



wwPDB EM Validation Summary Report ⓘ

Feb 4, 2023 – 08:50 AM EST

PDB ID : 8DD0
EMDB ID : EMD-27331
Title : The structure of the native cardiac thin filament junction region
Authors : Galkin, V.E.; Risi, C.M.
Deposited on : 2022-06-17
Resolution : 3.50 Å (reported)
Based on initial model : 7KO4

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.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.32.1

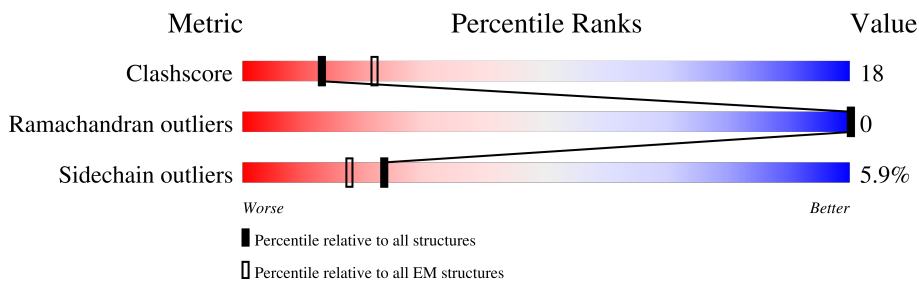
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.











Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	377	79% 17% ..
1	B	377	80% 16% ..
1	C	377	79% 17% ..
1	D	377	79% 17% ..
1	E	377	80% 16% ..
1	F	377	81% 15% ..
2	G	284	12% 7% . 81%
2	H	284	11% 7% . 81%

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Mol	Chain	Length	Quality of chain
2	I	284	 14% 10% • 75%
2	J	284	 16% 8% • 75%
2	L	284	 11% 8% • 81%
2	M	284	 10% 9% • 81%
2	N	284	 17% 7% • 75%
2	O	284	 17% 7% • 75%
3	K	295	 13% 11% • 75%
3	P	295	 11% 11% • 75%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 22892 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 Actin, alpha cardiac muscle 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	371	2898	1836	489	553	20	0	0
1	B	371	2898	1836	489	553	20	0	0
1	C	371	2898	1836	489	553	20	0	0
1	D	371	2898	1836	489	553	20	0	0
1	E	371	2898	1836	489	553	20	0	0
1	F	371	2898	1836	489	553	20	0	0

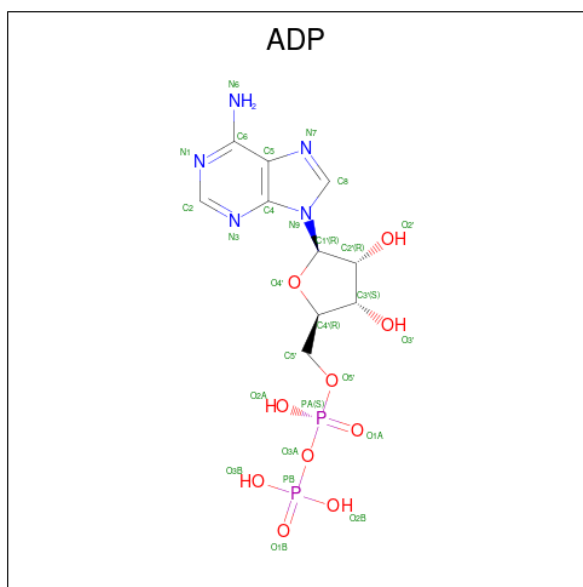
- Molecule 2 is a protein called Tropomyosin alpha-1 chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	G	55	447	277	72	97	1	0	0
2	H	55	447	277	72	97	1	0	0
2	I	70	562	343	98	118	3	0	0
2	J	70	562	343	98	118	3	0	0
2	L	55	447	277	72	97	1	0	0
2	M	55	447	277	72	97	1	0	0
2	N	70	562	343	98	118	3	0	0
2	O	70	562	343	98	118	3	0	0

- Molecule 3 is a protein called Troponin T2, cardiac type.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	K	75	Total	C	N	O	S	0	0
			650	394	132	123	1		
3	P	75	Total	C	N	O	S	0	0
			650	394	132	123	1		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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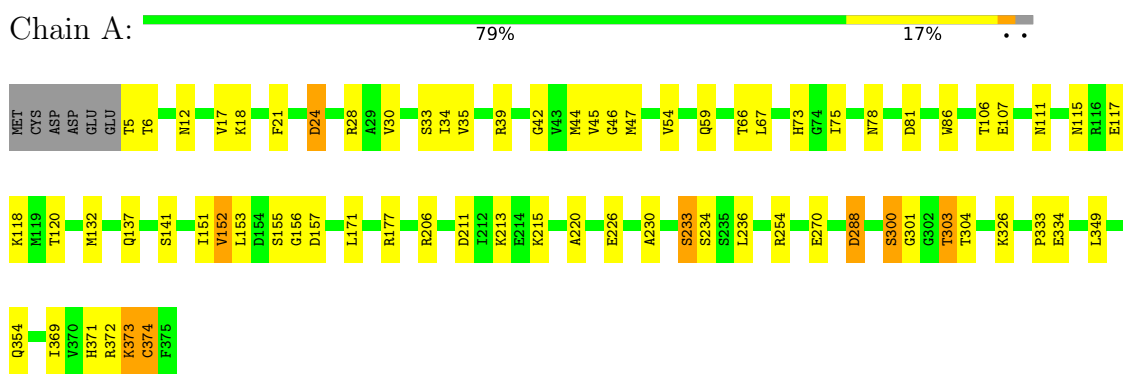
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Mol	Chain	Residues	Atoms		AltConf
5	B	1	Total 1	Mg 1	0
5	C	1	Total 1	Mg 1	0
5	D	1	Total 1	Mg 1	0
5	E	1	Total 1	Mg 1	0
5	F	1	Total 1	Mg 1	0

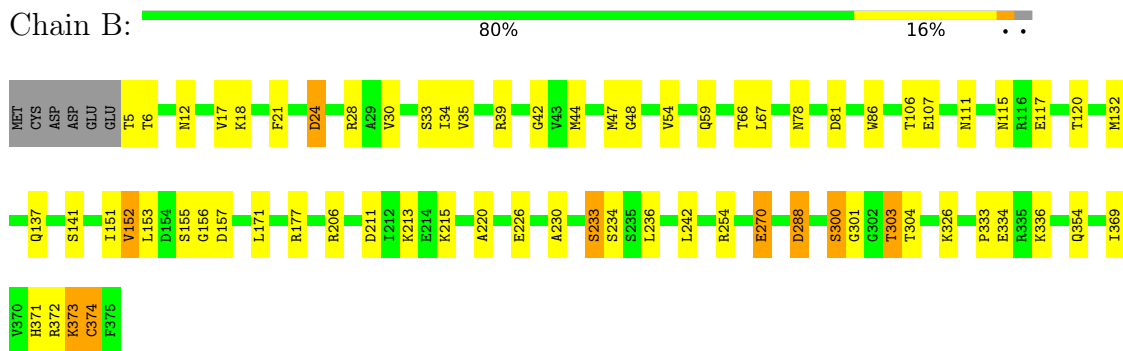
3 Residue-property plots

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.

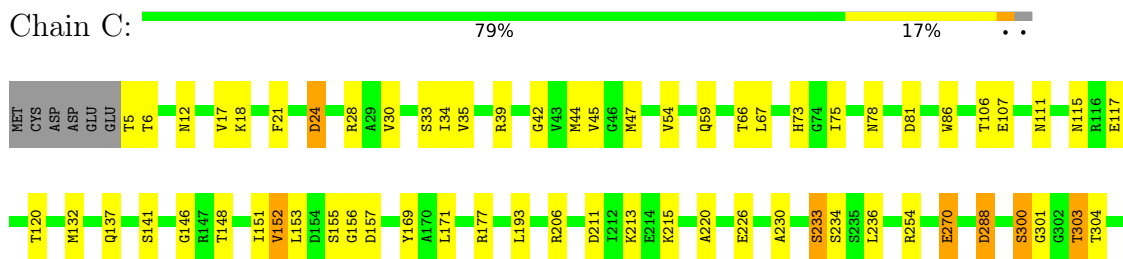
- Molecule 1: Actin, alpha cardiac muscle 1

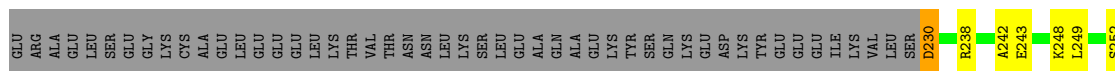


- Molecule 1: Actin, alpha cardiac muscle 1

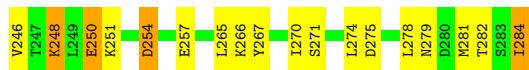
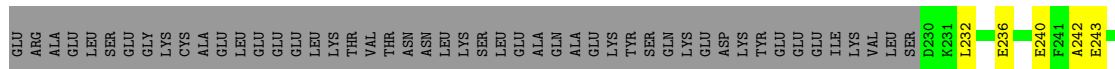
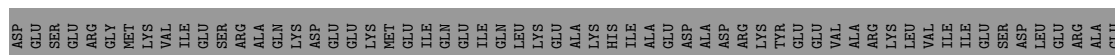
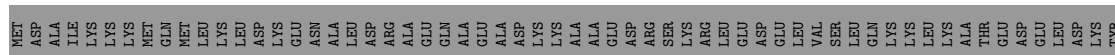


- Molecule 1: Actin, alpha cardiac muscle 1

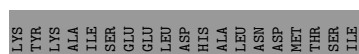
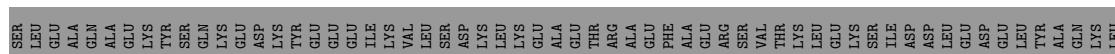
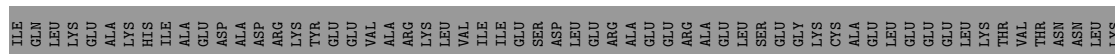
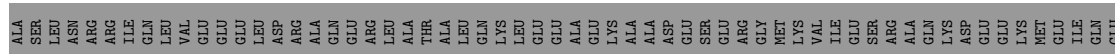
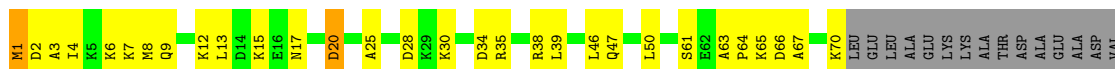




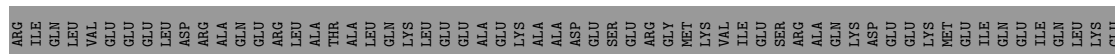
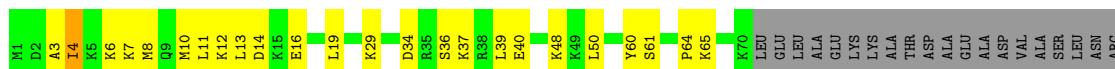
• Molecule 2: Tropomyosin alpha-1 chain

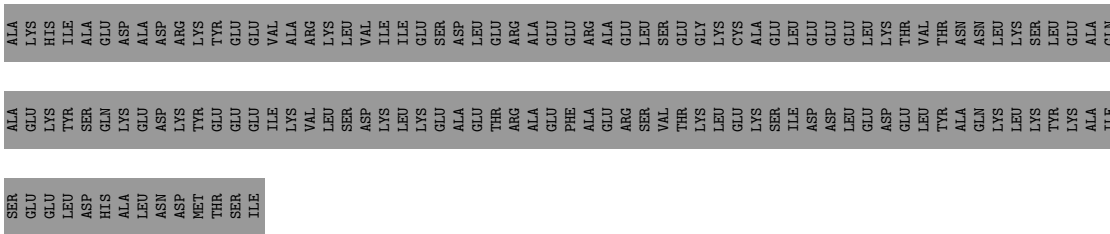


• Molecule 2: Tropomyosin alpha-1 chain



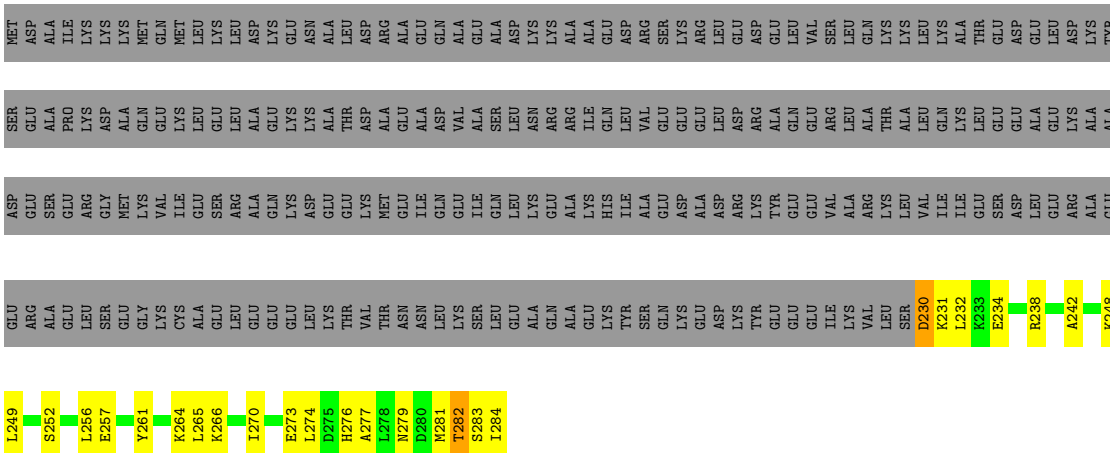
• Molecule 2: Tropomyosin alpha-1 chain





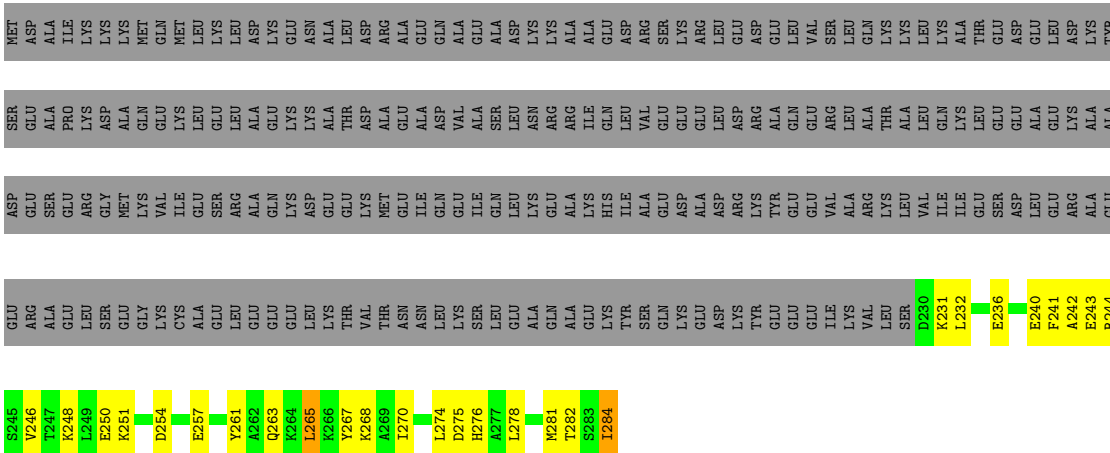
• Molecule 2: Tropomyosin alpha-1 chain

Chain L: 11% 8% . 81%



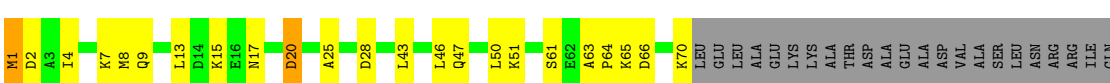
• Molecule 2: Tropomyosin alpha-1 chain

Chain M: 10% 9% . 81%



• Molecule 2: Tropomyosin alpha-1 chain

Chain N: 17% 7% . 75%



GLY	ARG	GLU	ALA	GLU	ASP	GLY	PRO	MET	GLU	SER	LYS	PRO	LYS	PRO	ARG	LEU	PHE	MET	PRO	ASN	LEU	V84	P85	P86	P89	E92	R93	V94	D95	F96	D97	D98	I99	H100	R101	K102	M103	M104	K105	D106	L107	L113	I114	E115	A116	H117	F118	E119	M120	A121	K122	K123	E124	E125	E126									
L127	V128	S129	L130	K131	D132	R133	I134	E135	R138	A139	E140	R141	Q145	R146	T149	E150	K153	E154	R155	Q156	T157	R158	LEU	ALA	GLU	GLU	ARG	ARG	ALA	ARG	VAL	LEU	ALA	ILE	ASP	HIS	GLU	ASN	ASN	ARG	ARG	LYS	LYS	ALA	ASP	GLN	LEU	LEU	ARG	GLU	LYS	ALA	ARG	LYS	GLU	TRP	LEU	SER	ILE	TYR	ASN	MET	HIS	PHE
ALA	GLU	LYS	PHE	ASP	LEU	GLN	GLU	LYS	PHE	GLN	LYS	GLN	ILE	ASN	VAL	LEU	ARG	ASN	ASN	ARG	ILE	ASN	ASP	ASN	GLN	LYS	VAL	SER	THR	ARG	GLY	LYS	ALA	LYS	VAL	THR	GLY	TRP	LYS																									

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	236359	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 ($e^-/\text{\AA}^2$)	34	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	10.091	Depositor
Minimum map value	-2.821	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	0.02	Depositor
Map size (Å)	97.631996, 108.479996, 225.096	wwPDB
Map dimensions	72, 80, 166	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.356, 1.356, 1.356	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, MG

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	0.43	0/2961	0.48	0/4011
1	B	0.43	0/2961	0.48	0/4011
1	C	0.43	0/2961	0.48	0/4011
1	D	0.43	0/2961	0.48	0/4011
1	E	0.43	0/2961	0.48	0/4011
1	F	0.43	0/2961	0.48	0/4011
2	G	0.38	0/450	0.65	0/600
2	H	0.40	0/450	0.71	0/600
2	I	0.34	0/563	0.59	0/744
2	J	0.34	0/563	0.61	0/744
2	L	0.39	0/450	0.68	0/600
2	M	0.40	0/450	0.71	0/600
2	N	0.36	0/563	0.65	0/744
2	O	0.35	0/563	0.63	0/744
3	K	0.37	0/656	0.86	1/871 (0.1%)
3	P	0.38	0/656	0.88	1/871 (0.1%)
All	All	0.42	0/23130	0.54	2/31184 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	93	ARG	C-N-CA	6.50	137.96	121.70
3	K	93	ARG	C-N-CA	5.99	136.68	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2898	0	2871	67	0
1	B	2898	0	2871	64	0
1	C	2898	0	2871	95	0
1	D	2898	0	2871	91	0
1	E	2898	0	2871	65	0
1	F	2898	0	2871	60	0
2	G	447	0	441	52	0
2	H	447	0	441	44	0
2	I	562	0	580	67	0
2	J	562	0	580	70	0
2	L	447	0	441	67	0
2	M	447	0	441	73	0
2	N	562	0	580	38	0
2	O	562	0	580	66	0
3	K	650	0	661	117	0
3	P	650	0	661	128	0
4	A	27	0	12	2	0
4	B	27	0	12	2	0
4	C	27	0	12	2	0
4	D	27	0	12	2	0
4	E	27	0	12	3	0
4	F	27	0	12	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
All	All	22892	0	22704	839	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 18.

The worst 5 of 839 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:G:284:ILE:CG2	2:J:12:LYS:CE	1.76	1.57
2:G:281:MET:SD	2:I:7:LYS:HE3	1.44	1.56
2:L:281:MET:SD	2:N:7:LYS:HE3	1.44	1.56
2:G:284:ILE:HG21	2:J:12:LYS:CE	1.35	1.45
2:G:284:ILE:CG2	2:J:12:LYS:HE3	0.98	1.43

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	369/377 (98%)	351 (95%)	18 (5%)	0	100	100
1	B	369/377 (98%)	350 (95%)	19 (5%)	0	100	100
1	C	369/377 (98%)	350 (95%)	19 (5%)	0	100	100
1	D	369/377 (98%)	351 (95%)	18 (5%)	0	100	100
1	E	369/377 (98%)	351 (95%)	18 (5%)	0	100	100
1	F	369/377 (98%)	350 (95%)	19 (5%)	0	100	100
2	G	53/284 (19%)	50 (94%)	3 (6%)	0	100	100
2	H	53/284 (19%)	51 (96%)	2 (4%)	0	100	100
2	I	68/284 (24%)	68 (100%)	0	0	100	100
2	J	68/284 (24%)	68 (100%)	0	0	100	100
2	L	53/284 (19%)	50 (94%)	3 (6%)	0	100	100
2	M	53/284 (19%)	51 (96%)	2 (4%)	0	100	100
2	N	68/284 (24%)	68 (100%)	0	0	100	100
2	O	68/284 (24%)	68 (100%)	0	0	100	100
3	K	73/295 (25%)	69 (94%)	4 (6%)	0	100	100
3	P	73/295 (25%)	70 (96%)	3 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2844/5124 (56%)	2716 (96%)	128 (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	A	314/320 (98%)	298 (95%)	16 (5%)	24	57
1	B	314/320 (98%)	298 (95%)	16 (5%)	24	57
1	C	314/320 (98%)	298 (95%)	16 (5%)	24	57
1	D	314/320 (98%)	298 (95%)	16 (5%)	24	57
1	E	314/320 (98%)	298 (95%)	16 (5%)	24	57
1	F	314/320 (98%)	297 (95%)	17 (5%)	22	55
2	G	49/245 (20%)	46 (94%)	3 (6%)	18	51
2	H	49/245 (20%)	42 (86%)	7 (14%)	3	19
2	I	60/245 (24%)	55 (92%)	5 (8%)	11	40
2	J	60/245 (24%)	58 (97%)	2 (3%)	38	68
2	L	49/245 (20%)	46 (94%)	3 (6%)	18	51
2	M	49/245 (20%)	44 (90%)	5 (10%)	7	32
2	N	60/245 (24%)	55 (92%)	5 (8%)	11	40
2	O	60/245 (24%)	58 (97%)	2 (3%)	38	68
3	K	71/257 (28%)	67 (94%)	4 (6%)	21	54
3	P	71/257 (28%)	59 (83%)	12 (17%)	2	12
All	All	2462/4394 (56%)	2317 (94%)	145 (6%)	23	53

5 of 145 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	K	106	ASP

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Mol	Chain	Res	Type
3	P	145	GLN
2	L	273	GLU
2	N	46	LEU
1	D	24	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
2	N	47	GLN
3	P	117	HIS
3	K	100	HIS
3	K	117	HIS
3	K	144	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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
4	ADP	A	401	5	24,29,29	0.97	1 (4%)	29,45,45	1.46	4 (13%)
4	ADP	B	401	5	24,29,29	0.97	1 (4%)	29,45,45	1.47	4 (13%)
4	ADP	D	401	5	24,29,29	0.97	1 (4%)	29,45,45	1.46	4 (13%)
4	ADP	E	401	5	24,29,29	0.96	1 (4%)	29,45,45	1.46	4 (13%)
4	ADP	F	401	5	24,29,29	0.96	1 (4%)	29,45,45	1.46	4 (13%)
4	ADP	C	401	5	24,29,29	0.96	1 (4%)	29,45,45	1.46	4 (13%)

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
4	ADP	A	401	5	-	3/12/32/32	0/3/3/3
4	ADP	B	401	5	-	3/12/32/32	0/3/3/3
4	ADP	D	401	5	-	3/12/32/32	0/3/3/3
4	ADP	E	401	5	-	3/12/32/32	0/3/3/3
4	ADP	F	401	5	-	3/12/32/32	0/3/3/3
4	ADP	C	401	5	-	3/12/32/32	0/3/3/3

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	401	ADP	C5-C4	2.18	1.46	1.40
4	D	401	ADP	C5-C4	2.18	1.46	1.40
4	E	401	ADP	C5-C4	2.17	1.46	1.40
4	C	401	ADP	C5-C4	2.17	1.46	1.40
4	B	401	ADP	C5-C4	2.16	1.46	1.40

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	401	ADP	PA-O3A-PB	-3.58	120.54	132.83
4	A	401	ADP	PA-O3A-PB	-3.58	120.56	132.83
4	C	401	ADP	PA-O3A-PB	-3.57	120.56	132.83
4	F	401	ADP	PA-O3A-PB	-3.57	120.59	132.83
4	B	401	ADP	PA-O3A-PB	-3.56	120.60	132.83

There are no chirality outliers.

5 of 18 torsion outliers are listed below:

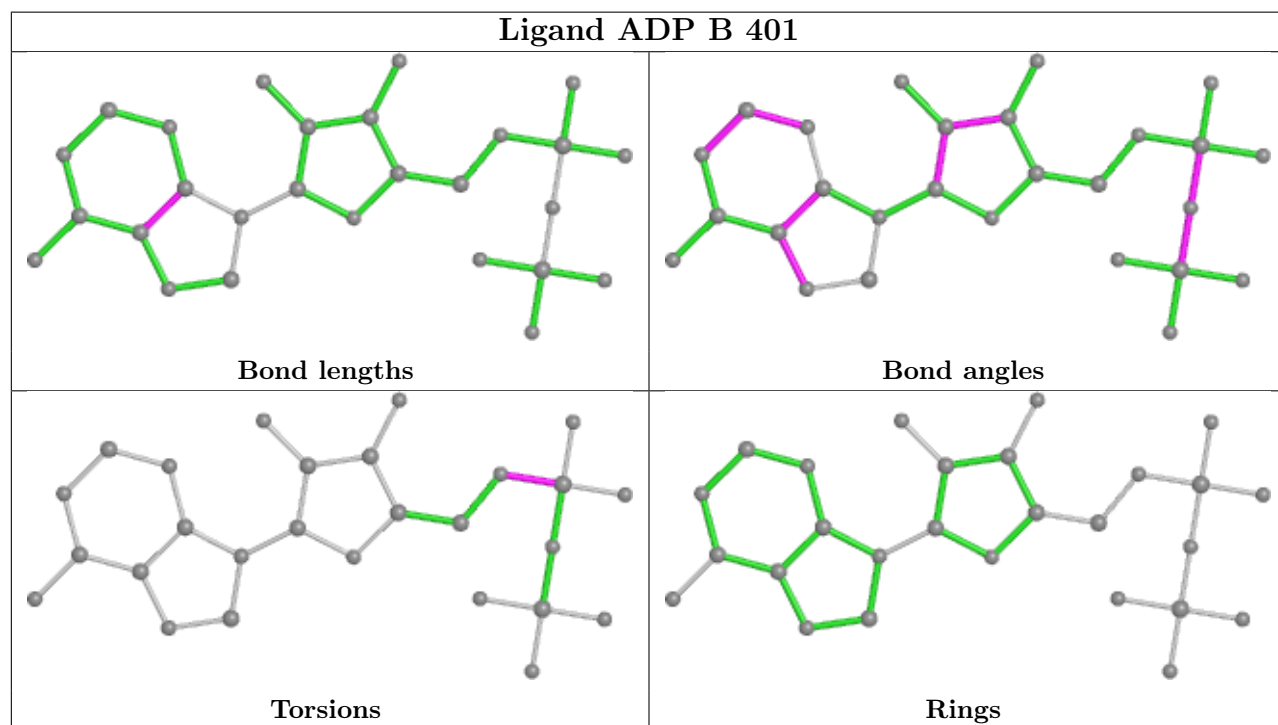
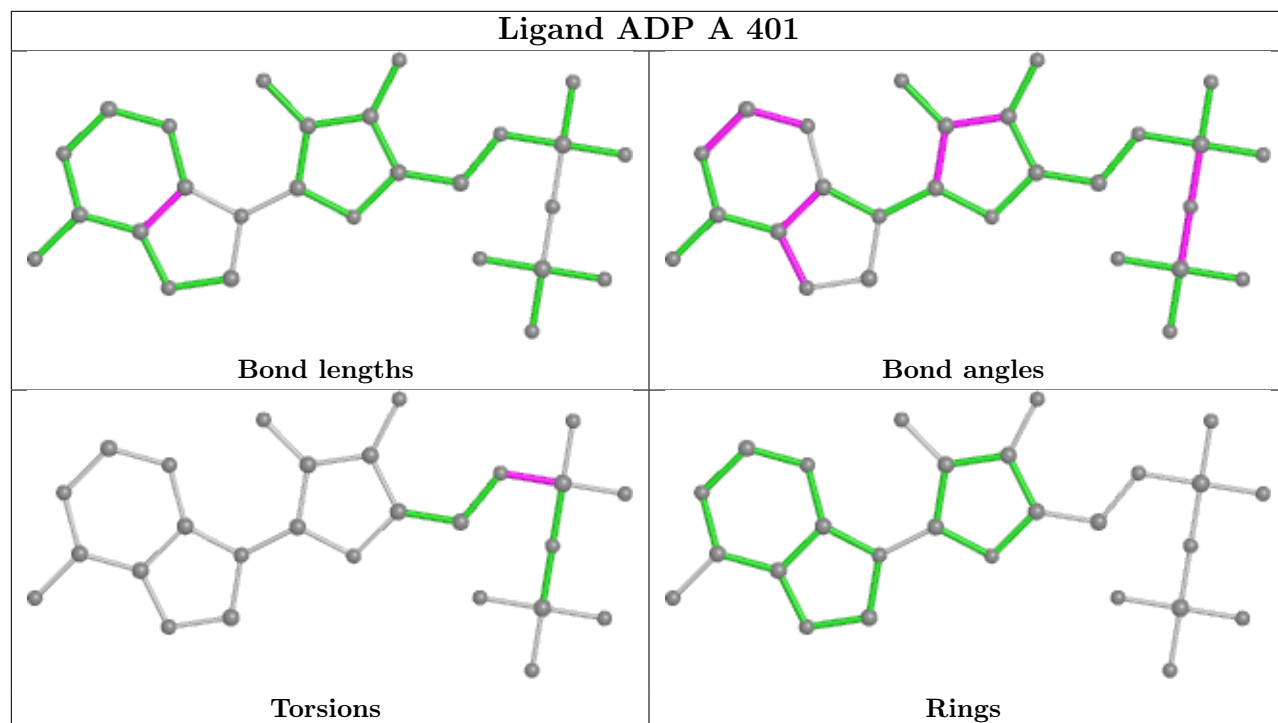
Mol	Chain	Res	Type	Atoms
4	A	401	ADP	C5'-O5'-PA-O1A
4	A	401	ADP	C5'-O5'-PA-O2A
4	B	401	ADP	C5'-O5'-PA-O1A
4	B	401	ADP	C5'-O5'-PA-O2A
4	C	401	ADP	C5'-O5'-PA-O1A

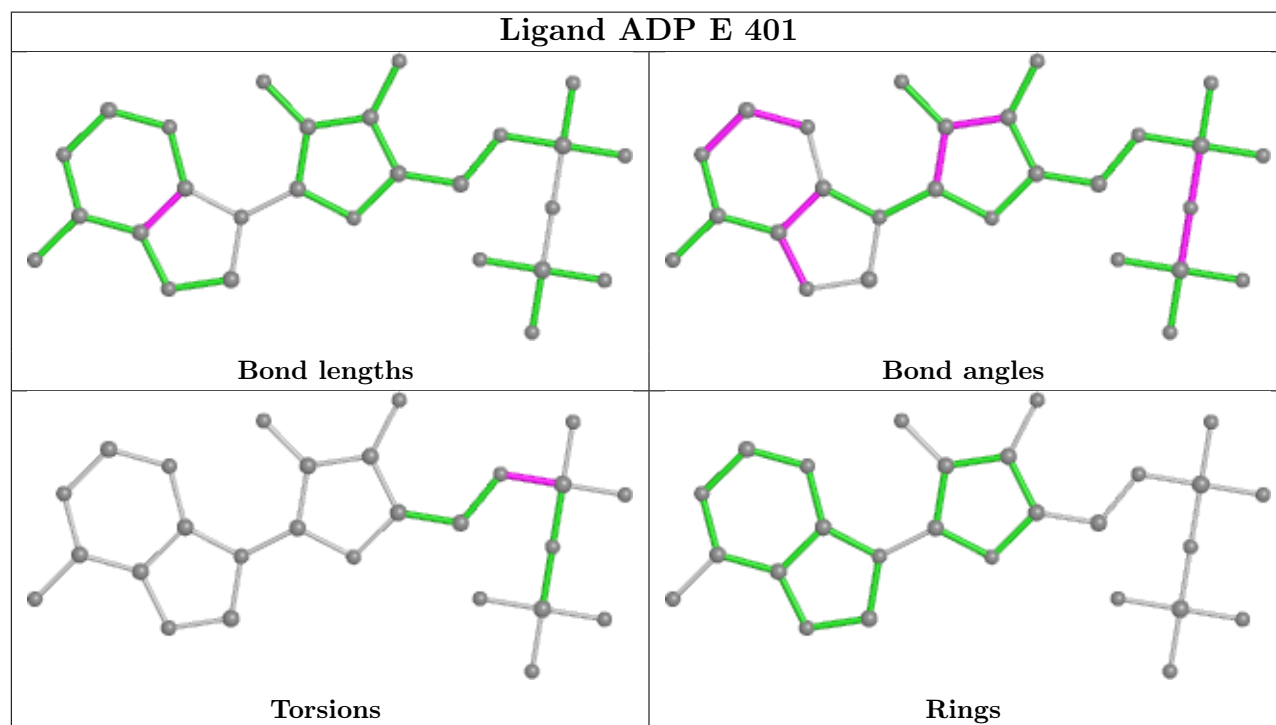
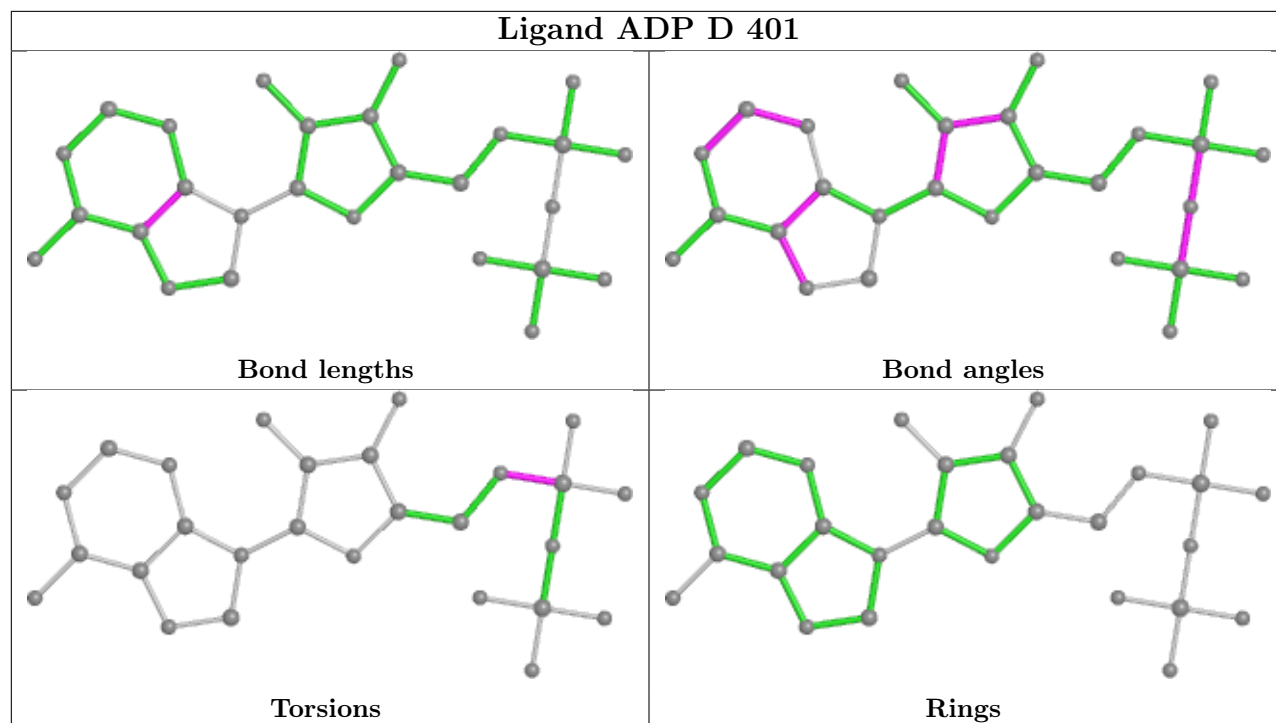
There are no ring outliers.

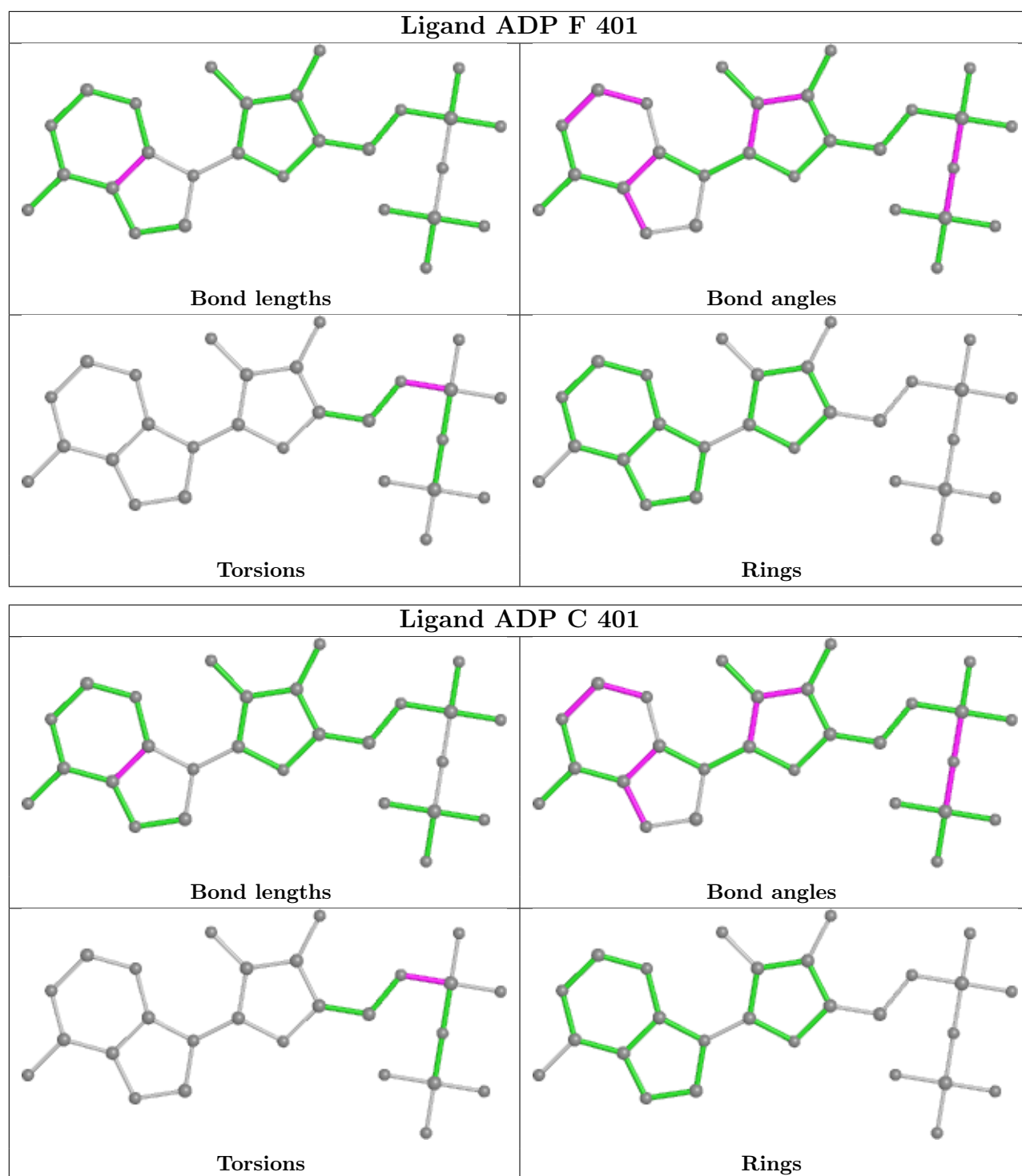
6 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	ADP	2	0
4	B	401	ADP	2	0
4	D	401	ADP	2	0
4	E	401	ADP	3	0
4	F	401	ADP	2	0
4	C	401	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

There are no chain breaks in this entry.

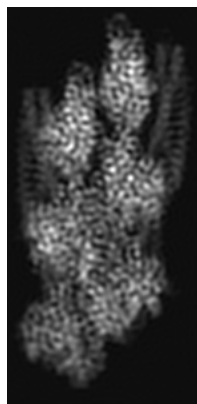
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27331. These allow visual inspection of the internal detail of the map and identification of artifacts.

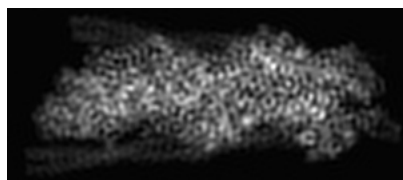
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

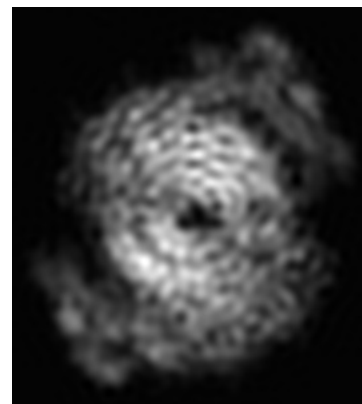
6.1.1 Primary map



X

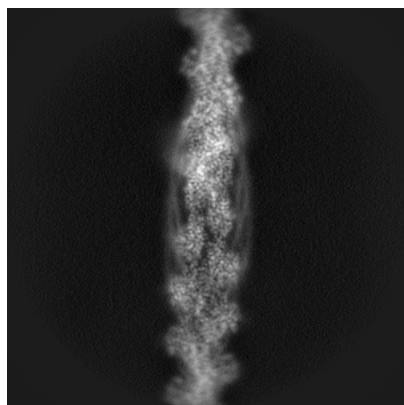


Y

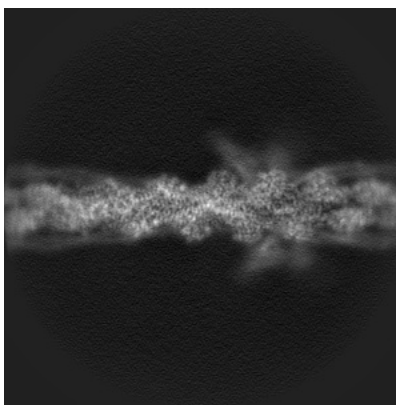


Z

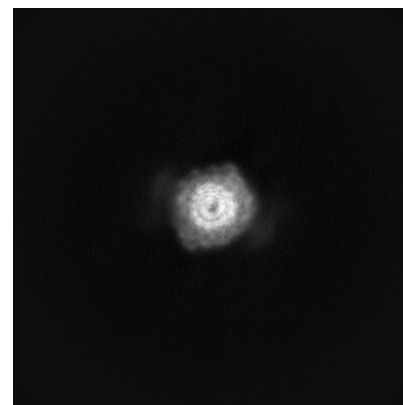
6.1.2 Raw map



X



Y

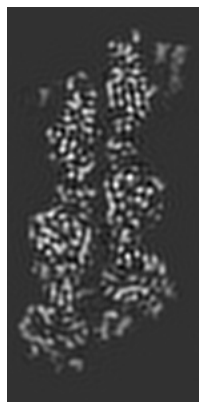


Z

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

6.2 Central slices [i](#)

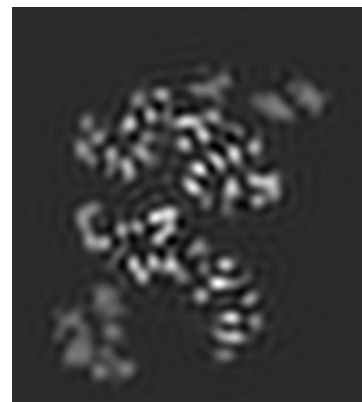
6.2.1 Primary map



X Index: 36

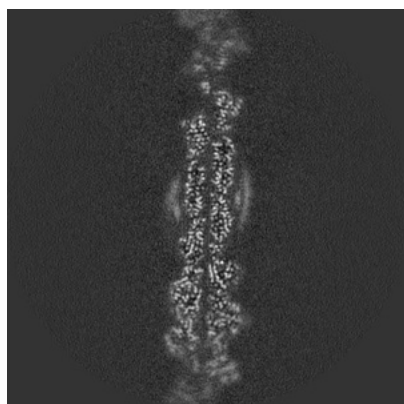


Y Index: 40

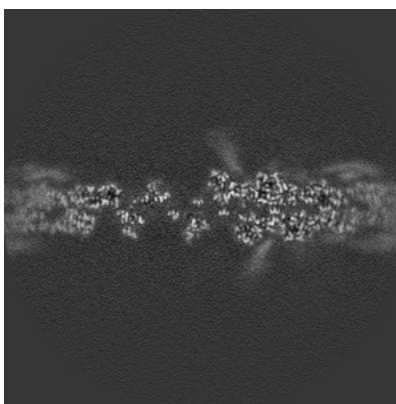


Z Index: 83

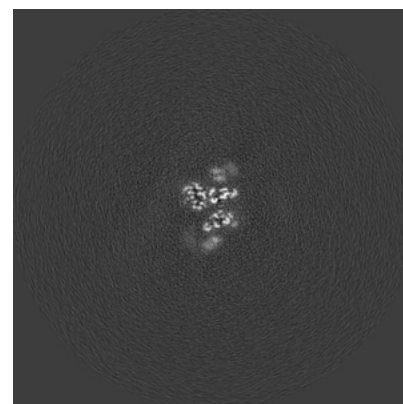
6.2.2 Raw map



X Index: 162



Y Index: 162

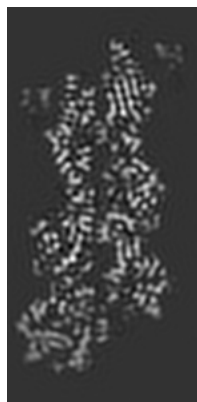


Z Index: 162

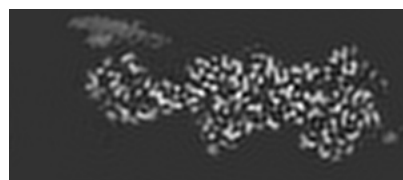
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

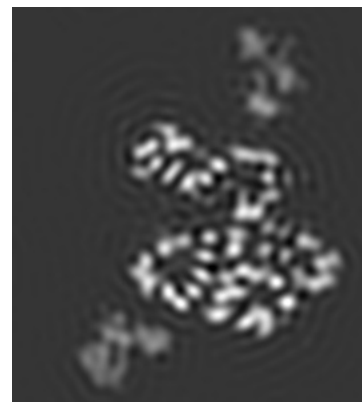
6.3.1 Primary map



X Index: 34

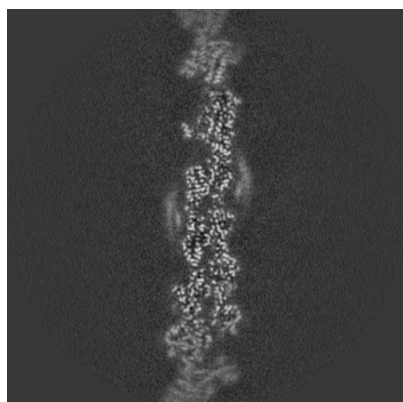


Y Index: 50

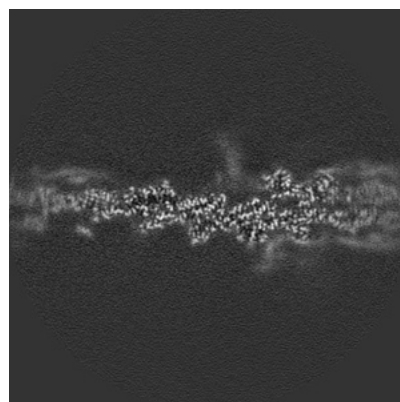


Z Index: 107

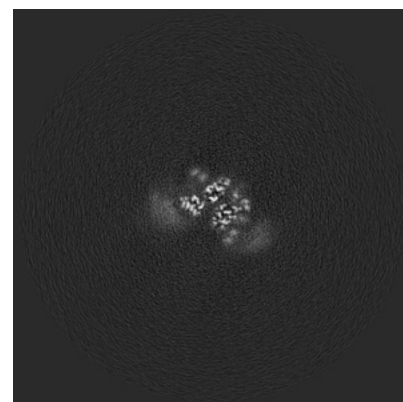
6.3.2 Raw map



X Index: 158



Y Index: 169



Z Index: 206

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

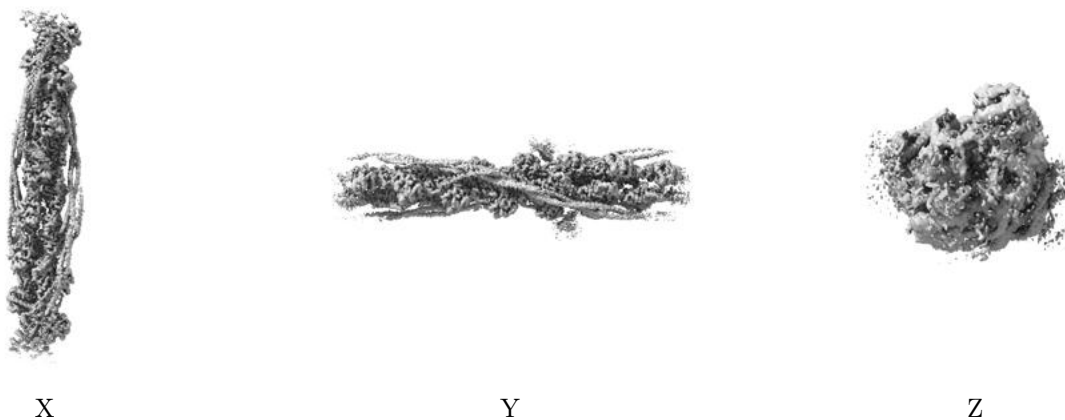
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.02. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 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.

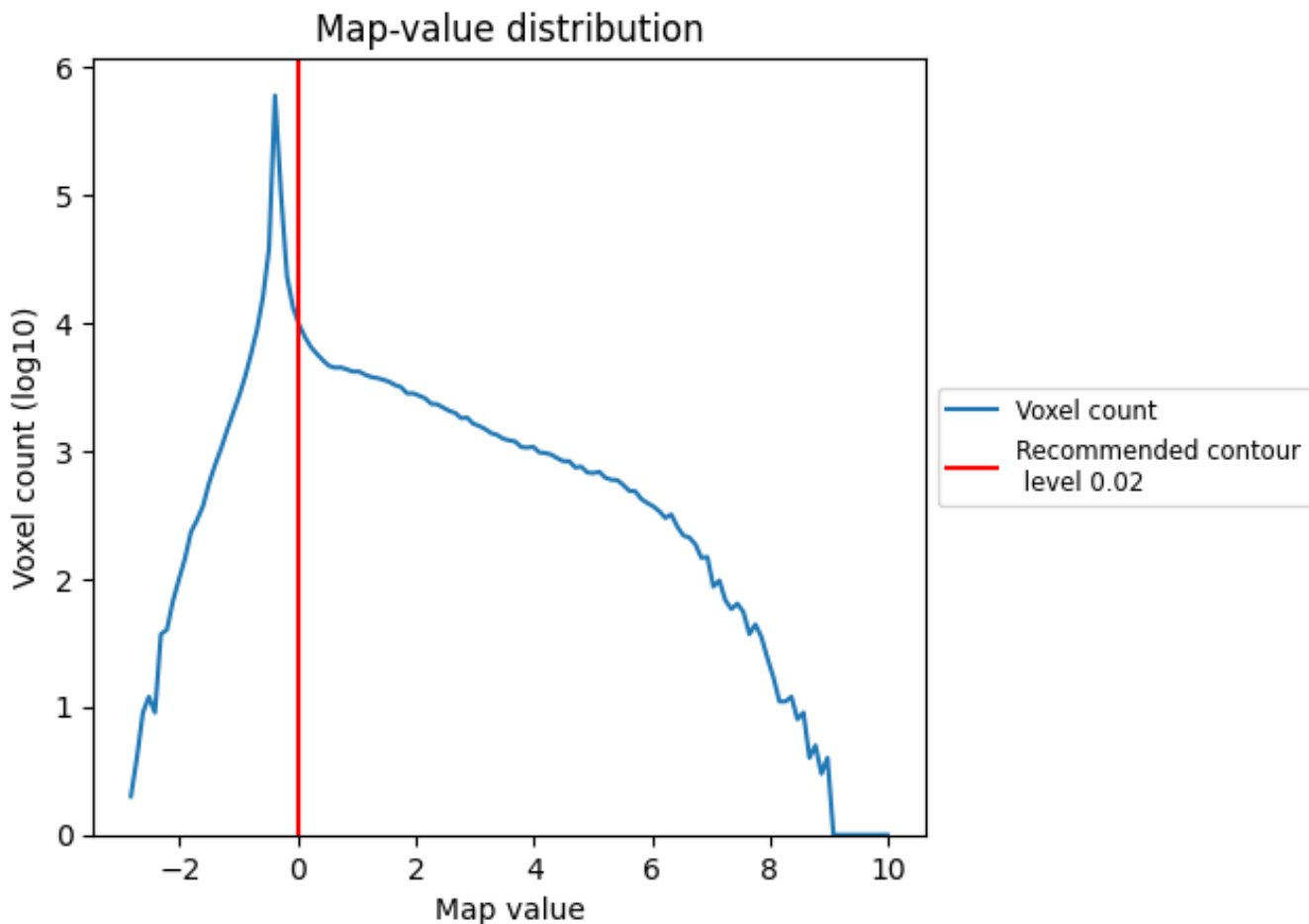
6.5 Mask visualisation [i](#)

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

7 Map analysis [i](#)

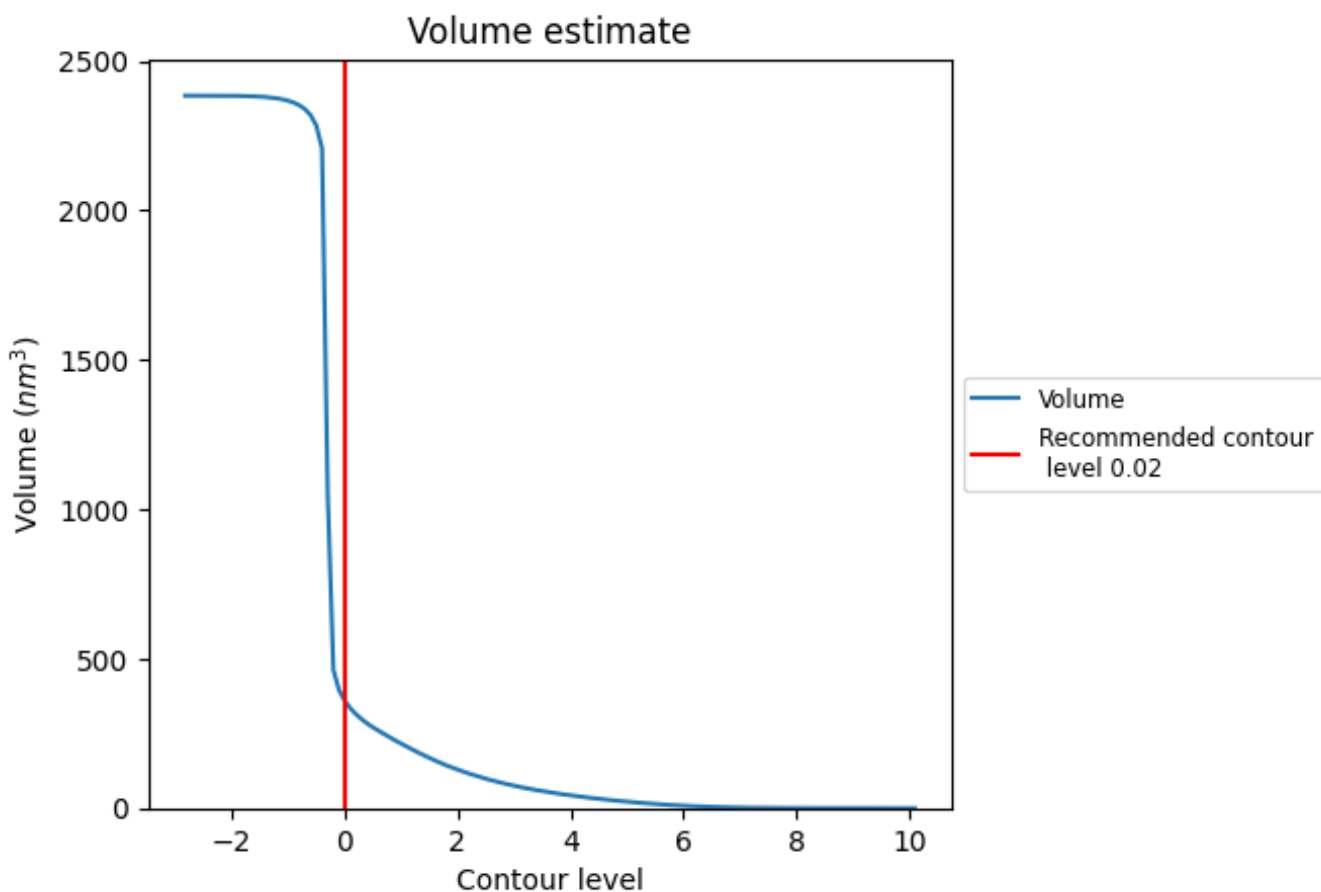
This section contains the results of statistical analysis of the map.

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 355 nm³; this corresponds to an approximate mass of 320 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.

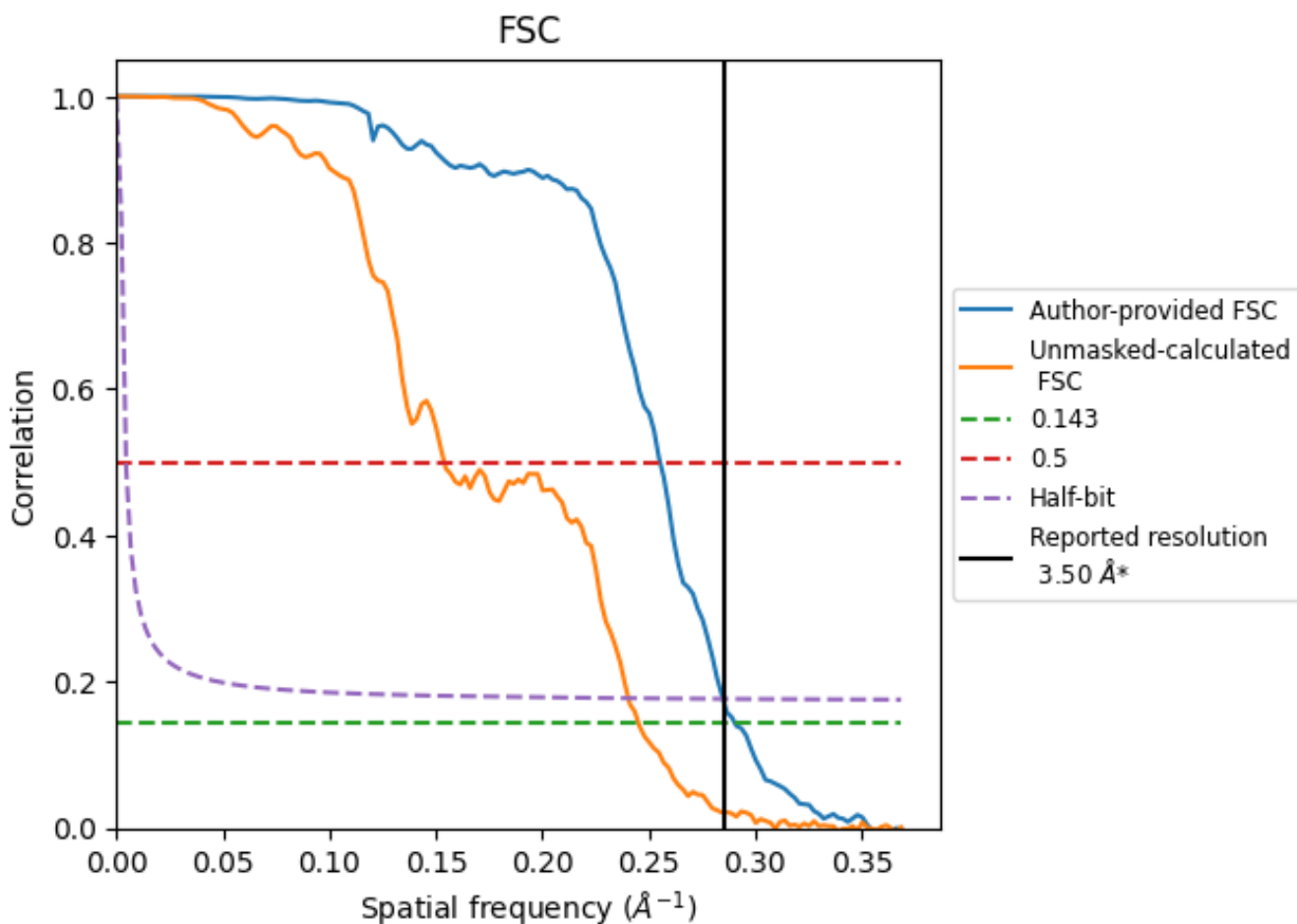
7.3 Rotationally averaged power spectrum [i](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

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 0.286 Å⁻¹

8.2 Resolution estimates [i](#)

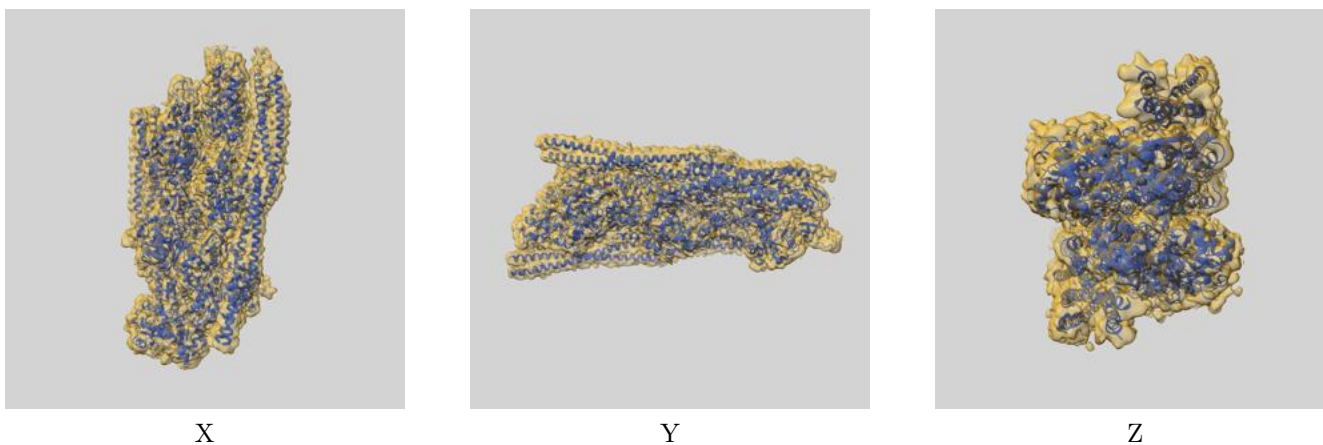
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	3.44	3.92	3.51
Unmasked-calculated*	4.08	6.49	4.16

*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 4.08 differs from the reported value 3.5 by more than 10 %

9 Map-model fit [i](#)

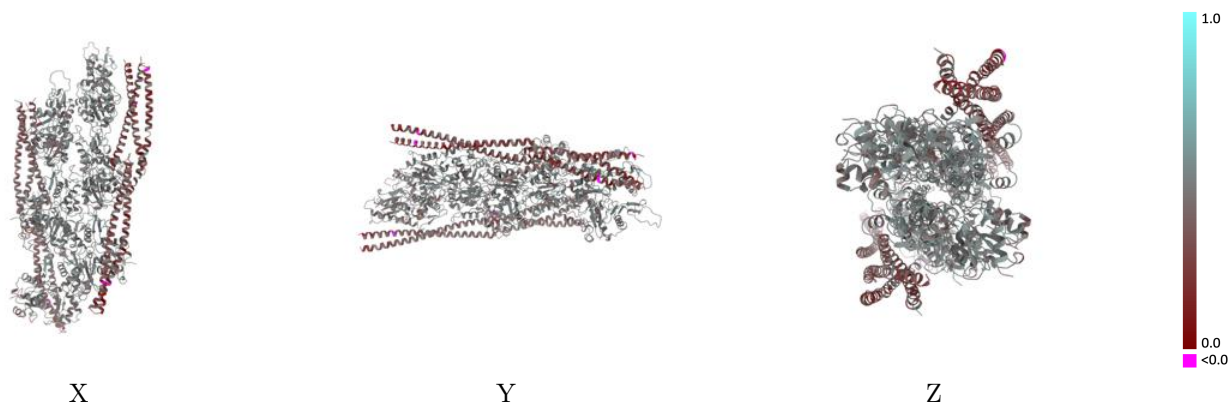
This section contains information regarding the fit between EMDB map EMD-27331 and PDB model 8DD0. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



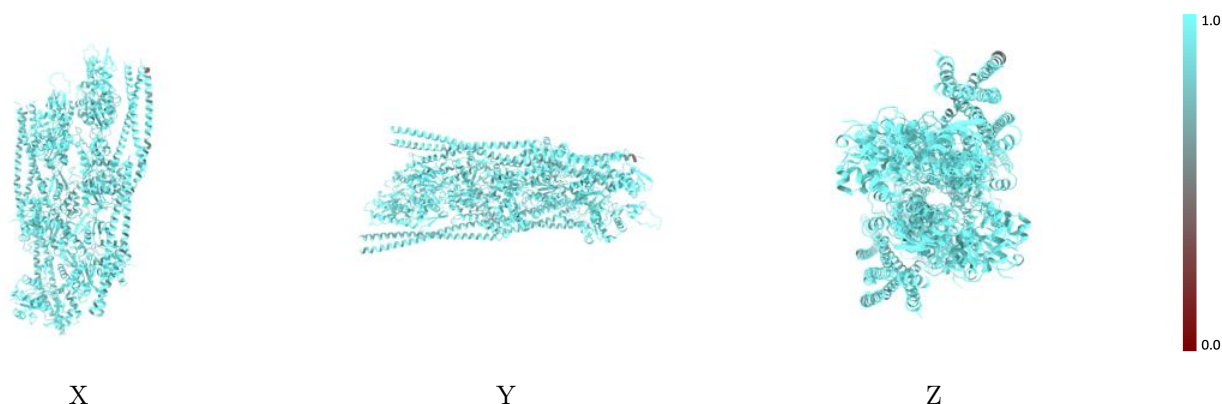
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



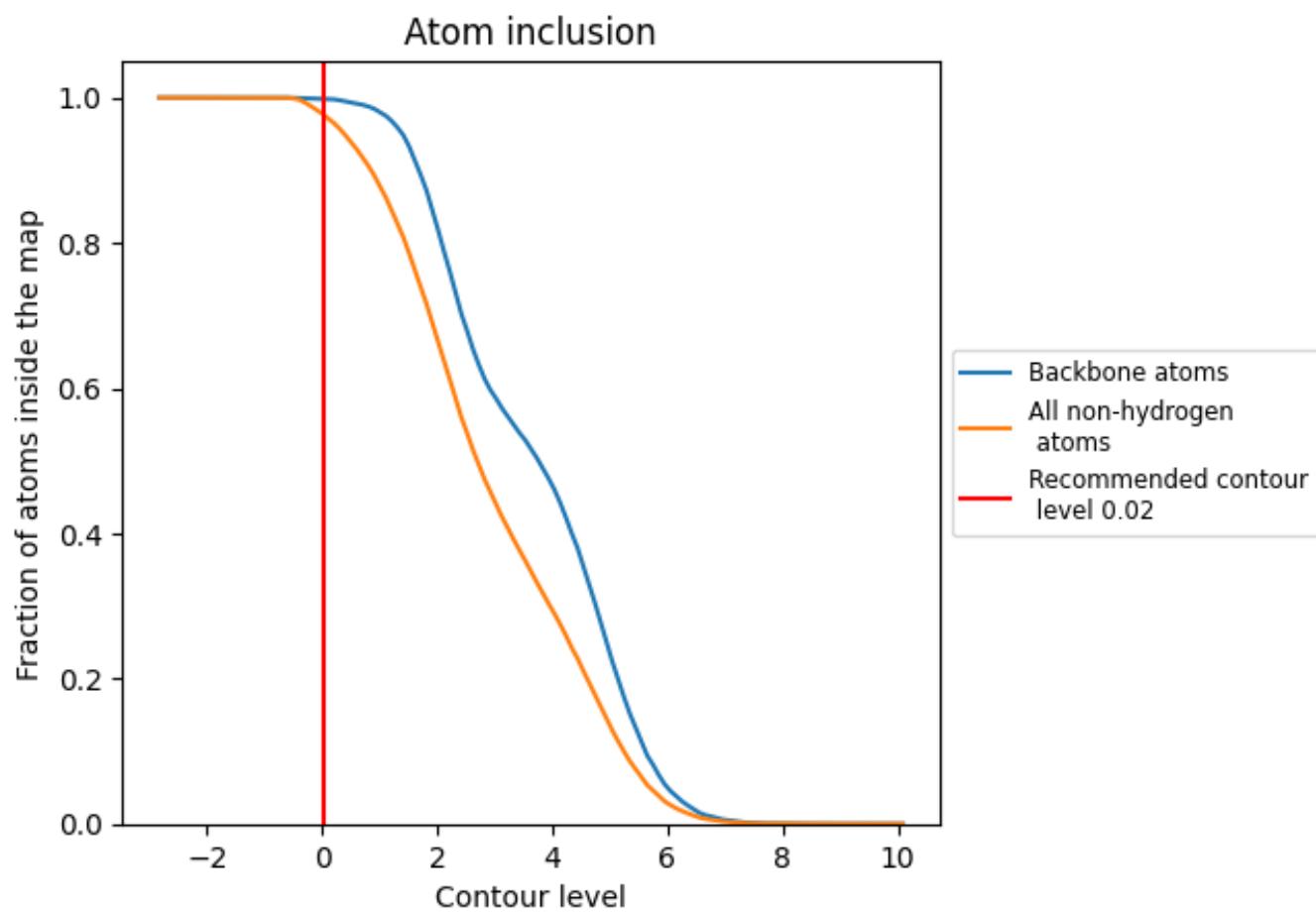
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.02).



































9.4 Atom inclusion [i](#)



At the recommended contour level, 100% of all backbone atoms, 98% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9772	 0.4430
A	 0.9875	 0.4650
B	 0.9895	 0.4820
C	 0.9941	 0.4900
D	 0.9920	 0.4910
E	 0.9934	 0.4900
F	 0.9902	 0.4840
G	 0.9592	 0.3200
H	 0.9841	 0.3420
I	 0.9207	 0.2980
J	 0.9441	 0.3300
K	 0.9215	 0.3360
L	 0.9342	 0.2880
M	 0.9660	 0.2870
N	 0.8919	 0.2860
O	 0.9441	 0.3120
P	 0.8798	 0.3050

