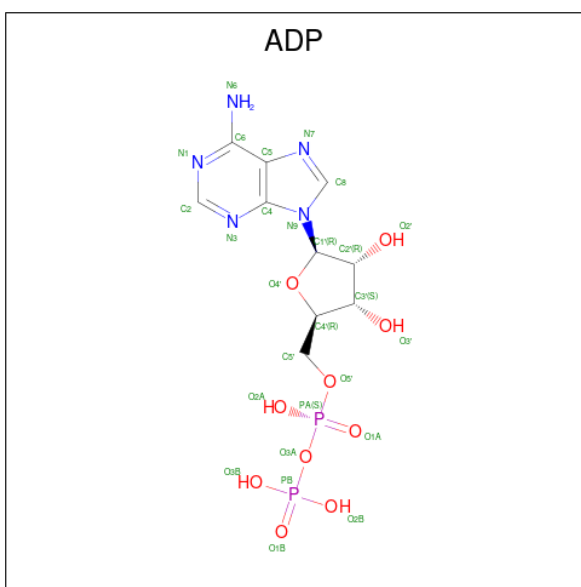
Chain	Residue	Modelled	Actual	Comment	Reference
E	0	LYS	-	expression tag	UNP P40340
F	-36	HIS	-	expression tag	UNP P40340
F	-35	HIS	-	expression tag	UNP P40340
F	-34	HIS	-	expression tag	UNP P40340
F	-33	HIS	-	expression tag	UNP P40340
F	-32	HIS	-	expression tag	UNP P40340
F	-31	HIS	-	expression tag	UNP P40340
F	-30	HIS	-	expression tag	UNP P40340
F	-29	HIS	-	expression tag	UNP P40340
F	-28	HIS	-	expression tag	UNP P40340
F	-27	HIS	-	expression tag	UNP P40340
F	-26	THR	-	expression tag	UNP P40340
F	-25	SER	-	expression tag	UNP P40340
F	-24	GLY	-	expression tag	UNP P40340
F	-23	SER	-	expression tag	UNP P40340
F	-22	MET	-	expression tag	UNP P40340
F	-21	ASP	-	expression tag	UNP P40340
F	-20	TYR	-	expression tag	UNP P40340
F	-19	LYS	-	expression tag	UNP P40340
F	-18	ASP	-	expression tag	UNP P40340
F	-17	HIS	-	expression tag	UNP P40340
F	-16	ASP	-	expression tag	UNP P40340
F	-15	GLY	-	expression tag	UNP P40340
F	-14	ASP	-	expression tag	UNP P40340
F	-13	TYR	-	expression tag	UNP P40340
F	-12	LYS	-	expression tag	UNP P40340
F	-11	ASP	-	expression tag	UNP P40340
F	-10	HIS	-	expression tag	UNP P40340
F	-9	ASP	-	expression tag	UNP P40340
F	-8	ILE	-	expression tag	UNP P40340
F	-7	ASP	-	expression tag	UNP P40340
F	-6	TYR	-	expression tag	UNP P40340
F	-5	LYS	-	expression tag	UNP P40340
F	-4	ASP	-	expression tag	UNP P40340
F	-3	ASP	-	expression tag	UNP P40340
F	-2	ASP	-	expression tag	UNP P40340
F	-1	ASP	-	expression tag	UNP P40340
F	0	LYS	-	expression tag	UNP P40340
G	-36	HIS	-	expression tag	UNP P40340
G	-35	HIS	-	expression tag	UNP P40340
G	-34	HIS	-	expression tag	UNP P40340
G	-33	HIS	-	expression tag	UNP P40340

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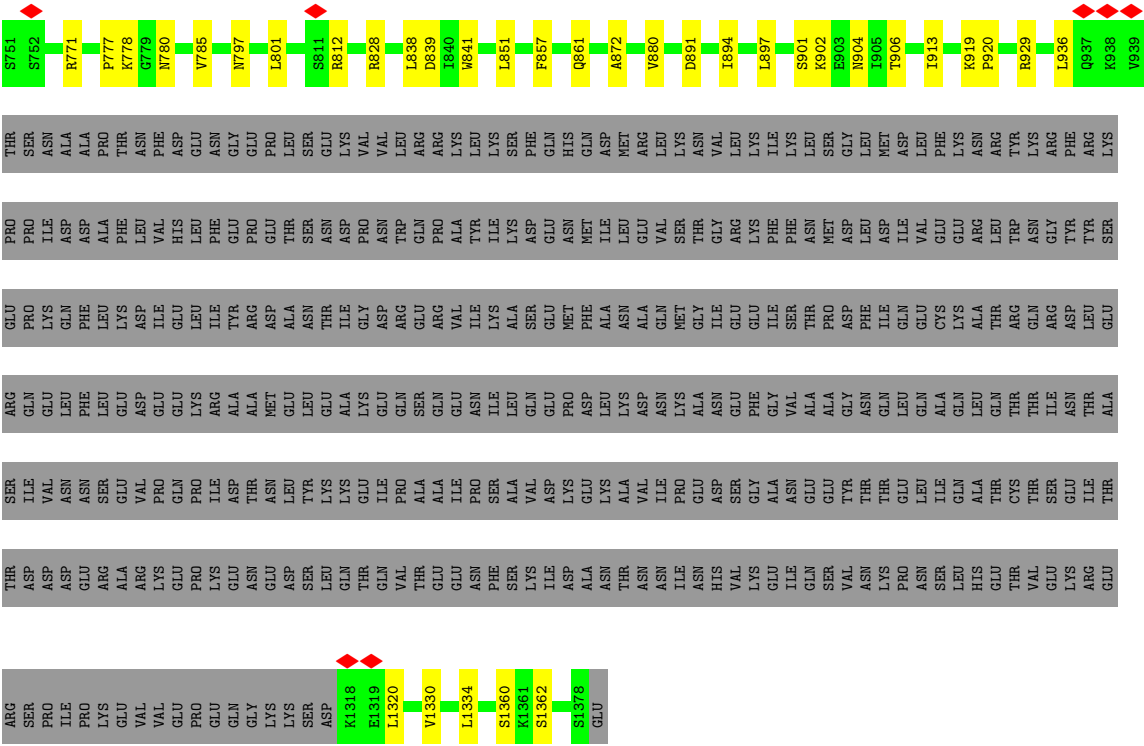
Chain	Residue	Modelled	Actual	Comment	Reference
G	-32	HIS	-	expression tag	UNP P40340
G	-31	HIS	-	expression tag	UNP P40340
G	-30	HIS	-	expression tag	UNP P40340
G	-29	HIS	-	expression tag	UNP P40340
G	-28	HIS	-	expression tag	UNP P40340
G	-27	HIS	-	expression tag	UNP P40340
G	-26	THR	-	expression tag	UNP P40340
G	-25	SER	-	expression tag	UNP P40340
G	-24	GLY	-	expression tag	UNP P40340
G	-23	SER	-	expression tag	UNP P40340
G	-22	MET	-	expression tag	UNP P40340
G	-21	ASP	-	expression tag	UNP P40340
G	-20	TYR	-	expression tag	UNP P40340
G	-19	LYS	-	expression tag	UNP P40340
G	-18	ASP	-	expression tag	UNP P40340
G	-17	HIS	-	expression tag	UNP P40340
G	-16	ASP	-	expression tag	UNP P40340
G	-15	GLY	-	expression tag	UNP P40340
G	-14	ASP	-	expression tag	UNP P40340
G	-13	TYR	-	expression tag	UNP P40340
G	-12	LYS	-	expression tag	UNP P40340
G	-11	ASP	-	expression tag	UNP P40340
G	-10	HIS	-	expression tag	UNP P40340
G	-9	ASP	-	expression tag	UNP P40340
G	-8	ILE	-	expression tag	UNP P40340
G	-7	ASP	-	expression tag	UNP P40340
G	-6	TYR	-	expression tag	UNP P40340
G	-5	LYS	-	expression tag	UNP P40340
G	-4	ASP	-	expression tag	UNP P40340
G	-3	ASP	-	expression tag	UNP P40340
G	-2	ASP	-	expression tag	UNP P40340
G	-1	ASP	-	expression tag	UNP P40340
G	0	LYS	-	expression tag	UNP P40340

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).

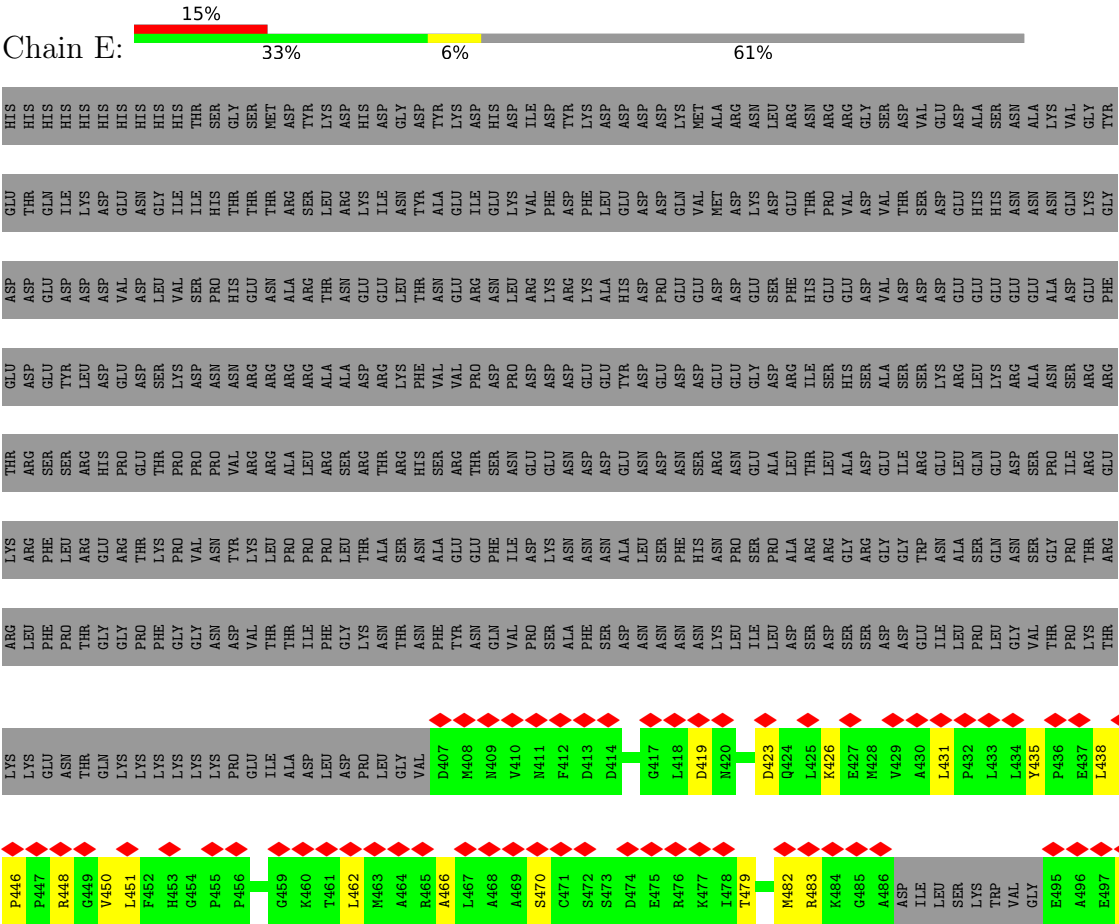


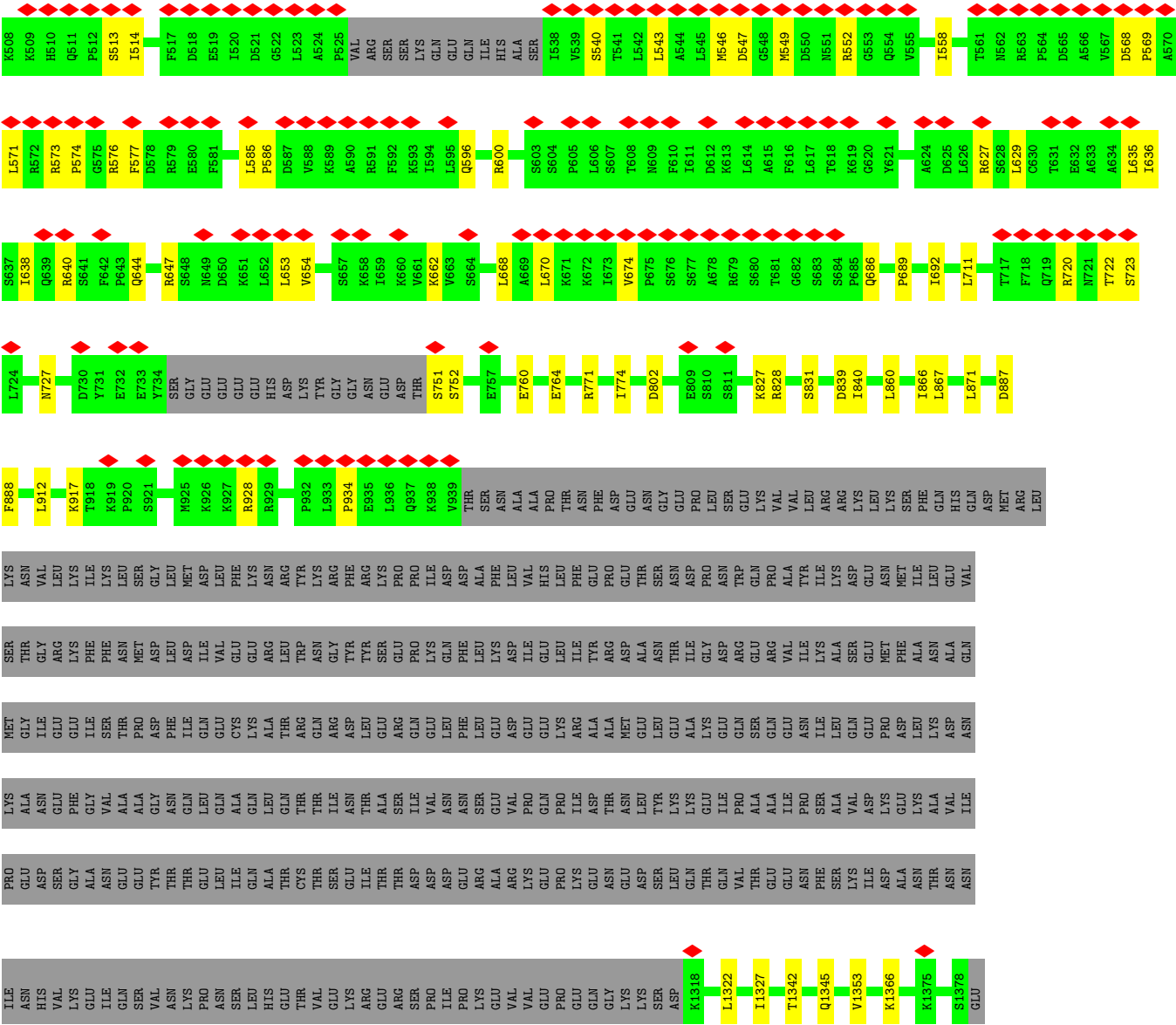
Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total 27	C 10	N 5	O 10	P 2	0
2	B	1	Total 27	C 10	N 5	O 10	P 2	0
2	C	1	Total 27	C 10	N 5	O 10	P 2	0
2	D	1	Total 27	C 10	N 5	O 10	P 2	0
2	E	1	Total 27	C 10	N 5	O 10	P 2	0



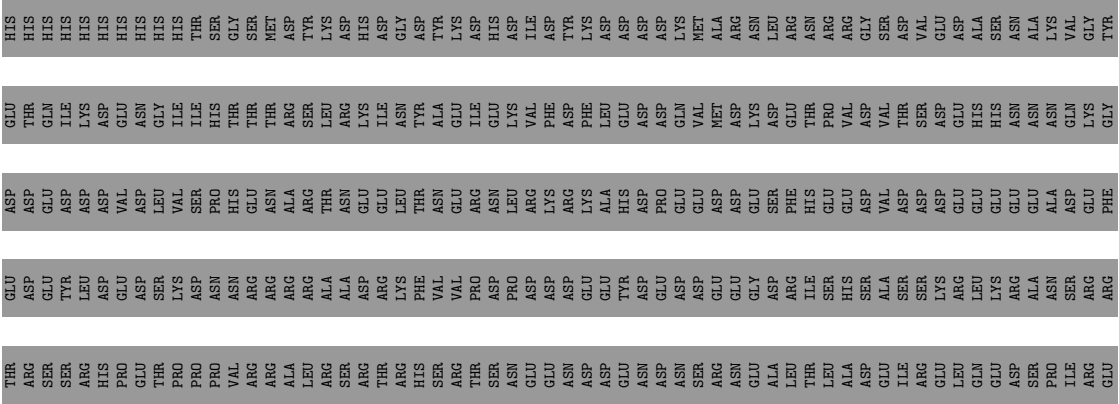


● Molecule 1: ATPase histone chaperone YTA7





• Molecule 1: ATPase histone chaperone YTA7







[illegible]

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	514878	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	65	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.111	Depositor
Minimum map value	-0.065	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.012	Depositor
Map size ($\text{\AA}$)	331.2, 331.2, 331.2	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.828, 0.828, 0.828	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: 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	A	0.23	0/4680	0.49	0/6323
1	B	0.32	2/4703 (0.0%)	0.52	2/6354 (0.0%)
1	C	0.23	0/4696	0.48	0/6344
1	D	0.22	0/4674	0.51	0/6314
1	E	0.23	0/4532	0.55	0/6120
1	F	0.27	1/4606 (0.0%)	0.57	2/6223 (0.0%)
1	G	0.31	0/1672	0.69	1/2237 (0.0%)
All	All	0.26	3/29563 (0.0%)	0.53	5/39915 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	2
1	E	0	1
1	G	0	1
All	All	0	7

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	577	PHE	CA-C	-7.80	1.46	1.53
1	B	410	VAL	C-O	-5.63	1.18	1.24
1	B	409	ASN	N-CA	-5.13	1.40	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	577	PHE	N-CA-C	-8.75	100.67	110.91
1	G	1118	LEU	N-CA-C	-7.22	103.34	111.14
1	B	408	MET	N-CA-C	6.39	119.75	109.79
1	F	550	ASP	N-CA-CB	-5.75	107.62	114.17
1	B	408	MET	CB-CG-SD	-5.08	97.46	112.70

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	533	GLU	Peptide
1	B	927	LYS	Peptide
1	C	511	GLN	Peptide
1	D	491	LYS	Peptide
1	D	653	LEU	Peptide
1	E	674	VAL	Peptide
1	G	346	ASN	Peptide

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	4591	0	4709	51	0
1	B	4614	0	4731	59	0
1	C	4607	0	4722	55	0
1	D	4586	0	4707	67	0
1	E	4448	0	4559	56	0
1	F	4519	0	4635	40	0
1	G	1642	0	1644	54	0
2	A	27	0	12	0	0
2	B	27	0	12	1	0
2	C	27	0	12	0	0
2	D	27	0	12	1	0
2	E	27	0	12	2	0
All	All	29142	0	29767	343	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 6.

All (343) 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:551:ASP:HB2	1:B:406:VAL:CG1	1.63	1.26
1:F:573:ARG:HB3	1:F:574:PRO:CD	1.75	1.16
1:F:573:ARG:HB3	1:F:574:PRO:HD2	1.15	1.10
1:A:551:ASP:HB2	1:B:406:VAL:HG13	1.12	1.08
1:B:449:GLY:HA2	1:B:557:VAL:O	1.67	0.95
1:A:551:ASP:CB	1:B:406:VAL:CG1	2.46	0.94
1:A:551:ASP:CB	1:B:406:VAL:HG13	1.98	0.93
1:D:449:GLY:HA2	1:D:557:VAL:O	1.69	0.91
1:G:998:ARG:HH21	1:G:1047:ASP:HB2	1.36	0.89
1:B:407:ASP:OD2	1:B:480:PHE:CD2	2.29	0.85
1:F:573:ARG:CB	1:F:574:PRO:HD2	2.06	0.84
1:A:551:ASP:CG	1:B:406:VAL:HG11	2.04	0.82
1:G:998:ARG:NH2	1:G:1047:ASP:HB2	1.96	0.79
1:A:551:ASP:O	1:B:406:VAL:HG13	1.83	0.79
1:F:573:ARG:CB	1:F:574:PRO:CD	2.58	0.78
1:G:326:PHE:CE2	1:G:1048:ILE:HD12	2.24	0.73
1:G:1048:ILE:O	1:G:1052:ARG:HD2	1.93	0.68
1:B:804:ALA:O	1:B:808:SER:HB3	1.95	0.67
1:G:1079:GLY:HA3	1:G:1083:ARG:HE	1.61	0.66
1:B:533:GLN:HE22	1:C:525:PRO:HG2	1.62	0.64
1:A:551:ASP:CB	1:B:406:VAL:HG11	2.21	0.64
1:F:448:ARG:O	1:F:576:ARG:HG2	1.98	0.64
1:F:449:GLY:HA2	1:F:557:VAL:O	1.97	0.64
1:G:1046:LEU:HD12	1:G:1046:LEU:C	2.23	0.63
1:D:731:TYR:HA	1:E:928:ARG:HE	1.64	0.63
1:E:640:ARG:HH22	1:E:662:LYS:HG2	1.65	0.62
1:B:406:VAL:HG12	1:B:406:VAL:O	1.99	0.61
1:G:1048:ILE:O	1:G:1048:ILE:HG22	2.00	0.61
1:G:1046:LEU:HD12	1:G:1046:LEU:O	2.01	0.61
1:C:445:THR:HG23	1:D:627:ARG:HH21	1.67	0.60
1:C:428:MET:SD	1:C:579:ARG:NH2	2.74	0.59
1:G:998:ARG:HH21	1:G:1047:ASP:CB	2.12	0.59
1:G:1117:ASP:HA	1:G:1120:ARG:HB3	1.83	0.59
1:G:1044:MET:HB2	1:G:1048:ILE:HG13	1.85	0.59
1:A:551:ASP:HB2	1:B:406:VAL:HG11	1.70	0.59
1:B:501:ARG:HD2	1:C:487:ASP:HB2	1.85	0.58
1:F:1342:THR:H	1:F:1345:GLN:HE21	1.50	0.58
1:E:860:LEU:HD21	1:E:866:ILE:HD12	1.85	0.58
1:A:450:GLY:HA2	1:A:558:VAL:O	2.03	0.58
1:B:574:PRO:HA	1:B:578:ASP:HB3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:596:GLN:O	1:E:600:ARG:NH1	2.35	0.58
1:B:462:LEU:HD22	2:B:1401:ADP:H2'	1.86	0.58
1:C:710:MET:HE2	1:C:770:PRO:HA	1.85	0.58
1:C:644:GLN:HE21	1:C:933:LEU:HD13	1.68	0.58
1:C:494:GLY:H	1:D:491:LYS:HD3	1.69	0.58
1:E:636:ILE:HG12	1:E:668:LEU:HD11	1.86	0.58
1:A:477:ARG:HE	1:A:513:PRO:HG3	1.69	0.57
1:D:797:ASN:HD21	1:D:828:ARG:HH21	1.51	0.57
1:E:479:THR:H	1:E:513:SER:HA	1.69	0.57
1:D:425:LEU:HD13	1:D:467:LEU:HD13	1.87	0.57
1:E:1342:THR:H	1:E:1345:GLN:HE21	1.53	0.57
1:E:540:SER:HA	1:E:543:LEU:HB2	1.84	0.57
1:A:853:VAL:HG22	1:B:803:LEU:HB3	1.86	0.57
1:A:704:LEU:HD13	1:A:790:ILE:HD11	1.86	0.57
1:G:326:PHE:CD2	1:G:1048:ILE:HD12	2.39	0.57
1:F:711:LEU:HD11	1:F:867:LEU:HD22	1.86	0.56
1:A:553:ARG:NH2	1:A:556:VAL:O	2.39	0.56
1:C:479:THR:HG23	1:C:512:PRO:HB2	1.87	0.56
1:E:439:TYR:HE2	1:E:446:PRO:HB3	1.70	0.56
1:E:448:ARG:HG3	1:E:549:MET:HG3	1.88	0.56
1:A:881:VAL:HG11	1:A:895:ILE:HD11	1.87	0.56
1:G:1047:ASP:C	1:G:1049:VAL:H	2.14	0.56
1:A:467:ALA:O	1:A:471:SER:HB3	2.06	0.55
1:B:703:LEU:HD13	1:B:789:ILE:HD11	1.87	0.55
1:B:492:TRP:HD1	1:B:494:GLY:H	1.53	0.55
1:D:475:GLU:HG2	1:D:476:ARG:HG3	1.88	0.55
1:G:1118:LEU:HD23	1:G:1121:GLN:HE21	1.71	0.55
1:C:436:PRO:O	1:C:552:ARG:NH2	2.40	0.55
1:E:839:ASP:OD1	1:E:839:ASP:N	2.39	0.55
1:D:514:ILE:HG22	1:D:556:ILE:HB	1.87	0.55
1:C:546:MET:HG3	1:C:576:ARG:HG2	1.89	0.55
1:F:506:GLU:OE1	1:F:510:HIS:ND1	2.39	0.55
1:B:711:LEU:HD11	1:B:867:LEU:HD22	1.89	0.54
1:D:462:LEU:HD22	2:D:1401:ADP:H2'	1.89	0.54
1:B:858:ARG:NH1	1:C:873:GLU:OE2	2.40	0.54
1:D:652:LEU:HD12	1:D:936:LEU:HB2	1.89	0.54
1:G:998:ARG:HA	1:G:1045:ASP:O	2.08	0.54
1:C:860:LEU:HD21	1:C:866:ILE:HD12	1.88	0.54
1:E:431:LEU:HB3	1:E:438:LEU:HD11	1.90	0.54
1:D:448:ARG:NH2	1:D:550:ASP:OD2	2.41	0.54
1:D:493:VAL:HG11	1:G:1092:ALA:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:591:ARG:HA	1:F:594:ILE:HG22	1.90	0.54
1:D:473:SER:HB3	1:D:477:LYS:HD2	1.88	0.53
1:C:919:LYS:HB2	1:C:922:ASP:HB2	1.90	0.53
1:B:754:ARG:NH1	1:C:914:GLU:OE2	2.40	0.53
1:G:973:MET:HA	1:G:976:LYS:HD2	1.90	0.53
1:A:775:ILE:HB	1:A:872:LEU:HG	1.91	0.53
1:D:880:VAL:HG11	1:D:894:ILE:HD11	1.89	0.53
1:B:815:GLU:HG3	1:B:849:VAL:HG22	1.91	0.53
1:D:427:GLU:HA	1:D:431:LEU:HD13	1.91	0.53
1:G:337:VAL:HG21	1:G:343:LYS:HZ2	1.73	0.53
1:E:435:TYR:HB2	1:E:438:LEU:HD22	1.89	0.53
1:G:326:PHE:CD2	1:G:1048:ILE:CD1	2.92	0.53
1:C:731:TYR:O	1:D:929:ARG:NH1	2.42	0.53
1:C:1318:LYS:NZ	1:C:1319:GLU:O	2.42	0.53
1:G:1092:ALA:HA	1:G:1095:GLN:HB2	1.91	0.53
1:A:429:MET:SD	1:A:580:ARG:NH2	2.83	0.52
1:B:455:PRO:O	1:B:458:THR:OG1	2.26	0.52
1:E:546:MET:HB3	1:E:576:ARG:HH12	1.74	0.52
1:B:689:PRO:HD2	1:B:692:ILE:HD11	1.90	0.52
1:A:587:PRO:O	1:A:592:ARG:NH1	2.43	0.52
1:B:891:ASP:OD1	1:B:891:ASP:N	2.40	0.52
1:E:917:LYS:HD3	1:E:1327:ILE:HG12	1.92	0.52
1:F:521:ASP:OD1	1:F:521:ASP:N	2.42	0.52
1:E:771:ARG:NH1	1:E:888:PHE:O	2.43	0.52
1:A:924:ILE:HG21	1:A:1362:LYS:HD2	1.91	0.52
1:D:429:VAL:HG11	1:D:467:LEU:HD11	1.92	0.52
1:D:678:ALA:HB1	1:D:682:GLY:HA2	1.92	0.52
1:C:475:GLU:OE2	1:D:649:ASN:ND2	2.41	0.52
1:D:451:LEU:HD11	1:D:561:THR:HG22	1.92	0.52
1:G:983:LEU:HD22	1:G:1050:GLU:OE1	2.10	0.52
1:G:987:MET:HE3	1:G:991:LYS:HB3	1.92	0.52
1:B:573:ARG:NH1	1:B:574:PRO:O	2.43	0.52
1:C:530:LYS:HD2	1:C:531:GLN:HB2	1.91	0.52
1:C:814:ILE:HG22	1:C:849:VAL:HG11	1.92	0.51
1:A:494:VAL:HB	1:B:534:ILE:HD13	1.92	0.51
1:A:778:PRO:HG2	1:A:781:ASN:HD22	1.76	0.51
1:C:550:ASP:OD1	1:C:550:ASP:N	2.44	0.51
1:E:720:ARG:HH21	1:E:727:ASN:HB2	1.76	0.51
1:G:1102:SER:HB3	1:G:1107:ILE:HD11	1.92	0.51
1:D:801:LEU:O	1:D:841:TRP:NE1	2.43	0.51
1:D:550:ASP:OD1	1:D:550:ASP:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:901:SER:OG	1:D:904:ASN:ND2	2.43	0.51
1:B:640:ARG:NH1	1:B:665:ASP:OD2	2.43	0.50
1:G:1044:MET:HG3	1:G:1046:LEU:H	1.75	0.50
1:G:326:PHE:CG	1:G:1048:ILE:HD13	2.46	0.50
1:C:543:LEU:HD21	1:C:570:ALA:HB1	1.93	0.50
1:F:448:ARG:HH12	1:F:551:ASN:H	1.58	0.50
1:A:699:ASP:N	1:A:699:ASP:OD1	2.44	0.50
1:A:502:ARG:HD3	1:B:487:ASP:HB2	1.93	0.50
1:D:689:PRO:HD2	1:D:692:ILE:HD11	1.92	0.50
1:E:483:ARG:HG3	1:E:503:LEU:HD22	1.93	0.50
1:A:786:VAL:HG21	1:A:898:LEU:HD11	1.92	0.50
1:D:491:LYS:NZ	1:G:1100:GLU:OE1	2.45	0.50
1:E:573:ARG:HD3	1:F:456:PRO:HB2	1.94	0.49
1:B:425:LEU:HD13	1:B:467:LEU:HD22	1.94	0.49
1:G:1068:ILE:HA	1:G:1071:ILE:HD12	1.94	0.49
1:A:529:SER:HB2	1:B:810:SER:HB3	1.94	0.49
1:F:434:LEU:HD13	1:F:476:ARG:HD3	1.94	0.49
1:G:1050:GLU:OE1	1:G:1050:GLU:HA	2.12	0.49
1:A:835:PHE:HE1	1:A:872:LEU:HD13	1.77	0.49
1:G:1110:CYS:O	1:G:1114:ARG:HB2	2.11	0.49
1:D:543:LEU:O	1:D:576:ARG:NE	2.45	0.49
1:F:458:THR:HG22	1:F:624:ALA:H	1.78	0.49
1:E:644:GLN:HA	1:E:647:ARG:HH12	1.78	0.49
1:B:410:VAL:O	1:B:410:VAL:HG23	2.11	0.48
1:C:683:SER:OG	1:C:684:SER:N	2.46	0.48
1:D:424:GLN:OE1	1:D:579:ARG:NH1	2.45	0.48
1:D:472:SER:OG	1:D:474:ASP:OD1	2.31	0.48
1:F:828:ARG:O	1:F:831:SER:OG	2.31	0.48
1:F:785:VAL:HG21	1:F:897:LEU:HD11	1.95	0.48
1:B:876:ASP:N	1:B:876:ASP:OD1	2.43	0.48
1:E:482:MET:O	1:E:483:ARG:NH2	2.44	0.48
1:C:528:SER:OG	1:C:530:LYS:O	2.29	0.48
1:E:543:LEU:O	1:E:576:ARG:NH1	2.47	0.48
1:G:1098:ILE:HD12	1:G:1101:ILE:HB	1.95	0.48
1:A:733:GLU:HB3	1:B:927:LYS:HG2	1.95	0.48
1:C:650:ASP:OD1	1:C:650:ASP:N	2.44	0.48
1:G:336:ASP:OD1	1:G:336:ASP:N	2.42	0.48
1:G:339:THR:OG1	1:G:340:ILE:N	2.45	0.48
1:G:961:LEU:HD21	1:G:1123:LEU:HD22	1.95	0.48
1:E:462:LEU:HD22	2:E:1401:ADP:H2'	1.96	0.48
1:G:1043:ASN:HD21	1:G:1078:ILE:HD12	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:ASP:OD1	1:A:577:ARG:NH1	2.45	0.47
1:G:995:LYS:HA	1:G:998:ARG:HG3	1.96	0.47
1:C:482:MET:O	1:C:483:ARG:NH2	2.46	0.47
1:D:715:ASP:OD1	1:D:715:ASP:N	2.44	0.47
1:D:771:ARG:HH21	1:D:857:PHE:HB3	1.78	0.47
1:E:751:SER:OG	1:E:752:SER:N	2.45	0.47
1:F:689:PRO:HD2	1:F:692:ILE:HD11	1.96	0.47
1:G:983:LEU:HD13	1:G:1053:LEU:HD12	1.95	0.47
1:C:828:ARG:O	1:C:831:SER:OG	2.33	0.47
1:D:595:LEU:HD21	1:D:626:LEU:HD22	1.94	0.47
1:A:656:ASP:OD1	1:A:656:ASP:N	2.47	0.47
1:B:429:VAL:HG21	1:B:467:LEU:HD21	1.96	0.47
1:B:851:LEU:HB3	1:C:840:ILE:HG23	1.96	0.47
1:G:340:ILE:HB	1:G:1027:LYS:HE3	1.95	0.47
1:B:799:GLN:HG3	1:B:825:ALA:HB2	1.97	0.47
1:C:648:SER:OG	1:C:650:ASP:OD1	2.33	0.47
1:G:332:PHE:HB2	1:G:973:MET:HE1	1.96	0.47
1:C:587:ASP:OD1	1:C:587:ASP:N	2.41	0.47
1:C:619:LYS:NZ	1:C:794:GLU:OE1	2.44	0.47
1:E:1322:LEU:HD13	1:E:1366:LYS:HB2	1.97	0.47
1:A:551:ASP:CG	1:B:406:VAL:CG1	2.77	0.47
1:C:531:GLN:HE21	1:C:533:GLN:HB2	1.80	0.47
1:E:828:ARG:O	1:E:831:SER:OG	2.33	0.47
1:F:472:SER:OG	1:F:473:SER:N	2.47	0.47
1:C:486:ALA:HA	1:C:523:LEU:HD13	1.97	0.47
1:D:586:PRO:HB2	1:D:591:ARG:HG3	1.97	0.47
1:E:887:ASP:OD1	1:E:887:ASP:N	2.41	0.47
1:A:467:ALA:O	1:A:471:SER:CB	2.63	0.46
1:F:720:ARG:HG3	1:F:722:THR:HG23	1.96	0.46
1:B:437:GLU:HG2	1:B:438:LEU:HD12	1.97	0.46
1:D:778:LYS:HA	1:D:778:LYS:HD2	1.72	0.46
1:E:514:ILE:HD12	1:E:514:ILE:HA	1.86	0.46
1:C:855:GLY:HA2	1:C:858:ARG:HG2	1.96	0.46
1:G:326:PHE:CG	1:G:1048:ILE:CD1	2.99	0.46
1:D:438:LEU:HD11	1:E:654:VAL:HG11	1.98	0.46
1:A:888:ASP:OD1	1:A:888:ASP:N	2.38	0.46
1:C:509:LYS:HB2	1:C:509:LYS:HE3	1.79	0.46
1:D:488:ILE:HB	1:D:523:LEU:HD21	1.98	0.46
1:E:585:LEU:HD12	1:E:586:PRO:HD2	1.98	0.46
1:E:547:ASP:N	1:E:547:ASP:OD1	2.49	0.46
1:D:433:LEU:HB3	1:D:476:ARG:HH11	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:574:PRO:HB3	1:D:628:SER:HB3	1.98	0.46
1:G:1053:LEU:HD23	1:G:1053:LEU:HA	1.85	0.46
1:A:678:SER:OG	1:A:685:SER:OG	2.34	0.46
1:G:965:LEU:O	1:G:969:GLN:CB	2.64	0.46
1:D:455:PRO:O	1:D:460:LYS:NZ	2.48	0.45
1:E:571:LEU:HD12	1:E:577:PHE:HB2	1.97	0.45
1:E:629:LEU:HD11	1:E:670:LEU:HD23	1.98	0.45
1:G:1034:GLU:OE1	1:G:1037:THR:OG1	2.34	0.45
1:D:919:LYS:HA	1:D:1320:LEU:HD13	1.98	0.45
1:F:446:PRO:HA	1:F:447:PRO:HD3	1.84	0.45
1:F:929:ARG:HD2	1:F:931:LYS:H	1.81	0.45
1:B:550:ASP:OD1	1:B:550:ASP:N	2.43	0.45
1:C:544:ALA:HB1	1:D:484:LYS:HB3	1.98	0.45
1:C:919:LYS:HZ1	1:C:1319:GLU:HA	1.80	0.45
1:E:827:LYS:HB2	1:E:827:LYS:HE3	1.77	0.45
1:E:912:LEU:HD21	1:E:1353:VAL:HB	1.98	0.45
1:B:923:ILE:HG21	1:B:1361:LYS:HD2	1.98	0.45
1:A:712:LEU:HD11	1:A:868:LEU:HD22	1.99	0.45
1:D:812:ARG:HE	1:D:812:ARG:HB2	1.65	0.45
1:E:689:PRO:HD2	1:E:692:ILE:HD11	1.97	0.45
1:C:450:VAL:HA	1:C:579:ARG:O	2.17	0.45
1:F:887:ASP:OD1	1:F:887:ASP:N	2.38	0.45
1:E:774:ILE:HB	1:E:871:LEU:HD23	1.99	0.45
1:G:1095:GLN:HA	1:G:1098:ILE:HG22	1.99	0.45
1:A:452:LEU:HB3	1:A:581:GLU:HG2	1.98	0.45
1:F:813:THR:HG23	1:F:816:ALA:H	1.81	0.45
1:B:916:LEU:HD13	1:B:1357:ILE:HG12	1.99	0.45
1:D:408:MET:HB3	1:D:465:ARG:HH21	1.82	0.45
1:D:617:LEU:HD12	1:D:618:THR:HG23	1.99	0.45
1:D:785:VAL:HG21	1:D:897:LEU:HD11	1.99	0.44
1:E:573:ARG:HB3	1:E:576:ARG:HB2	1.98	0.44
1:A:645:GLN:HE21	1:A:934:LEU:HG	1.81	0.44
1:F:563:ARG:HD2	1:F:565:ASP:H	1.80	0.44
1:G:954:PRO:HB2	1:G:955:LEU:H	1.59	0.44
1:G:1104:PRO:HA	1:G:1107:ILE:HB	1.98	0.44
1:F:588:VAL:HA	1:F:591:ARG:HG2	1.97	0.44
1:D:487:ASP:N	1:D:487:ASP:OD1	2.50	0.44
1:A:454:HIS:HA	1:A:562:THR:O	2.17	0.44
1:E:711:LEU:HD11	1:E:867:LEU:HD22	1.99	0.44
1:A:916:LEU:HD21	1:A:1359:TRP:HE3	1.83	0.44
1:B:710:MET:HE3	1:B:710:MET:HB2	1.87	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:851:LEU:HB3	1:E:840:ILE:HG23	1.98	0.43
1:G:965:LEU:O	1:G:969:GLN:HB3	2.17	0.43
1:B:637:SER:OG	1:B:665:ASP:OD2	2.33	0.43
1:B:926:LYS:H	1:B:926:LYS:HG2	1.55	0.43
1:C:444:ILE:HB	1:D:631:THR:HG22	1.98	0.43
1:D:423:ASP:OD1	1:D:423:ASP:N	2.49	0.43
1:G:1088:SER:HA	1:G:1091:PHE:HB3	2.00	0.43
1:G:976:LYS:HE2	1:G:1057:TYR:HA	2.01	0.43
1:C:1329:LYS:HA	1:C:1329:LYS:HD2	1.83	0.43
1:D:589:LYS:HB3	1:D:589:LYS:HE3	1.86	0.43
1:E:451:LEU:HD23	1:E:451:LEU:HA	1.93	0.43
1:E:627:ARG:NH1	2:E:1401:ADP:O3'	2.52	0.43
1:B:448:ARG:NH1	1:B:553:GLY:O	2.51	0.43
1:B:497:GLU:OE1	1:B:537:SER:OG	2.30	0.43
1:C:531:GLN:HE22	1:C:535:HIS:HB2	1.84	0.43
1:C:1322:LEU:HD12	1:C:1322:LEU:HA	1.93	0.43
1:F:654:VAL:HG12	1:F:656:PRO:HD3	2.00	0.43
1:F:419:ASP:HA	1:F:422:ILE:HB	2.00	0.43
1:G:1060:GLU:HG3	1:G:1062:LYS:HB2	2.01	0.43
1:B:542:LEU:HG	1:B:546:MET:HE3	2.01	0.43
1:D:587:ASP:OD1	1:D:587:ASP:N	2.41	0.43
1:F:412:PHE:HD2	1:F:422:ILE:HG23	1.83	0.43
1:C:411:ASN:OD1	1:C:411:ASN:N	2.50	0.42
1:F:651:LYS:HA	1:F:651:LYS:HD3	1.75	0.42
1:A:776:ASN:ND2	1:A:897:GLN:OE1	2.44	0.42
1:E:419:ASP:OD1	1:E:419:ASP:N	2.50	0.42
1:B:427:GLU:O	1:B:439:TYR:OH	2.34	0.42
1:C:600:ARG:HG2	1:C:601:LYS:HD2	2.01	0.42
1:C:527:ARG:HD3	1:C:527:ARG:HA	1.86	0.42
1:F:640:ARG:HH12	1:F:662:LYS:HD3	1.85	0.42
1:G:1110:CYS:HA	1:G:1113:THR:HG22	2.01	0.42
1:C:427:GLU:HB2	1:D:646:TYR:HE2	1.84	0.42
1:D:777:PRO:HD2	1:D:780:ASN:HD22	1.84	0.42
1:E:466:ALA:O	1:E:470:SER:OG	2.32	0.42
1:A:551:ASP:OD2	1:B:406:VAL:HG11	2.17	0.42
1:D:441:ASN:OD1	1:D:441:ASN:N	2.52	0.42
1:D:1360:SER:O	1:D:1360:SER:OG	2.33	0.42
1:F:712:ASN:O	1:F:716:THR:OG1	2.31	0.42
1:A:429:MET:HE1	1:A:450:GLY:H	1.85	0.42
1:A:682:THR:OG1	1:A:683:GLY:N	2.53	0.42
1:C:543:LEU:HD12	1:C:571:LEU:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:913:ILE:HD13	1:D:913:ILE:HA	1.95	0.42
1:C:801:LEU:HB2	1:C:835:ILE:HG12	2.02	0.42
1:C:877:ILE:HG22	1:C:881:LYS:HE3	2.02	0.42
1:D:891:ASP:N	1:D:891:ASP:OD1	2.48	0.42
1:D:902:LYS:O	1:D:906:THR:OG1	2.31	0.42
1:E:722:THR:OG1	1:E:723:SER:N	2.53	0.42
1:A:499:ARG:HH12	1:G:1104:PRO:HD2	1.84	0.41
1:D:474:ASP:OD1	1:D:474:ASP:N	2.35	0.41
1:E:450:VAL:HB	1:E:558:ILE:HG23	2.01	0.41
1:D:838:LEU:HB3	1:D:872:ALA:HB2	2.01	0.41
1:A:727:GLN:HE22	1:B:647:ARG:HH12	1.68	0.41
1:D:839:ASP:OD1	1:D:839:ASP:N	2.50	0.41
1:D:861:GLN:NE2	1:E:686:GLN:OE1	2.52	0.41
1:D:920:PRO:HB3	1:D:1362:SER:HA	2.02	0.41
1:C:637:SER:OG	1:C:665:ASP:OD2	2.37	0.41
1:E:635:LEU:HA	1:E:638:ILE:HG22	2.02	0.41
1:E:760:GLU:O	1:E:764:GLU:HG3	2.21	0.41
1:C:846:PRO:HG2	1:C:849:VAL:HG23	2.02	0.41
1:C:1353:VAL:HG13	1:C:1373:ILE:HG12	2.02	0.41
1:E:498:ARG:NH2	1:G:1076:ASN:O	2.53	0.41
1:E:653:LEU:HD13	1:E:934:PRO:HB2	2.02	0.41
1:B:521:ASP:OD1	1:B:521:ASP:N	2.52	0.41
1:B:826:LYS:HA	1:B:866:ILE:HD11	2.01	0.41
1:G:999:LYS:HD3	1:G:1000:PRO:HD2	2.03	0.41
1:A:533:GLU:HG2	1:B:531:GLN:HE21	1.86	0.41
1:A:802:LEU:HB2	1:A:836:ILE:HG12	2.03	0.41
1:D:573:ARG:HD3	1:D:576:ARG:HD3	2.03	0.41
1:D:646:TYR:O	1:D:647:ARG:NE	2.54	0.41
1:F:660:LYS:HA	1:F:660:LYS:HD3	1.81	0.41
1:C:538:ILE:HD13	1:C:538:ILE:HA	1.96	0.41
1:C:891:ASP:N	1:C:891:ASP:OD1	2.49	0.41
1:F:876:ASP:N	1:F:876:ASP:OD1	2.51	0.41
1:B:407:ASP:OD2	1:B:480:PHE:CE2	2.73	0.41
1:B:672:LYS:HE2	1:B:672:LYS:HB3	1.91	0.41
1:C:407:ASP:HA	1:C:465:ARG:HD3	2.02	0.41
1:D:575:GLY:N	1:D:578:ASP:OD1	2.43	0.41
1:E:568:ASP:HA	1:E:569:PRO:HD3	1.94	0.41
1:E:802:ASP:N	1:E:802:ASP:OD1	2.54	0.41
1:F:437:GLU:HA	1:F:440:GLN:HE21	1.86	0.41
1:B:492:TRP:HE1	1:B:495:GLU:HB2	1.86	0.41
1:F:563:ARG:HE	1:F:566:ALA:HB2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:549:MET:O	1:E:552:ARG:NH1	2.54	0.40
1:B:435:TYR:HB3	1:B:438:LEU:HD13	2.03	0.40
1:D:446:PRO:HA	1:D:447:PRO:HD3	1.96	0.40
1:F:587:ASP:O	1:F:591:ARG:N	2.39	0.40
1:F:802:ASP:HB3	1:F:836:PRO:HG2	2.03	0.40
1:D:1330:VAL:O	1:D:1334:LEU:HB2	2.22	0.40
1:A:484:ARG:HA	1:A:484:ARG:HD3	1.97	0.40
1:A:705:LYS:O	1:A:709:ASP:N	2.52	0.40
1:A:709:ASP:O	1:A:715:LYS:NZ	2.55	0.40
1:A:839:LEU:HB3	1:A:873:ALA:HB2	2.03	0.40
1:E:573:ARG:HH11	1:E:574:PRO:HD2	1.86	0.40
1:F:561:THR:OG1	1:F:562:ASN:N	2.54	0.40
1:F:839:ASP:OD1	1:F:839:ASP:N	2.54	0.40
1:A:457:PRO:HG3	1:A:563:ASN:HD21	1.87	0.40
1:E:423:ASP:HA	1:E:426:LYS:HB3	2.03	0.40
1:G:1025:TYR:HA	1:G:1033:LEU:O	2.22	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	570/1416 (40%)	527 (92%)	43 (8%)	0	100	100
1	B	573/1416 (40%)	543 (95%)	30 (5%)	0	100	100
1	C	572/1416 (40%)	531 (93%)	41 (7%)	0	100	100
1	D	570/1416 (40%)	519 (91%)	51 (9%)	0	100	100
1	E	548/1416 (39%)	496 (90%)	52 (10%)	0	100	100
1	F	559/1416 (40%)	509 (91%)	49 (9%)	1 (0%)	43	73
1	G	191/1416 (14%)	182 (95%)	7 (4%)	2 (1%)	12	41
All	All	3583/9912 (36%)	3307 (92%)	273 (8%)	3 (0%)	49	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	1048	ILE
1	G	1047	ASP
1	F	573	ARG

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	518/1281 (40%)	518 (100%)	0	100	100
1	B	521/1281 (41%)	519 (100%)	2 (0%)	84	86
1	C	520/1281 (41%)	519 (100%)	1 (0%)	87	89
1	D	518/1281 (40%)	517 (100%)	1 (0%)	87	89
1	E	502/1281 (39%)	502 (100%)	0	100	100
1	F	510/1281 (40%)	508 (100%)	2 (0%)	84	86
1	G	179/1281 (14%)	173 (97%)	6 (3%)	32	63
All	All	3268/8967 (36%)	3256 (100%)	12 (0%)	81	86

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	411	ASN
1	B	674	VAL
1	C	627	ARG
1	D	474	ASP
1	F	576	ARG
1	F	599	THR
1	G	340	ILE
1	G	1046	LEU
1	G	1047	ASP
1	G	1048	ILE
1	G	1050	GLU
1	G	1118	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (81) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	442	ASN
1	A	555	GLN
1	A	563	ASN
1	A	645	GLN
1	A	687	GLN
1	A	703	ASN
1	A	713	ASN
1	A	727	GLN
1	A	781	ASN
1	A	792	ASN
1	A	800	GLN
1	A	801	ASN
1	A	830	GLN
1	A	849	ASN
1	A	883	ASN
1	A	1351	HIS
1	B	420	ASN
1	B	441	ASN
1	B	453	HIS
1	B	510	HIS
1	B	531	GLN
1	B	533	GLN
1	B	639	GLN
1	B	702	ASN
1	B	719	GLN
1	B	726	GLN
1	B	797	ASN
1	B	863	ASN
1	B	893	ASN
1	B	911	ASN
1	B	1337	HIS
1	C	443	ASN
1	C	499	GLN
1	C	531	GLN
1	C	535	HIS
1	C	562	ASN
1	C	644	GLN
1	C	726	GLN
1	C	797	ASN
1	C	800	ASN
1	C	893	ASN

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Mol	Chain	Res	Type
1	C	1350	HIS
1	D	791	ASN
1	D	799	GLN
1	D	848	ASN
1	D	863	ASN
1	D	904	ASN
1	D	911	ASN
1	D	1350	HIS
1	E	409	ASN
1	E	424	GLN
1	E	443	ASN
1	E	551	ASN
1	E	596	GLN
1	E	639	GLN
1	E	719	GLN
1	E	727	ASN
1	E	775	ASN
1	E	791	ASN
1	E	820	GLN
1	E	837	ASN
1	E	861	GLN
1	E	863	ASN
1	E	896	GLN
1	E	1337	HIS
1	E	1340	ASN
1	E	1345	GLN
1	F	511	GLN
1	F	596	GLN
1	F	797	ASN
1	F	829	GLN
1	F	848	ASN
1	F	863	ASN
1	F	896	GLN
1	F	904	ASN
1	F	1339	GLN
1	F	1345	GLN
1	G	344	ASN
1	G	977	ASN
1	G	1095	GLN
1	G	1121	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](#)

5 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$
2	ADP	D	1401	-	28,29,29	1.36	4 (14%)	43,45,45	1.83	8 (18%)
2	ADP	B	1401	-	28,29,29	1.35	4 (14%)	43,45,45	1.86	10 (23%)
2	ADP	E	1401	-	28,29,29	1.41	5 (17%)	43,45,45	1.88	8 (18%)
2	ADP	C	1401	-	28,29,29	1.34	4 (14%)	43,45,45	1.89	8 (18%)
2	ADP	A	1401	-	28,29,29	1.37	4 (14%)	43,45,45	1.90	8 (18%)

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
2	ADP	D	1401	-	-	1/16/32/32	0/3/3/3
2	ADP	B	1401	-	-	2/16/32/32	0/3/3/3
2	ADP	E	1401	-	-	3/16/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	C	1401	-	-	4/16/32/32	0/3/3/3
2	ADP	A	1401	-	-	4/16/32/32	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1401	ADP	C5-C4	4.65	1.47	1.39
2	A	1401	ADP	C5-C4	4.47	1.47	1.39
2	D	1401	ADP	C5-C4	4.46	1.47	1.39
2	B	1401	ADP	C5-C4	4.28	1.46	1.39
2	C	1401	ADP	C5-C4	4.28	1.46	1.39
2	C	1401	ADP	C5-N7	-2.75	1.34	1.39
2	A	1401	ADP	C5-N7	-2.72	1.34	1.39
2	B	1401	ADP	C5-N7	-2.63	1.34	1.39
2	D	1401	ADP	C5-N7	-2.58	1.34	1.39
2	E	1401	ADP	C5-C6	2.57	1.48	1.41
2	E	1401	ADP	C5-N7	-2.53	1.34	1.39
2	D	1401	ADP	C5-C6	2.46	1.47	1.41
2	A	1401	ADP	C5-C6	2.43	1.47	1.41
2	B	1401	ADP	C5-C6	2.31	1.47	1.41
2	C	1401	ADP	C5-C6	2.18	1.47	1.41
2	D	1401	ADP	C8-N7	2.16	1.35	1.31
2	B	1401	ADP	C4-N9	-2.13	1.33	1.37
2	C	1401	ADP	C4-N9	-2.11	1.33	1.37
2	E	1401	ADP	C8-N7	2.09	1.35	1.31
2	A	1401	ADP	C8-N7	2.05	1.35	1.31
2	E	1401	ADP	PA-O3A	2.02	1.61	1.59

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1401	ADP	C5-C4-N3	-6.30	118.04	126.72
2	E	1401	ADP	C5-C4-N3	-6.18	118.21	126.72
2	D	1401	ADP	C5-C4-N3	-5.94	118.54	126.72
2	B	1401	ADP	C5-C4-N3	-5.85	118.66	126.72
2	C	1401	ADP	C5-C4-N3	-5.76	118.78	126.72
2	A	1401	ADP	N3-C4-N9	5.08	135.81	127.17
2	E	1401	ADP	N3-C4-N9	4.96	135.61	127.17
2	C	1401	ADP	N3-C4-N9	4.94	135.56	127.17
2	B	1401	ADP	N3-C4-N9	4.73	135.22	127.17
2	D	1401	ADP	N3-C4-N9	4.70	135.16	127.17
2	A	1401	ADP	C2-N3-C4	3.76	121.03	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1401	ADP	C2-N3-C4	3.71	120.89	111.83
2	E	1401	ADP	C2-N3-C4	3.66	120.76	111.83
2	B	1401	ADP	C2-N3-C4	3.64	120.72	111.83
2	C	1401	ADP	C2-N3-C4	3.58	120.58	111.83
2	E	1401	ADP	C4-C5-N7	-3.48	106.60	110.58
2	B	1401	ADP	C4-C5-N7	-3.36	106.75	110.58
2	C	1401	ADP	N3-C2-N1	-3.33	123.54	128.58
2	D	1401	ADP	C4-C5-N7	-3.32	106.78	110.58
2	D	1401	ADP	N3-C2-N1	-3.26	123.65	128.58
2	A	1401	ADP	C4-C5-N7	-3.23	106.88	110.58
2	B	1401	ADP	N3-C2-N1	-3.17	123.78	128.58
2	A	1401	ADP	N3-C2-N1	-3.13	123.85	128.58
2	E	1401	ADP	N3-C2-N1	-2.92	124.16	128.58
2	C	1401	ADP	C4-N9-C8	2.91	108.80	105.74
2	C	1401	ADP	C4-C5-N7	-2.88	107.28	110.58
2	B	1401	ADP	C4-N9-C8	2.81	108.69	105.74
2	D	1401	ADP	C3'-C2'-C1'	2.74	106.65	101.46
2	A	1401	ADP	C3'-C2'-C1'	2.74	106.64	101.46
2	B	1401	ADP	C3'-C2'-C1'	2.72	106.61	101.46
2	E	1401	ADP	C3'-C2'-C1'	2.65	106.47	101.46
2	E	1401	ADP	C5-N7-C8	2.63	107.59	103.45
2	C	1401	ADP	C3'-C2'-C1'	2.58	106.34	101.46
2	E	1401	ADP	C4-N9-C8	2.55	108.42	105.74
2	B	1401	ADP	C5-N7-C8	2.52	107.42	103.45
2	D	1401	ADP	C4-N9-C8	2.46	108.32	105.74
2	D	1401	ADP	C5-N7-C8	2.42	107.26	103.45
2	A	1401	ADP	C5-N7-C8	2.42	107.25	103.45
2	A	1401	ADP	C4-N9-C8	2.37	108.23	105.74
2	C	1401	ADP	C5-N7-C8	2.31	107.09	103.45
2	B	1401	ADP	C2'-C1'-N9	-2.22	107.80	113.30
2	B	1401	ADP	N9-C8-N7	-2.04	111.04	113.94

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1401	ADP	C5'-O5'-PA-O1A
2	A	1401	ADP	C5'-O5'-PA-O2A
2	A	1401	ADP	C5'-O5'-PA-O3A
2	B	1401	ADP	C5'-O5'-PA-O3A
2	C	1401	ADP	C5'-O5'-PA-O2A
2	C	1401	ADP	C5'-O5'-PA-O3A

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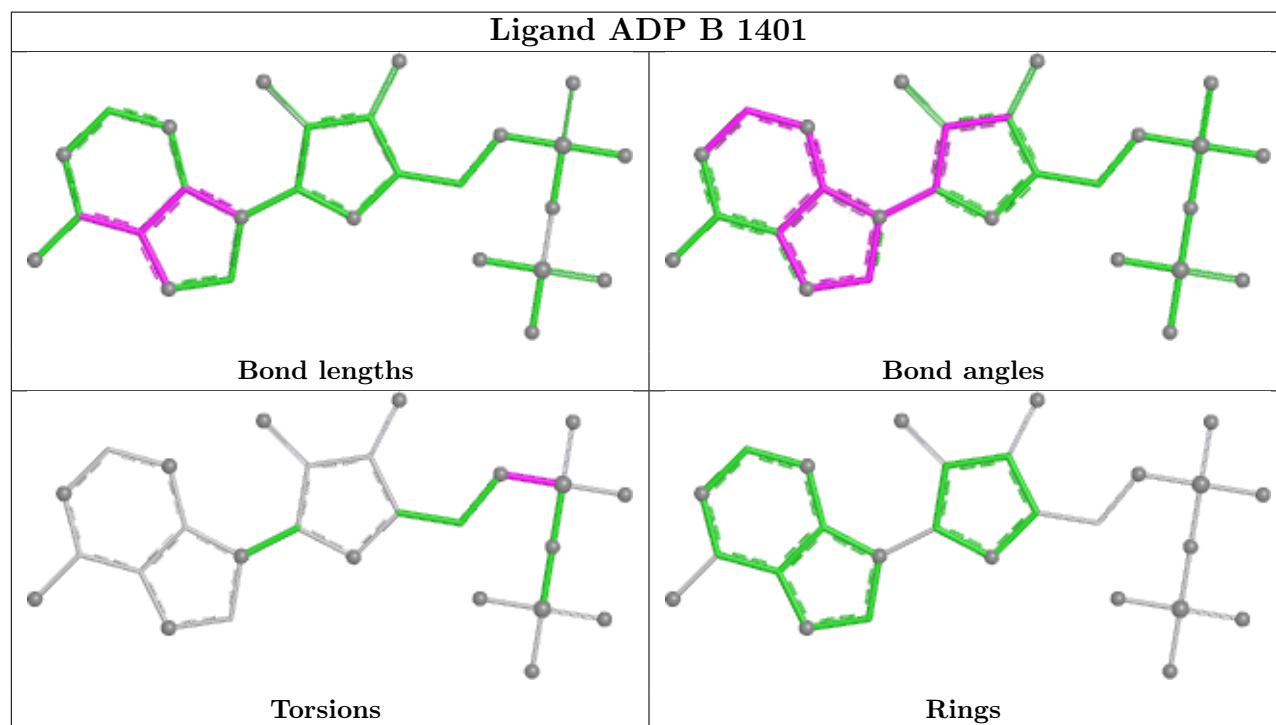
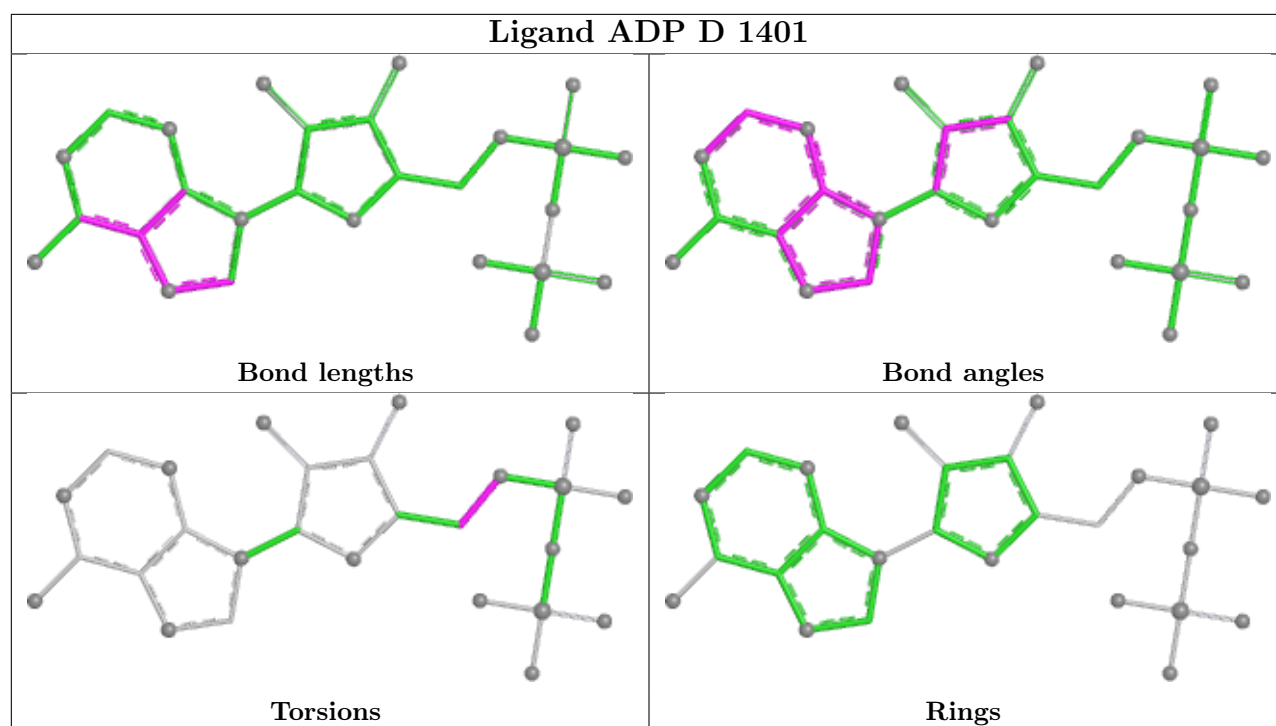
Mol	Chain	Res	Type	Atoms
2	E	1401	ADP	C5'-O5'-PA-O1A
2	E	1401	ADP	C5'-O5'-PA-O2A
2	E	1401	ADP	C5'-O5'-PA-O3A
2	C	1401	ADP	C3'-C4'-C5'-O5'
2	C	1401	ADP	O4'-C4'-C5'-O5'
2	B	1401	ADP	C5'-O5'-PA-O1A
2	D	1401	ADP	C4'-C5'-O5'-PA
2	A	1401	ADP	C3'-C4'-C5'-O5'

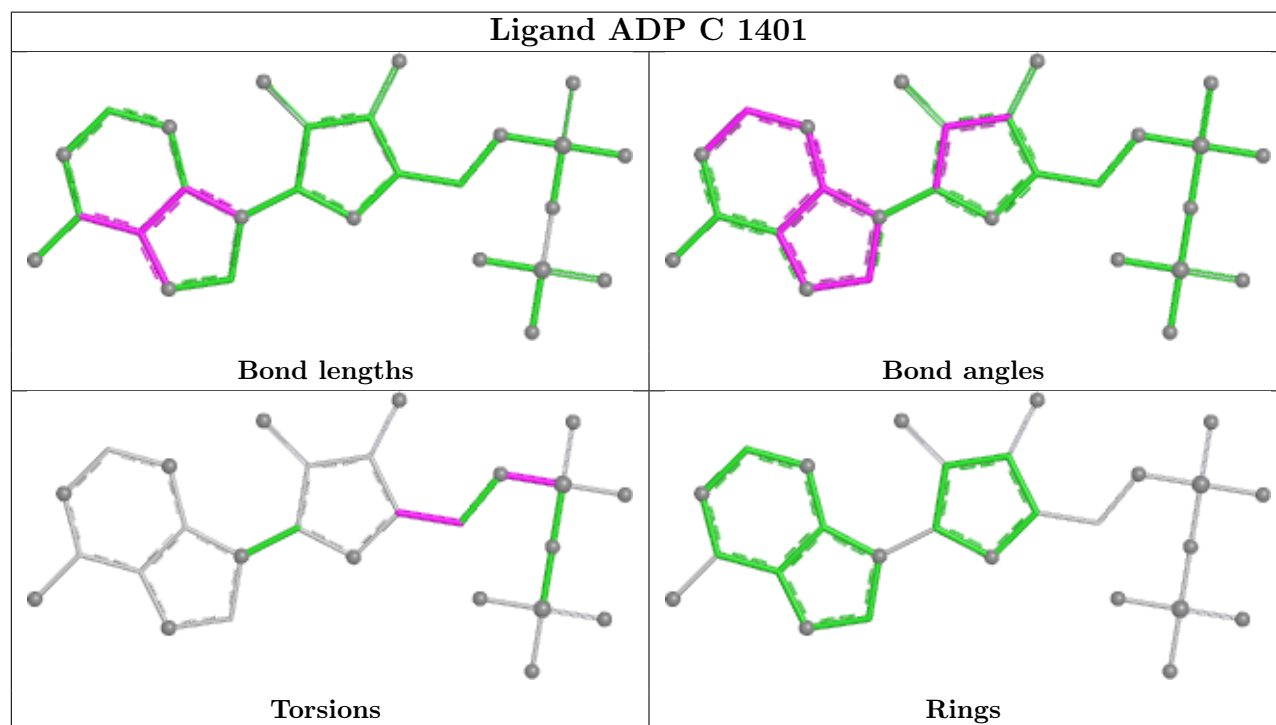
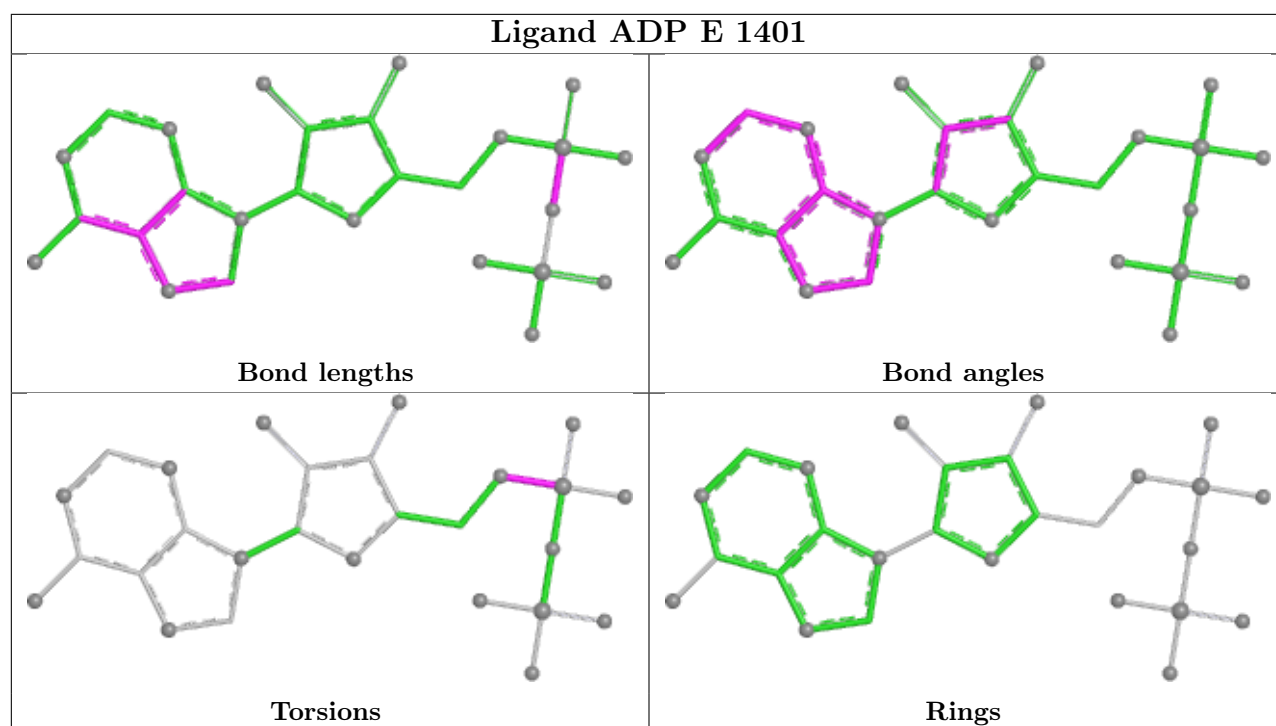
There are no ring outliers.

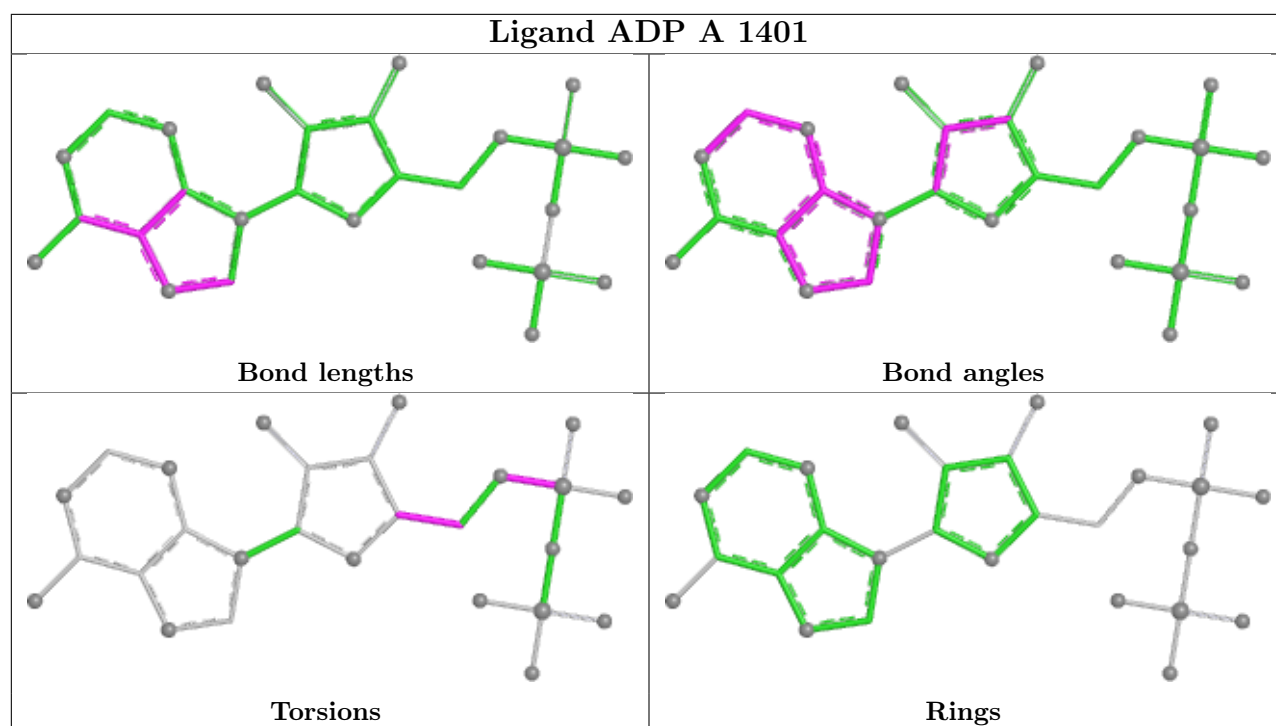
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1401	ADP	1	0
2	B	1401	ADP	1	0
2	E	1401	ADP	2	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.

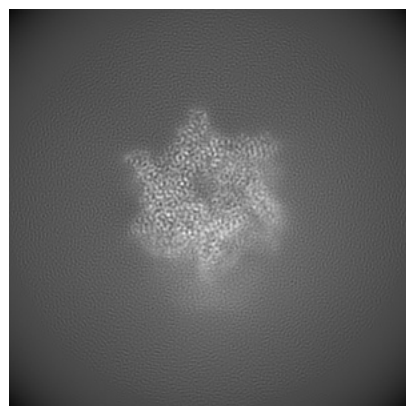
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26697. These allow visual inspection of the internal detail of the map and identification of artifacts.

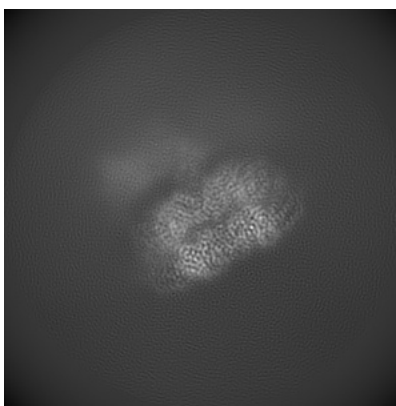
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

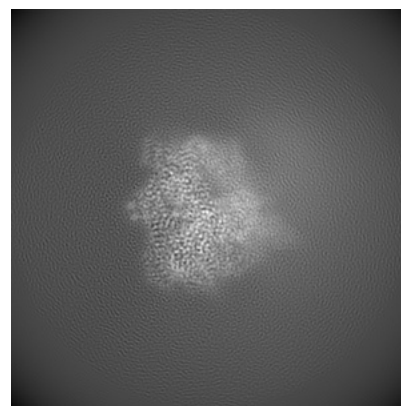
6.1.1 Primary map



X

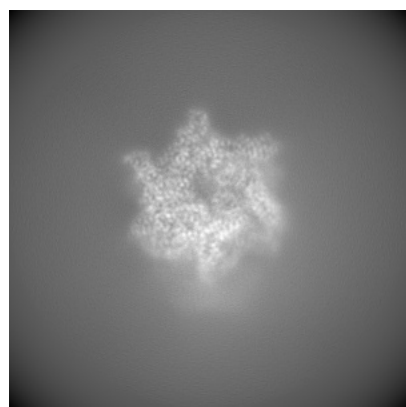


Y

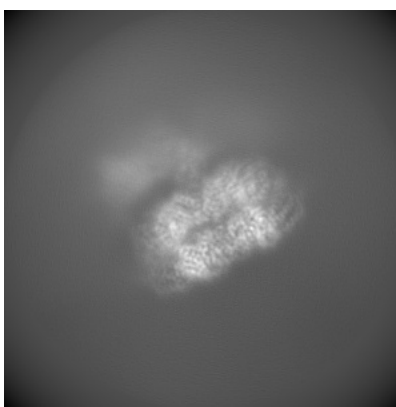


Z

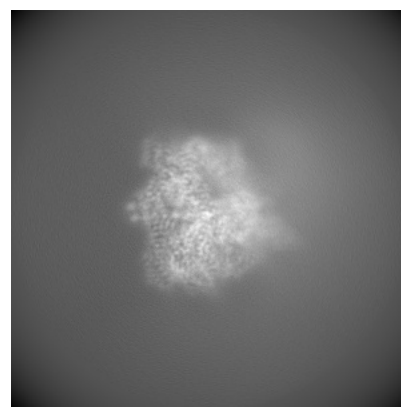
6.1.2 Raw map



X



Y

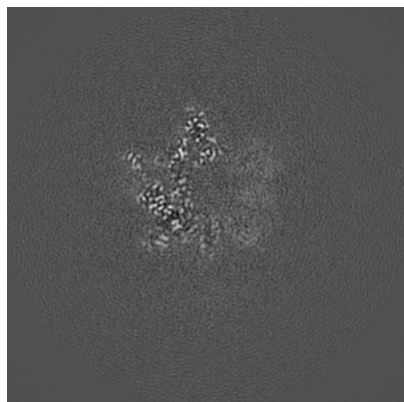


Z

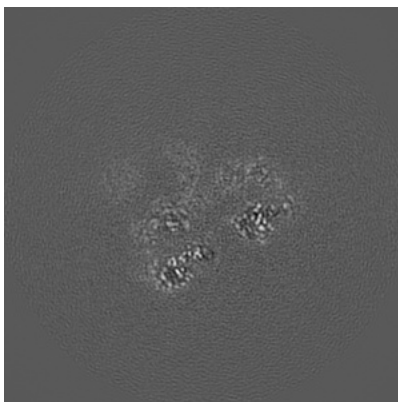
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

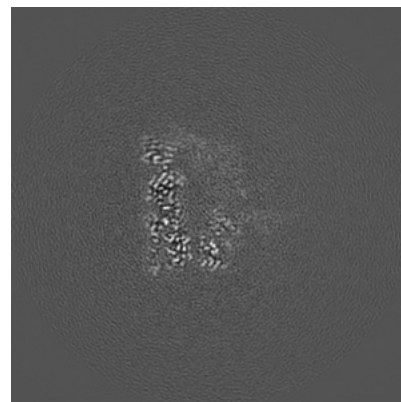
6.2.1 Primary map



X Index: 200

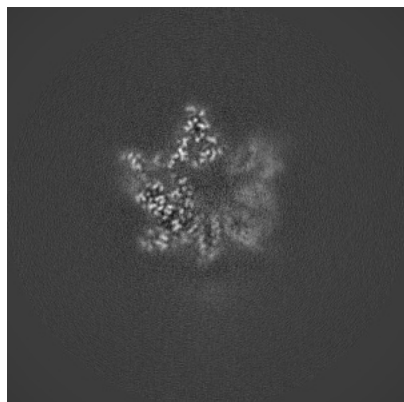


Y Index: 200

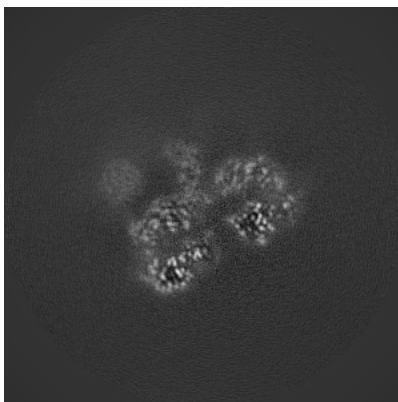


Z Index: 200

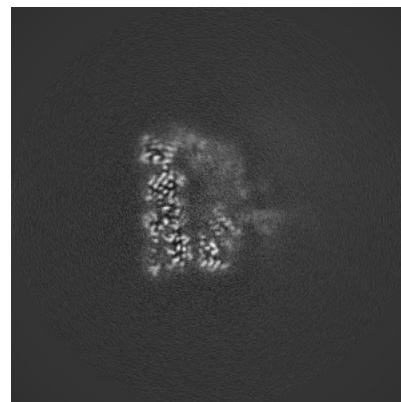
6.2.2 Raw map



X Index: 200



Y Index: 200

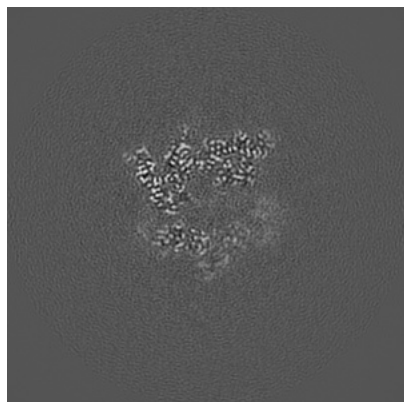


Z Index: 200

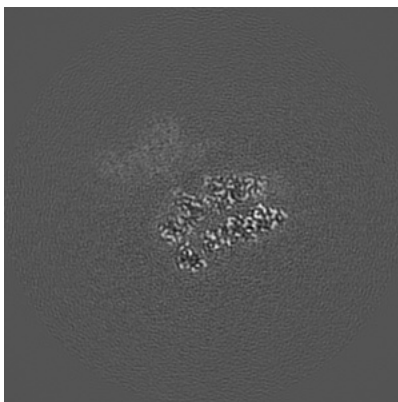
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

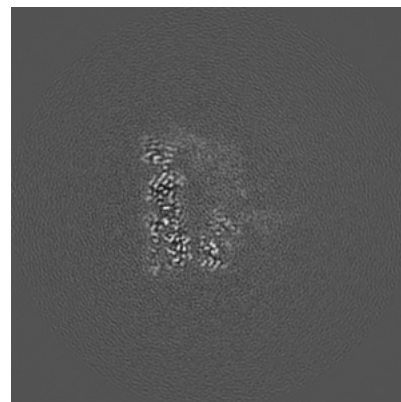
6.3.1 Primary map



X Index: 176

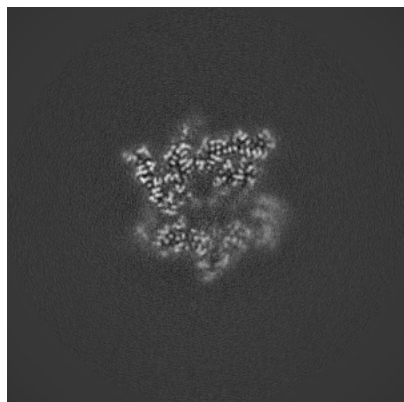


Y Index: 173

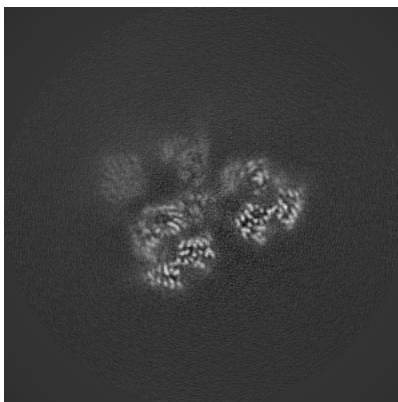


Z Index: 200

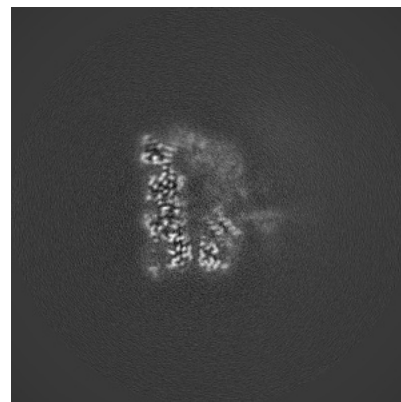
6.3.2 Raw map



X Index: 176



Y Index: 194

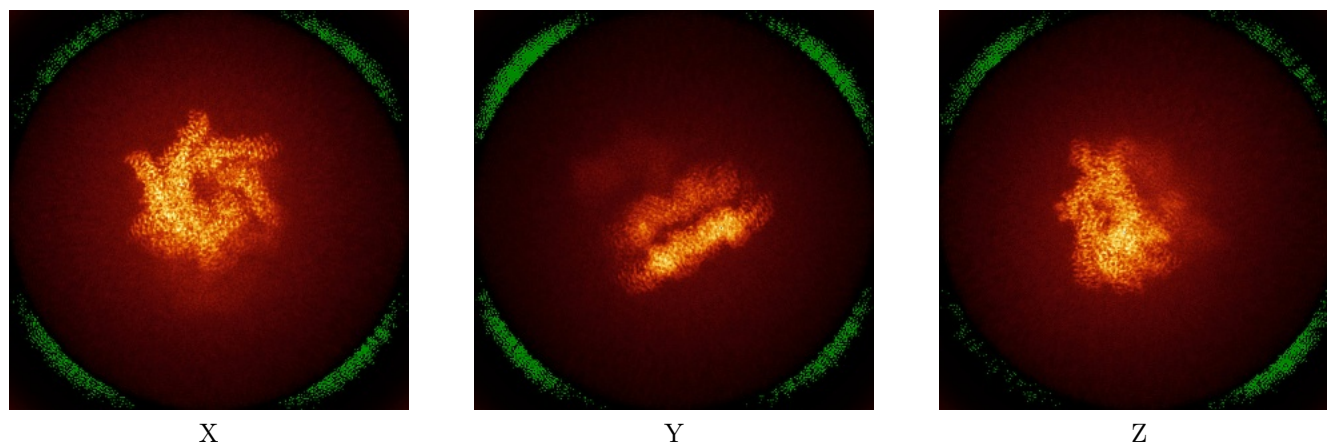


Z Index: 201

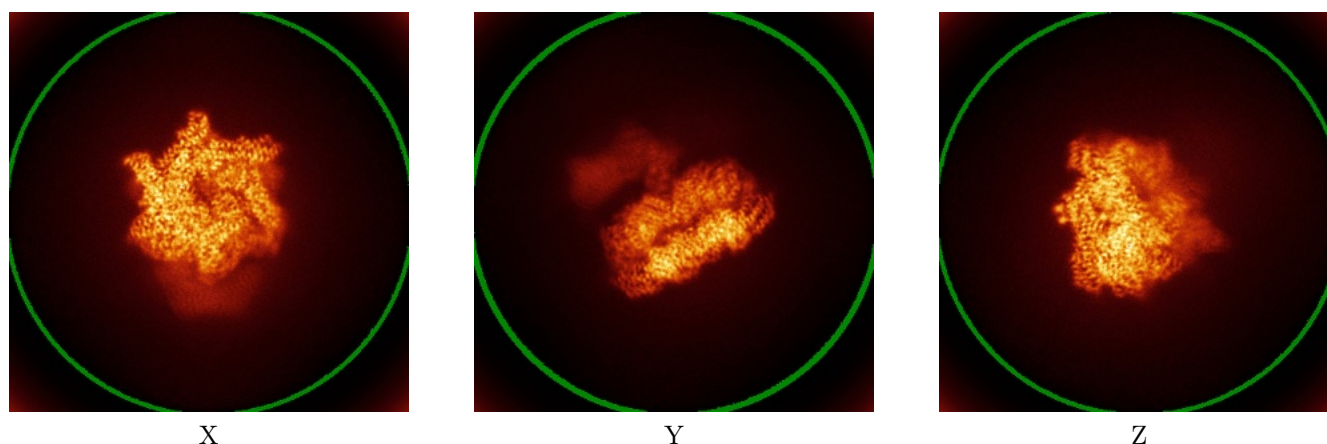
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

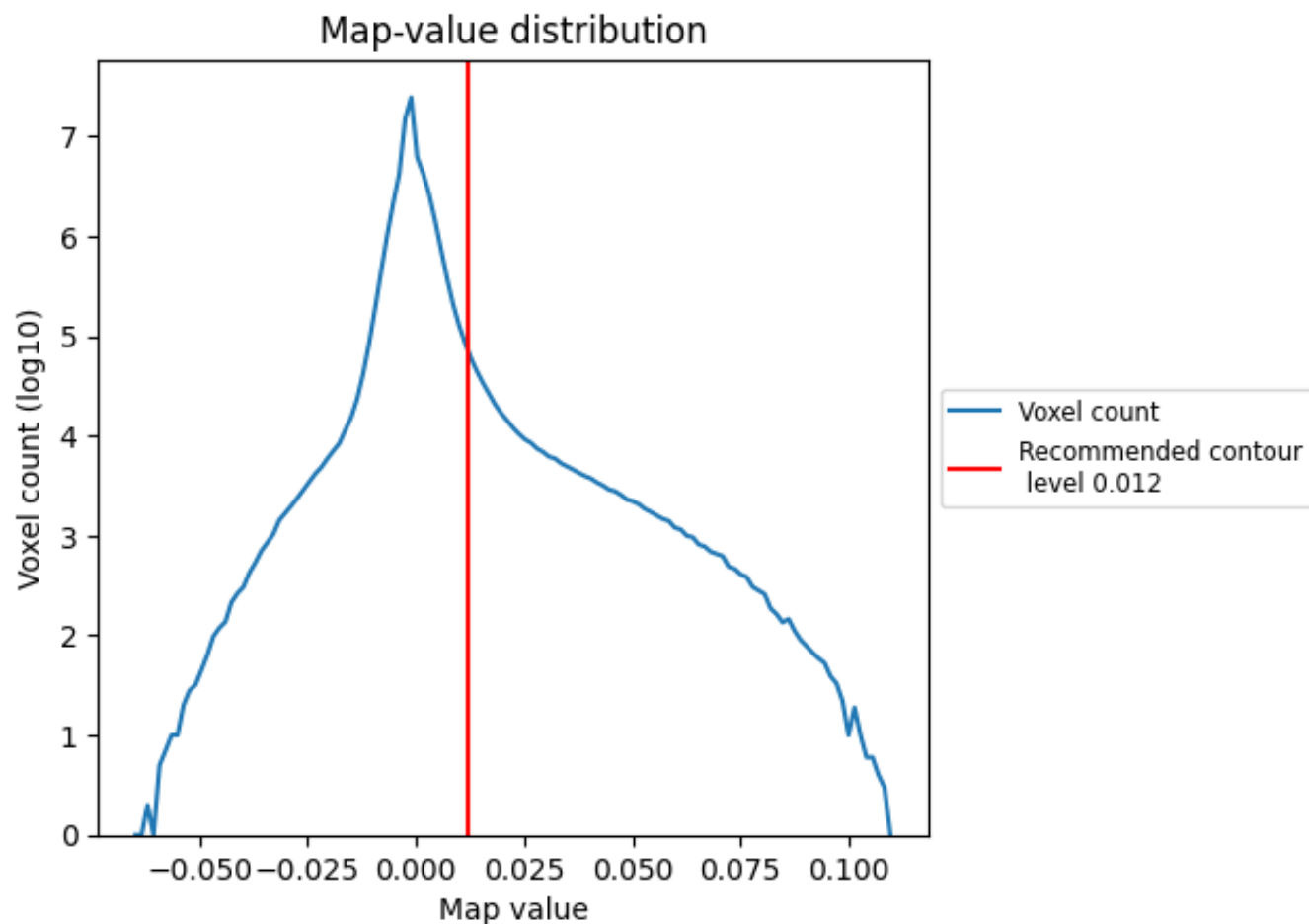
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

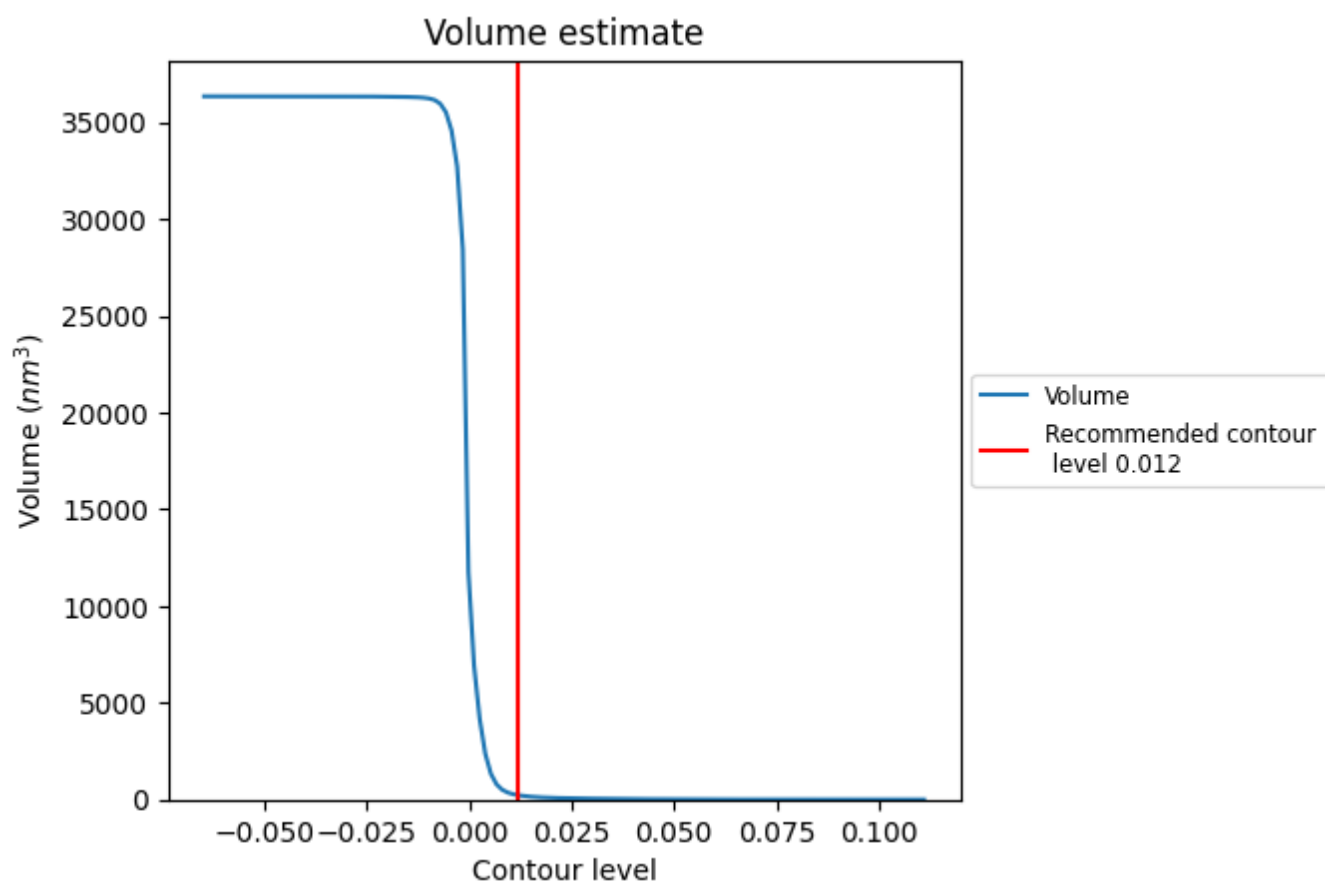
This section contains the results of statistical analysis of the map.

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.

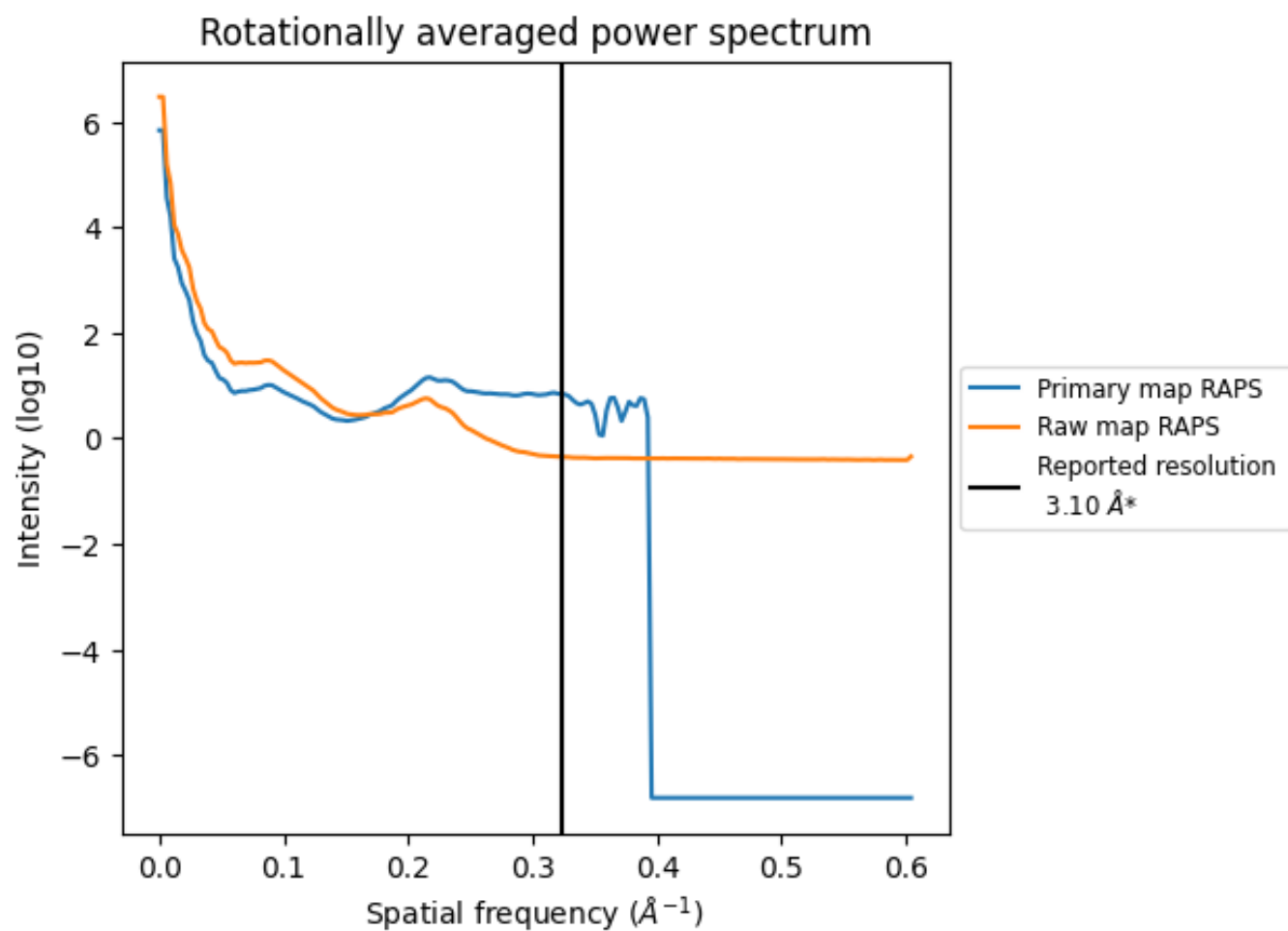
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 224 nm³; this corresponds to an approximate mass of 202 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 ⓘ

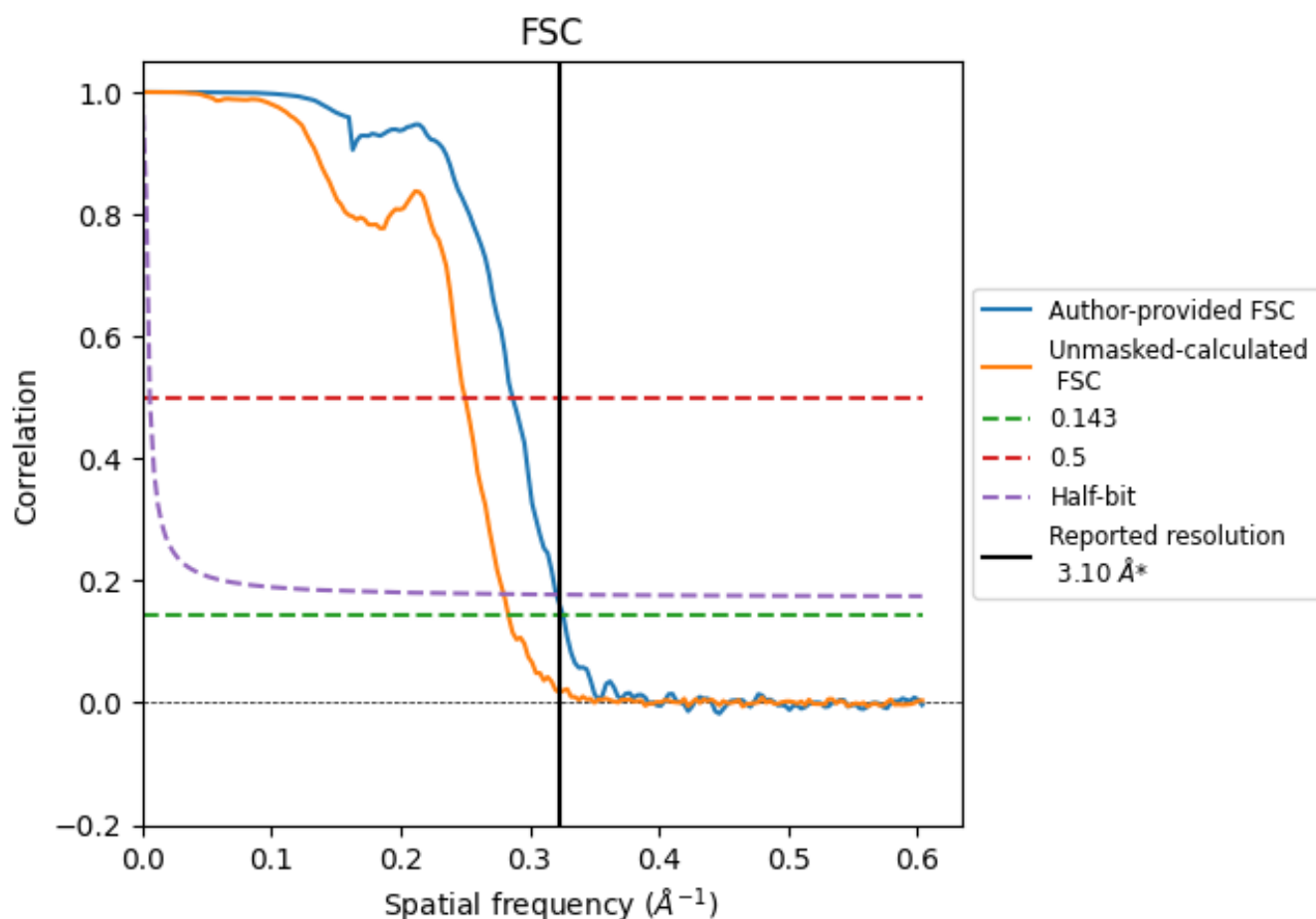


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 0.323 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

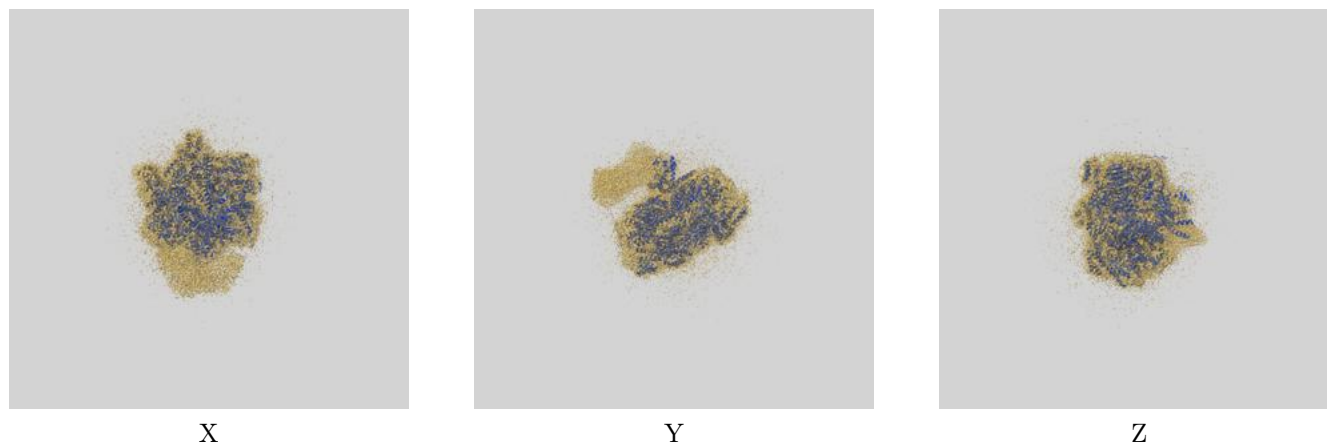
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.07	3.49	3.12
Unmasked-calculated*	3.53	4.00	3.57

*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 3.53 differs from the reported value 3.1 by more than 10 %

9 Map-model fit [i](#)

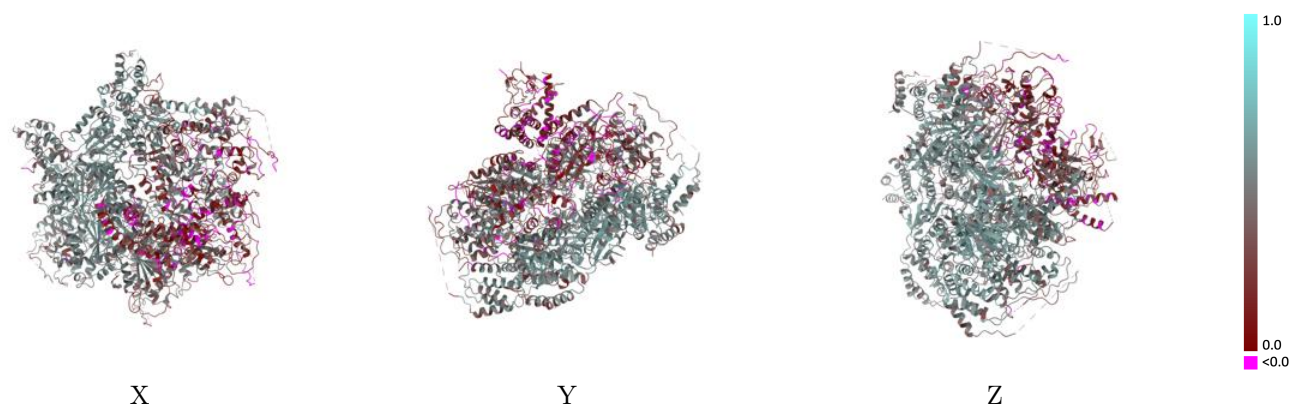
This section contains information regarding the fit between EMDB map EMD-26697 and PDB model 7UQK. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



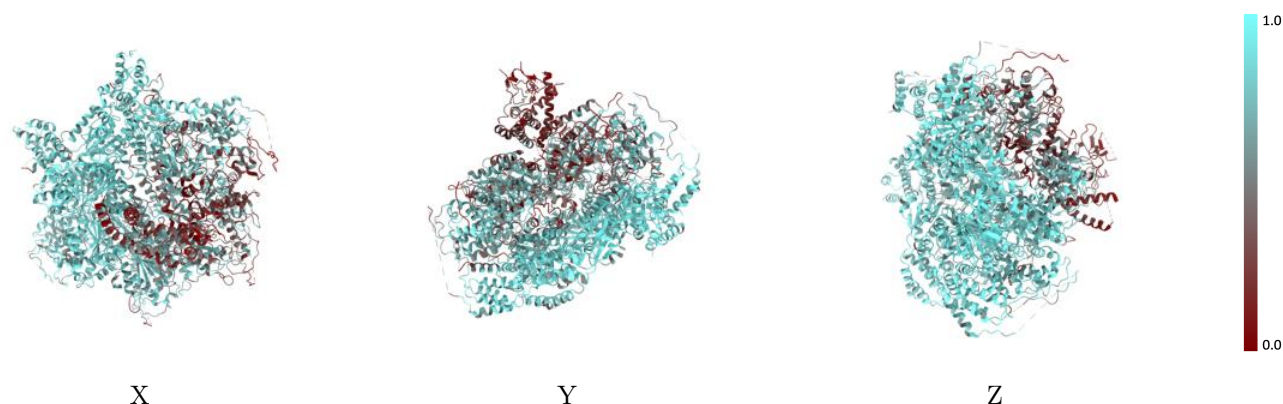
The images above show the 3D surface view of the map at the recommended contour level 0.012 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



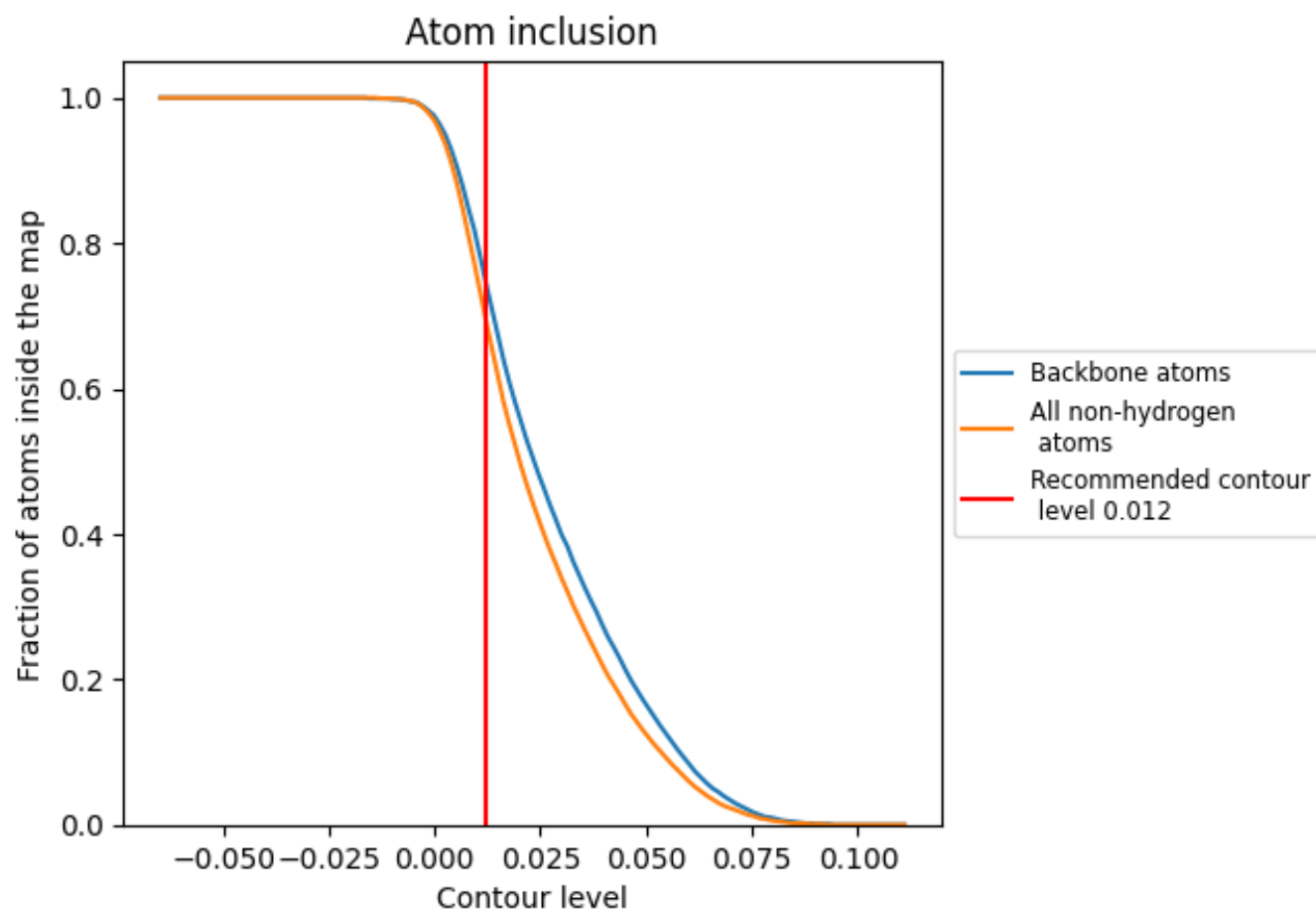
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



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 (0.012).

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 70% 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 (0.012) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6970	0.4170
A	0.8140	0.5040
B	0.8310	0.5080
C	0.7950	0.4810
D	0.6960	0.4080
E	0.5480	0.3120
F	0.6530	0.3760
G	0.2390	0.1540

1.0

0.0

<0.0