



wwPDB EM Validation Summary Report ⓘ

Mar 6, 2026 – 11:46 PM UTC

PDB ID : 6ZNO / pdb_00006zno
EMDB ID : EMD-11319
Title : The pointed end complex of dynactin with the p150 projection docked
Authors : Lau, C.K.; Lacey, S.E.; Carter, A.P.
Deposited on : 2020-07-06
Resolution : 6.80 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:

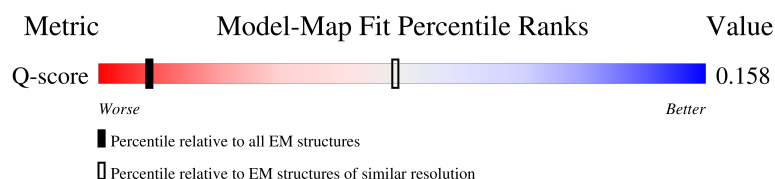
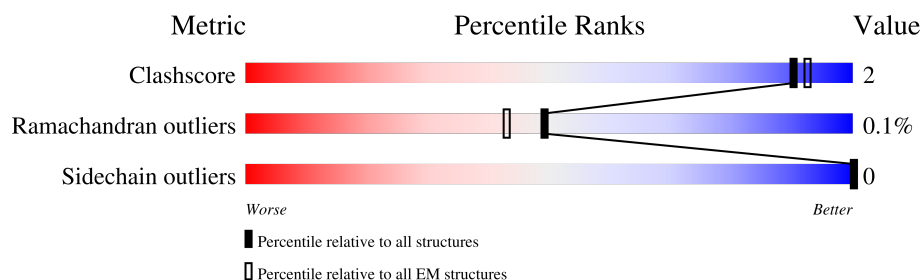
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 6.80 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.



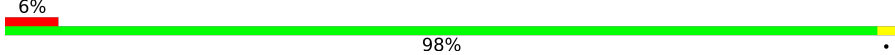
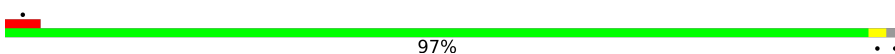
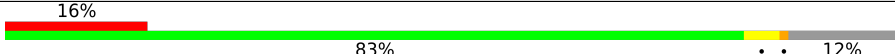
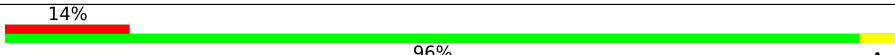
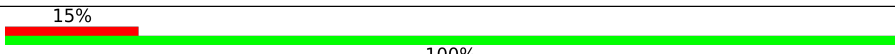
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	503 (6.30 - 7.30)

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	G	376	
1	I	376	
2	H	375	
3	J	417	

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Mol	Chain	Length	Quality of chain
4	M	405	
5	U	167	
6	V	182	
7	Y	467	
8	Z	189	
9	z	188	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 19737 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 ARP1 actin related protein 1 homolog A.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	G	370	Total	C	N	O	S	0	0
			2956	1892	509	545	10		
1	I	370	Total	C	N	O	S	0	0
			2941	1885	509	537	10		

- Molecule 2 is a protein called Actin, cytoplasmic 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	370	Total	C	N	O	S	0	0
			2885	1827	486	550	22		

- Molecule 3 is a protein called Arp11.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	379	Total	C	N	O	S	0	0
			2932	1888	496	532	16		

- Molecule 4 is a protein called Dynactin subunit 2.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	M	75	Total	C	N	O	0	0
			583	363	99	121		

- Molecule 5 is a protein called Dynactin 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	U	167	Total	C	N	O	S	0	0
			1224	771	212	231	10		

- Molecule 6 is a protein called Dynactin subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	V	179	Total	C	N	O	S	0	0
			1260	818	222	211	9		

- Molecule 7 is a protein called Dynactin subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	Y	410	Total	C	N	O	S	0	0
			2956	1866	543	527	20		

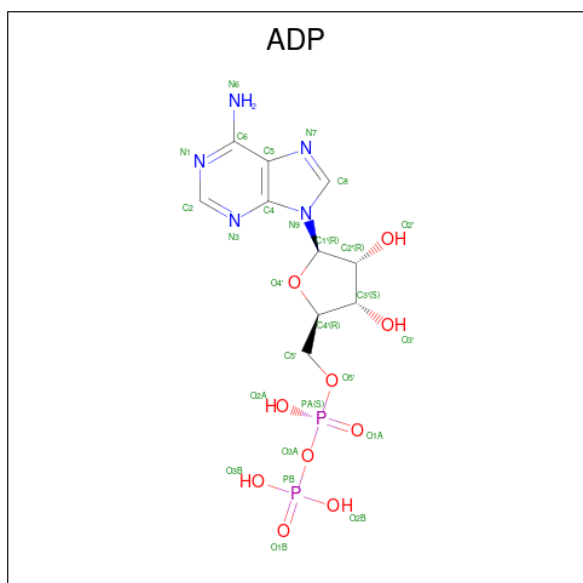
- Molecule 8 is a protein called Dynactin subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	Z	189	Total	C	N	O		0	0
			945	567	189	189			

- Molecule 9 is a protein called Dynactin subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	z	188	Total	C	N	O		0	0
			940	564	188	188			

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



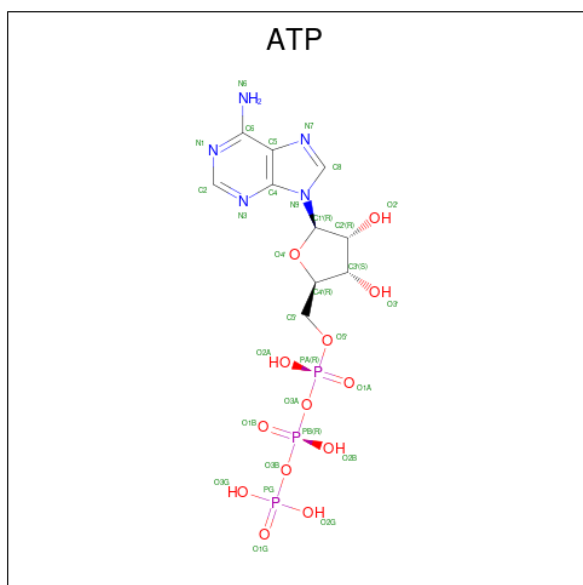
Mol	Chain	Residues	Atoms					AltConf
10	G	1	Total	C	N	O	P	0
			27	10	5	10	2	

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

- Molecule 11 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

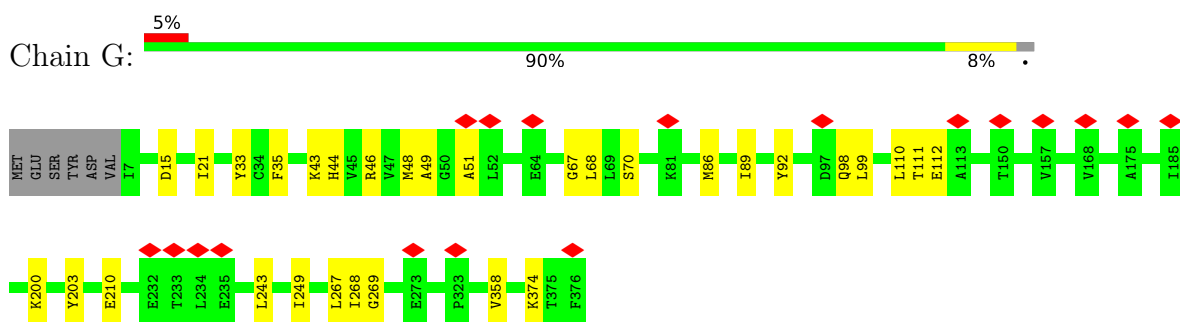
- Molecule 12 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
12	Y	3	Total	Zn	0
			3	3	

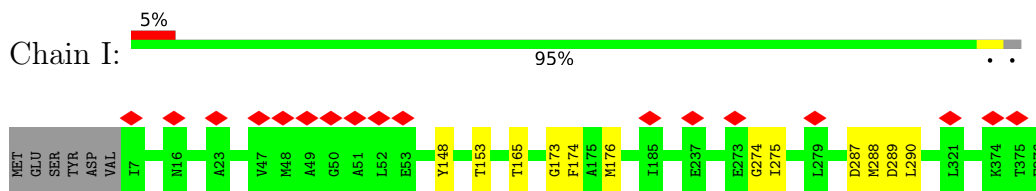
3 Residue-property plots

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

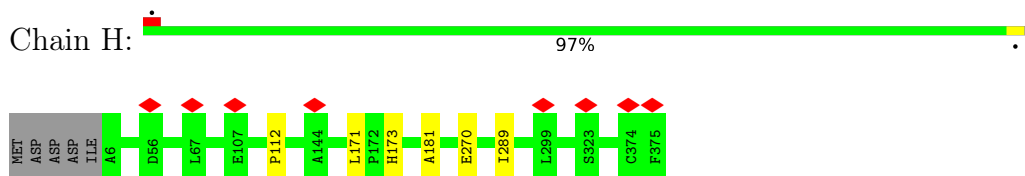
- Molecule 1: ARP1 actin related protein 1 homolog A



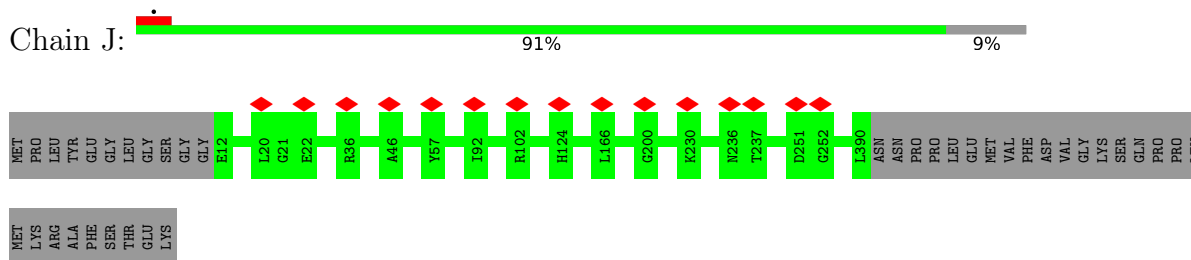
- Molecule 1: ARP1 actin related protein 1 homolog A



- Molecule 2: Actin, cytoplasmic 1



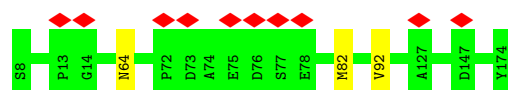
- Molecule 3: Arp11



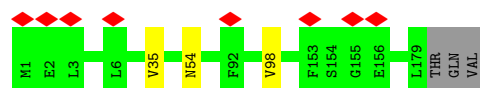
- Molecule 4: Dynactin subunit 2

GLN	LEU	LEU	VAL	GLY	GLU	ASP	VAL	MET
LEU	LEU	VAL	ARG	ALA	ARG	GLY	GLN	ALA
THR	GLN	VAL	LEU	ALA	LEU	LEU	GLU	D3
HIS	THR	ASN	THR	THR	ALA	ALA	THR	P4
LEU	ILE	ILE	GLU	LYS	LYS	THR	THR	K5
ASP	ALA	ALA	LEU	ARG	ARG	GLU	GLU	Y6
THR	LYS	LYS	GLU	LEU	LEU	VAL	VAL	A7
THR	HIS	HIS	ALA	LEU	LEU	GLU	GLU	D8
GLN	LYS	LYS	THR	LEU	LEU	LYS	LYS	L9
GLN	ALA	ALA	VAL	GLN	GLN	ILE	ILE	
MET	SER	SER	ARG	LEU	LEU	LYS	LYS	E14
ILE	VAL	CYS	VAL	GLU	GLU	MET	MET	M15
ALA	GLU	ASP	ALA	ALA	THR	THR	THR	E16
CYS	GLN	GLN	GLN	THR	THR	VAL	VAL	
SER	SER	ALA	ASP	ASP	LYS	LYS	LYS	
LEU	ASP	ALA	ALA	ASN	ASN	GLU	GLU	L25
LYS	THR	THR	GLN	THR	THR	SER	SER	P26
ASP	GLN	GLN	ASN	LYS	LYS	ALA	ALA	GLU
ASN	SER	PRO	PRO	GLY	GLY	THR	THR	ASP
ALA	LEU	LEU	LEU	ALA	ALA	GLU	GLU	ASP
THR	VAL	SER	SER	GLY	GLY	GLU	GLU	GLN
LEU	HIS	ALA	ALA	SER	LYS	LYS	ALA	ALA
LEU	GLN	GLY	GLY	GLY	LEU	LEU	LEU	GLU
THR	LEU	LEU	LEU	GLY	THR	THR	THR	L37
VAL	TYR	GLN	GLN	LYS	THR	PRO	GLN	E38
GLN	THR	ALA	ALA	THR	THR	VAL	VAL	E39
THR	ILE	CYS	CYS	SER	SER	LEU	LEU	
THR	GLN	LEU	LEU	GLY	GLY	ALA	ALA	
MET	ARG	MET	MET	SER	SER	LYS	LYS	
THR	TRP	GLU	GLU	PRO	PRO	GLN	GLN	
GLU	SER	THR	THR	THR	THR	ALA	ALA	
ASN	PRO	VAL	VAL	ASP	ASP	LEU	LEU	
LEU	ILE	GLU	GLU	SER	SER	ALA	ALA	
SER	ALA	LEU	LEU	SER	SER	LYS	LYS	I75
THR	SER	LEU	LEU	LEU	LEU	LYS	LYS	
VAL	THR	GLN	GLN	VAL	VAL	GLN	GLN	
GLU	LEU	LEU	ALA	ALA	THR	GLN	GLN	G86
GLY	GLY	PRO	LYS	LYS	TYR	LEU	LEU	E87
ASN	GLU	VAL	VAL	GLU	GLU	VAL	VAL	TYR
PHE	LEU	SER	SER	LEU	LEU	ALA	ALA	GLU
ALA	VAL	ALA	ALA	HIS	HIS	SER	SER	MET
ASN	GLN	LEU	LEU	SER	SER	HIS	HIS	LEU
ILE	ARG	ASP	ASP	ASP	ASP	PRO	PRO	GLY
ASP	LEU	LEU	LEU	PRO	PRO	GLU	GLU	LEU
GLU	VAL	ALA	ALA	GLN	GLN	LYS	LYS	GLY
ARG	THR	VAL	VAL	GLN	GLN	LEU	LEU	VAL
MET	ILE	LEU	LEU	ASP	ASP	LEU	LEU	VAL
LYS	LYS	LYS	ASP	LYS	LYS	GLY	GLY	VAL
LYS	GLN	GLN	GLN	PHE	PHE	PRO	PRO	LYS
LEU	LEU	LEU	VAL	SER	SER	ASP	ASP	GLU
GLY	HIS	HIS	GLU	GLU	GLN	ALA	ALA	THR
LYS	GLU	GLU	ALA	ALA	ALA	ALA	ALA	PRO
	GLN	ARG	ARG	ALA	ALA	ILE	ILE	GLN
	ALA	LEU	LEU	LYS	LYS	ASN	ASN	GLN
	ALA	ALA	GLN	VAL	VAL	THR	THR	LYS
	MET	MET	GLN	VAL	VAL	ALA	ALA	TYR
	GLN	GLN	SER	ALA	ALA	THR	THR	GLN
	THR	PHE	THR	GLU	GLU	ASP	ASP	ARG
	GLY	GLY	LEU	LEU	LEU	PRO	PRO	LEU
								LEU
								HIS

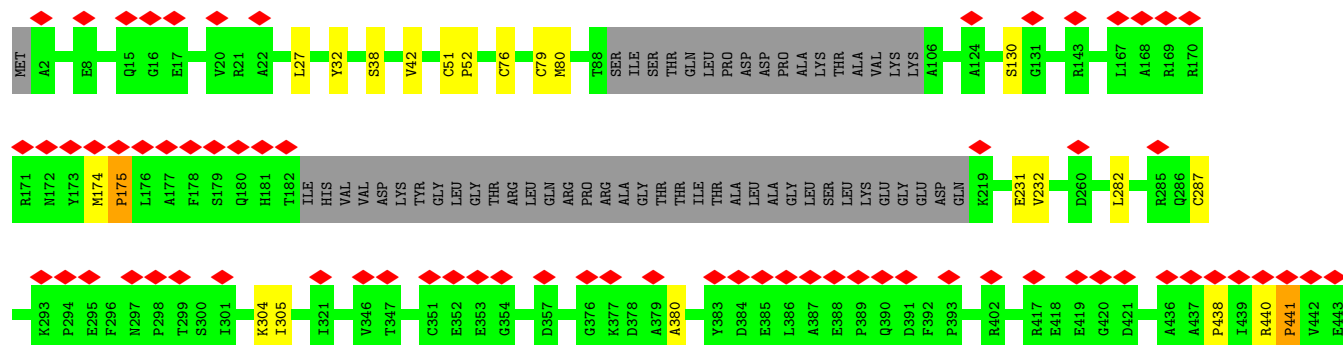
Chain U: 98% 6%

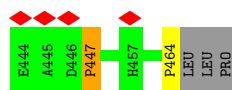


Chain V: 97%



Chain Y: 16% 83% 12%

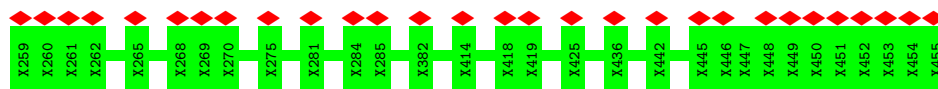




- Molecule 8: Dynactin subunit 1



- Molecule 9: Dynactin subunit 1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	20037	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 ($e^-/\text{\AA}^2$)	52, 40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k), GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.177	Depositor
Minimum map value	-0.065	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.0492	Depositor
Map size ($\text{\AA}$)	348.4, 348.4, 348.4	wwPDB
Map dimensions	260, 260, 260	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.34, 1.34, 1.34	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, ZN

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	G	1.03	0/3025	1.54	0/4085
1	I	1.03	0/3010	1.51	1/4066 (0.0%)
2	H	1.04	0/2948	1.54	2/3991 (0.1%)
3	J	1.04	0/2994	1.57	0/4066
4	M	1.03	0/593	1.44	0/800
5	U	1.11	0/1241	1.45	0/1691
6	V	1.12	0/1286	1.49	0/1757
7	Y	1.10	0/3016	1.53	13/4114 (0.3%)
All	All	1.06	0/18113	1.53	16/24570 (0.1%)

There are no bond length outliers.

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	Y	438	PRO	N-CA-CB	8.46	110.06	103.30
7	Y	447	PRO	N-CA-CB	6.54	110.11	103.25
7	Y	441	PRO	N-CA-CB	6.48	110.05	103.25
7	Y	464	PRO	N-CA-CB	6.36	109.99	103.00
1	I	274	GLY	CA-C-O	-6.27	118.15	122.22

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	2956	0	2950	43	0
1	I	2941	0	2936	26	0
2	H	2885	0	2856	9	0
3	J	2932	0	3003	0	0
4	M	583	0	558	0	0
5	U	1224	0	1215	2	0
6	V	1260	0	1213	4	0
7	Y	2956	0	2714	10	0
8	Z	945	0	192	6	0
9	z	940	0	192	0	0
10	G	27	0	12	0	0
10	I	27	0	12	0	0
10	J	27	0	12	0	0
11	H	31	0	12	0	0
12	Y	3	0	0	0	0
All	All	19737	0	17877	63	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 2.

The worst 5 of 63 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:G:44:HIS:CD2	1:I:176:MET:HG3	1.58	1.36
1:G:44:HIS:CG	1:I:176:MET:HG3	1.78	1.17
1:G:51:ALA:HB2	1:I:173:GLY:HA3	1.30	1.07
1:G:44:HIS:CD2	1:I:176:MET:CG	2.40	1.05
1:G:70:SER:HB2	2:H:270:GLU:OE2	1.68	0.94

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	G	368/376 (98%)	353 (96%)	15 (4%)	0	100	100
1	I	368/376 (98%)	347 (94%)	21 (6%)	0	100	100
2	H	368/375 (98%)	352 (96%)	16 (4%)	0	100	100
3	J	377/417 (90%)	347 (92%)	30 (8%)	0	100	100
4	M	71/405 (18%)	62 (87%)	9 (13%)	0	100	100
5	U	165/167 (99%)	141 (86%)	24 (14%)	0	100	100
6	V	177/182 (97%)	161 (91%)	16 (9%)	0	100	100
7	Y	404/467 (86%)	355 (88%)	46 (11%)	3 (1%)	18	56
All	All	2298/2765 (83%)	2118 (92%)	177 (8%)	3 (0%)	49	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	Y	175	PRO
7	Y	441	PRO
7	Y	447	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	318/324 (98%)	318 (100%)	0	100	100
1	I	314/324 (97%)	314 (100%)	0	100	100
2	H	313/318 (98%)	313 (100%)	0	100	100
3	J	323/363 (89%)	323 (100%)	0	100	100
4	M	63/346 (18%)	63 (100%)	0	100	100
5	U	129/142 (91%)	129 (100%)	0	100	100
6	V	121/163 (74%)	121 (100%)	0	100	100
7	Y	273/416 (66%)	273 (100%)	0	100	100
All	All	1854/2396 (77%)	1854 (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. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
6	V	100	ASN
7	Y	457	HIS
7	Y	142	GLN
3	J	232	ASN
6	V	20	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
10	ADP	I	800	-	28,29,29	0.46	0	43,45,45	0.44	0
10	ADP	G	800	-	28,29,29	0.43	0	43,45,45	0.44	0
11	ATP	H	401	-	32,33,33	0.53	0	48,52,52	0.56	0
10	ADP	J	800	-	28,29,29	0.50	0	43,45,45	0.51	0

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
10	ADP	I	800	-	-	4/16/32/32	0/3/3/3
10	ADP	G	800	-	-	4/16/32/32	0/3/3/3
11	ATP	H	401	-	-	6/22/38/38	0/3/3/3
10	ADP	J	800	-	-	3/16/32/32	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

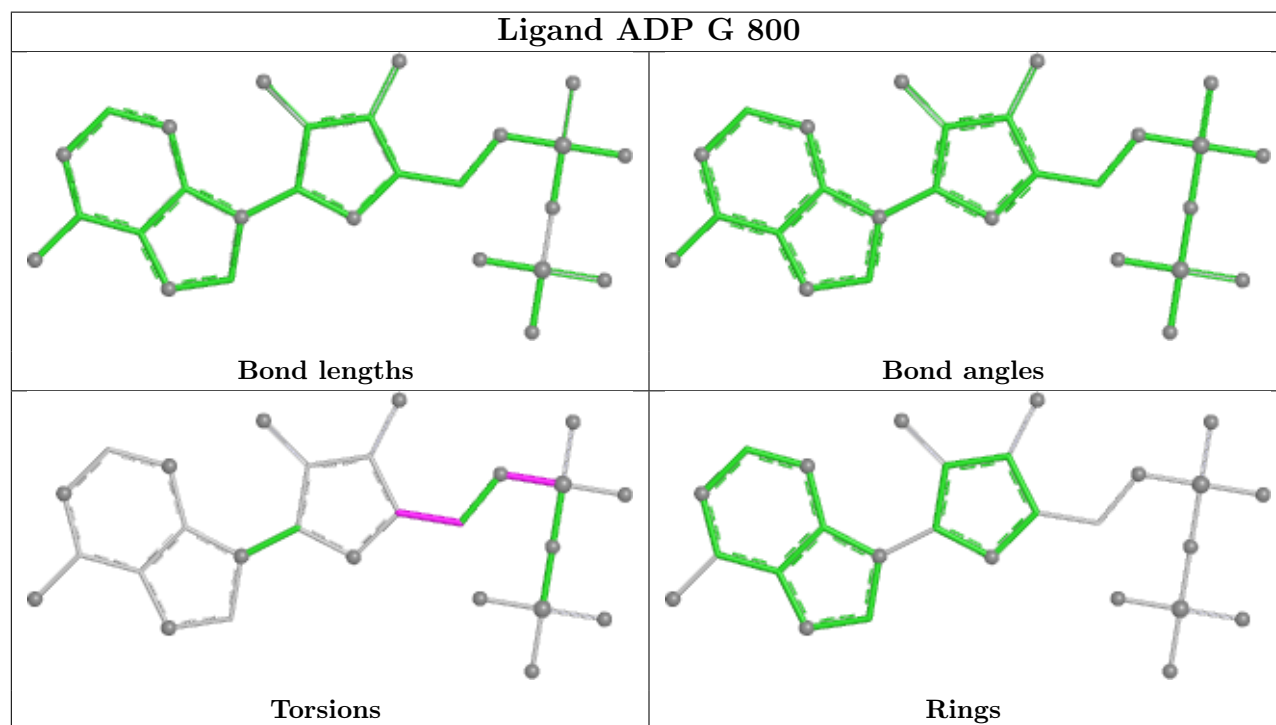
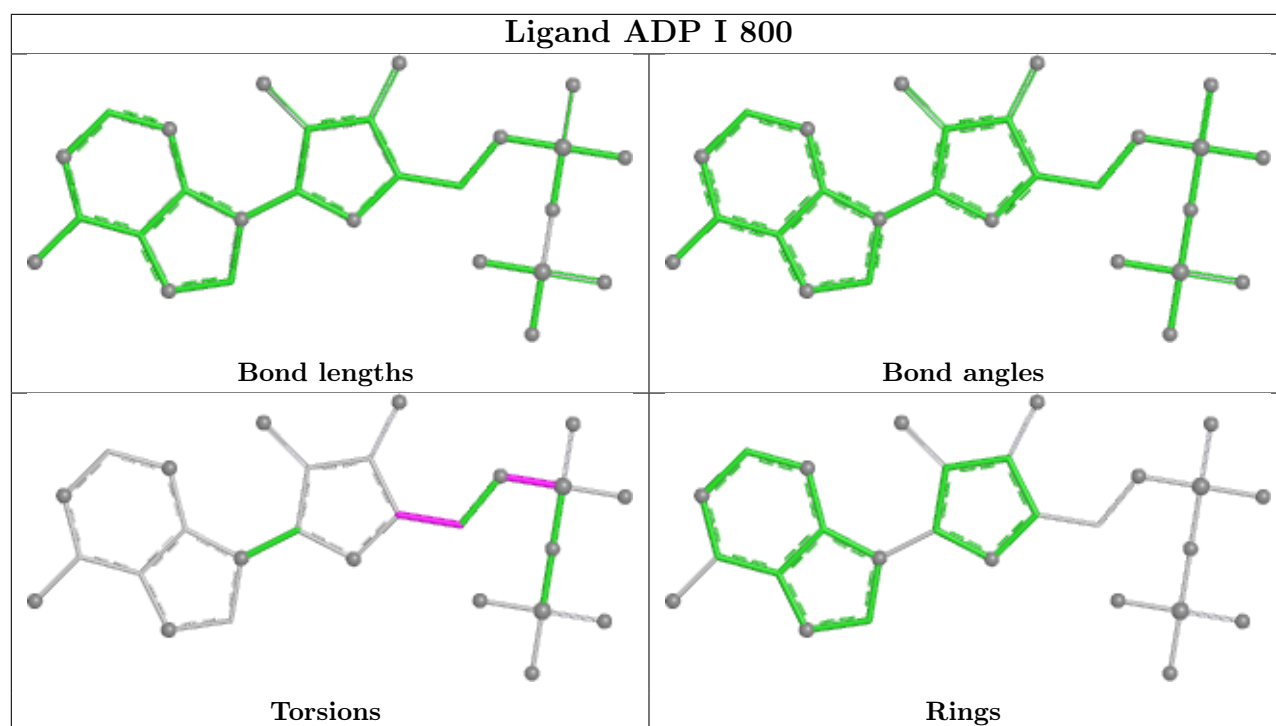
5 of 17 torsion outliers are listed below:

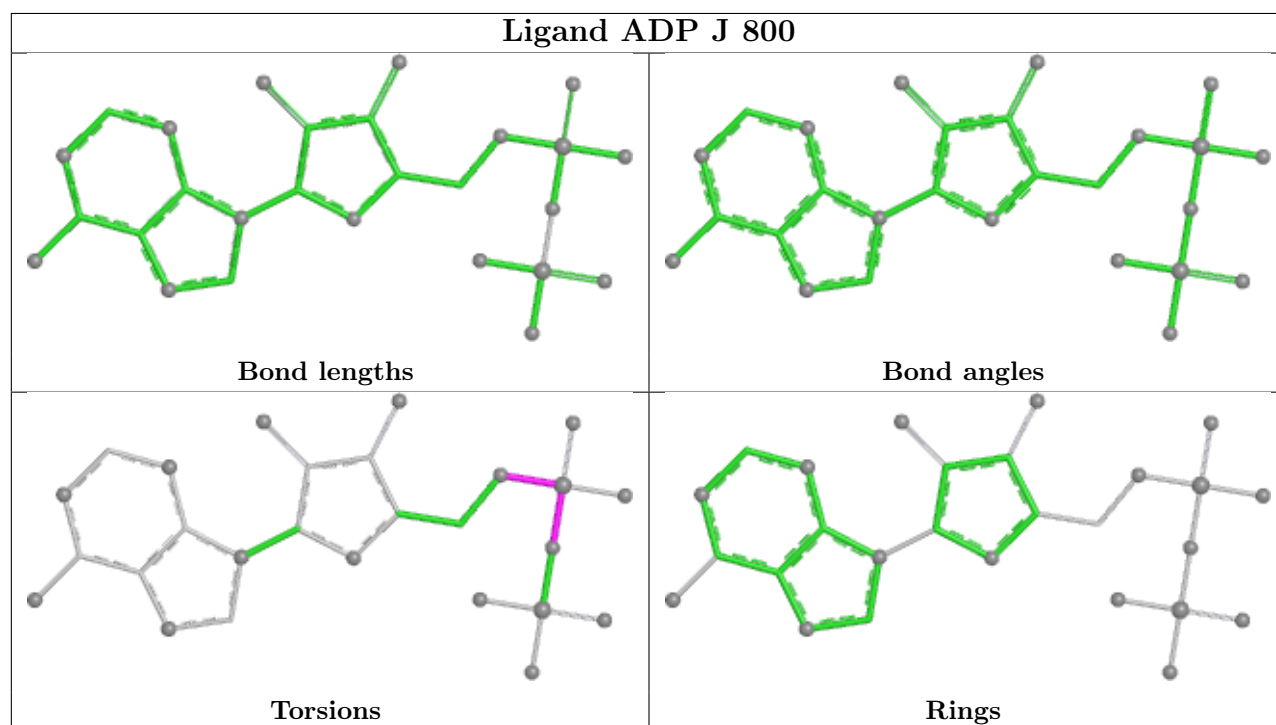
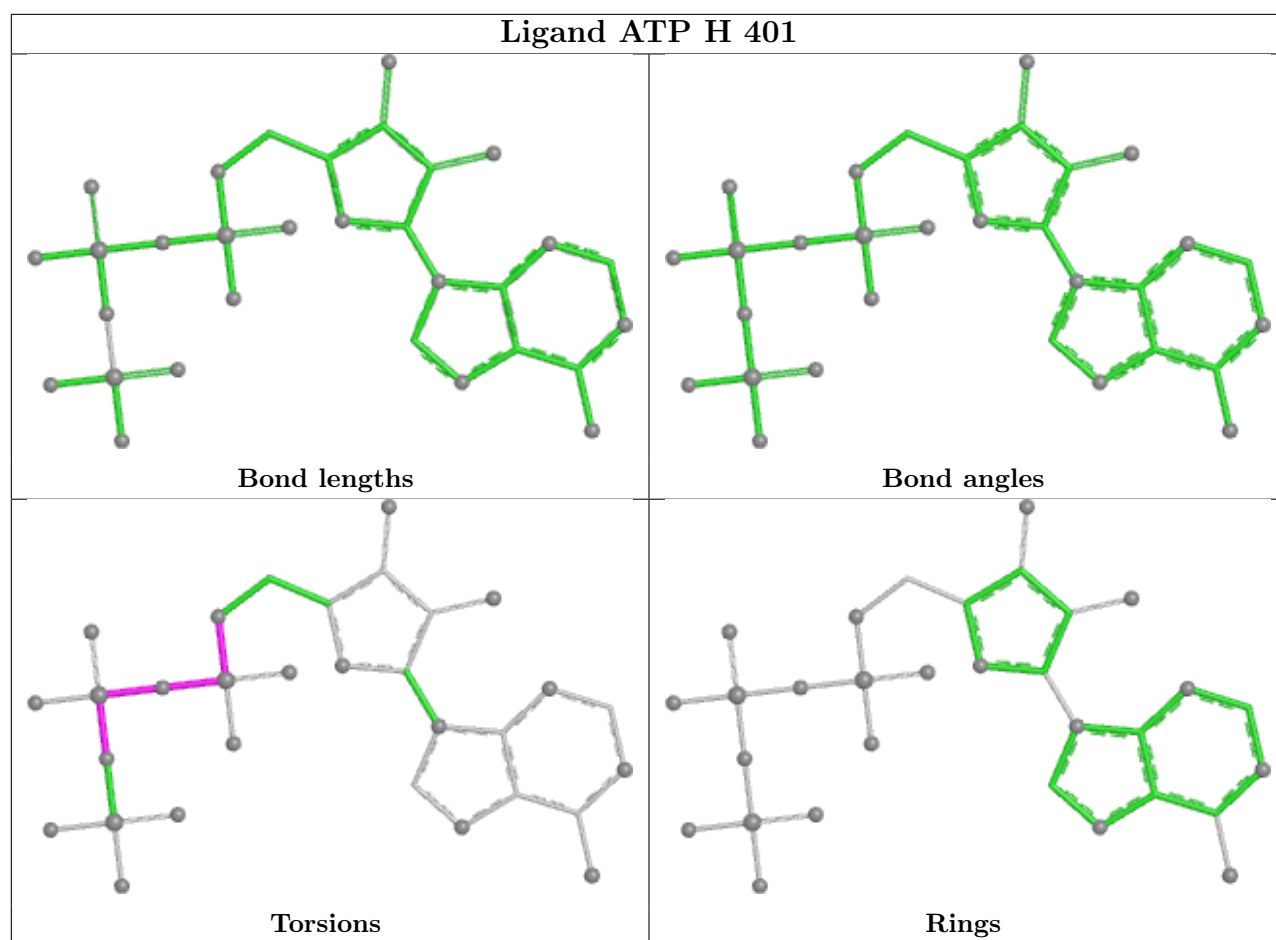
Mol	Chain	Res	Type	Atoms
10	G	800	ADP	C5'-O5'-PA-O2A
10	G	800	ADP	C5'-O5'-PA-O3A
10	I	800	ADP	C5'-O5'-PA-O2A
10	I	800	ADP	C5'-O5'-PA-O3A
11	H	401	ATP	PB-O3A-PA-O5'

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
9	<i>z</i>	1
8	<i>Z</i>	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	<i>z</i>	344:UNK	C	354:UNK	N	16.34
1	<i>Z</i>	342:UNK	C	355:UNK	N	12.59

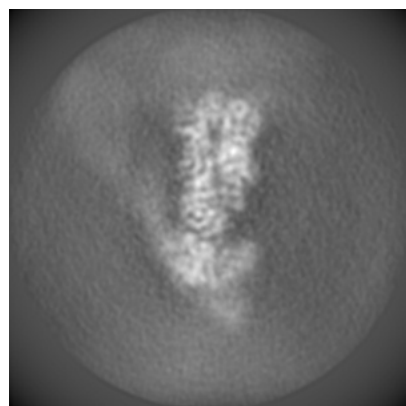
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11319. These allow visual inspection of the internal detail of the map and identification of artifacts.

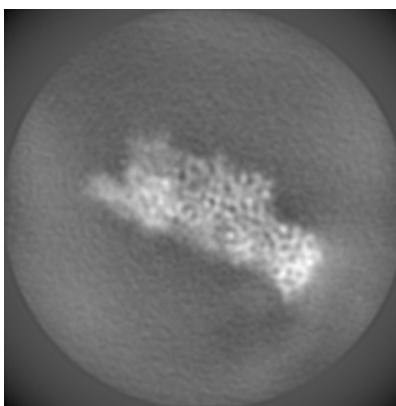
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

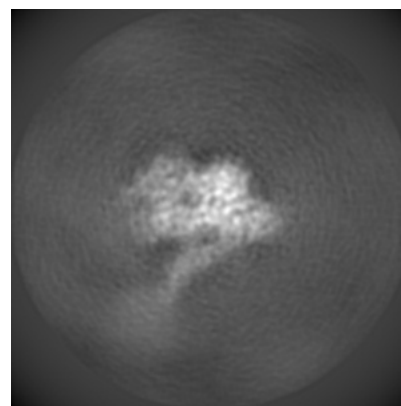
6.1.1 Primary map



X

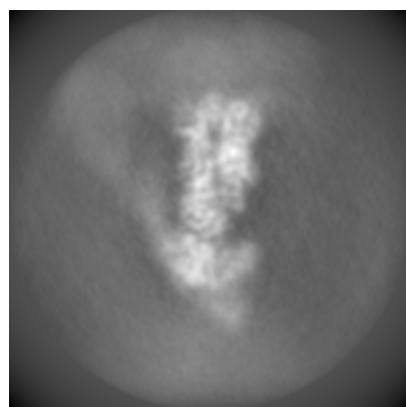


Y

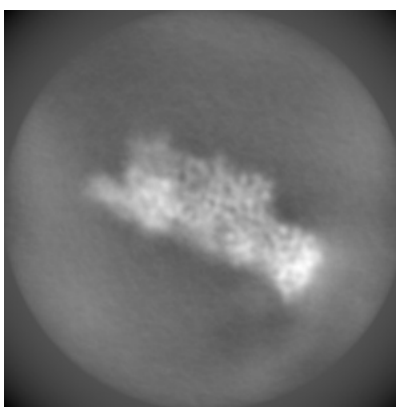


Z

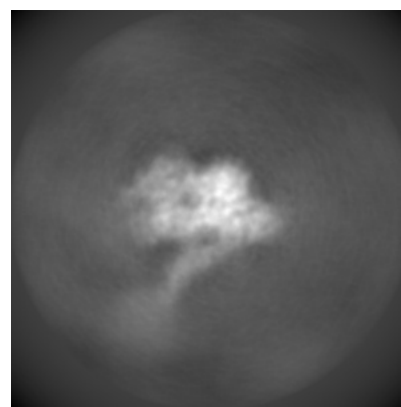
6.1.2 Raw map



X



Y

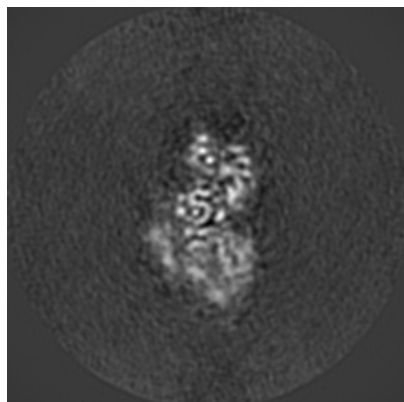


Z

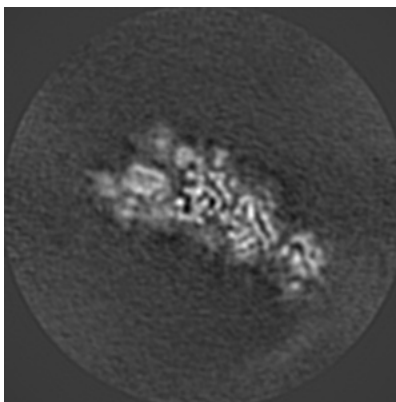
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

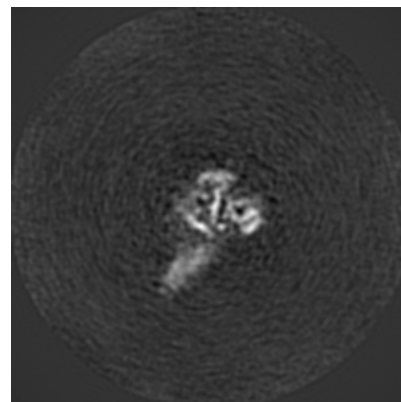
6.2.1 Primary map



X Index: 130

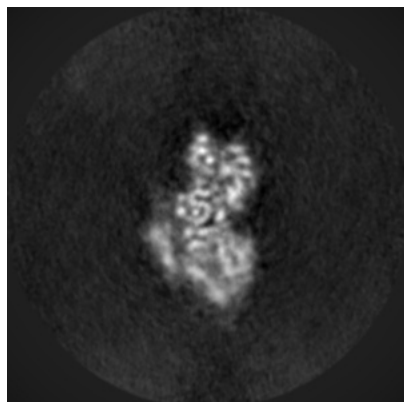


Y Index: 130

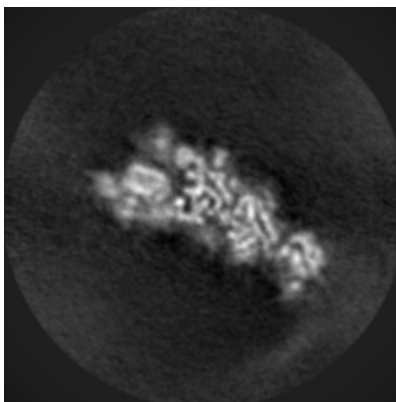


Z Index: 130

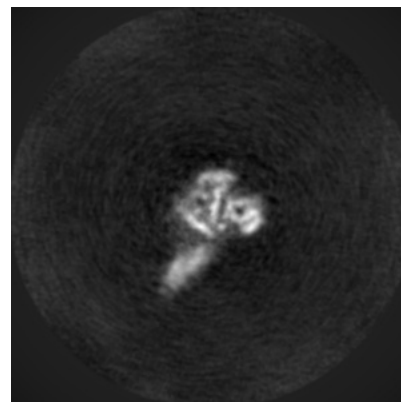
6.2.2 Raw map



X Index: 130



Y Index: 130

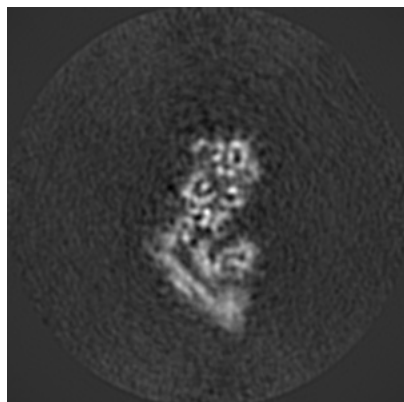


Z Index: 130

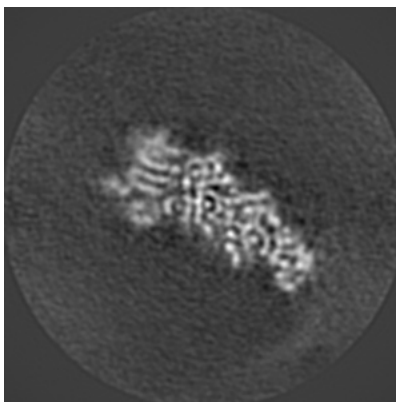
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

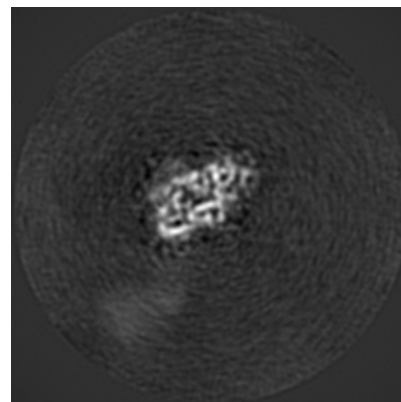
6.3.1 Primary map



X Index: 139

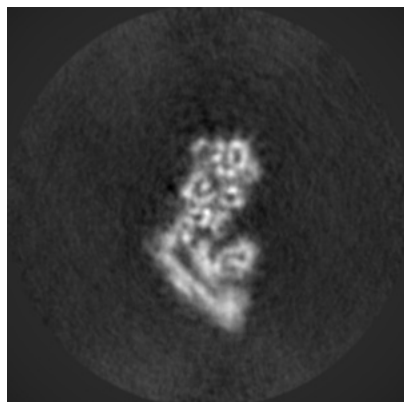


Y Index: 123

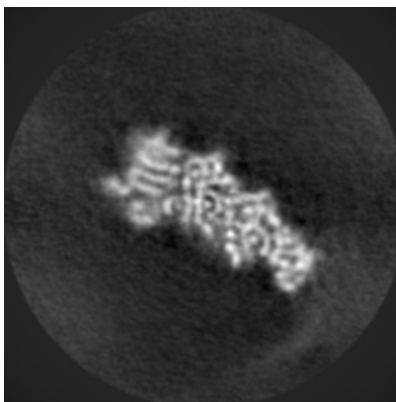


Z Index: 159

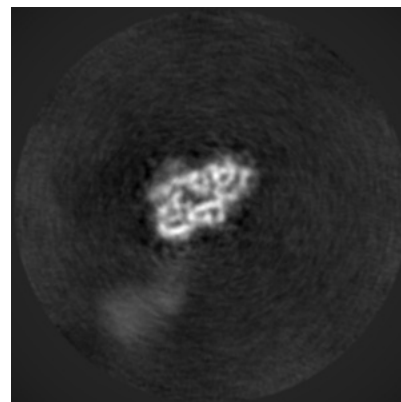
6.3.2 Raw map



X Index: 139



Y Index: 123

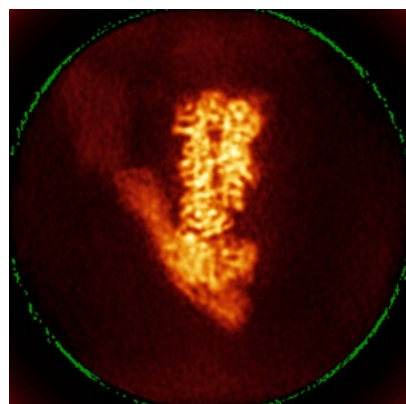


Z Index: 159

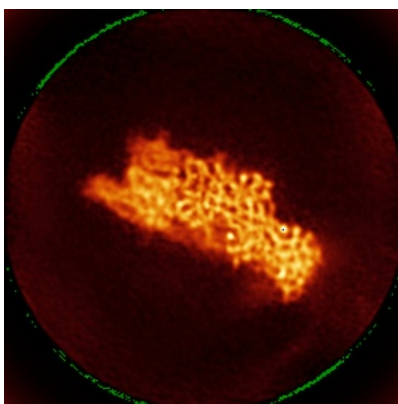
The images above show the largest variance slices of the map in three orthogonal directions.

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

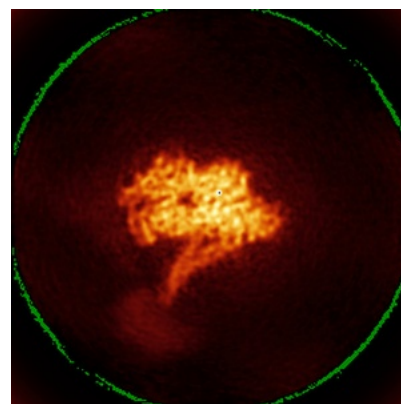
6.4.1 Primary map



X

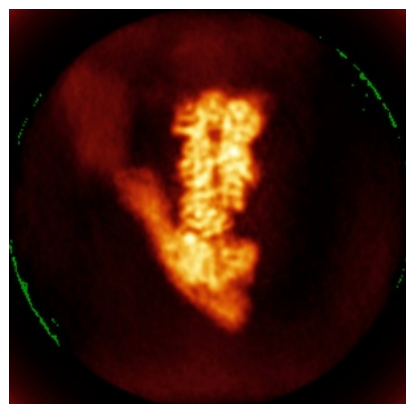


Y

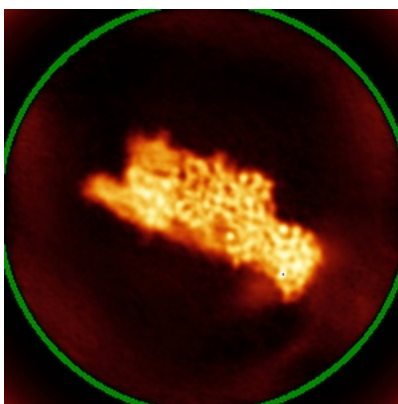


Z

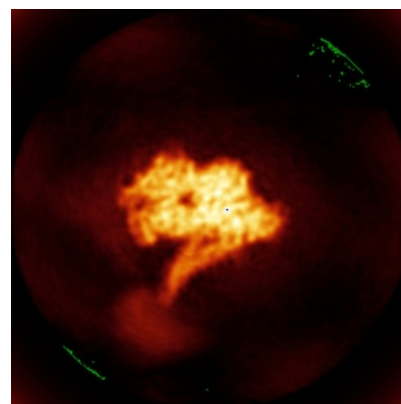
6.4.2 Raw map



X



Y

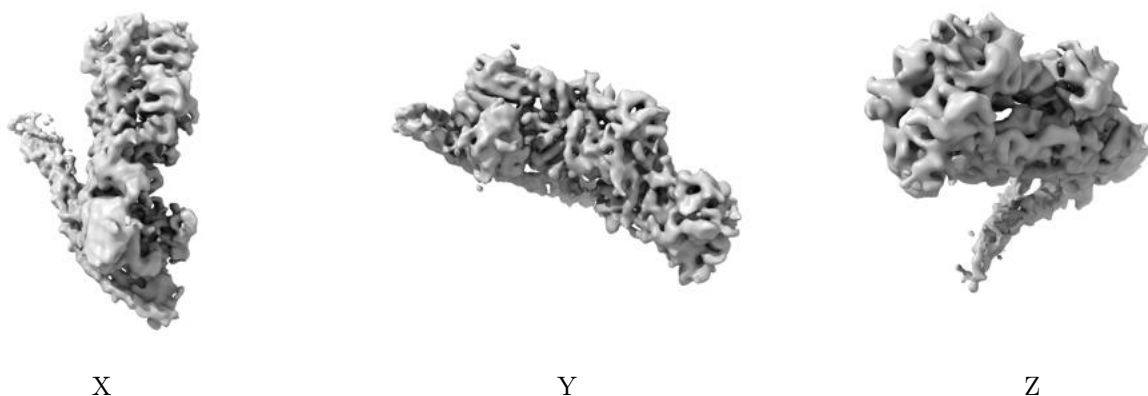


Z

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.

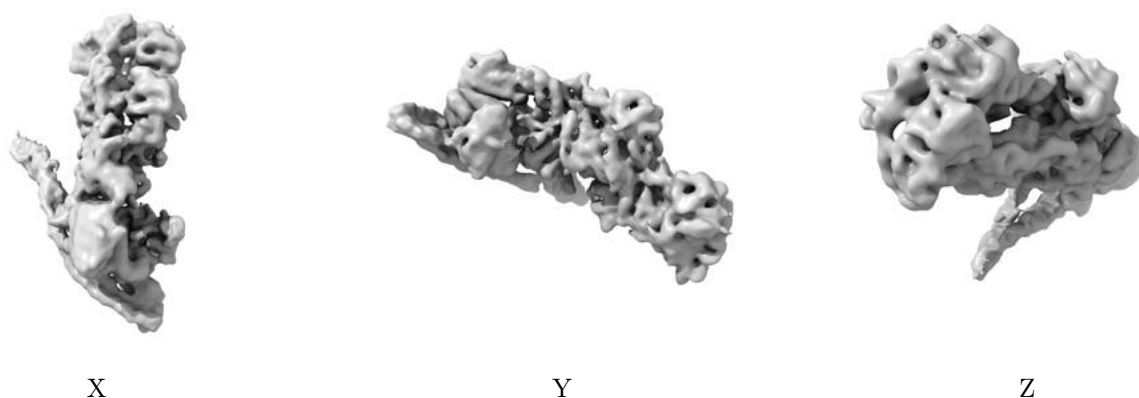
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0492. 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.

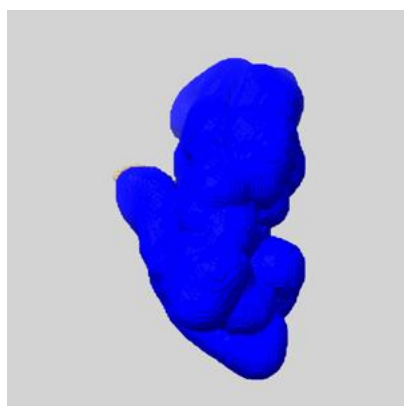
6.6 Mask visualisation [i](#)

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

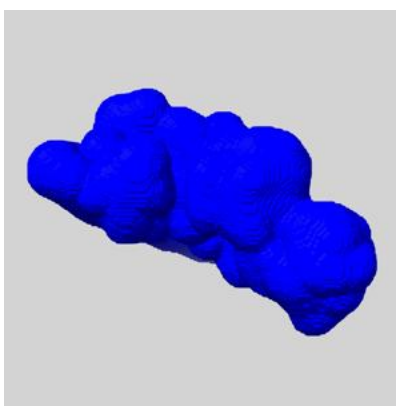
A mask typically either:

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

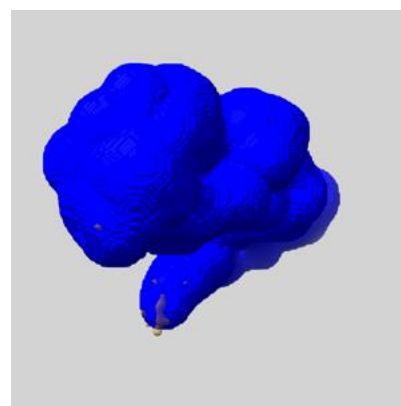
6.6.1 emd_11319_msk_1.map [i](#)



X



Y

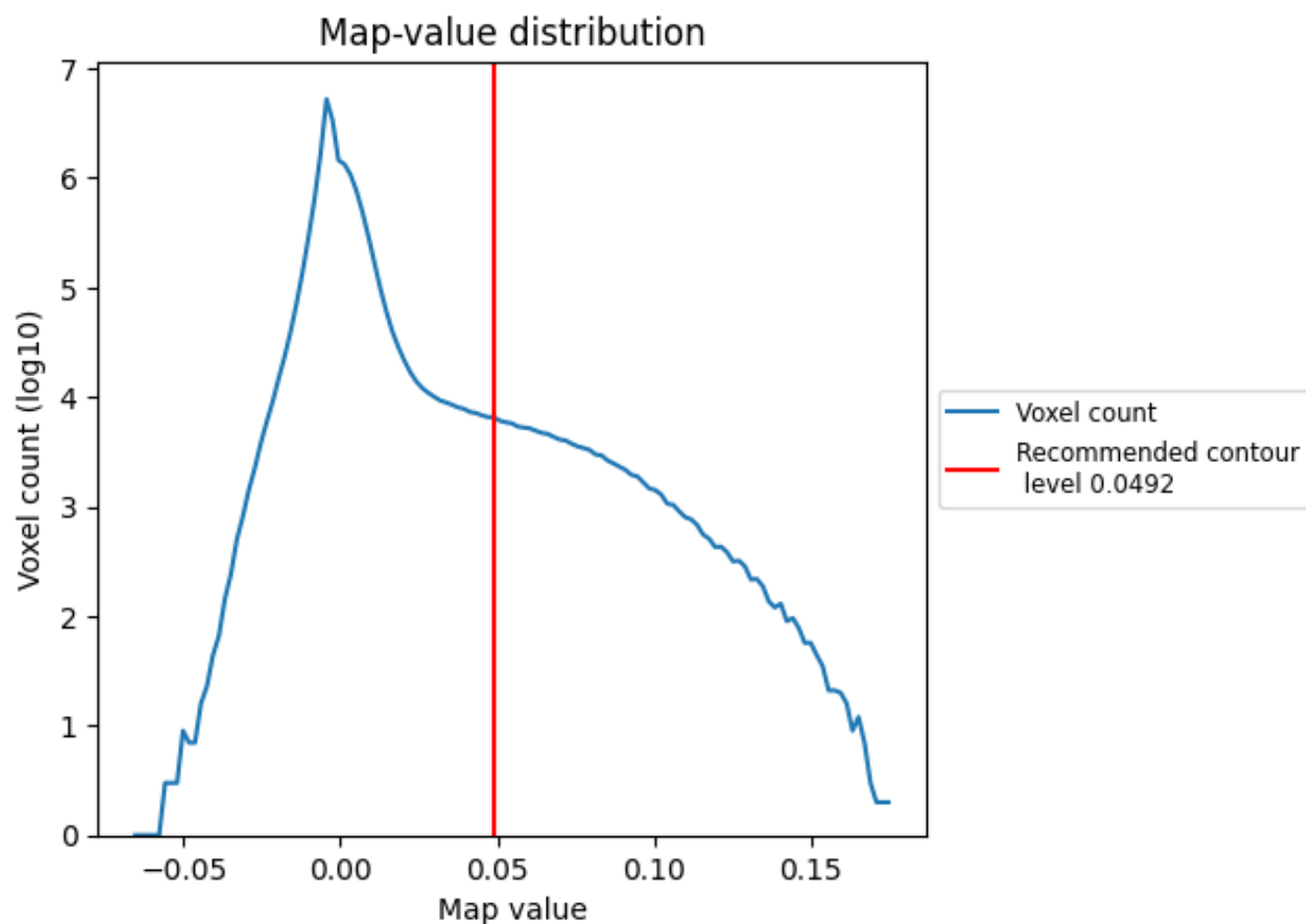


Z

7 Map analysis [i](#)

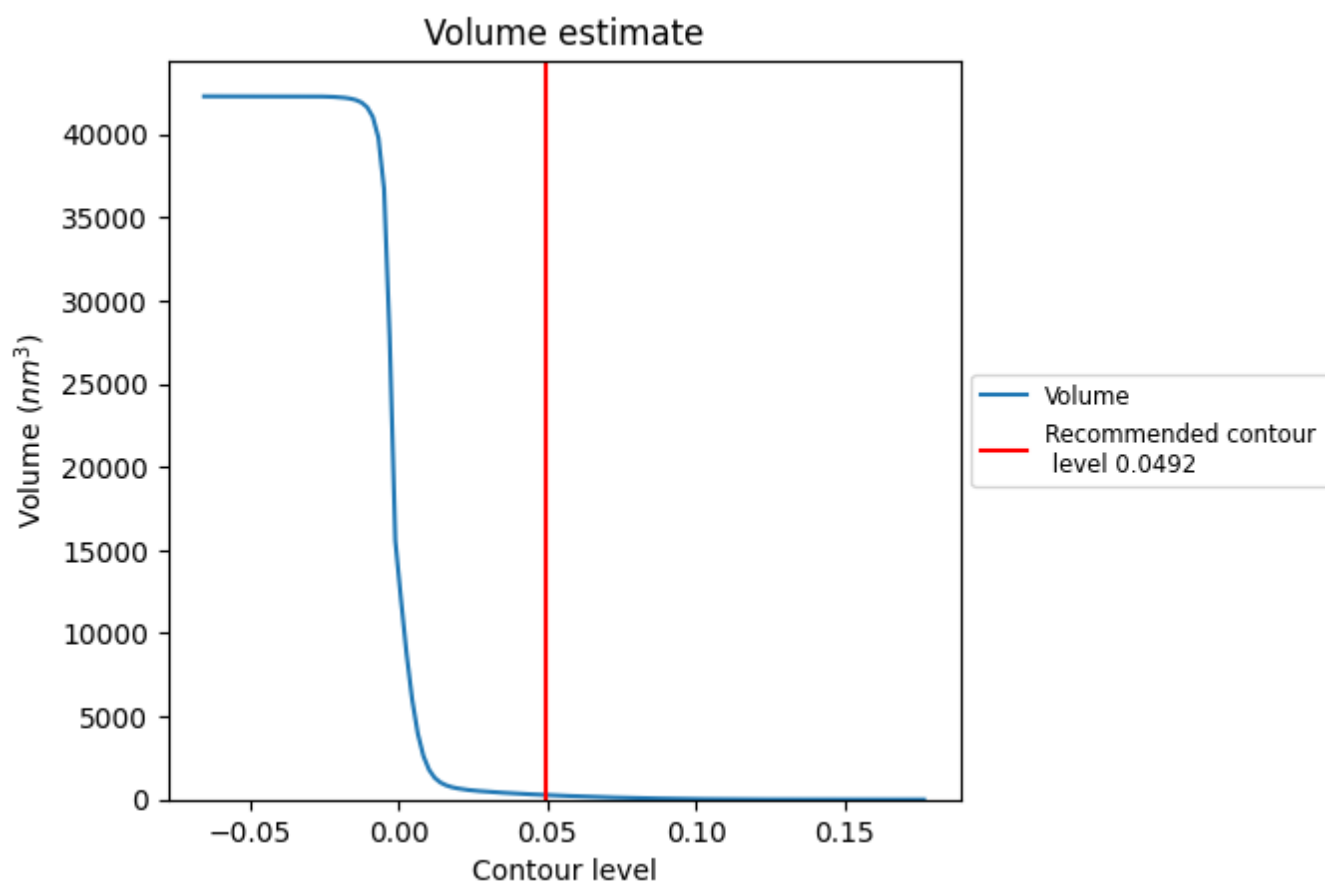
This section contains the results of statistical analysis of the map.

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.

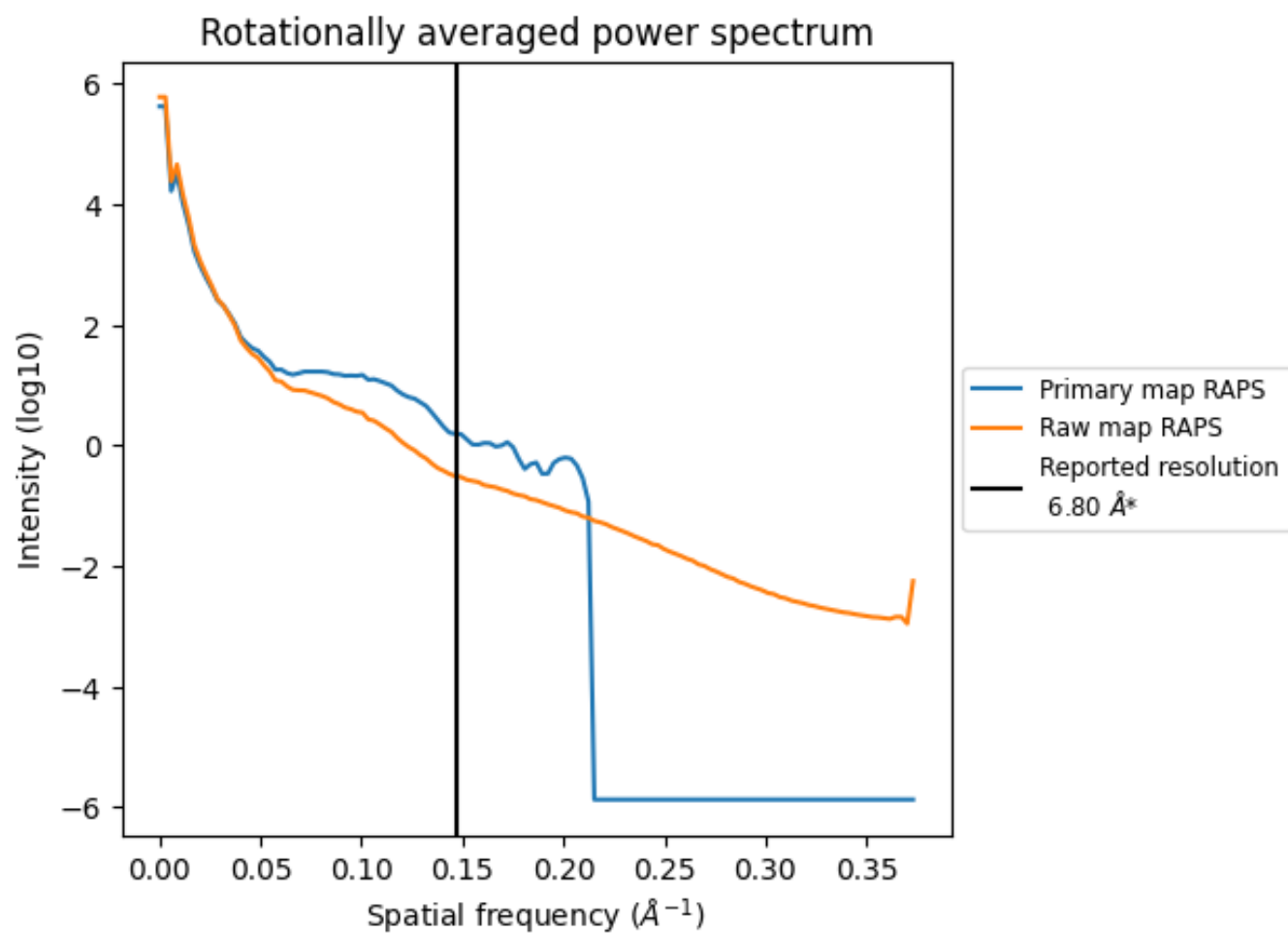
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 277 nm³; this corresponds to an approximate mass of 250 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 ⓘ

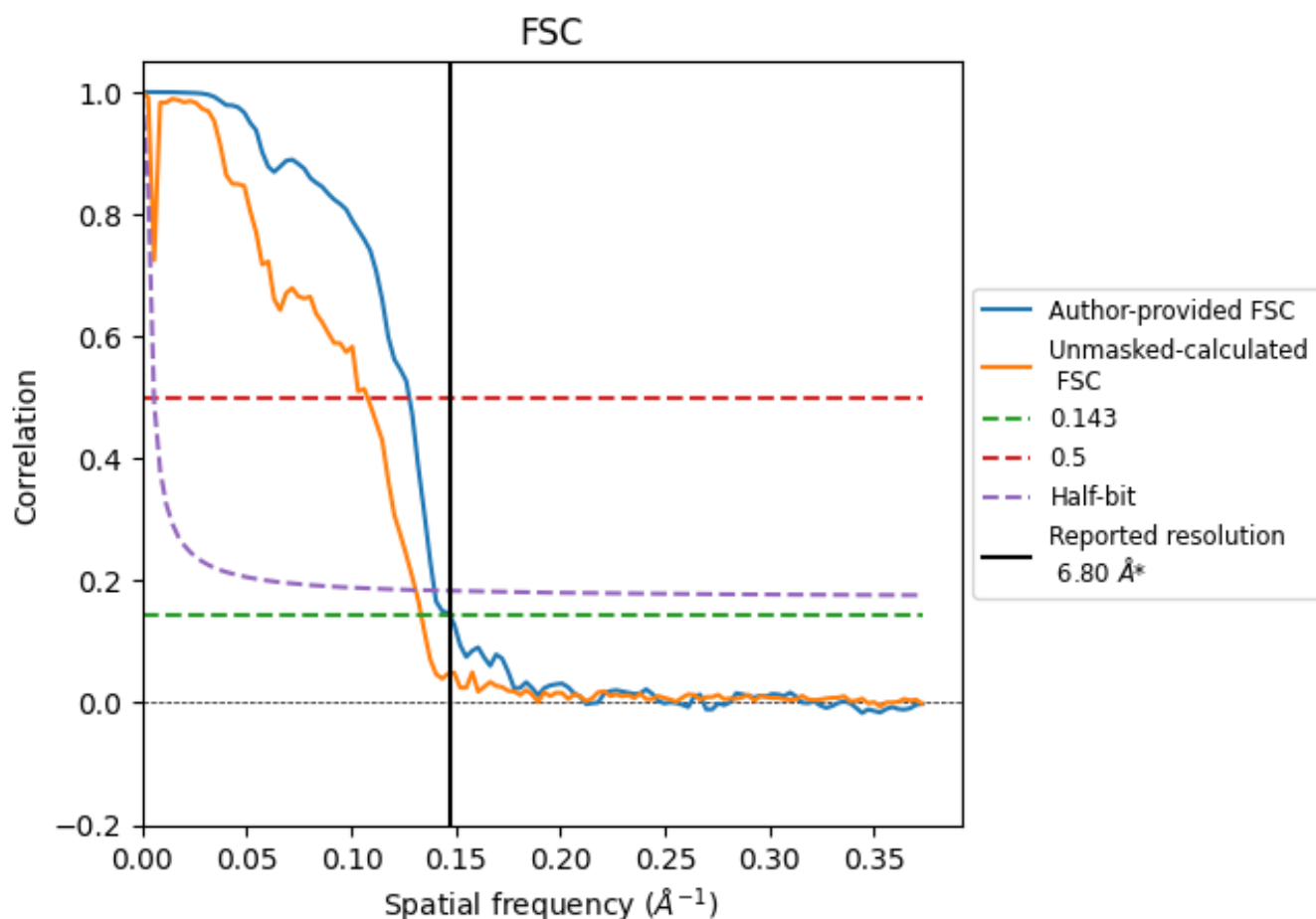


*Reported resolution corresponds to spatial frequency of 0.147 $\text{\AA}^{-1}$

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 $0.147 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

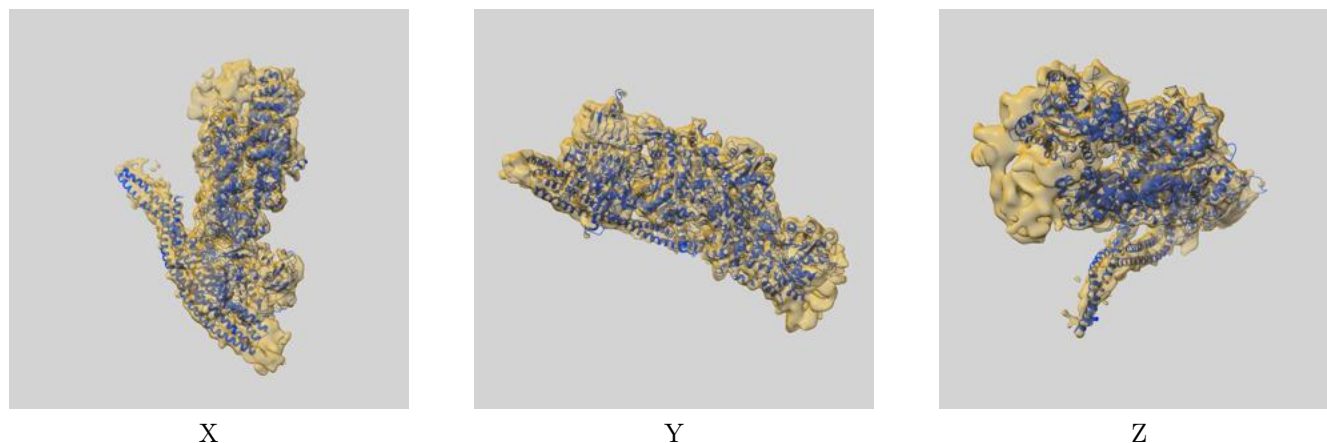
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	6.80	-	-
Author-provided FSC curve	6.80	7.83	7.15
Unmasked-calculated*	7.49	9.29	7.65

*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 7.49 differs from the reported value 6.8 by more than 10 %

9 Map-model fit [i](#)

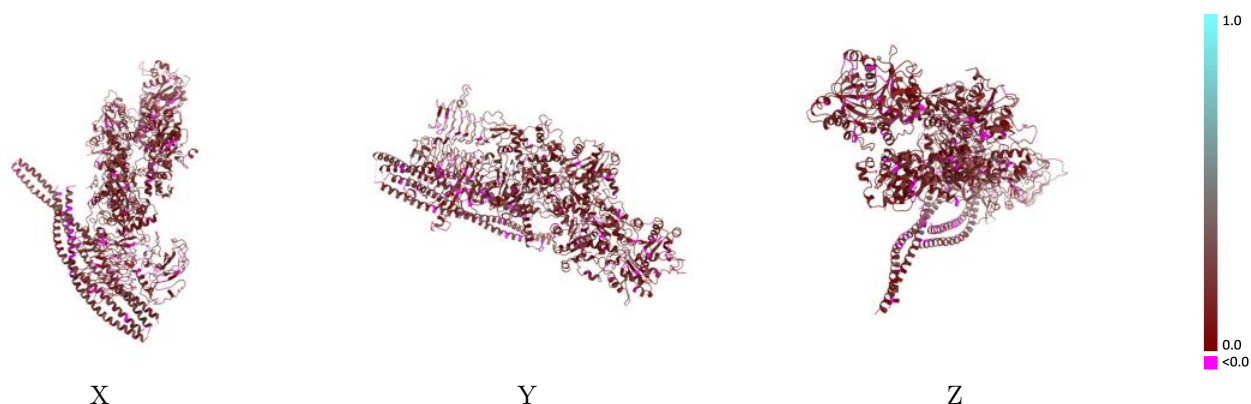
This section contains information regarding the fit between EMDB map EMD-11319 and PDB model 6ZNO. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



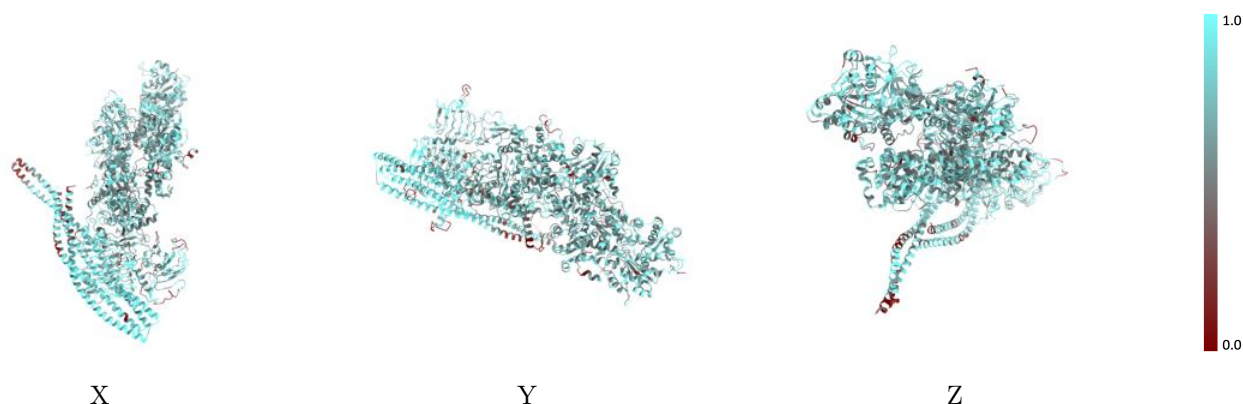
The images above show the 3D surface view of the map at the recommended contour level 0.0492 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



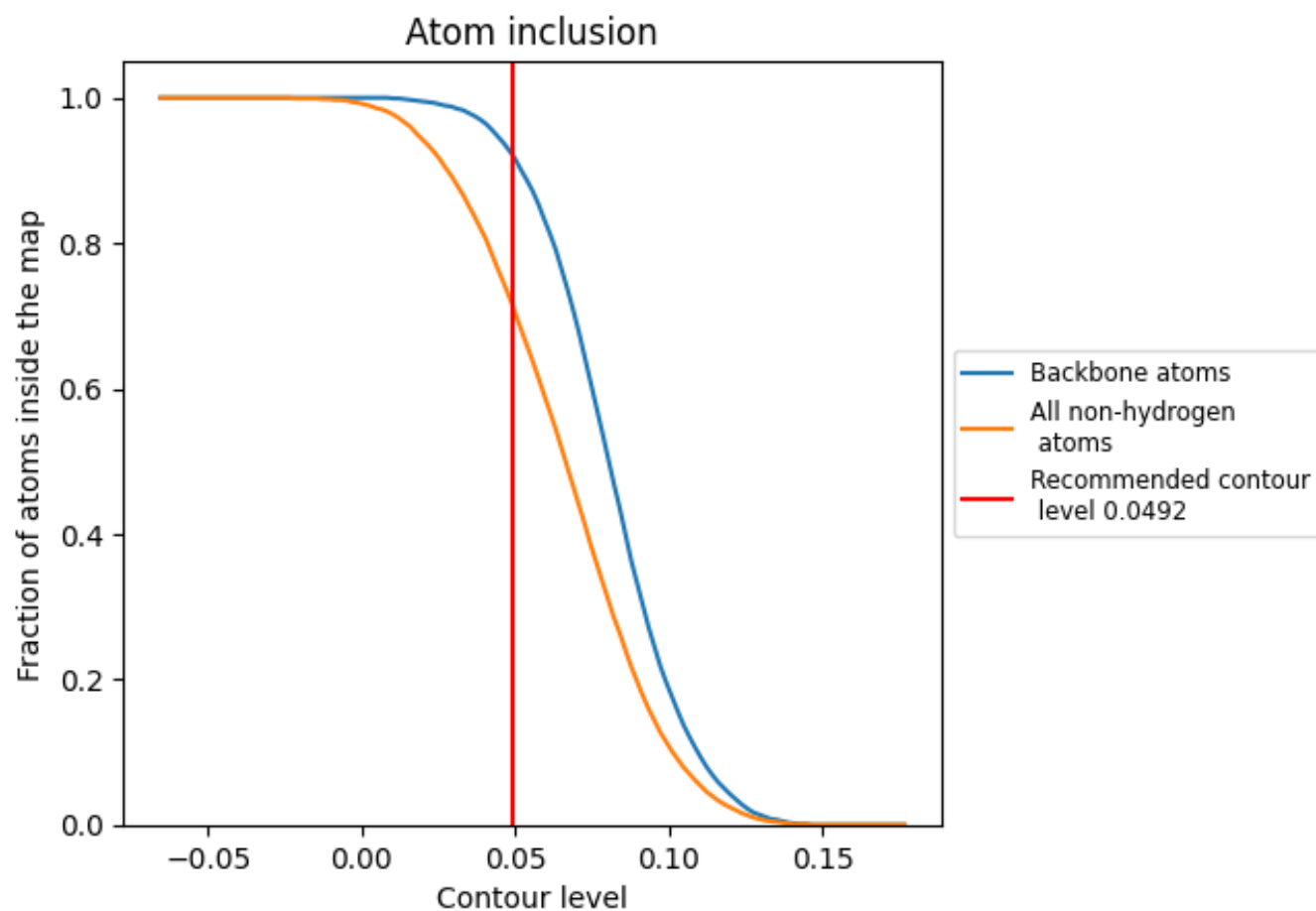
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0492).

9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 71% 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 (0.0492) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7140	0.1580
G	0.7010	0.1430
H	0.7160	0.1600
I	0.7050	0.1550
J	0.6910	0.1720
M	0.6430	0.1710
U	0.7360	0.1400
V	0.7610	0.1630
Y	0.6670	0.1480
Z	0.8330	0.1900
z	0.8220	0.1760

1.0

0.0

<0.0