



## wwPDB EM Validation Summary Report ⓘ

Nov 19, 2022 – 02:05 PM EST

PDB ID : 3J94  
EMDB ID : EMD-6204  
Title : Structure of ATP-bound N-ethylmaleimide sensitive factor determined by single particle cryoelectron microscopy  
Authors : Zhao, M.; Wu, S.; Cheng, Y.; Brunger, A.T.  
Deposited on : 2014-12-05  
Resolution : 4.20 Å(reported)  
Based on initial model : 1NSF

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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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) : 2.31.3

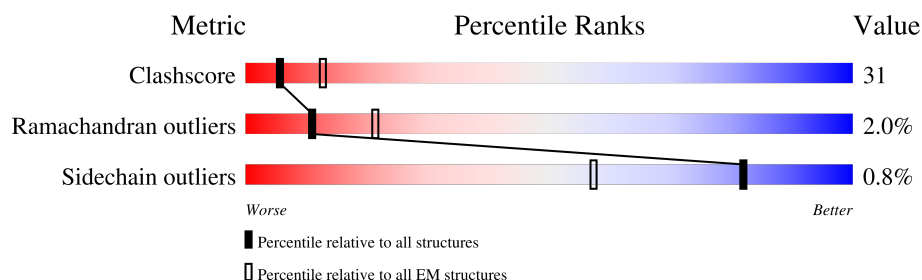
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 4.20 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  $\geq 3$ , 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	747	<div> <div>5%</div> <div>32%</div> <div>32%</div> <div>34%</div> </div>
1	B	747	<div> <div>32%</div> <div>31%</div> <div>35%</div> </div>
1	C	747	<div> <div>34%</div> <div>30%</div> <div>35%</div> </div>
1	D	747	<div> <div>5%</div> <div>33%</div> <div>32%</div> <div>35%</div> </div>
1	E	747	<div> <div>14%</div> <div>28%</div> <div>36%</div> <div>35%</div> </div>
1	F	747	<div> <div>31%</div> <div>29%</div> <div>32%</div> <div>38%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ATP	D	801	-	-	X	-

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21509 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.

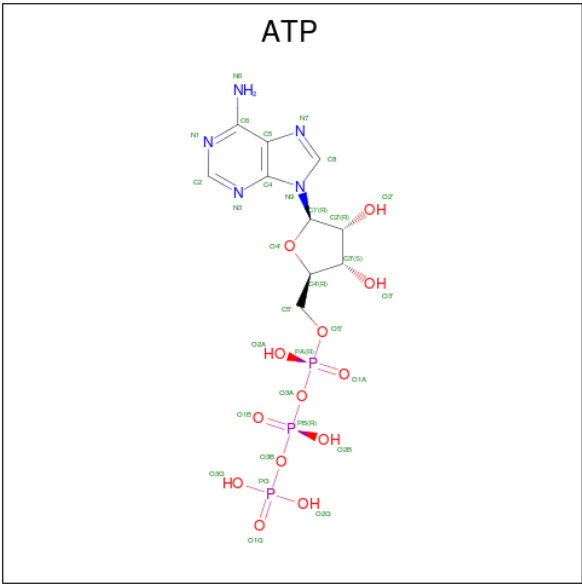
- Molecule 1 is a protein called Vesicle-fusing ATPase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	490	Total	C	N	O	S	0	0
			3559	2266	613	665	15		
1	B	484	Total	C	N	O	S	0	0
			3563	2264	618	665	16		
1	C	488	Total	C	N	O	S	0	0
			3573	2268	621	669	15		
1	D	485	Total	C	N	O	S	0	0
			3526	2244	603	663	16		
1	E	482	Total	C	N	O	S	0	0
			3529	2242	611	661	15		
1	F	466	Total	C	N	O	S	0	0
			3418	2164	598	640	16		

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P18708
A	-1	ALA	-	EXPRESSION TAG	UNP P18708
A	0	HIS	-	EXPRESSION TAG	UNP P18708
B	-2	GLY	-	EXPRESSION TAG	UNP P18708
B	-1	ALA	-	EXPRESSION TAG	UNP P18708
B	0	HIS	-	EXPRESSION TAG	UNP P18708
C	-2	GLY	-	EXPRESSION TAG	UNP P18708
C	-1	ALA	-	EXPRESSION TAG	UNP P18708
C	0	HIS	-	EXPRESSION TAG	UNP P18708
D	-2	GLY	-	EXPRESSION TAG	UNP P18708
D	-1	ALA	-	EXPRESSION TAG	UNP P18708
D	0	HIS	-	EXPRESSION TAG	UNP P18708
E	-2	GLY	-	EXPRESSION TAG	UNP P18708
E	-1	ALA	-	EXPRESSION TAG	UNP P18708
E	0	HIS	-	EXPRESSION TAG	UNP P18708
F	-2	GLY	-	EXPRESSION TAG	UNP P18708
F	-1	ALA	-	EXPRESSION TAG	UNP P18708
F	0	HIS	-	EXPRESSION TAG	UNP P18708

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).

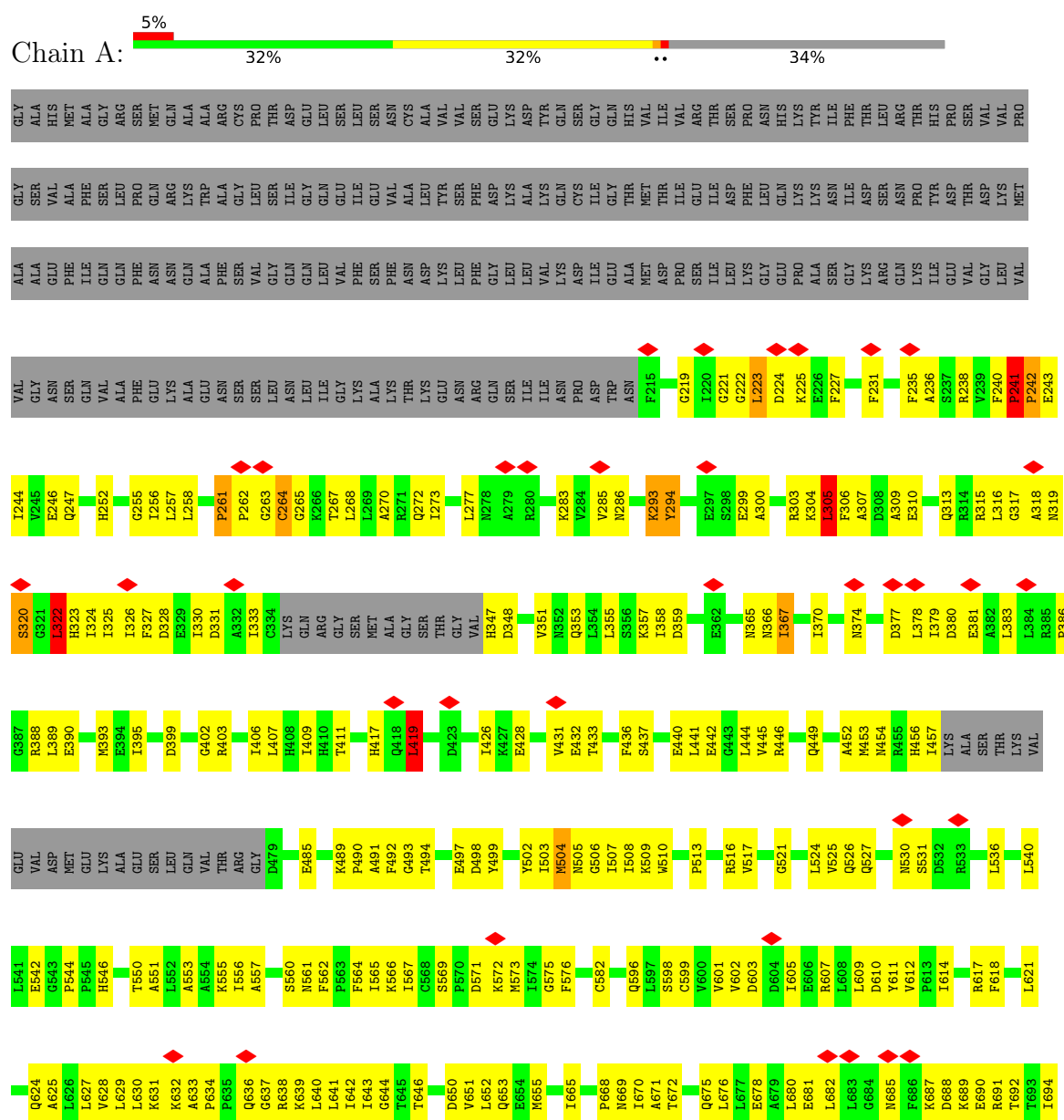


Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	A	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	B	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	B	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	C	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	C	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	D	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	D	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	E	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	E	1	Total	C	N	O	P	0
			62	20	10	26	6	
2	F	1	Total	C	N	O	P	0
			31	10	5	13	3	

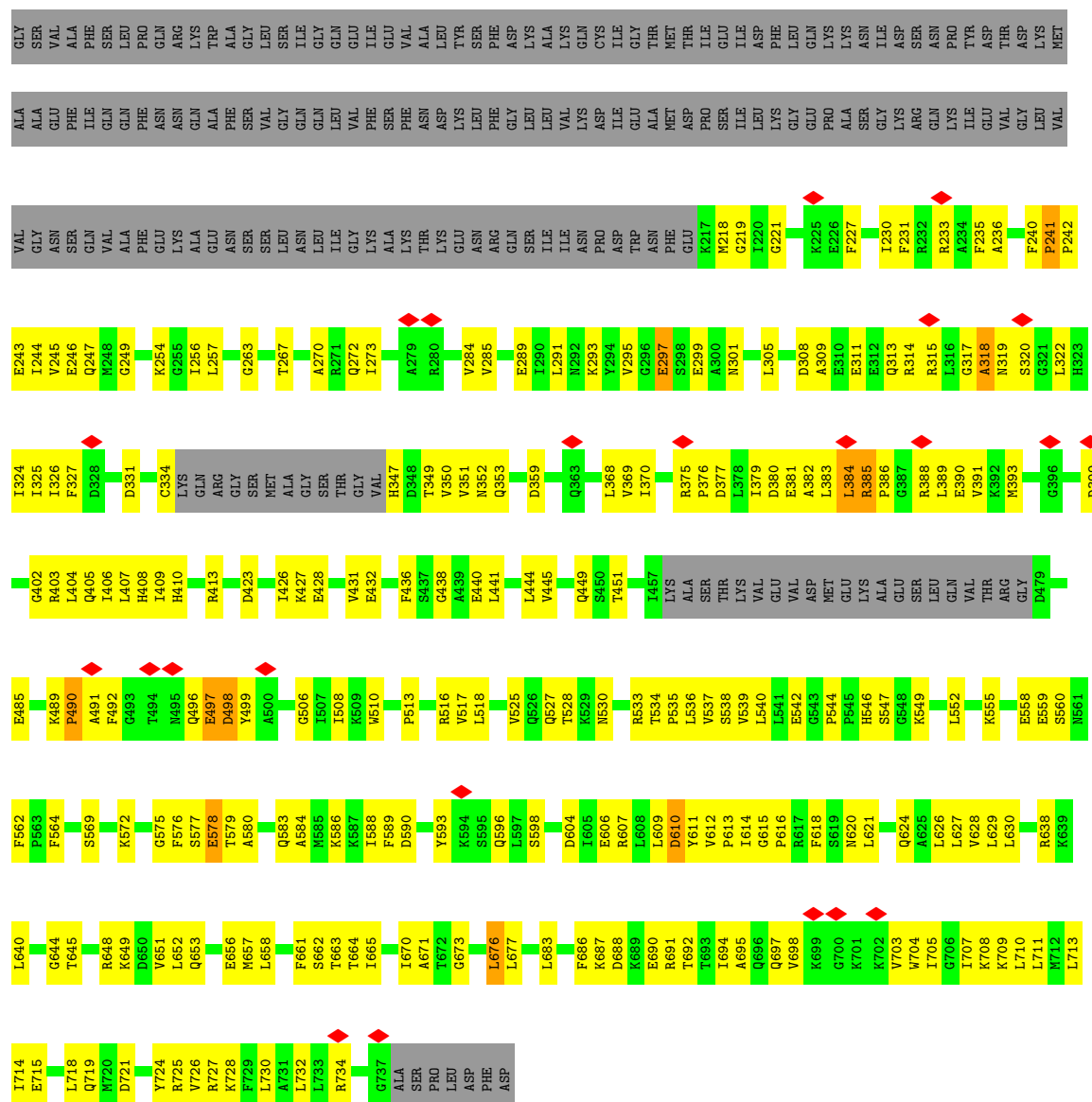
### 3 Residue-property plots

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.

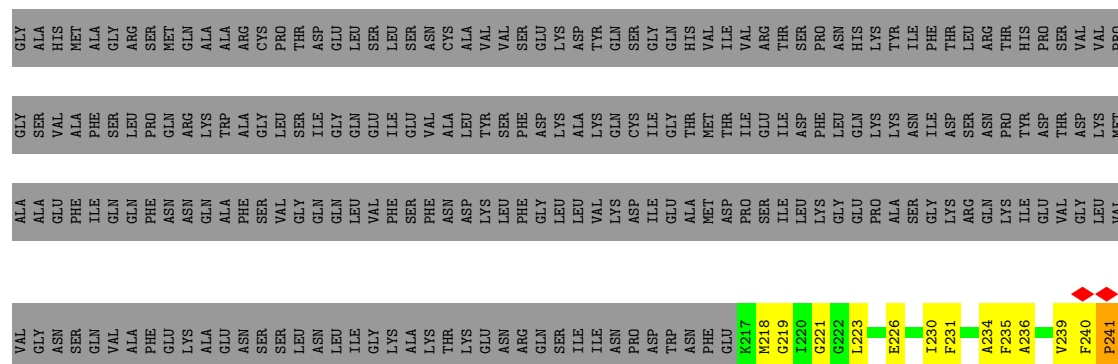
#### • Molecule 1: Vesicle-fusing ATPase



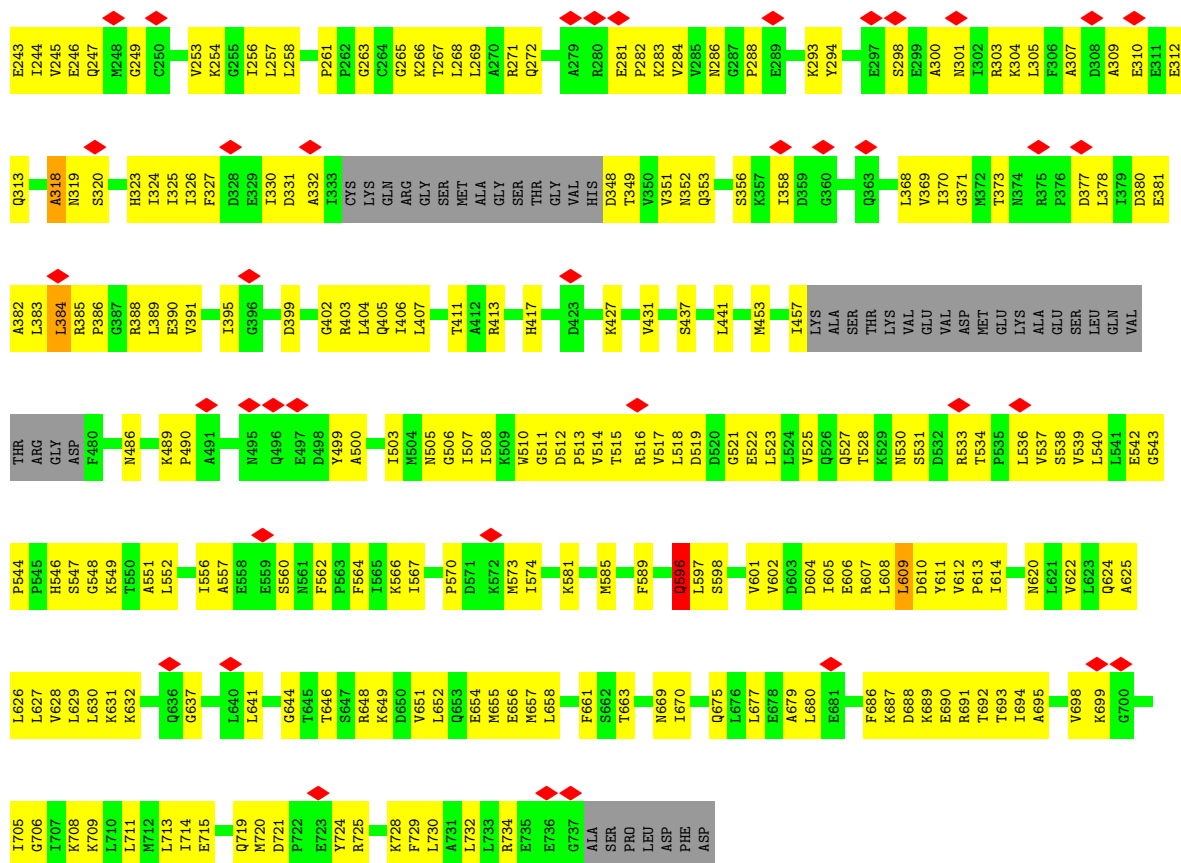




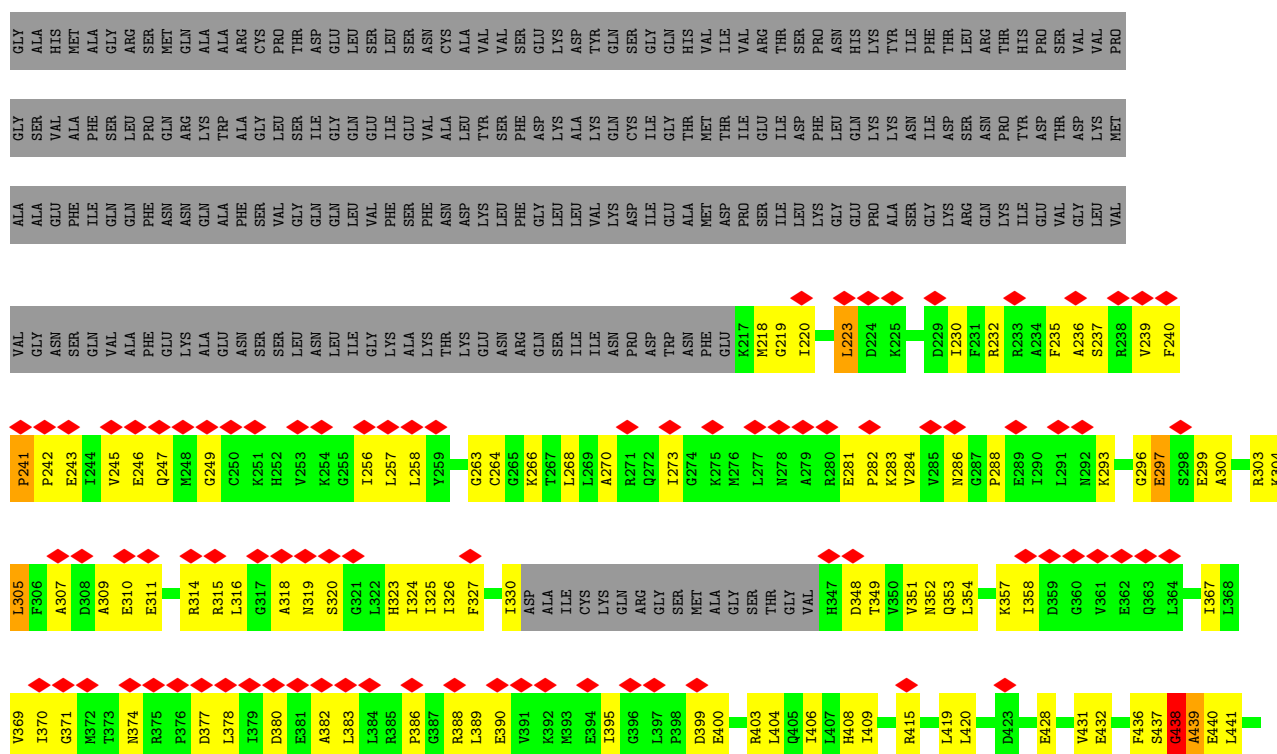
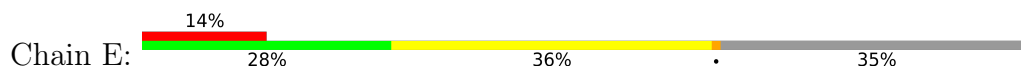
### • Molecule 1: Vesicle-fusing ATPase

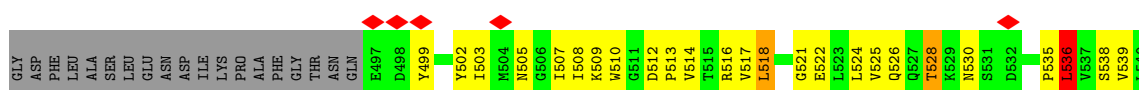






● Molecule 1: Vesicle-fusing ATPase





L541	E542	G543	P544	H546	S547	G548	K549	T550	A551	L552	A553	A554	K555	I556	A557	E558	E559	S560	F564	I565	K566	I567	C568	S569	P570	D571	K572	M573	I574	S577	E578	T579	A580	K581	C582	Q583	K586	F589	Y593	L597	S598	C599	V600	V601	V602	D603	D604	I605	E606	R607	L608	L609
D610	Y611	V612	P613	I614	G615	P616	R617	F618	S619	N620	Q624	V628	L629	K632	A633	P634	P635	R638	K639	L640	L641	I642	I643	G644	T645	T646	S647	R648	K649	D650	Q653	E654	M655	E656	M657	L658	F661	S662	T663	T664	I665	H666	V667	P668	M669	I670	G673	E674	Q675	L676	L677	E678
E681	L682	L683	G684	N685	F686	E690	I694	A695	V698	K699	G700	K701	K702	V703	W704	I707	K708	L710	L711	M712	L713	I714	E715	M716	S717	L718	Q719	M720	D721	Y724	R725	V726	R727	K728	L732	L733	R734	E735	E736	G737	ALA	SER	PRO	LEU	ASP	PHE	ASP					

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	50781	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	Each particle	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	26.4	Depositor
Minimum defocus (nm)	-1800	Depositor
Maximum defocus (nm)	-2800	Depositor
Magnification	31000	Depositor
Image detector	GATAN K2 (4k x 4k)	Depositor
Maximum map value	17.301	Depositor
Minimum map value	-6.891	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	6.0	Depositor
Map size ( $\text{\AA}$ )	311.1936, 311.1936, 311.1936	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.2156, 1.2156, 1.2156	Depositor

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ATP

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	0.37	0/3612	0.81	4/4904 (0.1%)
1	B	0.37	0/3615	0.77	1/4898 (0.0%)
1	C	0.38	0/3625	0.77	3/4915 (0.1%)
1	D	0.36	0/3577	0.77	3/4856 (0.1%)
1	E	0.35	0/3580	0.79	2/4854 (0.0%)
1	F	0.39	0/3467	0.82	3/4697 (0.1%)
All	All	0.37	0/21476	0.79	16/29124 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	8

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(°)
1	E	223	LEU	CA-CB-CG	6.62	130.52	115.30
1	F	536	LEU	CA-CB-CG	6.52	130.31	115.30
1	F	518	LEU	CA-CB-CG	6.32	129.84	115.30
1	C	385	ARG	NE-CZ-NH2	6.20	123.40	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	609	LEU	CA-CB-CG	6.07	129.27	115.30

There are no chirality outliers.

5 of 8 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	241	PRO	Peptide
1	A	261	PRO	Peptide
1	B	438	GLY	Peptide
1	B	546	HIS	Sidechain
1	C	530	ASN	Sidechain

## 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	A	3559	0	3451	224	0
1	B	3563	0	3497	215	0
1	C	3573	0	3487	212	0
1	D	3526	0	3429	251	0
1	E	3529	0	3439	263	0
1	F	3418	0	3332	228	0
2	A	62	0	23	11	0
2	B	62	0	21	11	0
2	C	62	0	22	9	0
2	D	62	0	20	18	0
2	E	62	0	17	9	0
2	F	31	0	9	5	0
All	All	21509	0	20747	1292	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 31.

The worst 5 of 1292 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:505:ASN:HB2	2:A:801:ATP:H1'	1.39	1.01
1:C:257:LEU:HB2	1:C:389:LEU:HD13	1.56	0.87
1:B:526:GLN:HE21	1:C:719:GLN:HB3	1.39	0.85
1:B:386:PRO:HA	1:B:390:GLU:HA	1.58	0.84
1:A:705:ILE:HD13	1:A:710:LEU:HD12	1.59	0.84

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	484/747 (65%)	452 (93%)	20 (4%)	12 (2%)	5	35
1	B	478/747 (64%)	436 (91%)	31 (6%)	11 (2%)	6	37
1	C	482/747 (64%)	454 (94%)	19 (4%)	9 (2%)	8	41
1	D	479/747 (64%)	448 (94%)	25 (5%)	6 (1%)	12	48
1	E	474/747 (64%)	443 (94%)	21 (4%)	10 (2%)	7	39
1	F	460/747 (62%)	425 (92%)	26 (6%)	9 (2%)	7	40
All	All	2857/4482 (64%)	2658 (93%)	142 (5%)	57 (2%)	11	40

5 of 57 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	LEU
1	A	293	LYS
1	A	294	TYR
1	A	318	ALA
1	A	320	SER

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	348/638 (54%)	344 (99%)	4 (1%)	73	84
1	B	356/638 (56%)	353 (99%)	3 (1%)	81	89
1	C	353/638 (55%)	351 (99%)	2 (1%)	86	92
1	D	348/638 (54%)	347 (100%)	1 (0%)	92	95
1	E	350/638 (55%)	349 (100%)	1 (0%)	92	95
1	F	339/638 (53%)	334 (98%)	5 (2%)	65	80
All	All	2094/3828 (55%)	2078 (99%)	16 (1%)	82	89

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	528	THR
1	F	365	ASN
1	C	676	LEU
1	F	322	LEU
1	C	305	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	527	GLN
1	F	546	HIS
1	D	272	GLN
1	D	319	ASN
1	D	352	ASN

### 5.3.3 RNA ⓘ

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](#)

11 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	ATP	B	802	-	26,33,33	0.94	1 (3%)	31,52,52	1.45	5 (16%)
2	ATP	C	801	-	26,33,33	0.91	1 (3%)	31,52,52	1.64	5 (16%)
2	ATP	E	801	-	26,33,33	1.01	3 (11%)	31,52,52	2.34	9 (29%)
2	ATP	B	801	-	26,33,33	0.98	2 (7%)	31,52,52	1.59	8 (25%)
2	ATP	A	802	-	26,33,33	0.91	1 (3%)	31,52,52	1.56	5 (16%)
2	ATP	F	801	-	26,33,33	0.99	1 (3%)	31,52,52	1.68	5 (16%)
2	ATP	E	802	-	26,33,33	0.96	1 (3%)	31,52,52	1.62	5 (16%)
2	ATP	C	802	-	26,33,33	1.03	1 (3%)	31,52,52	1.53	5 (16%)
2	ATP	D	802	-	26,33,33	1.01	1 (3%)	31,52,52	1.85	9 (29%)
2	ATP	D	801	-	26,33,33	0.93	1 (3%)	31,52,52	1.52	6 (19%)
2	ATP	A	801	-	26,33,33	1.09	2 (7%)	31,52,52	1.56	5 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATP	B	802	-	-	6/18/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATP	C	801	-	-	2/18/38/38	0/3/3/3
2	ATP	E	801	-	-	5/18/38/38	0/3/3/3
2	ATP	B	801	-	-	2/18/38/38	0/3/3/3
2	ATP	A	802	-	-	1/18/38/38	0/3/3/3
2	ATP	F	801	-	-	4/18/38/38	0/3/3/3
2	ATP	E	802	-	-	5/18/38/38	0/3/3/3
2	ATP	C	802	-	-	6/18/38/38	0/3/3/3
2	ATP	D	802	-	-	4/18/38/38	0/3/3/3
2	ATP	D	801	-	-	4/18/38/38	0/3/3/3
2	ATP	A	801	-	-	3/18/38/38	0/3/3/3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	802	ATP	C5-C4	2.92	1.48	1.40
2	C	802	ATP	C5-C4	2.90	1.48	1.40
2	A	801	ATP	C5-C4	2.65	1.48	1.40
2	E	802	ATP	C5-C4	2.64	1.47	1.40
2	B	801	ATP	C5-C4	2.63	1.47	1.40

The worst 5 of 67 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	801	ATP	N3-C2-N1	-5.16	120.62	128.68
2	E	801	ATP	C5-C6-N6	5.05	128.03	120.35
2	F	801	ATP	PA-O3A-PB	-4.78	116.42	132.83
2	D	802	ATP	N3-C2-N1	-4.62	121.46	128.68
2	E	801	ATP	PB-O3B-PG	-4.51	117.35	132.83

There are no chirality outliers.

5 of 42 torsion outliers are listed below:

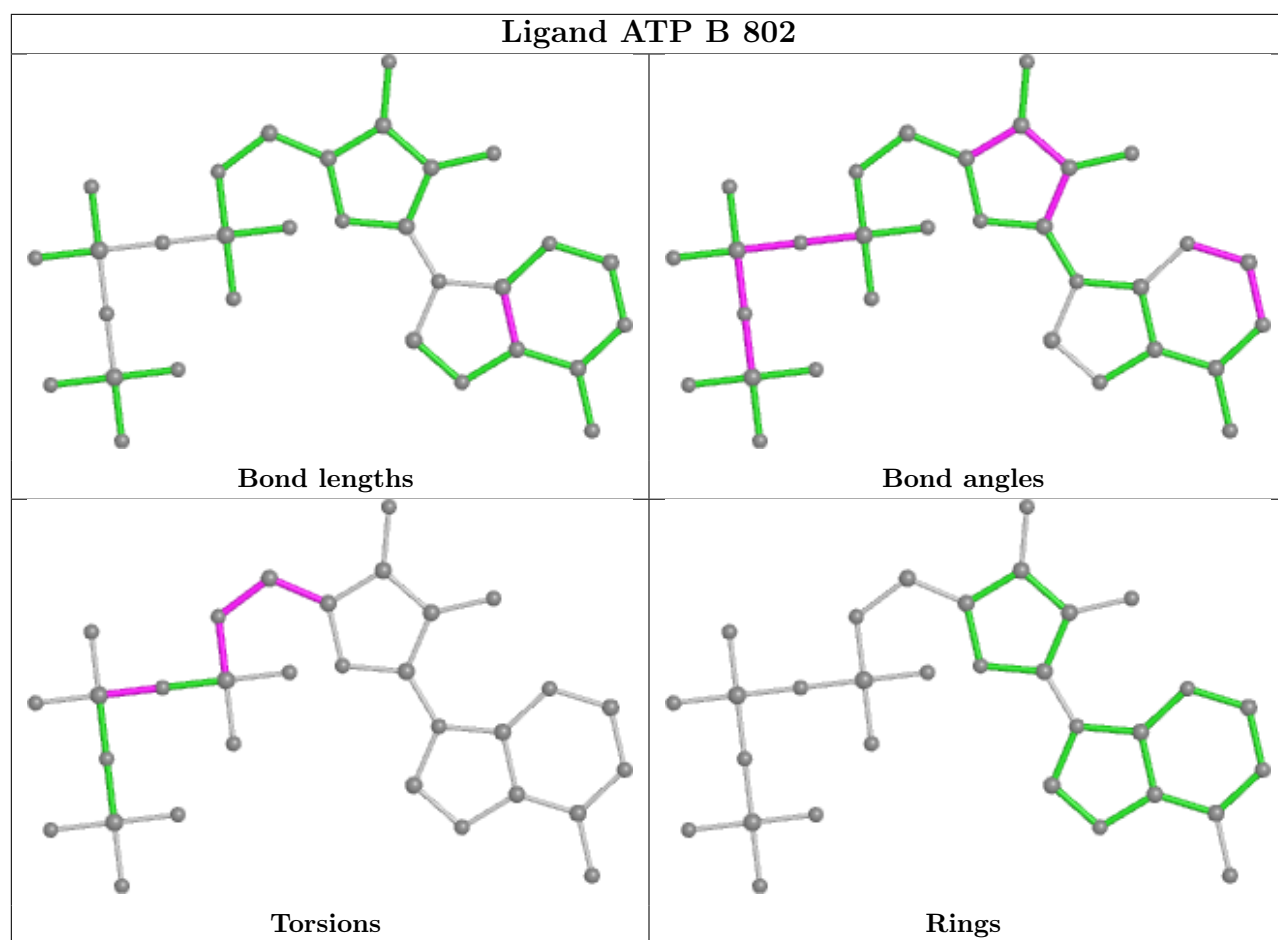
Mol	Chain	Res	Type	Atoms
2	B	802	ATP	C5'-O5'-PA-O3A
2	B	802	ATP	C4'-C5'-O5'-PA
2	C	802	ATP	PB-O3B-PG-O3G
2	C	802	ATP	C5'-O5'-PA-O3A
2	D	801	ATP	C5'-O5'-PA-O1A

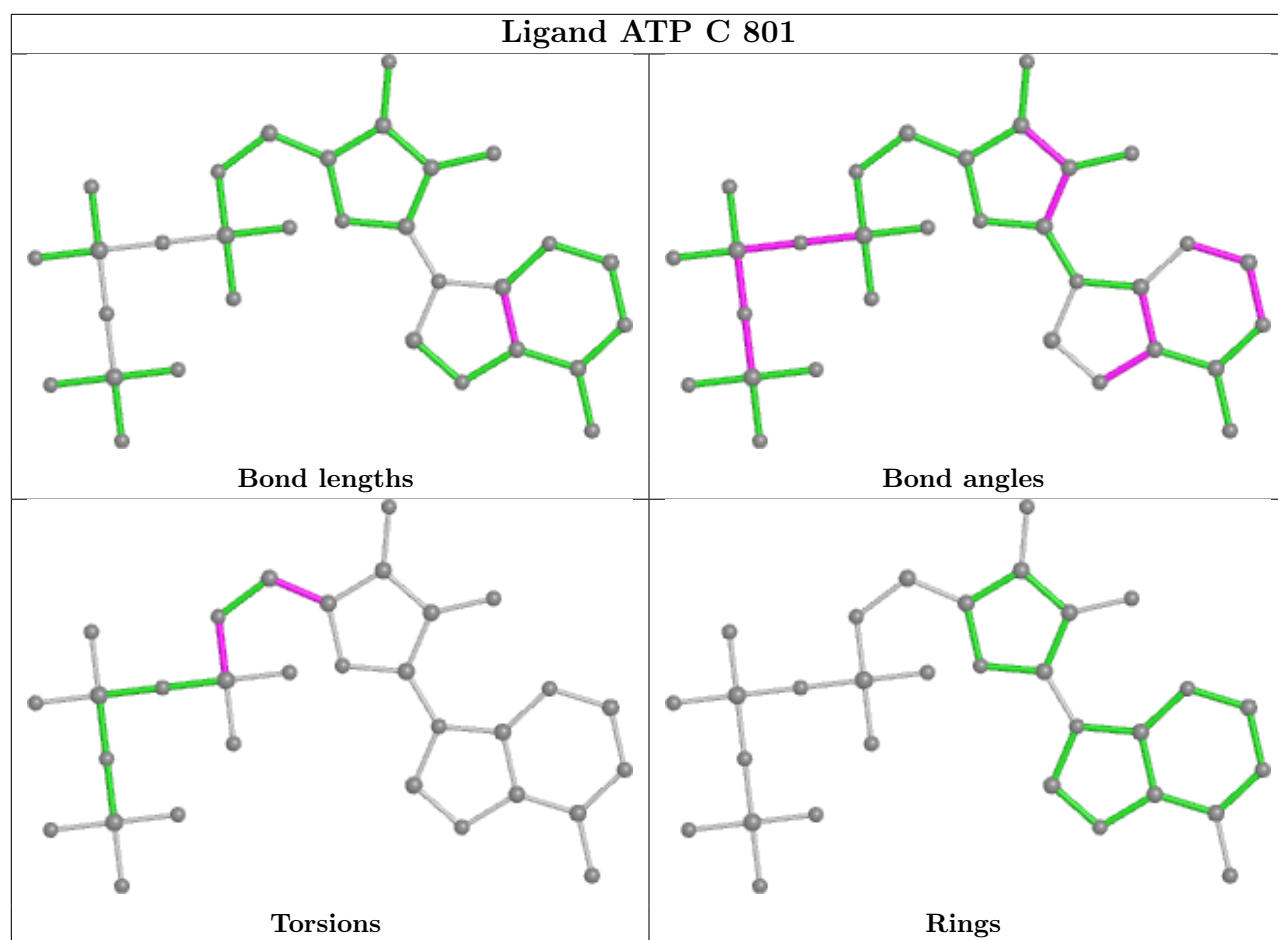
There are no ring outliers.

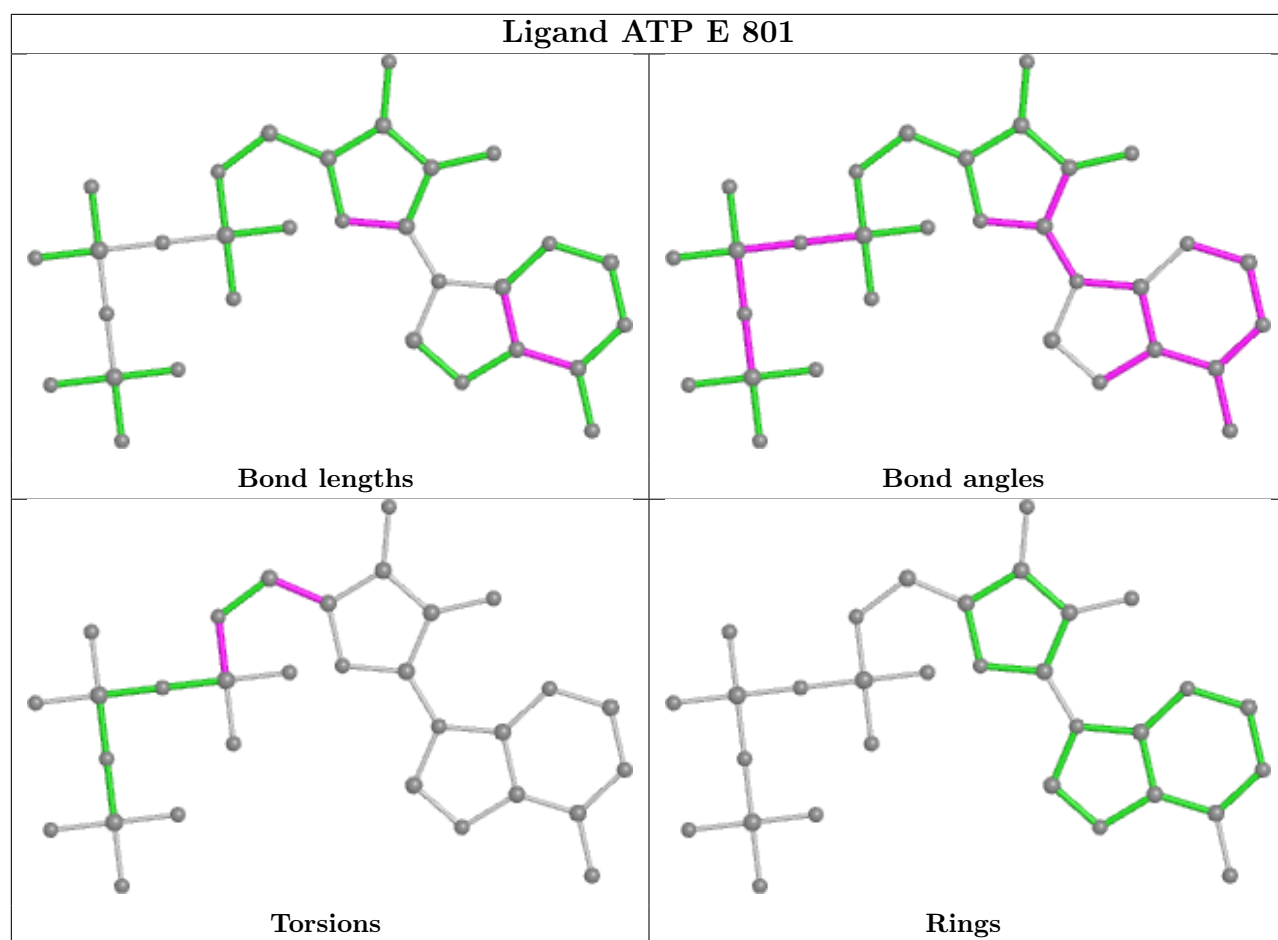
11 monomers are involved in 63 short contacts:

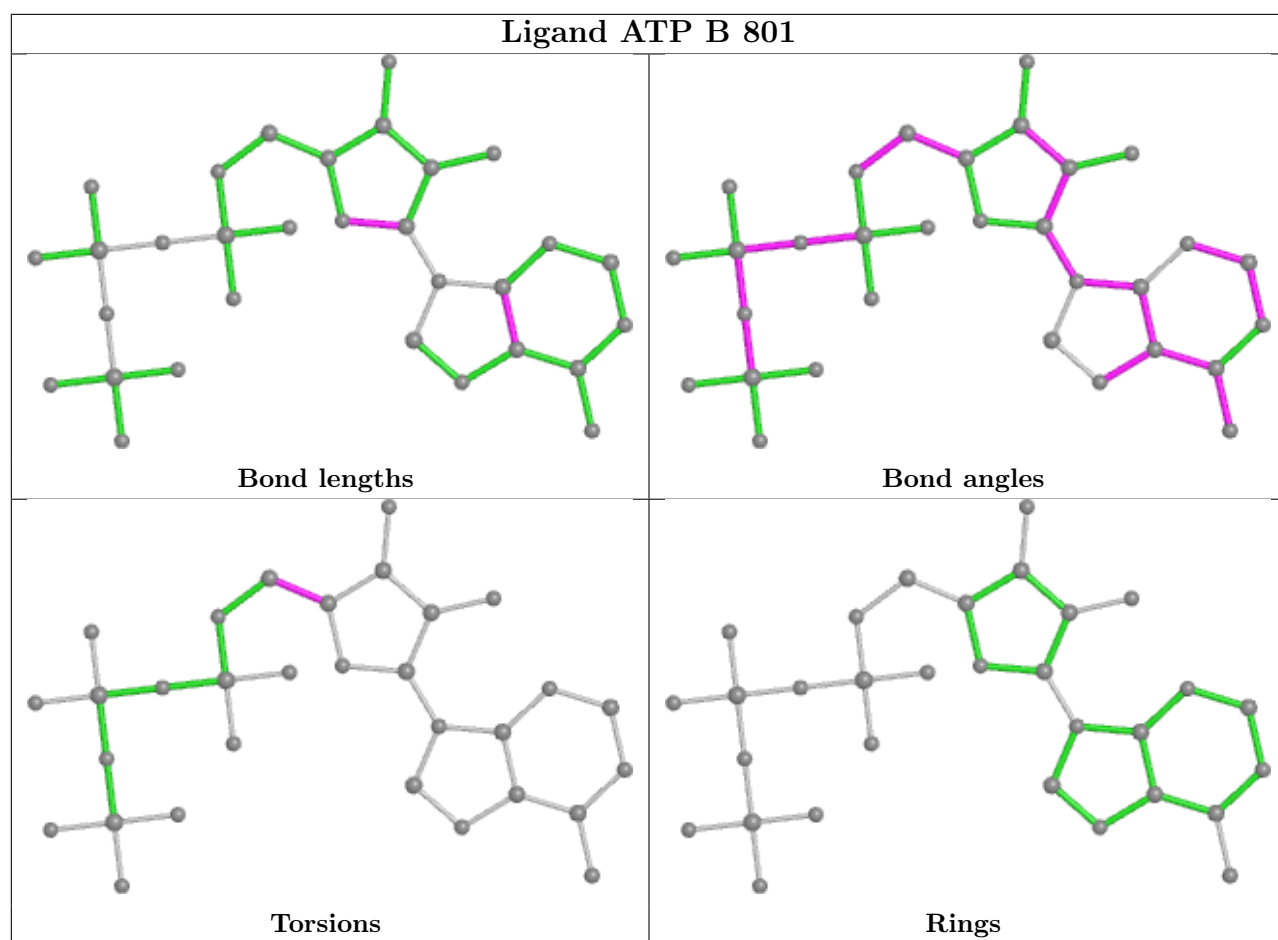
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	802	ATP	4	0
2	C	801	ATP	1	0
2	E	801	ATP	5	0
2	B	801	ATP	7	0
2	A	802	ATP	3	0
2	F	801	ATP	5	0
2	E	802	ATP	4	0
2	C	802	ATP	8	0
2	D	802	ATP	5	0
2	D	801	ATP	13	0
2	A	801	ATP	8	0

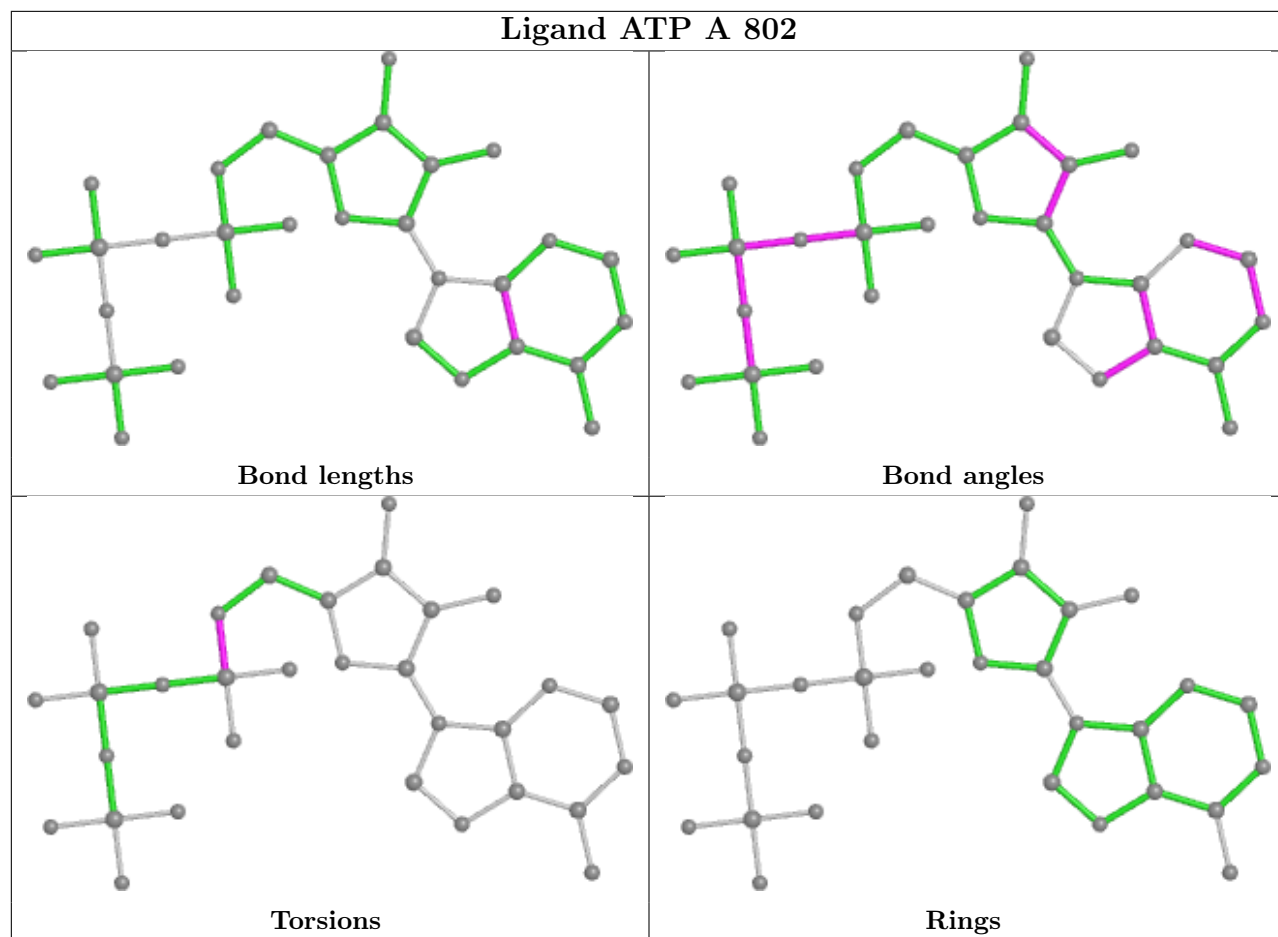
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



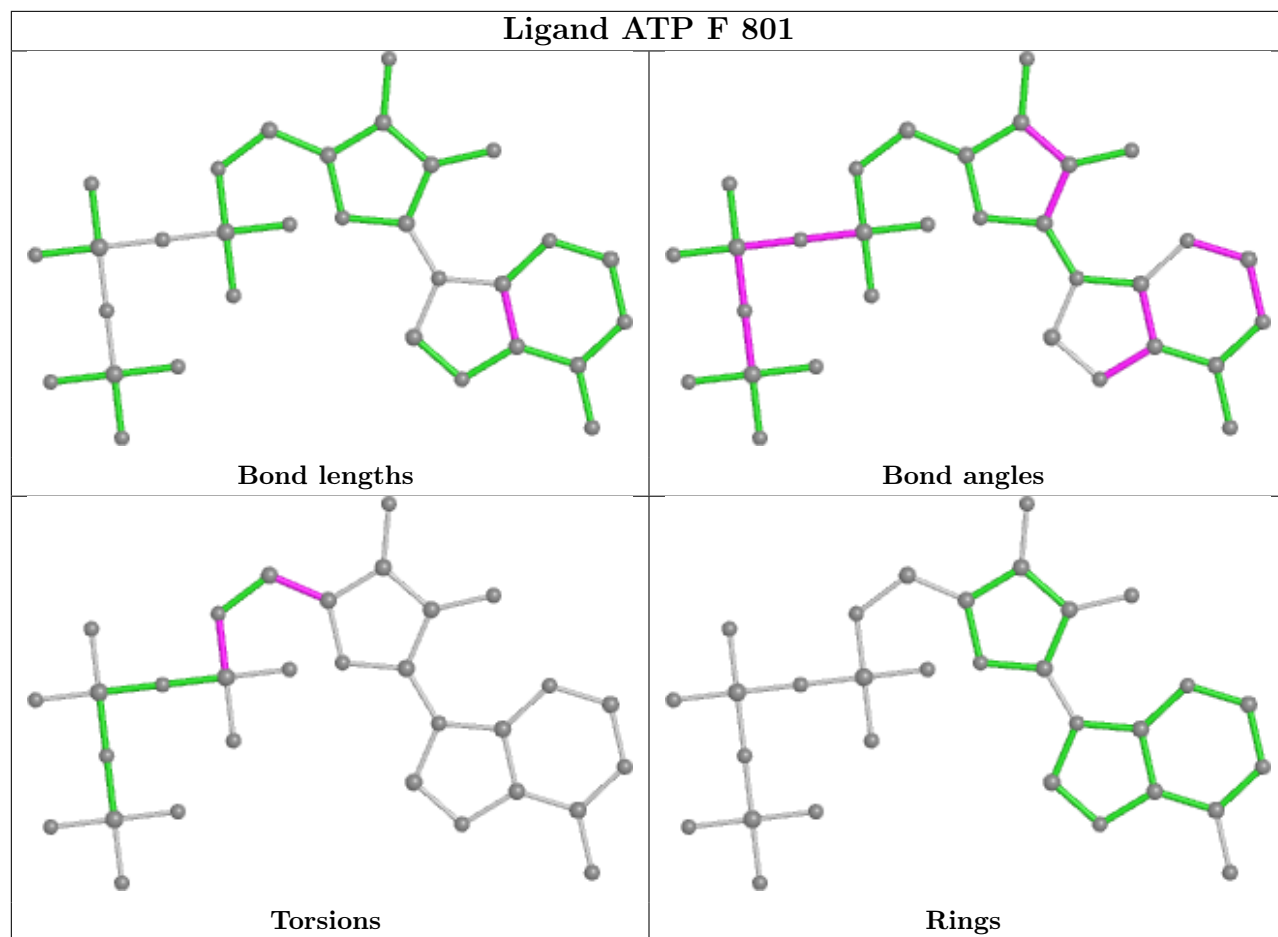


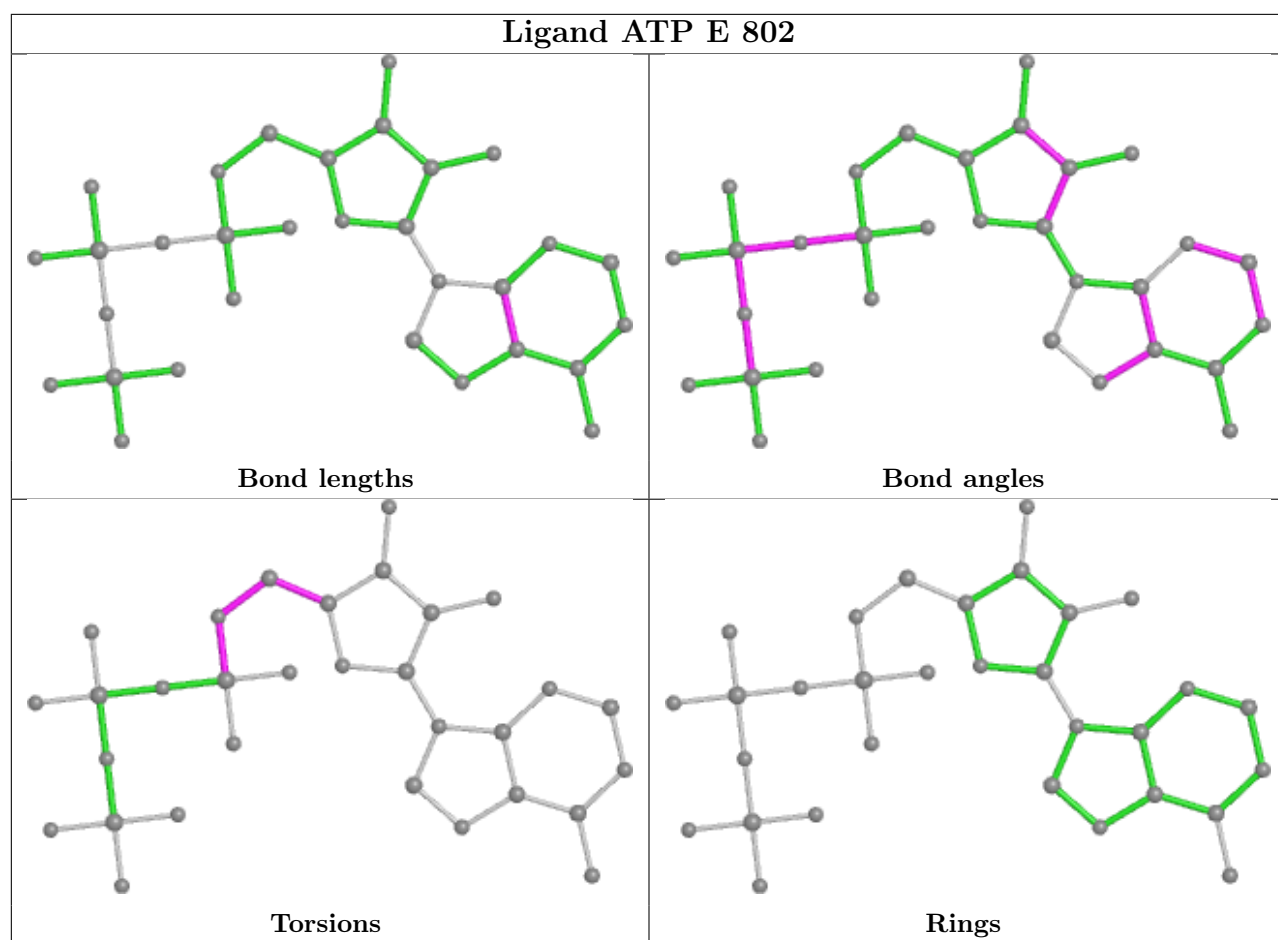


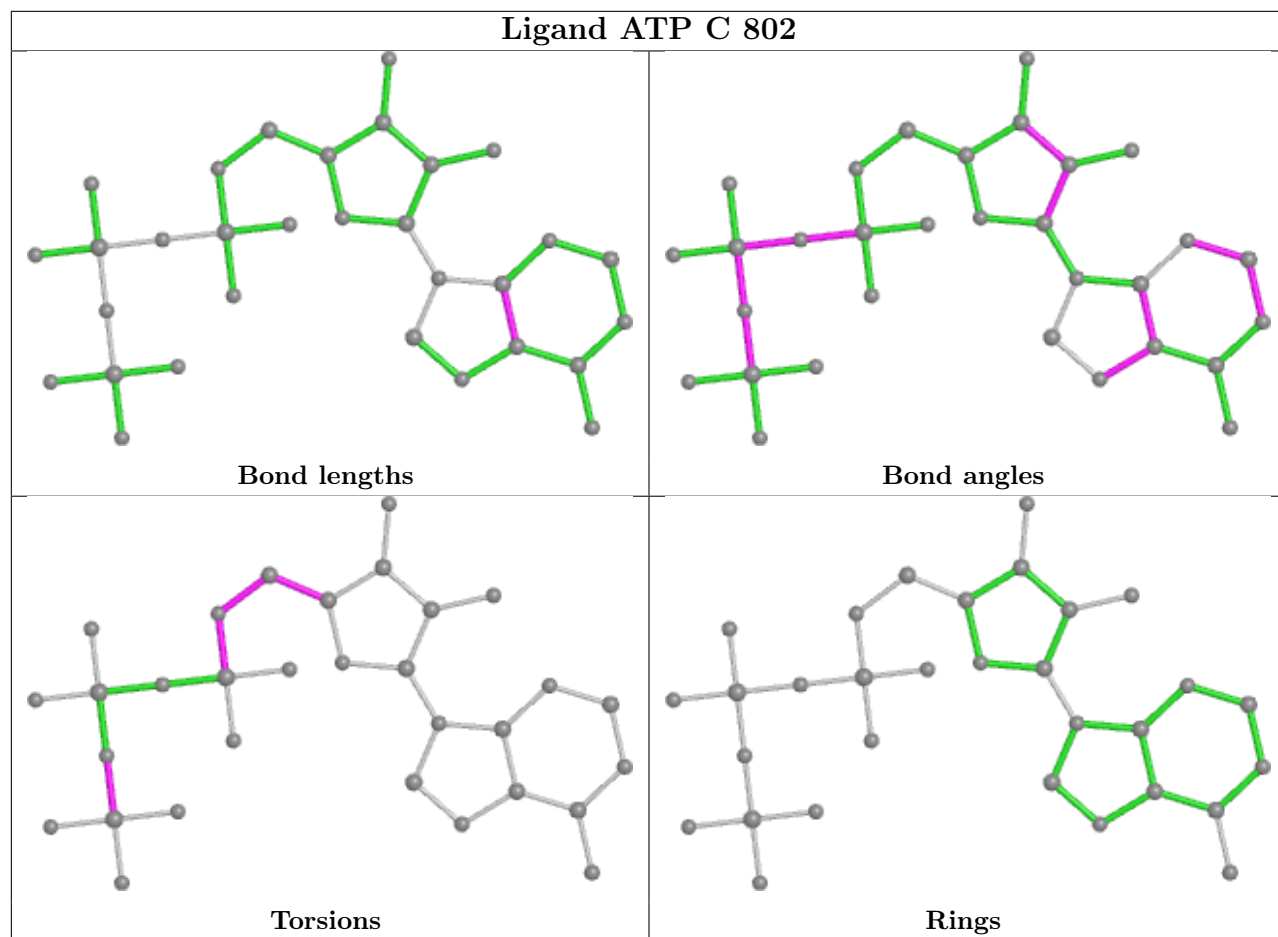


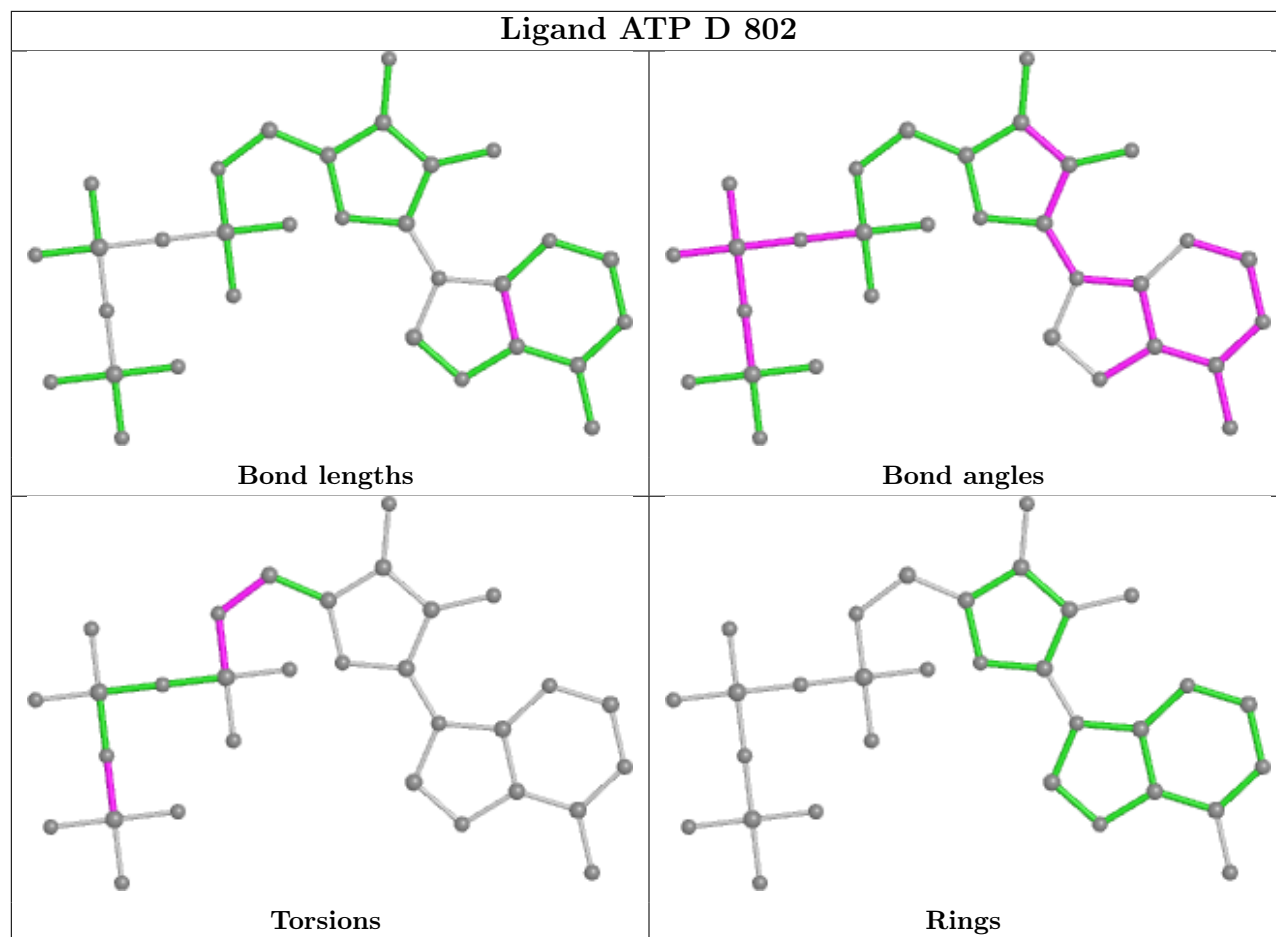


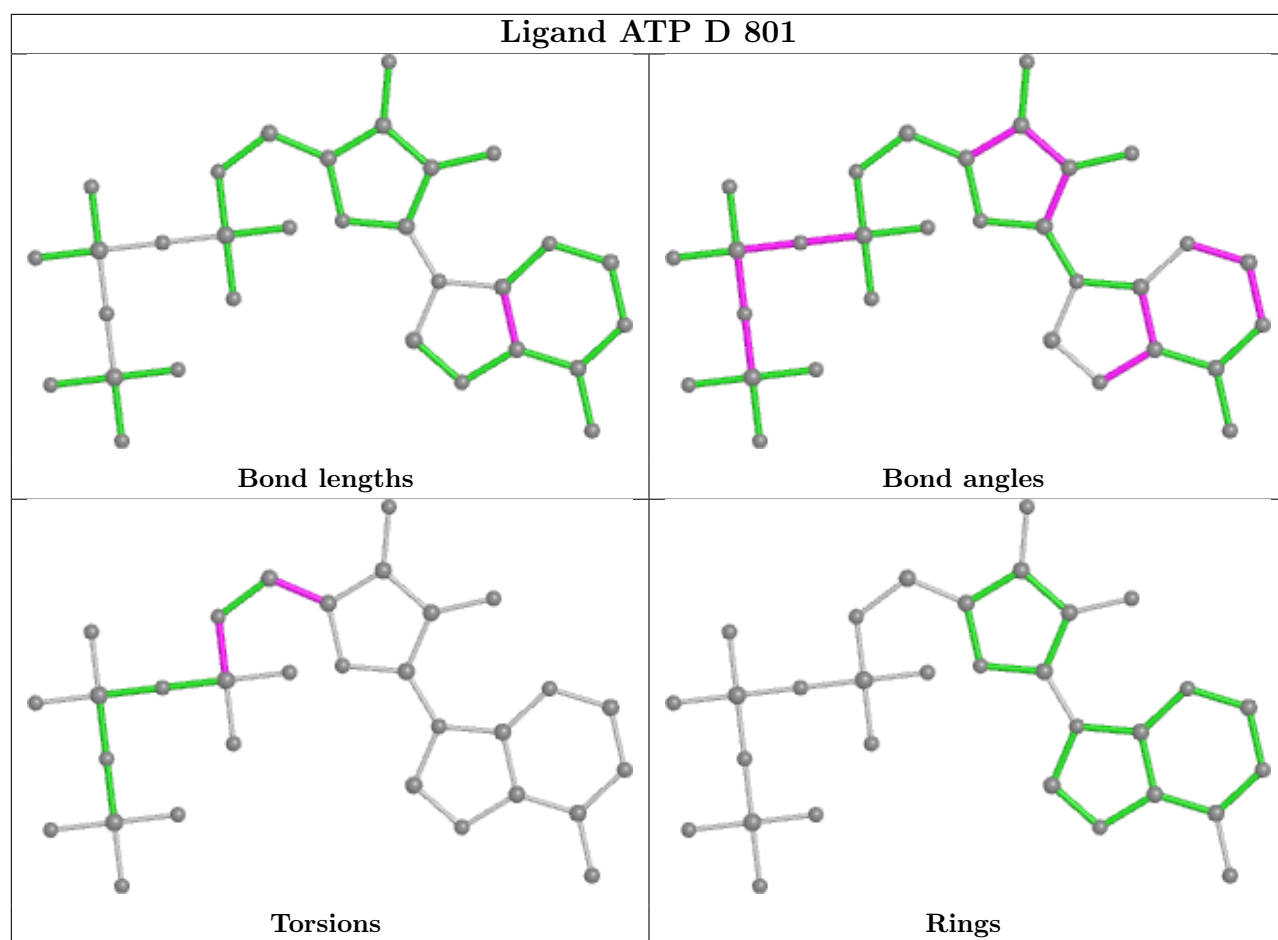


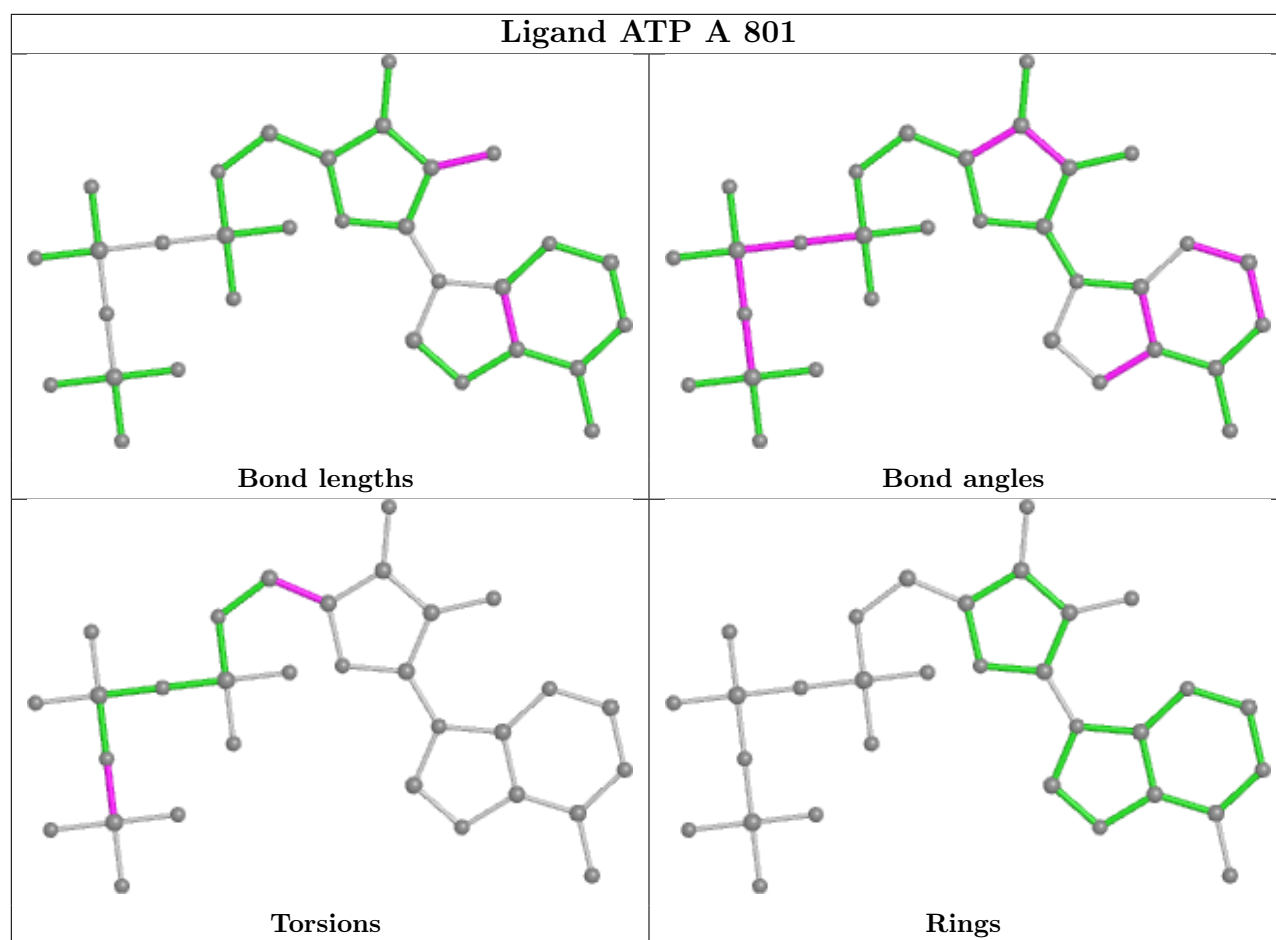












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

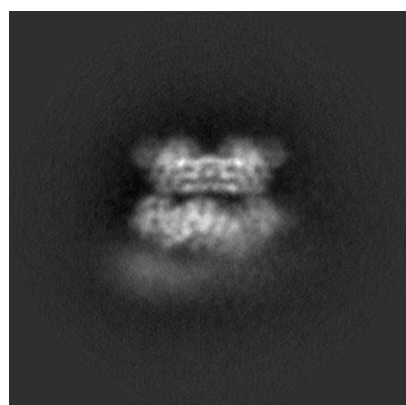
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-6204. These allow visual inspection of the internal detail of the map and identification of artifacts.

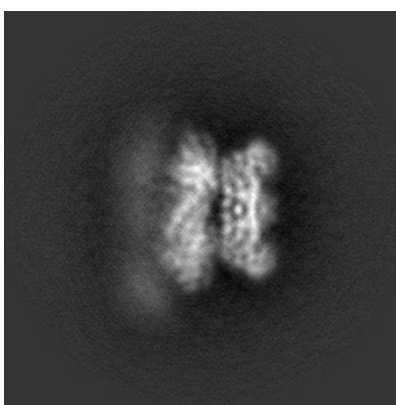
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

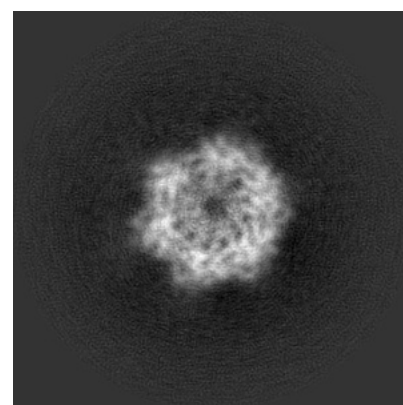
#### 6.1.1 Primary map



X



Y

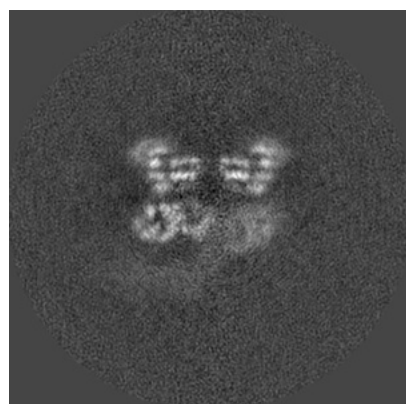


Z

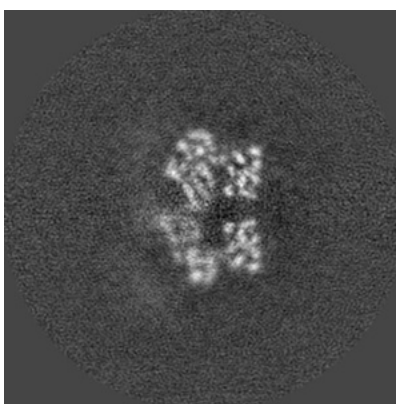
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

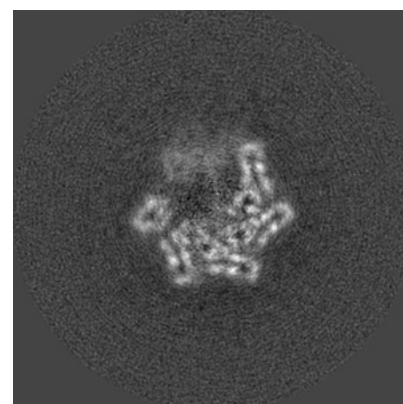
#### 6.2.1 Primary map



X Index: 128



Y Index: 128

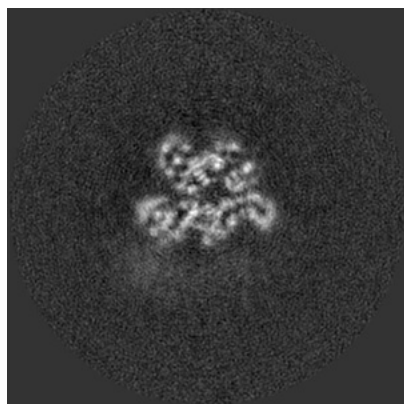


Z Index: 128

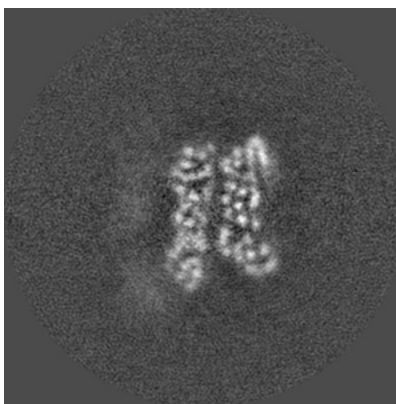
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

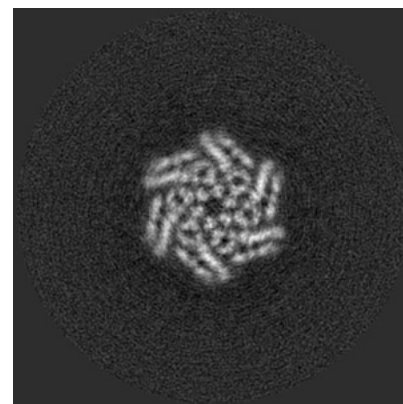
### 6.3.1 Primary map



X Index: 154



Y Index: 112



Z Index: 158

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



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 6.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.



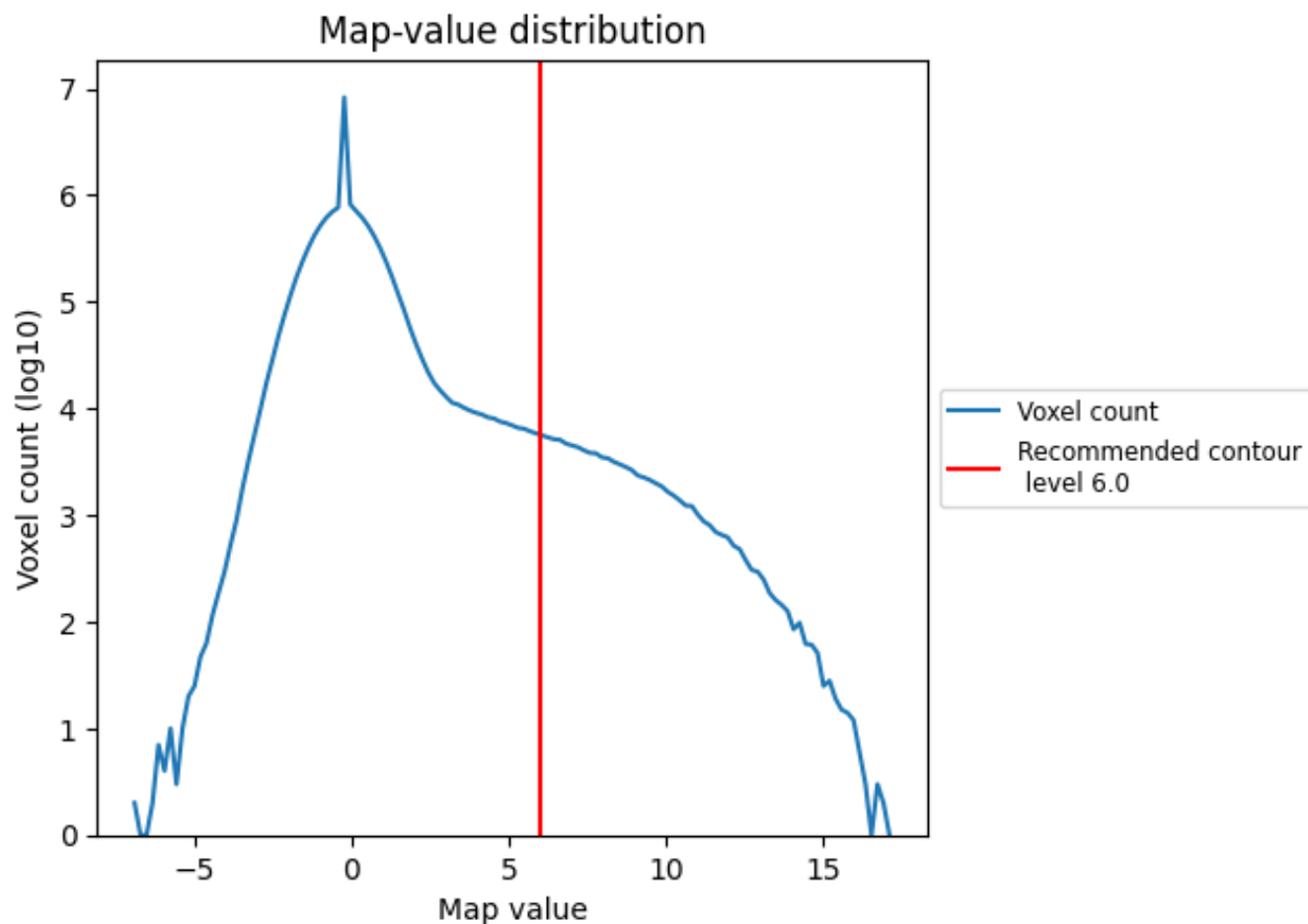
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

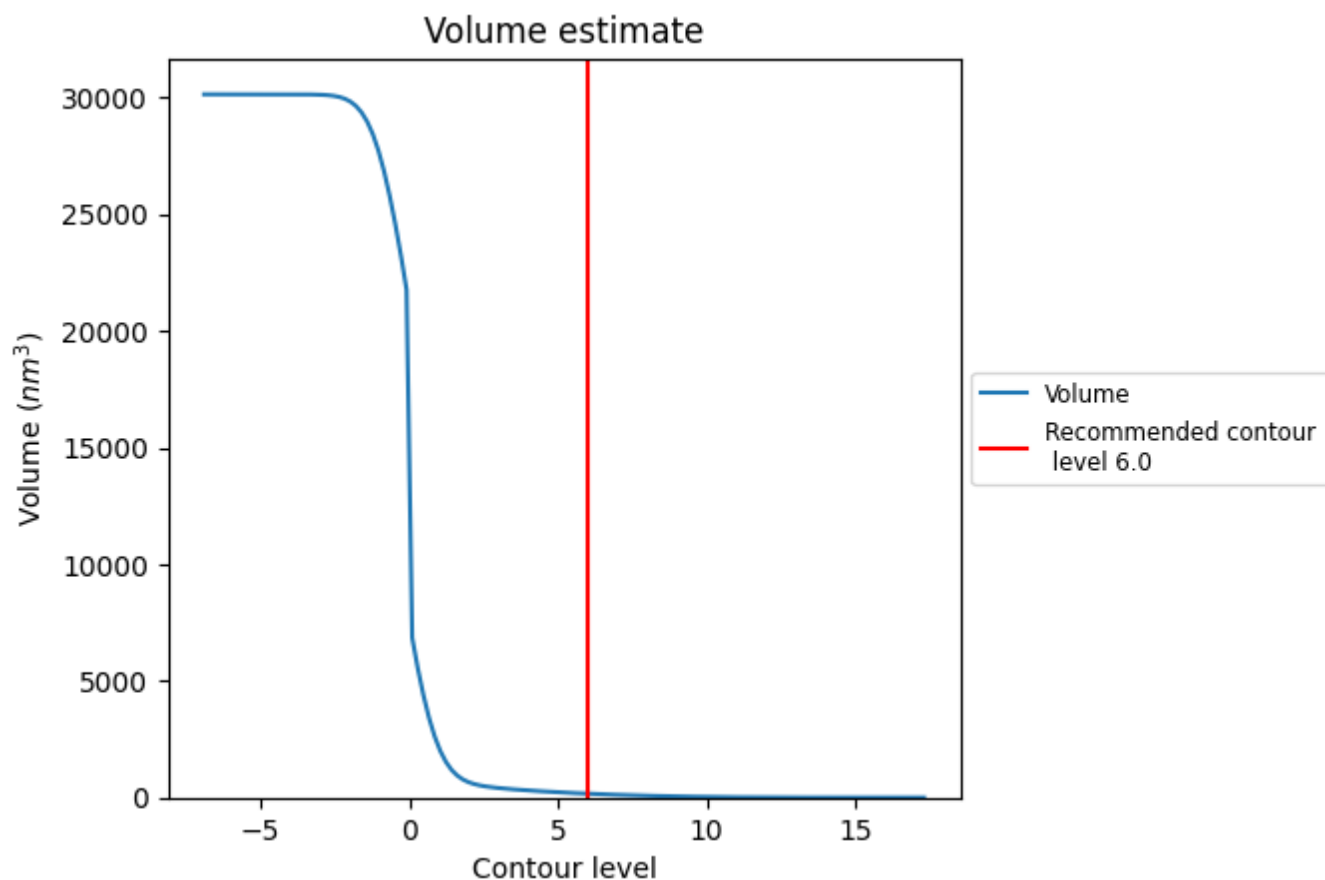
This section contains the results of statistical analysis of the map.

### 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.

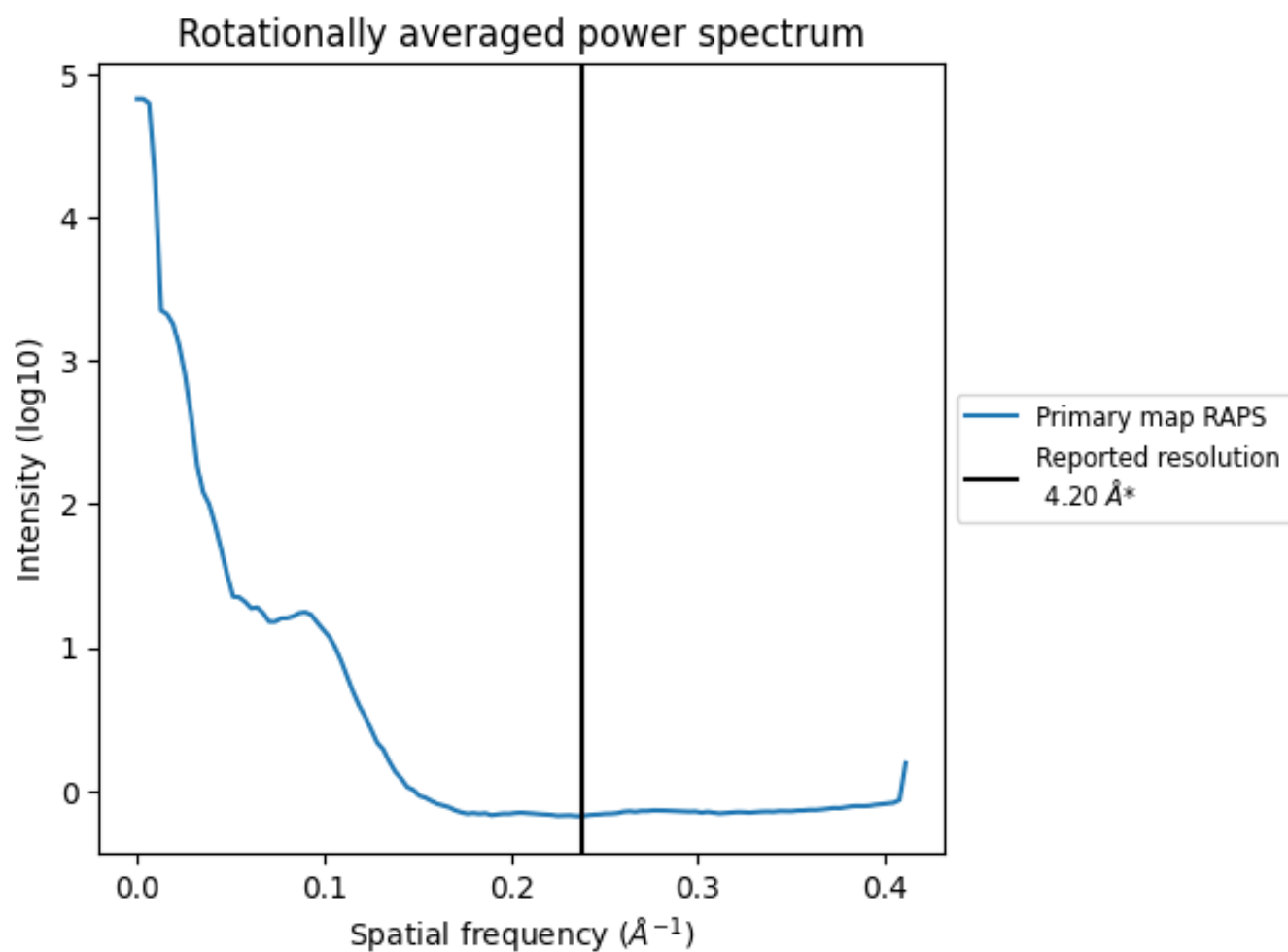
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 166 nm<sup>3</sup>; this corresponds to an approximate mass of 150 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 0.238 Å<sup>-1</sup>

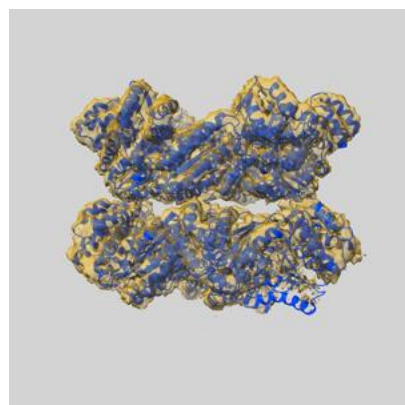
## 8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

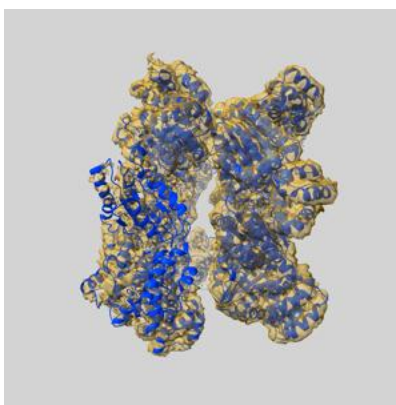
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-6204 and PDB model 3J94. Per-residue inclusion information can be found in section [3](#) on page [6](#).

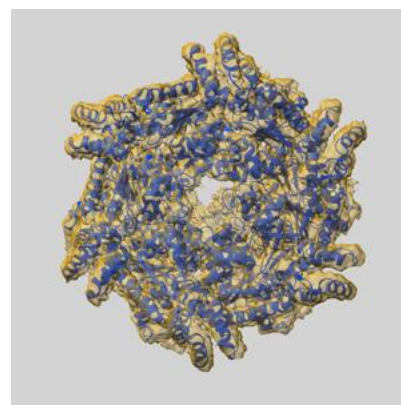
### 9.1 Map-model overlay [i](#)



X



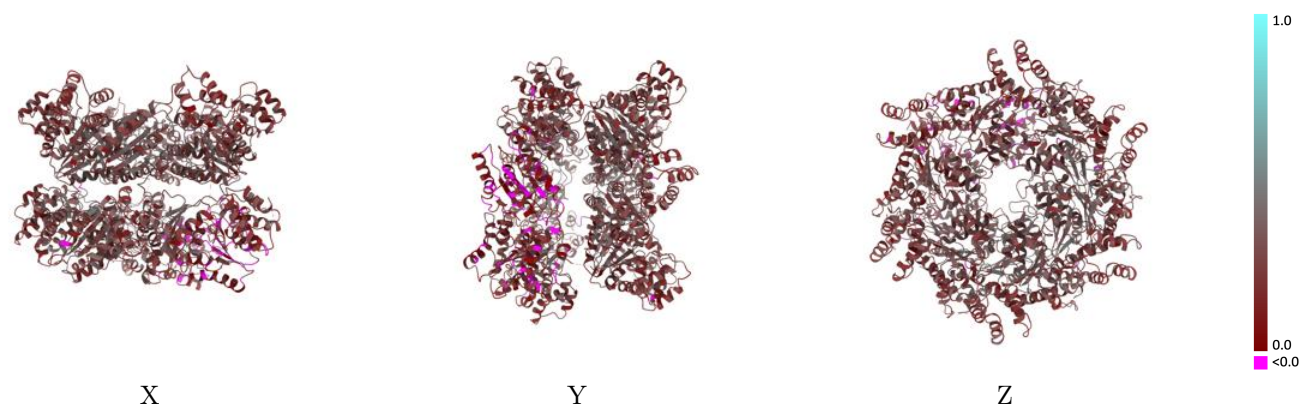
Y



Z

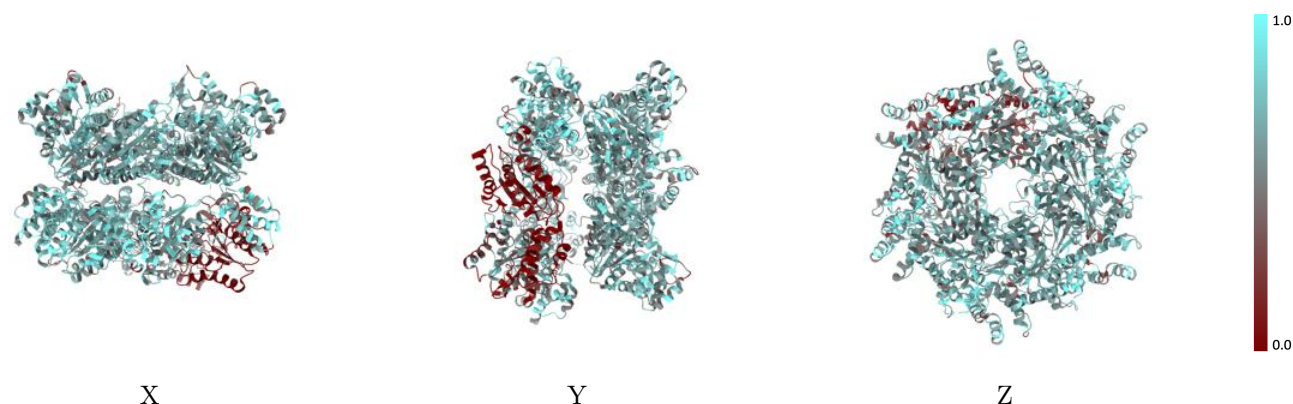
The images above show the 3D surface view of the map at the recommended contour level 6.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



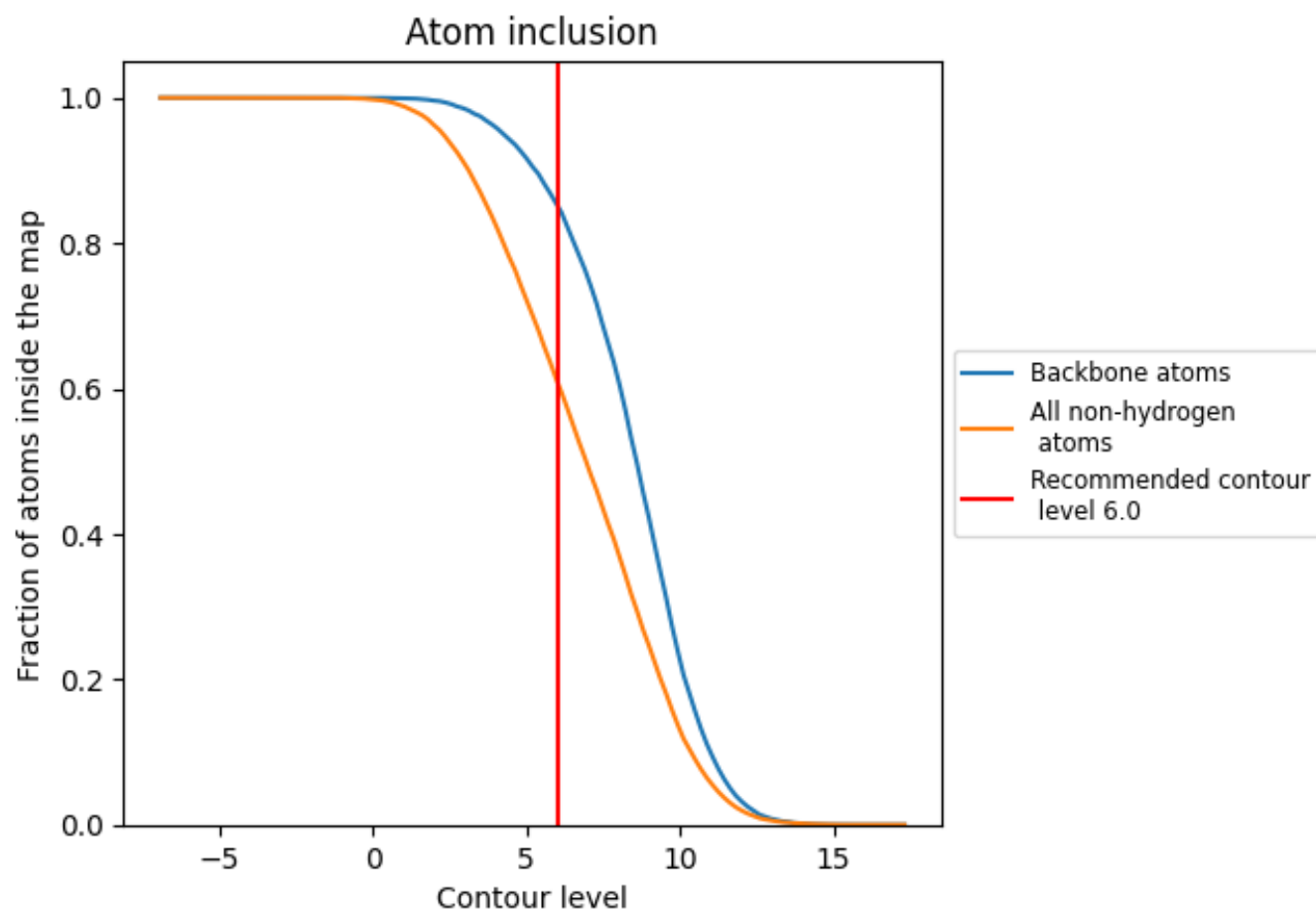
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (6.0).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 61% 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 (6.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6103	<div></div> 0.2550
A	<div></div> 0.6659	<div></div> 0.2720
B	<div></div> 0.6871	<div></div> 0.2850
C	<div></div> 0.6739	<div></div> 0.2830
D	<div></div> 0.6540	<div></div> 0.2640
E	<div></div> 0.5849	<div></div> 0.2330
F	<div></div> 0.3846	<div></div> 0.1860

1.0

0.0

<0.0