



wwPDB EM Validation Summary Report ⓘ

Mar 24, 2026 – 07:54 AM UTC

PDB ID : 8EUF / pdb_00008euf
EMDB ID : EMD-28613
Title : Class2 of the INO80-Nucleosome complex
Authors : Wu, H.; Munoz, E.; Gourdet, M.; Narlikar, G.; Cheng, Y.F.
Deposited on : 2022-10-18
Resolution : 3.41 Å(reported)

This is a wwPDB EM Validation Summary 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:

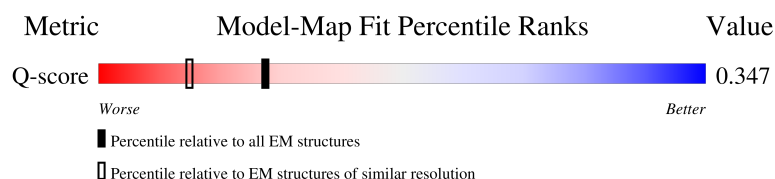
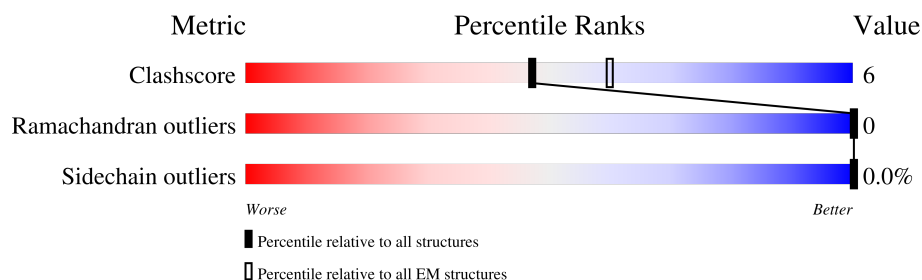
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.41 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.



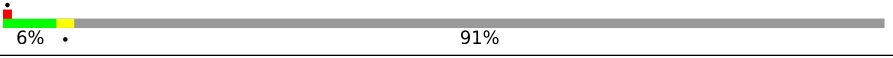
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13997 (2.91 - 3.91)

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	Q	1489	
2	R	755	
3	S	166	
4	T	463	

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Mol	Chain	Length	Quality of chain
4	V	463	 5% 84% 12% 5%
4	X	463	 9% 83% 13% 5%
5	U	460	 8% 83% 12% 5%
5	W	460	 10% 83% 13% 5%
5	Y	460	 7% 83% 13% 5%
6	Z	320	 6% 91% 6% 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 28707 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 Chromatin-remodeling ATPase INO80.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	Q	421	Total	C	N	O	S	0	0
			3409	2167	575	652	15		

- Molecule 2 is a protein called Actin-related protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	R	442	Total	C	N	O	S	0	0
			3536	2252	595	678	11		

- Molecule 3 is a protein called Chromatin-remodeling complex subunit IES6.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	S	128	Total	C	N	O	S	0	0
			1025	658	188	177	2		

- Molecule 4 is a protein called RuvB-like protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	T	437	Total	C	N	O	S	0	0
			3353	2113	578	652	10		
4	V	443	Total	C	N	O	S	0	0
			3404	2149	585	660	10		
4	X	442	Total	C	N	O	S	0	0
			3397	2144	584	659	10		

- Molecule 5 is a protein called RuvB-like protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	U	436	Total	C	N	O	S	0	0
			3354	2100	583	660	11		
5	W	442	Total	C	N	O	S	0	0
			3398	2123	590	673	12		

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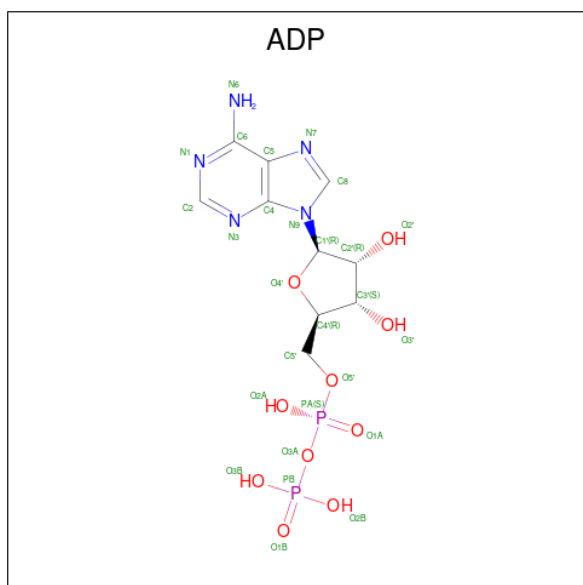
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Mol	Chain	Residues	Atoms					AltConf	Trace
5	Y	445	Total	C	N	O	S	0	0
			3421	2138	594	677	12		

- Molecule 6 is a protein called Ino eighty subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	Z	28	Total	C	N	O	S	0	0
			248	158	45	43	2		

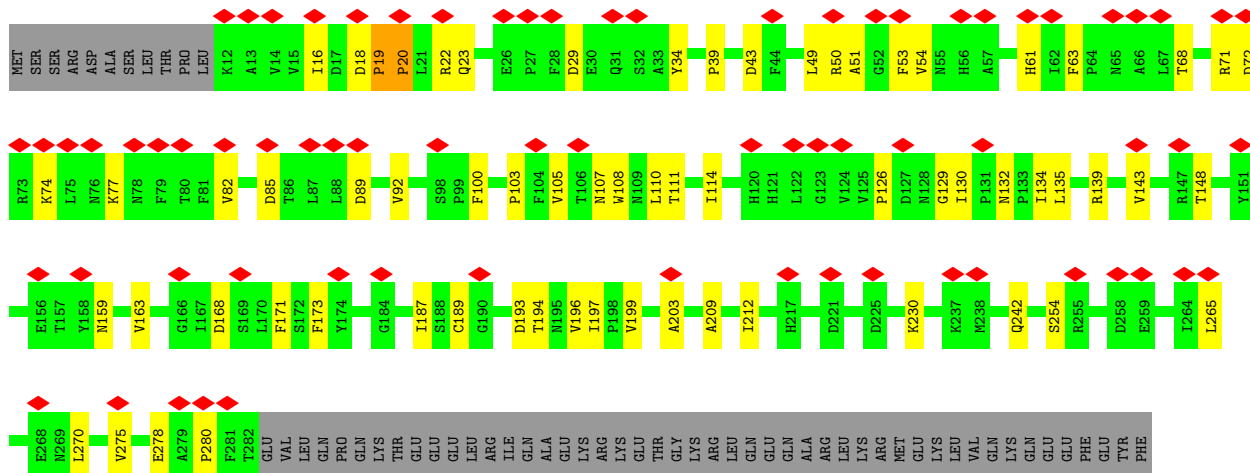
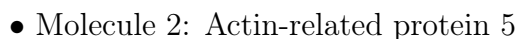
- Molecule 7 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).

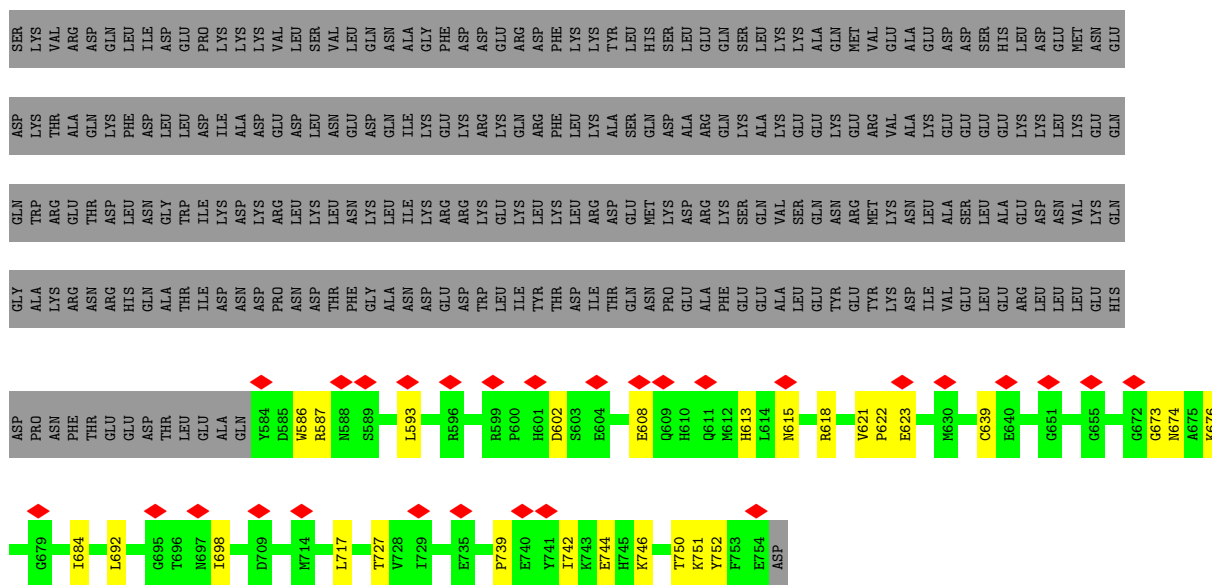


These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

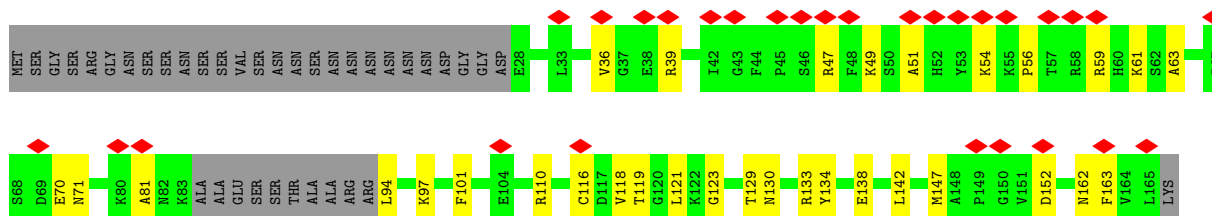
Chain Q: 11% 23% 5% 72%

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

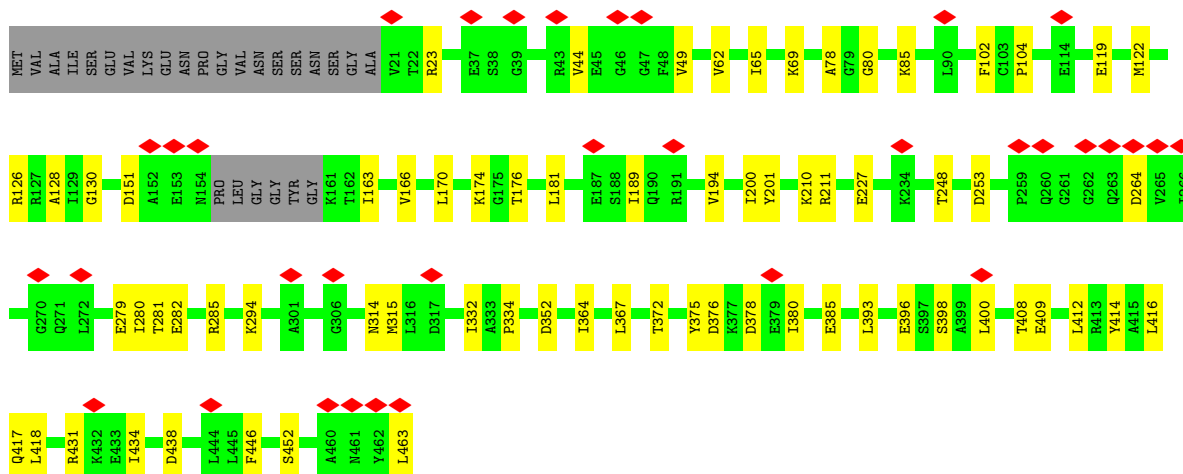
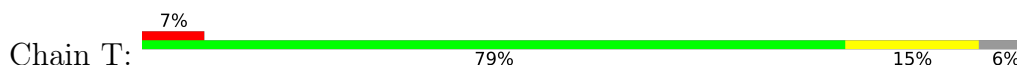




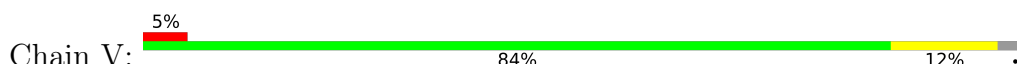
- Molecule 3: Chromatin-remodeling complex subunit IES6

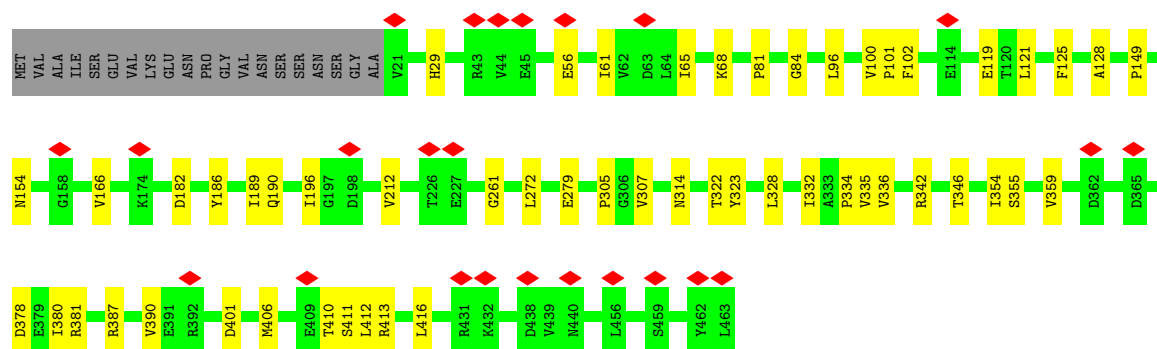


- Molecule 4: RuvB-like protein 1

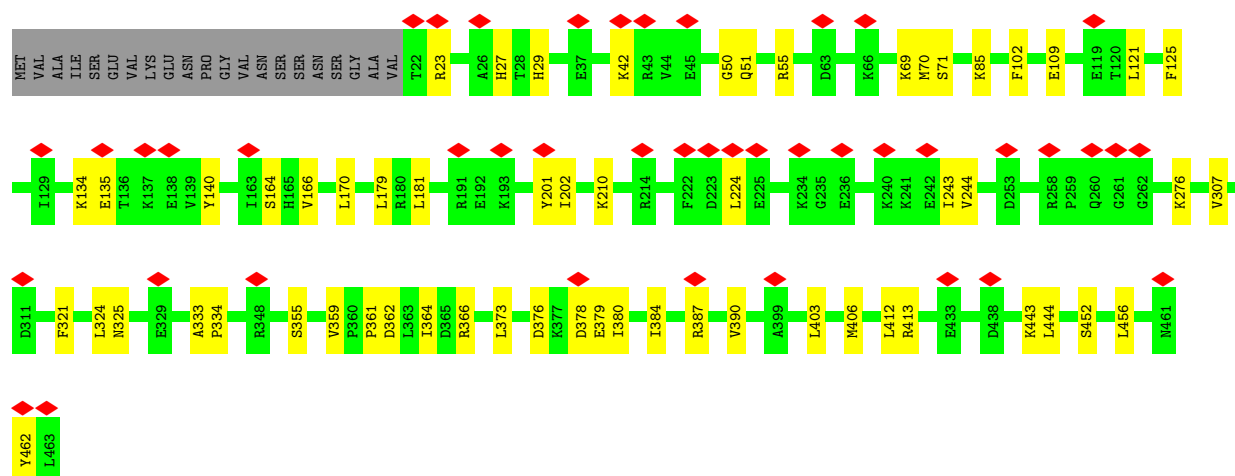
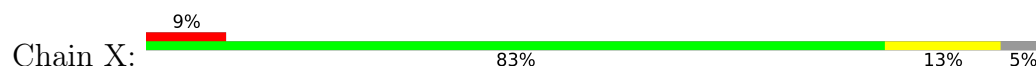


- Molecule 4: RuvB-like protein 1

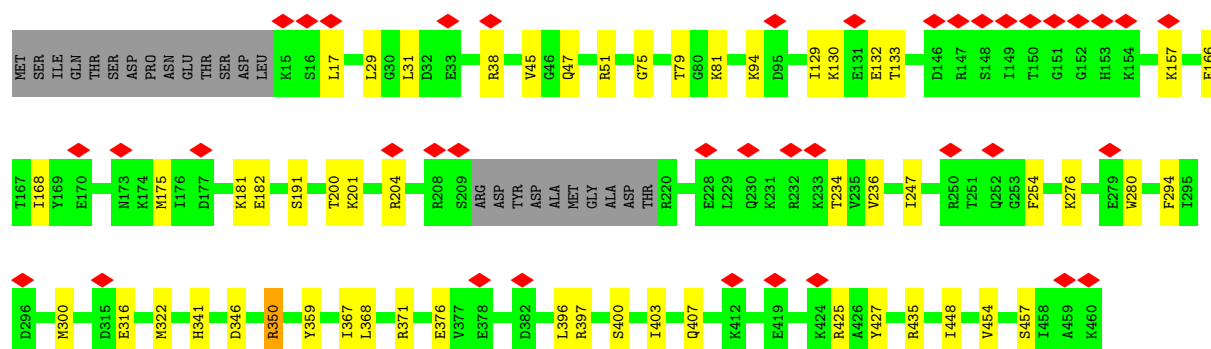
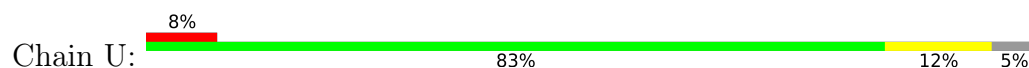




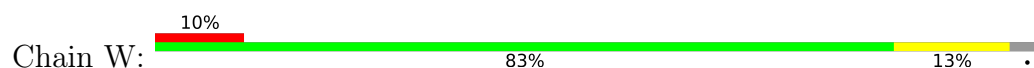
• Molecule 4: RuvB-like protein 1



• Molecule 5: RuvB-like protein 2



• Molecule 5: RuvB-like protein 2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	44535	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	43	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.782	Depositor
Minimum map value	-0.024	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.024	Depositor
Recommended contour level	0.174	Depositor
Map size ($\text{\AA}$)	400.8, 400.8, 400.8	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.835, 0.835, 0.835	Depositor

5 Model quality

5.1 Standard geometry

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	Q	0.28	0/3475	0.65	3/4711 (0.1%)
2	R	0.30	1/3627 (0.0%)	0.67	4/4934 (0.1%)
3	S	0.31	0/1047	0.70	1/1408 (0.1%)
4	T	0.27	0/3392	0.54	0/4588
4	V	0.27	0/3446	0.50	0/4662
4	X	0.25	0/3439	0.59	4/4652 (0.1%)
5	U	0.25	0/3390	0.52	1/4566 (0.0%)
5	W	0.24	0/3436	0.51	1/4632 (0.0%)
5	Y	0.25	0/3459	0.50	0/4662
6	Z	0.31	0/254	0.71	1/338 (0.3%)
All	All	0.27	1/28965 (0.0%)	0.57	15/39153 (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
5	U	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	R	20	PRO	N-CD	-6.73	1.38	1.47

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	1182	PRO	CA-N-CD	-8.32	100.35	112.00
2	R	20	PRO	CA-N-CD	7.92	123.09	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	1014	PRO	CA-N-CD	-7.85	101.01	112.00
2	R	20	PRO	N-CA-CB	-7.56	95.82	103.15
3	S	56	PRO	CA-N-CD	-6.93	102.29	112.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	U	350	ARG	Sidechain

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	Q	3409	0	3389	55	0
2	R	3536	0	3416	67	0
3	S	1025	0	1055	24	0
4	T	3353	0	3471	46	0
4	V	3404	0	3542	41	0
4	X	3397	0	3533	37	0
5	U	3354	0	3448	42	0
5	W	3398	0	3470	41	0
5	Y	3421	0	3499	39	0
6	Z	248	0	242	5	0
7	T	27	0	12	1	0
7	U	27	0	12	4	0
7	V	27	0	12	2	0
7	W	27	0	12	1	0
7	X	27	0	12	1	0
7	Y	27	0	12	1	0
All	All	28707	0	29137	348	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.

The worst 5 of 348 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:18:ASP:C	2:R:20:PRO:HD2	2.07	0.79
2:R:19:PRO:N	2:R:20:PRO:CD	2.45	0.79
4:T:44:VAL:HG22	4:T:49:VAL:HG23	1.65	0.78
2:R:265:LEU:HD12	3:S:121:LEU:HD11	1.74	0.70
5:Y:325:THR:HG22	5:Y:327:ARG:H	1.59	0.67

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	Q	409/1489 (28%)	390 (95%)	19 (5%)	0	100	100
2	R	438/755 (58%)	418 (95%)	20 (5%)	0	100	100
3	S	124/166 (75%)	114 (92%)	10 (8%)	0	100	100
4	T	433/463 (94%)	417 (96%)	16 (4%)	0	100	100
4	V	441/463 (95%)	428 (97%)	13 (3%)	0	100	100
4	X	440/463 (95%)	420 (96%)	20 (4%)	0	100	100
5	U	432/460 (94%)	415 (96%)	17 (4%)	0	100	100
5	W	440/460 (96%)	426 (97%)	14 (3%)	0	100	100
5	Y	443/460 (96%)	428 (97%)	15 (3%)	0	100	100
6	Z	26/320 (8%)	26 (100%)	0	0	100	100
All	All	3626/5499 (66%)	3482 (96%)	144 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	Q	392/1350 (29%)	391 (100%)	1 (0%)	86	84
2	R	391/682 (57%)	391 (100%)	0	100	100
3	S	108/142 (76%)	108 (100%)	0	100	100
4	T	368/391 (94%)	368 (100%)	0	100	100
4	V	375/391 (96%)	375 (100%)	0	100	100
4	X	374/391 (96%)	374 (100%)	0	100	100
5	U	372/394 (94%)	372 (100%)	0	100	100
5	W	375/394 (95%)	375 (100%)	0	100	100
5	Y	378/394 (96%)	378 (100%)	0	100	100
6	Z	27/285 (10%)	27 (100%)	0	100	100
All	All	3160/4814 (66%)	3159 (100%)	1 (0%)	100	100

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

Mol	Chain	Res	Type
1	Q	1432	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
5	U	408	GLN
5	W	63	ASN
4	V	94	GLN
5	W	415	ASN
2	R	90	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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
7	ADP	Y	501	-	28,29,29	1.38	4 (14%)	43,45,45	1.82	8 (18%)
7	ADP	W	501	-	28,29,29	1.39	5 (17%)	43,45,45	1.81	7 (16%)
7	ADP	V	501	-	28,29,29	1.37	5 (17%)	43,45,45	1.78	7 (16%)
7	ADP	U	501	-	28,29,29	1.38	5 (17%)	43,45,45	1.84	8 (18%)
7	ADP	X	501	-	28,29,29	1.33	5 (17%)	43,45,45	1.85	8 (18%)
7	ADP	T	501	-	28,29,29	1.36	6 (21%)	43,45,45	1.92	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
7	ADP	Y	501	-	-	4/16/32/32	0/3/3/3
7	ADP	W	501	-	-	2/16/32/32	0/3/3/3
7	ADP	V	501	-	-	0/16/32/32	0/3/3/3
7	ADP	U	501	-	-	3/16/32/32	0/3/3/3
7	ADP	X	501	-	-	3/16/32/32	0/3/3/3
7	ADP	T	501	-	-	3/16/32/32	0/3/3/3

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	Y	501	ADP	C5-C4	4.53	1.47	1.39

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	W	501	ADP	C5-C4	4.49	1.47	1.39
7	U	501	ADP	C5-C4	4.43	1.47	1.39
7	V	501	ADP	C5-C4	4.36	1.46	1.39
7	X	501	ADP	C5-C4	4.17	1.46	1.39

The worst 5 of 48 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	V	501	ADP	C5-C4-N3	-5.98	118.48	126.72
7	Y	501	ADP	C5-C4-N3	-5.96	118.50	126.72
7	X	501	ADP	C5-C4-N3	-5.87	118.64	126.72
7	U	501	ADP	C5-C4-N3	-5.85	118.66	126.72
7	T	501	ADP	C5-C4-N3	-5.78	118.75	126.72

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	T	501	ADP	C5'-O5'-PA-O1A
7	T	501	ADP	C5'-O5'-PA-O2A
7	T	501	ADP	C5'-O5'-PA-O3A
7	U	501	ADP	C5'-O5'-PA-O1A
7	U	501	ADP	C5'-O5'-PA-O2A

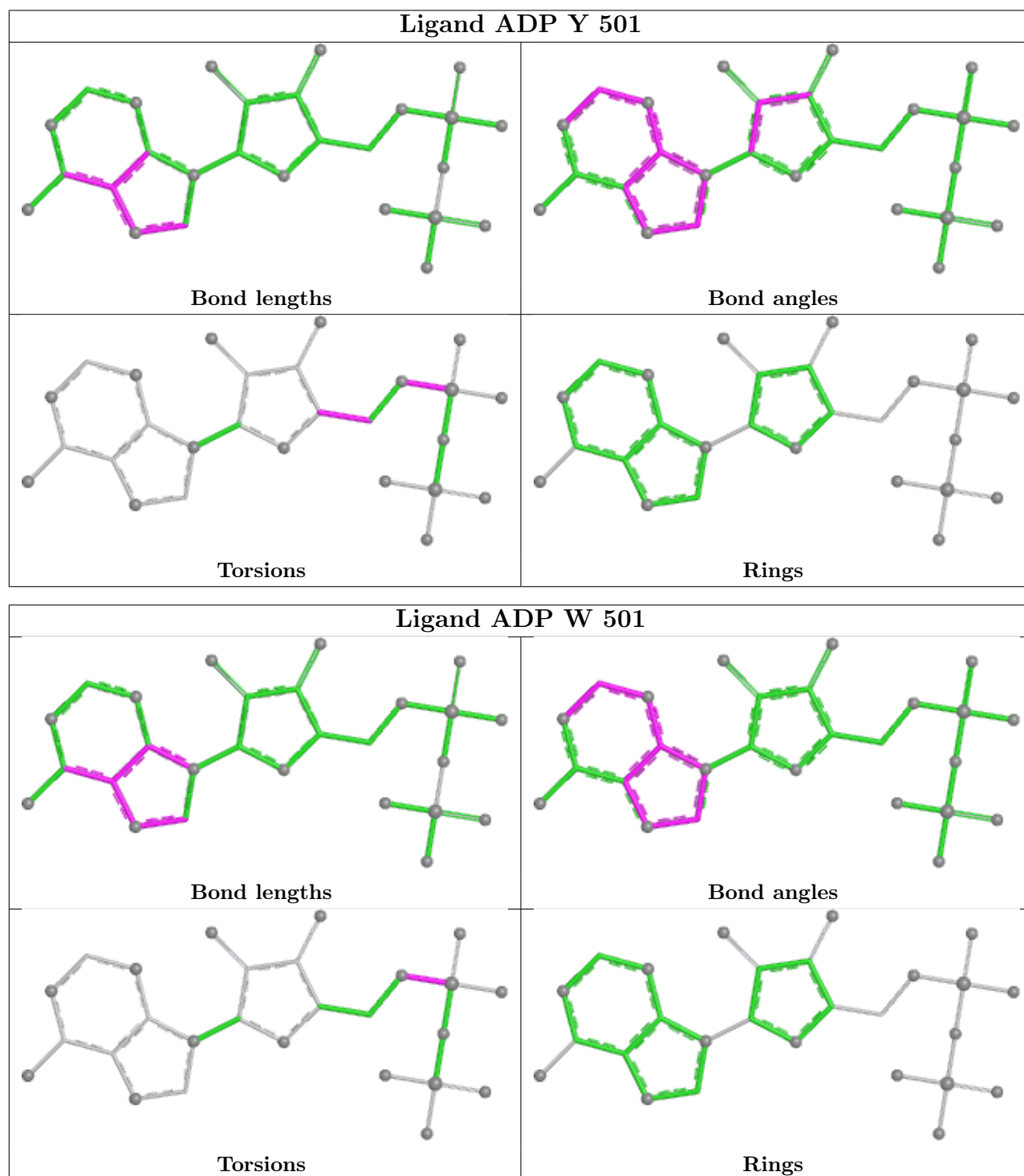
There are no ring outliers.

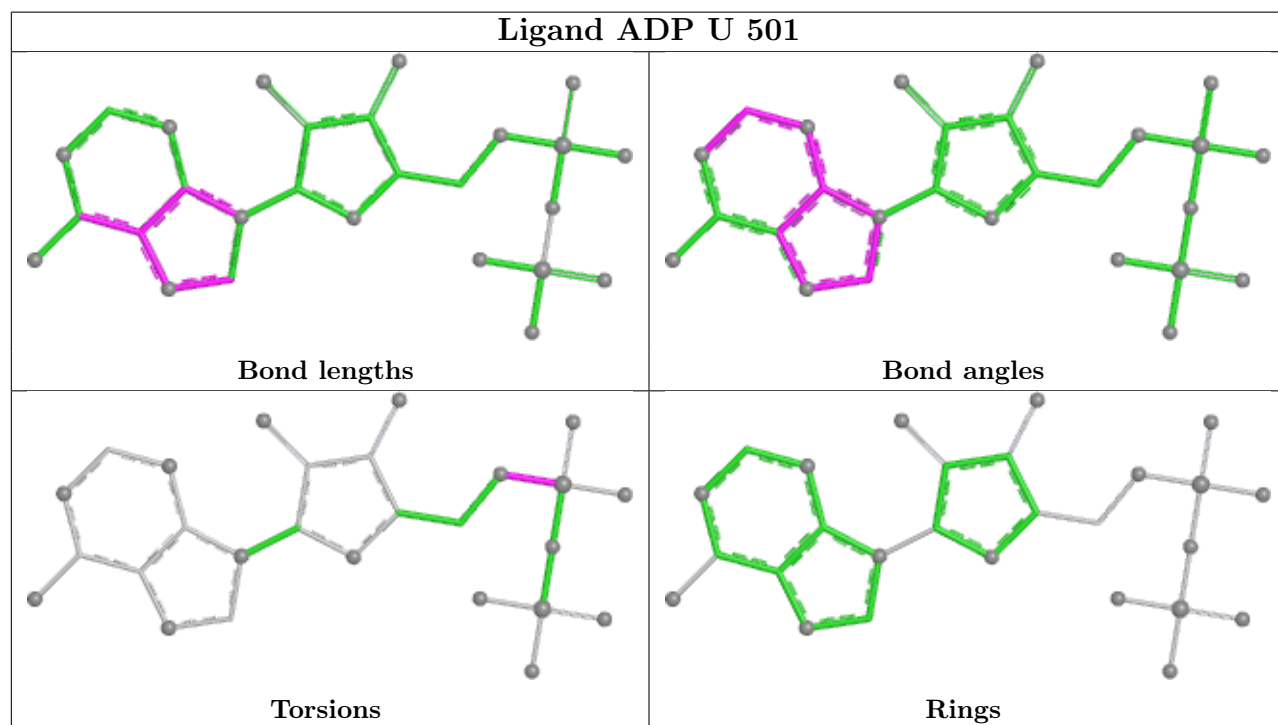
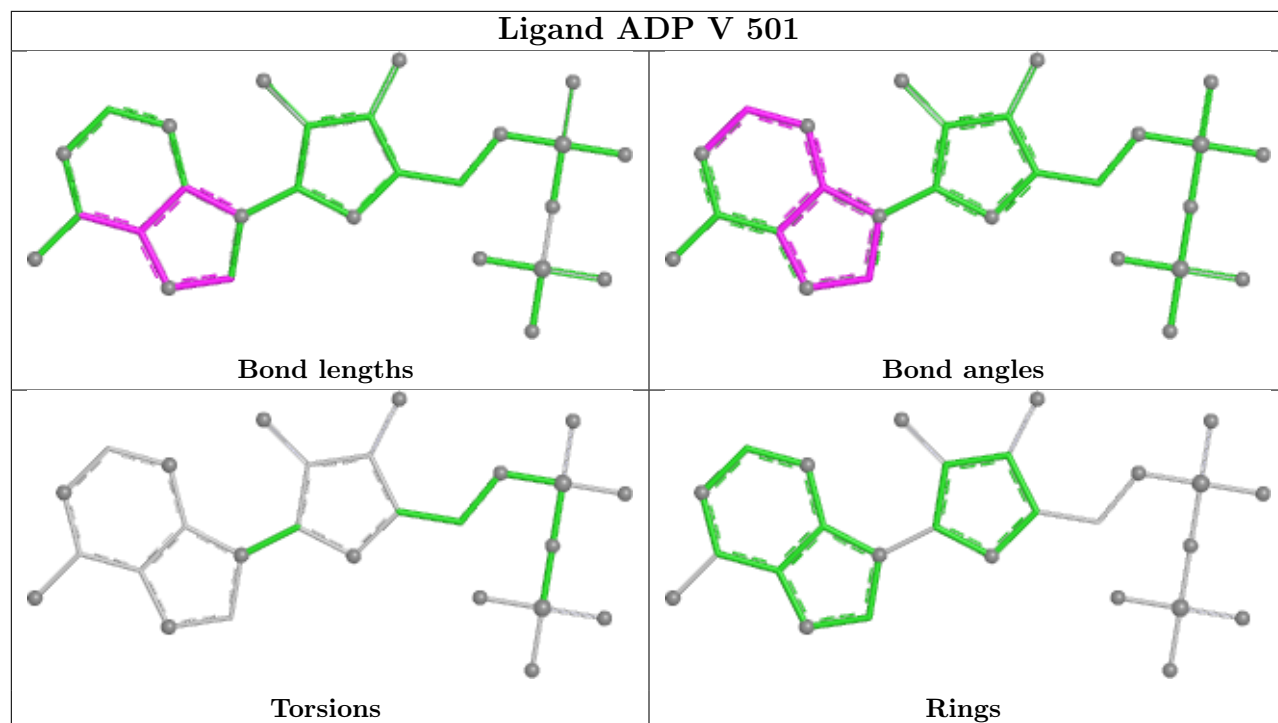
6 monomers are involved in 10 short contacts:

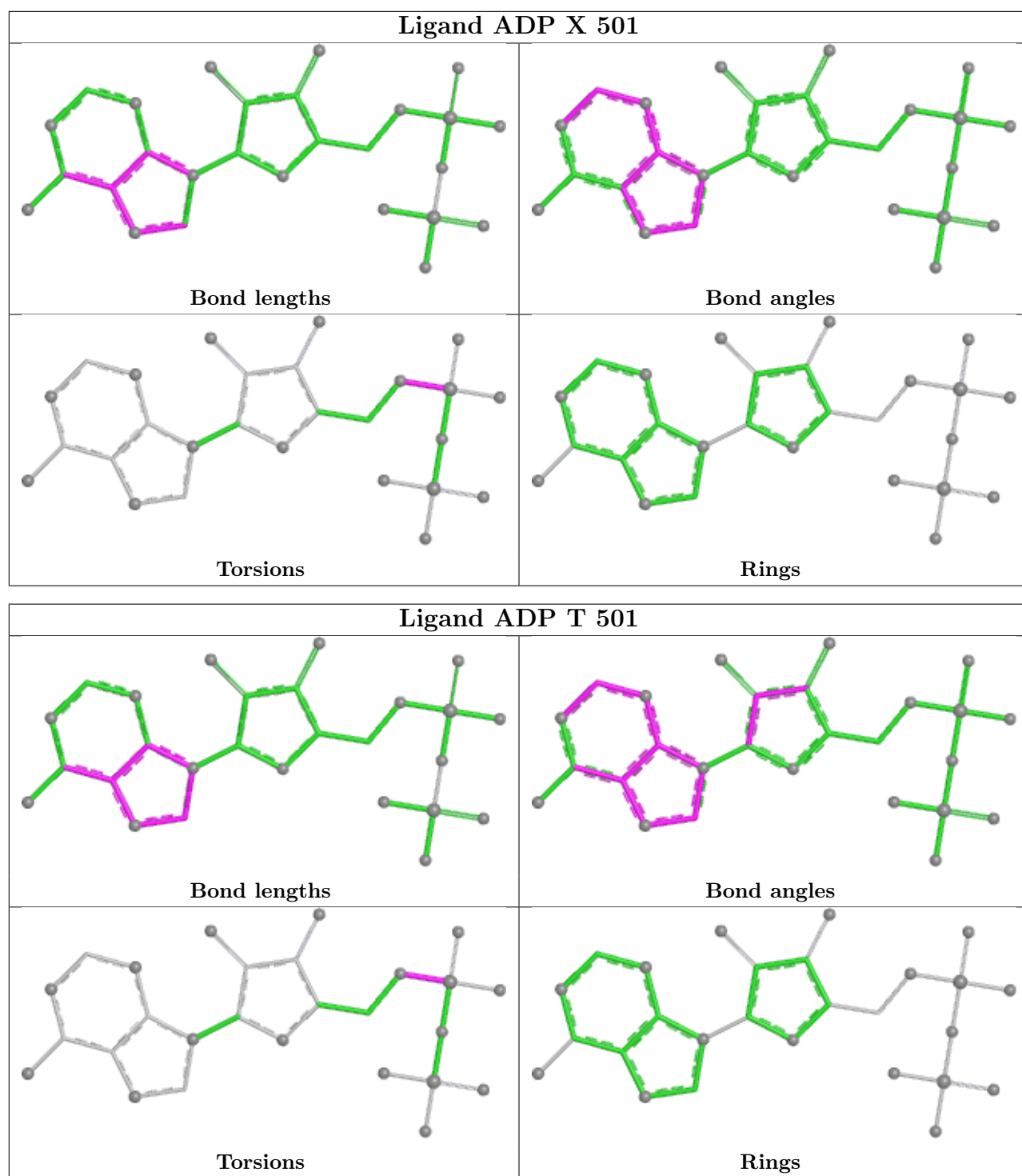
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	Y	501	ADP	1	0
7	W	501	ADP	1	0
7	V	501	ADP	2	0
7	U	501	ADP	4	0
7	X	501	ADP	1	0
7	T	501	ADP	1	0

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 ⓘ

There are no chain breaks in this entry.

6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-28613. These allow visual inspection of the internal detail of the map and identification of artifacts.

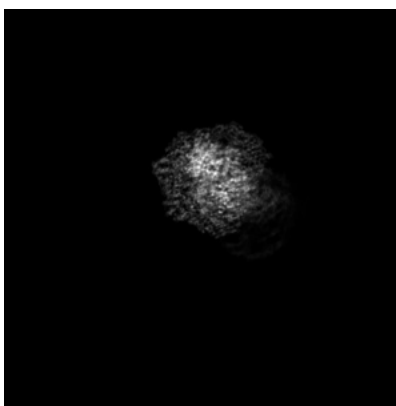
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

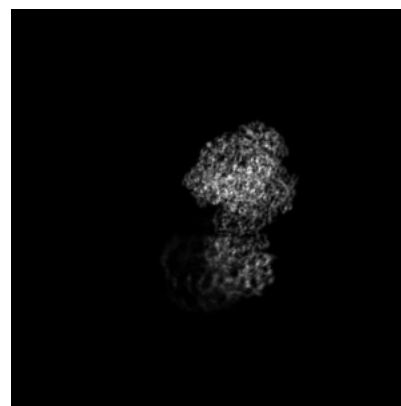
6.1.1 Primary map



X

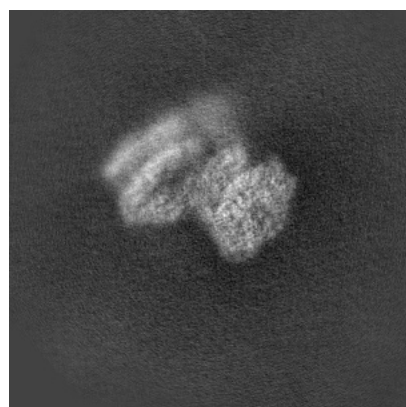


Y

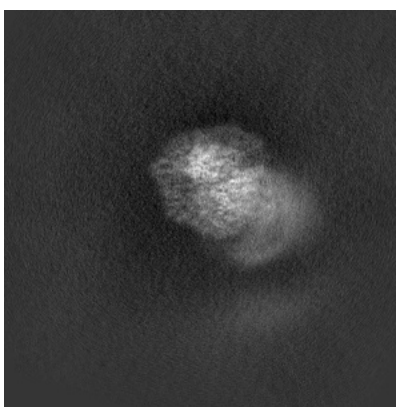


Z

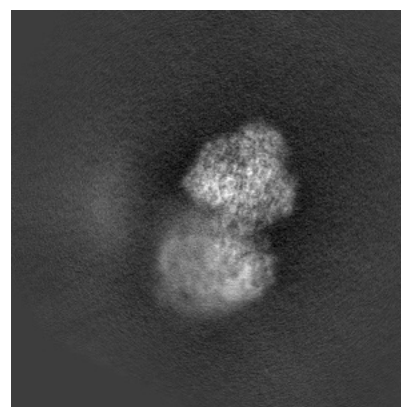
6.1.2 Raw map



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: 240

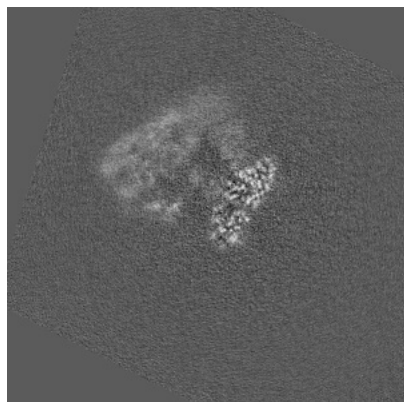


Y Index: 240

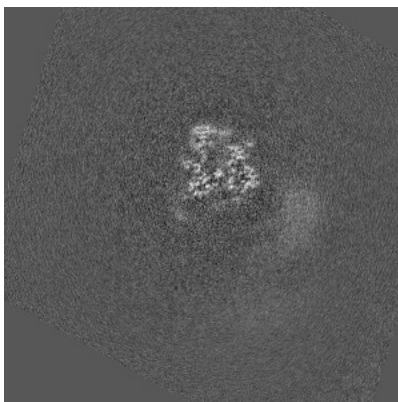


Z Index: 240

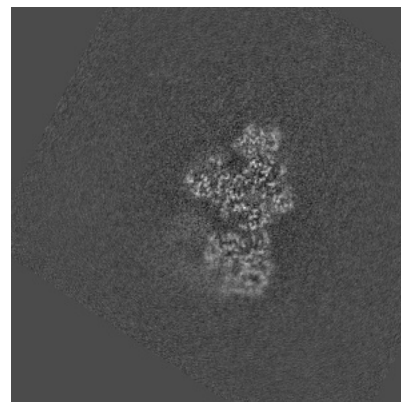
6.2.2 Raw map



X Index: 240



Y Index: 240



Z Index: 240

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

6.3 Largest variance slices [i](#)

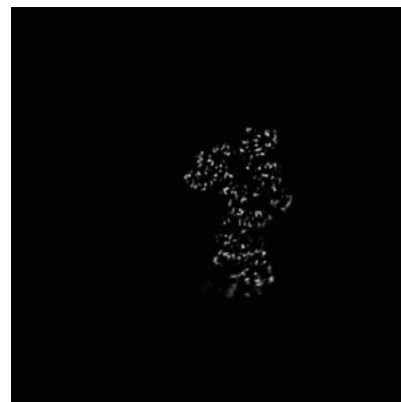
6.3.1 Primary map



X Index: 284

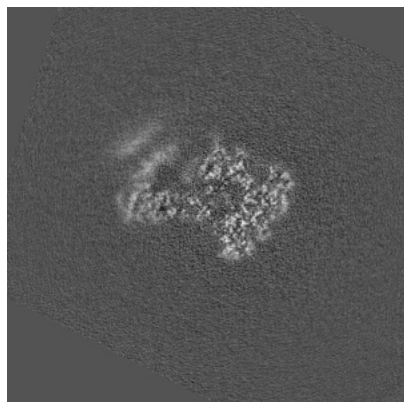


Y Index: 276

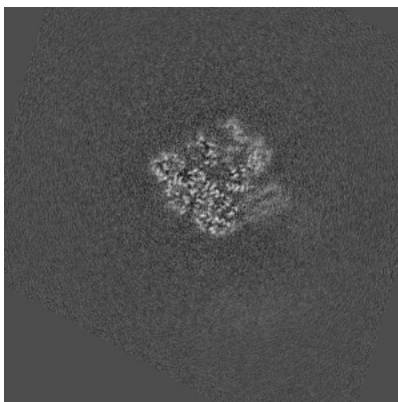


Z Index: 250

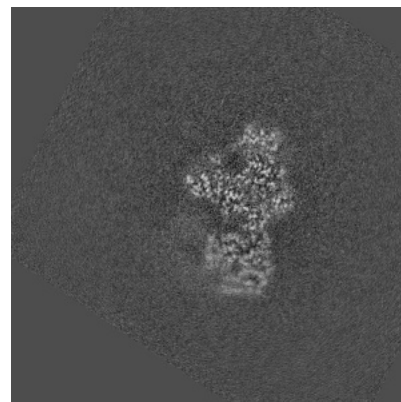
6.3.2 Raw map



X Index: 284



Y Index: 276

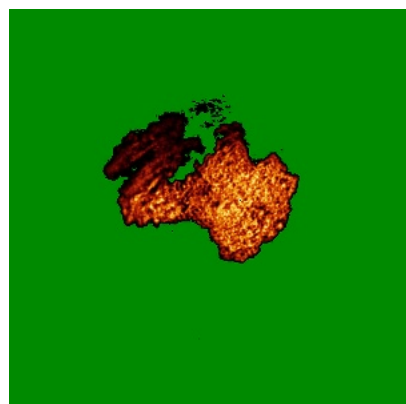


Z Index: 239

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

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

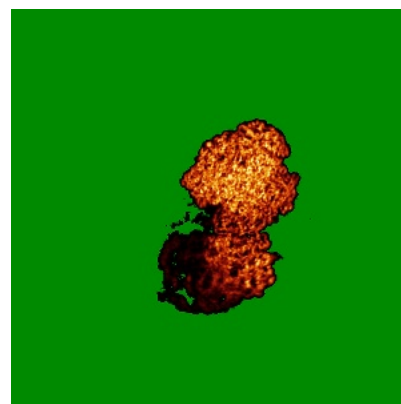
6.4.1 Primary map



X

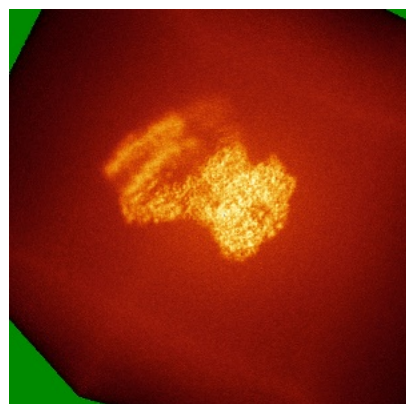


Y

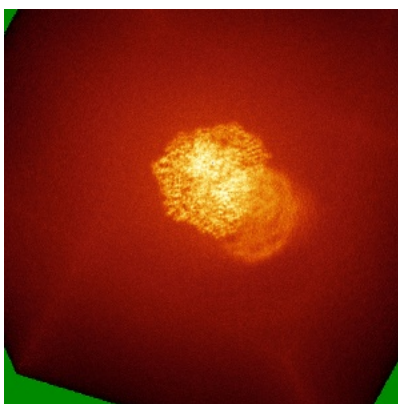


Z

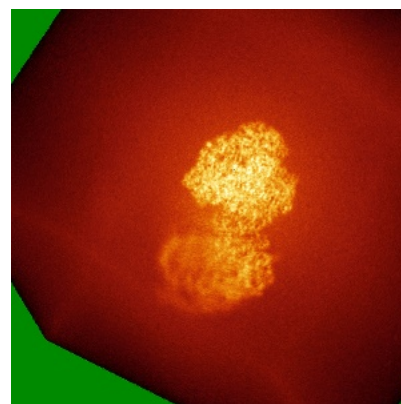
6.4.2 Raw map



X



Y

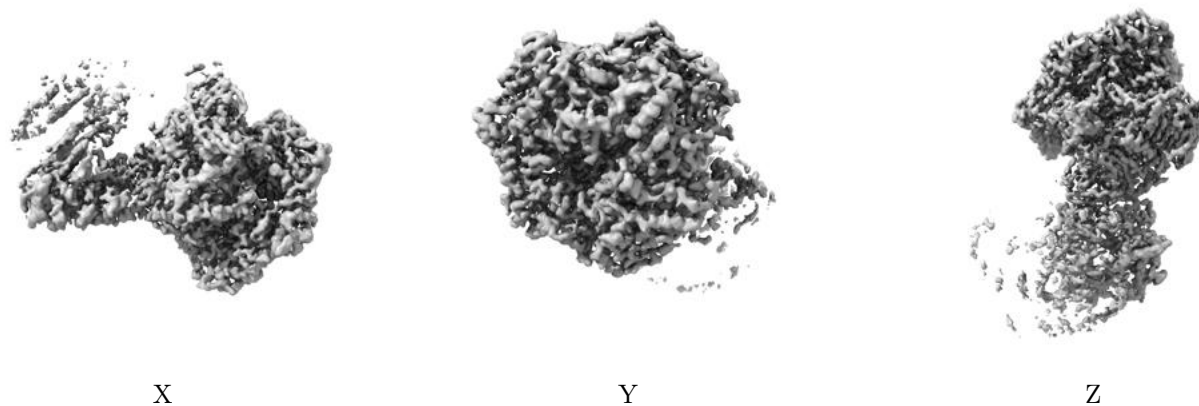


Z

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.

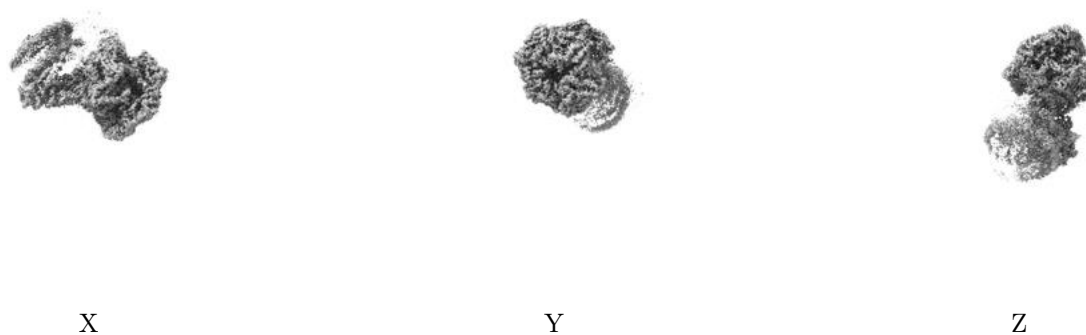
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.174. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

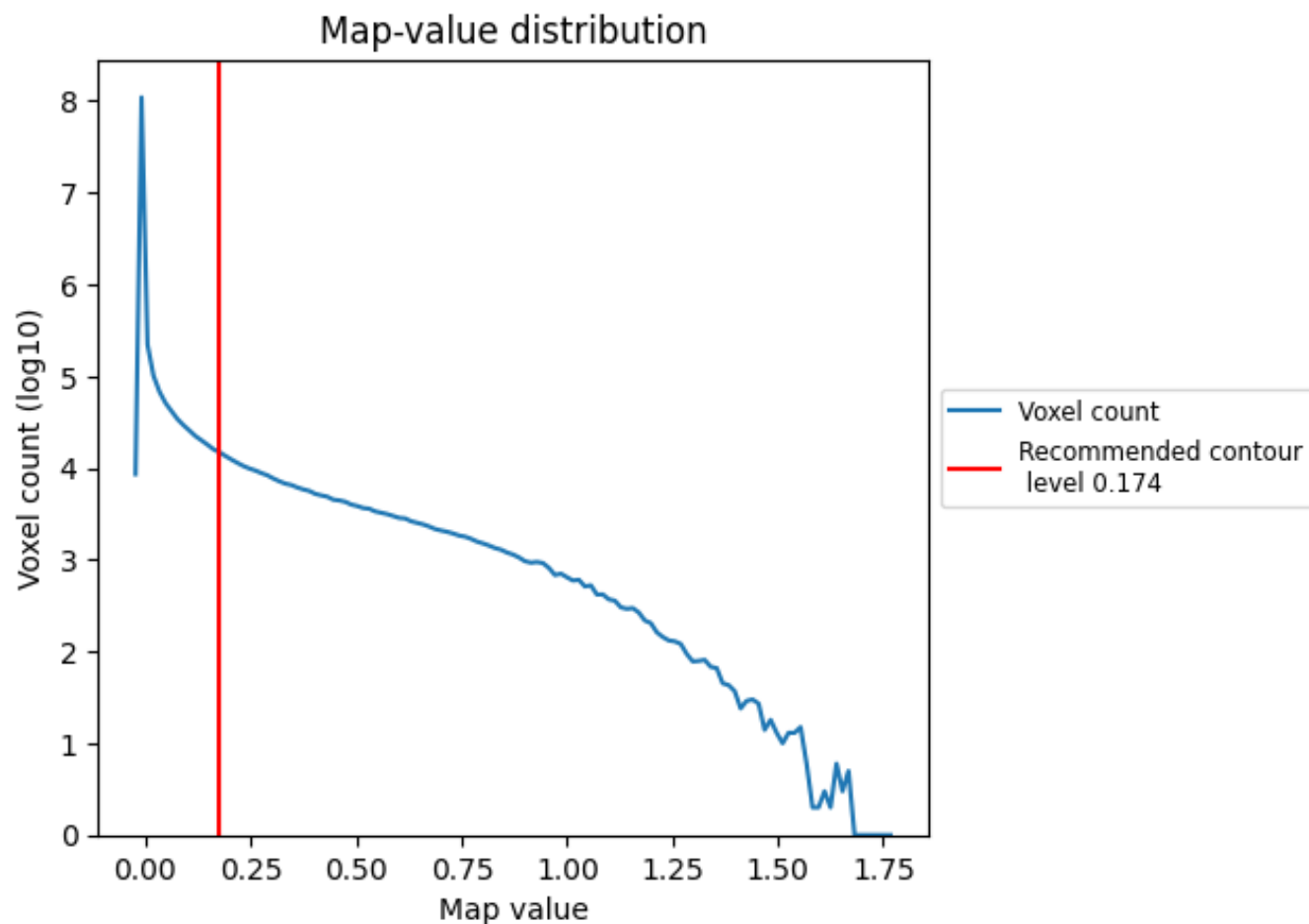
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

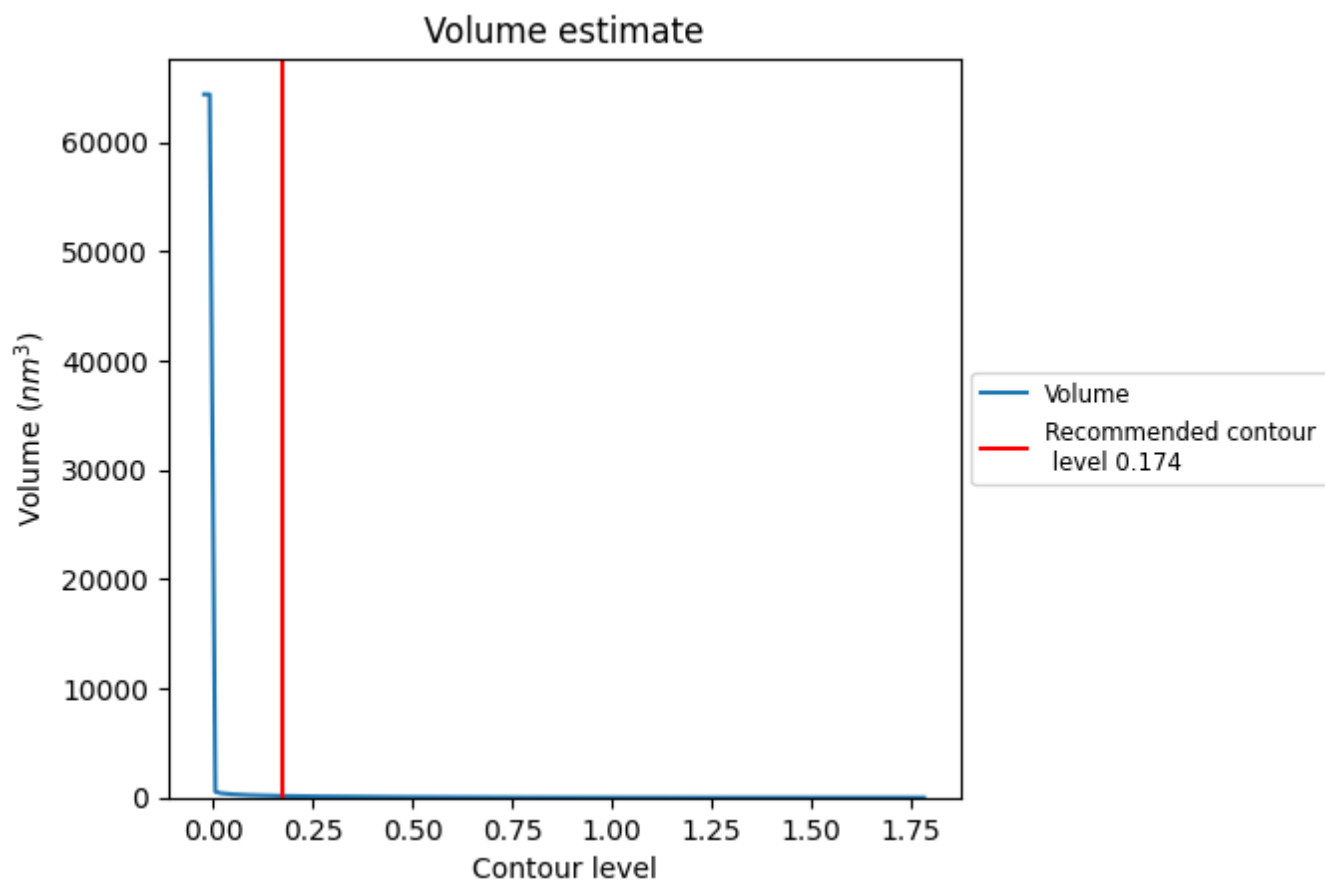
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

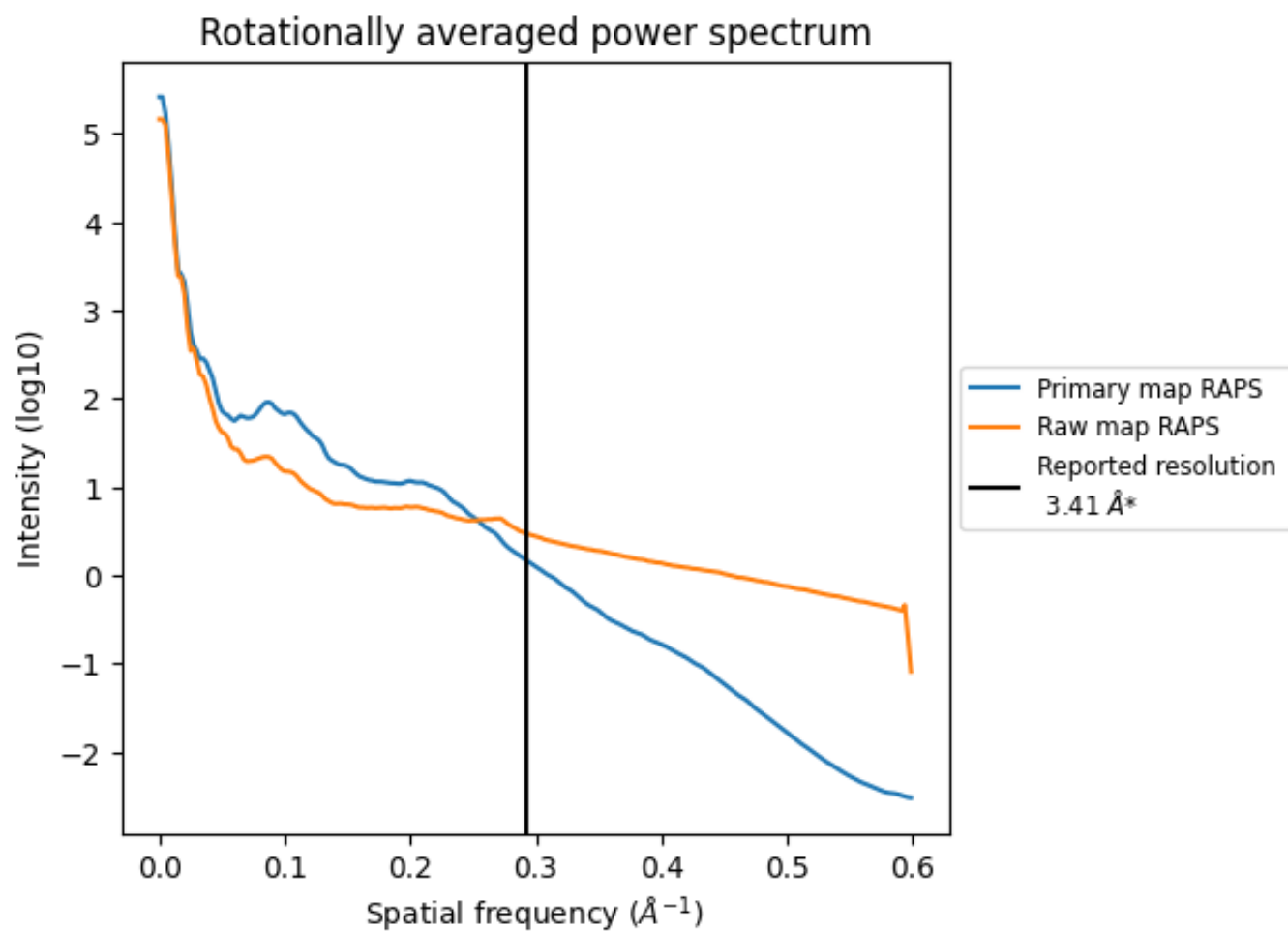
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 149 nm³; this corresponds to an approximate mass of 134 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 ⓘ

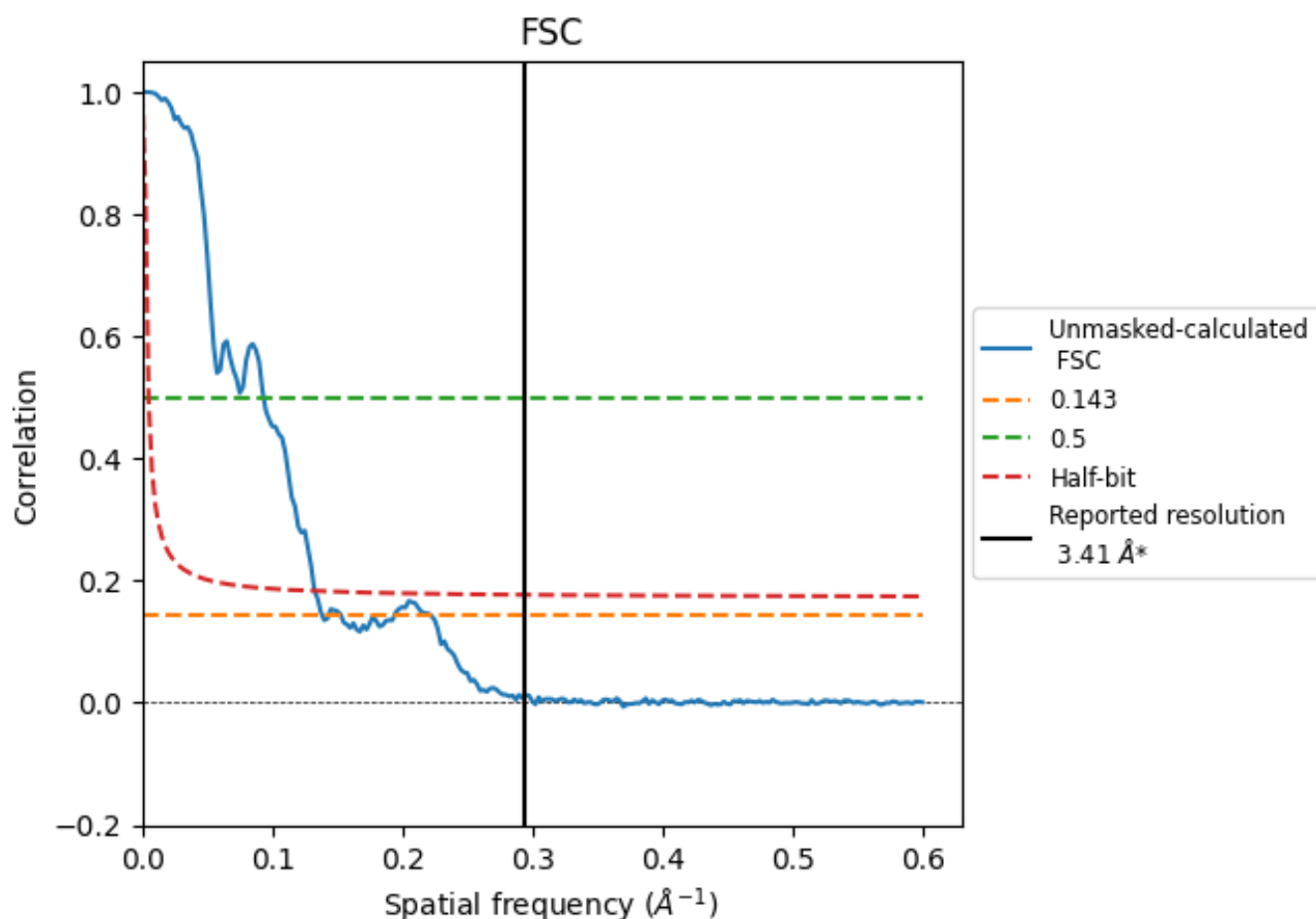


*Reported resolution corresponds to spatial frequency of 0.293 Å⁻¹

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 0.293 Å⁻¹

8.2 Resolution estimates [i](#)

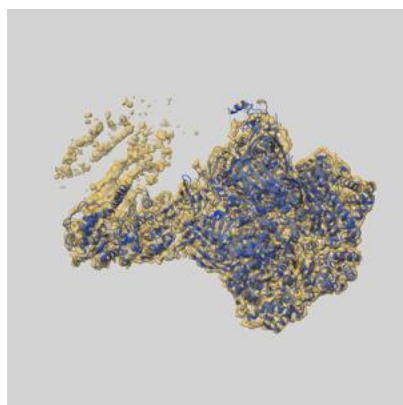
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.41	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	7.24	10.73	7.55

*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 7.24 differs from the reported value 3.41 by more than 10 %

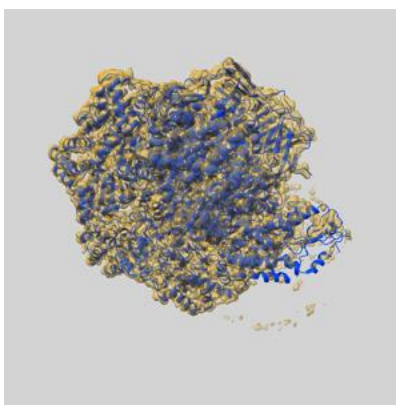
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-28613 and PDB model 8EUF. Per-residue inclusion information can be found in section 3 on page 6.

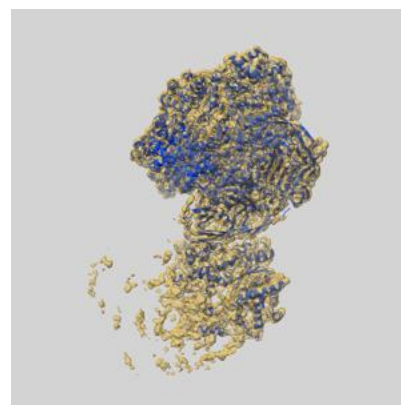
9.1 Map-model overlay [i](#)



X



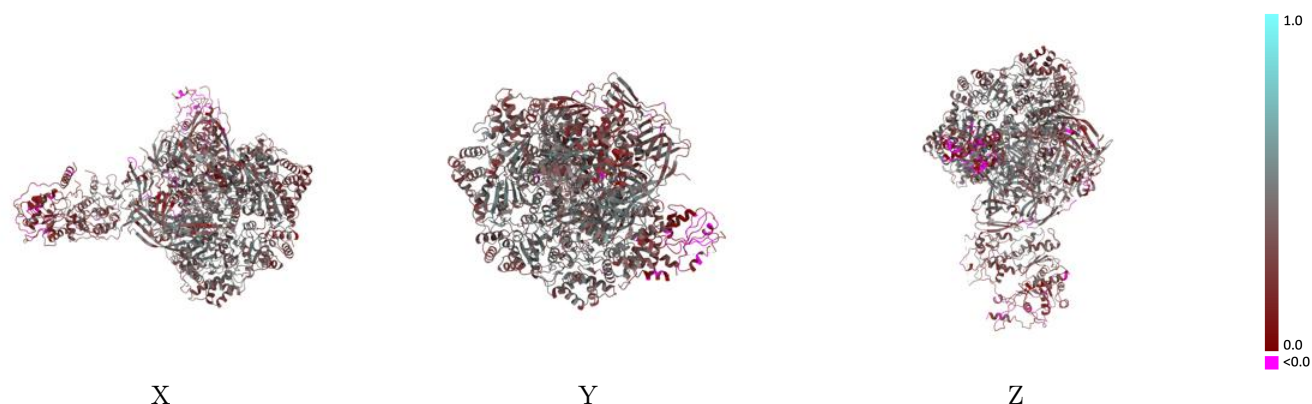
Y



Z

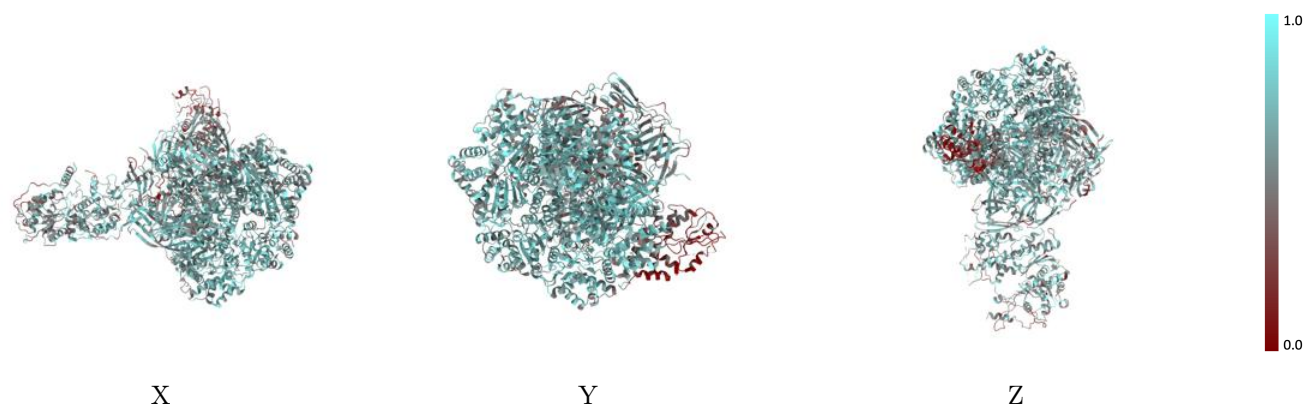
The images above show the 3D surface view of the map at the recommended contour level 0.174 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



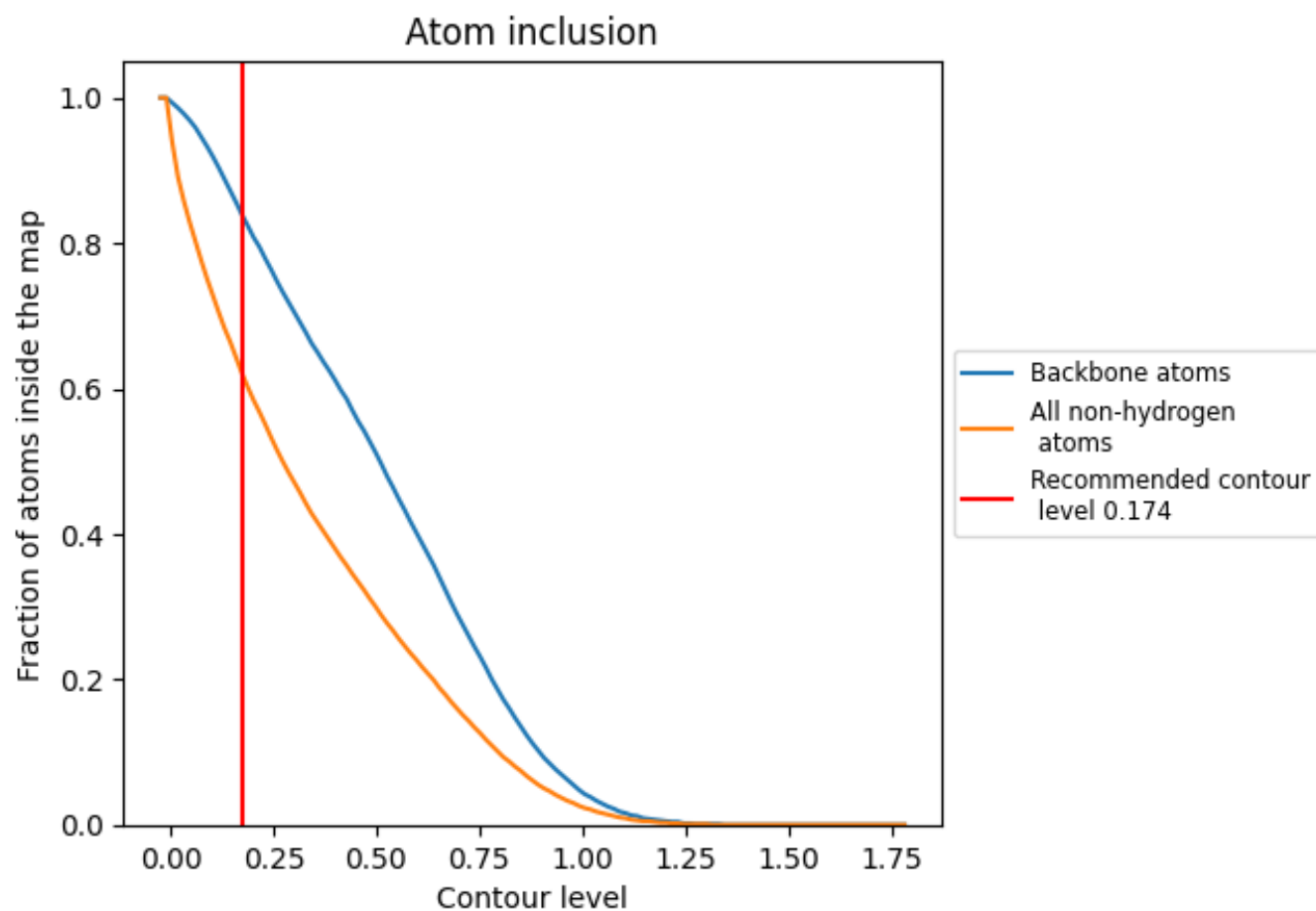
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.174).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 62% 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 (0.174) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6210	0.3470
Q	0.4670	0.2620
R	0.5580	0.2350
S	0.5750	0.2690
T	0.6650	0.3990
U	0.6610	0.4010
V	0.6890	0.4150
W	0.6510	0.3720
X	0.6350	0.3390
Y	0.6590	0.3800
Z	0.6360	0.3600

1.0

0.0

<0.0