



## wwPDB EM Validation Summary Report ⓘ

Jul 8, 2024 – 01:47 am BST

PDB ID : 7QV9  
EMDB ID : EMD-14171  
Title : CryoEM structure of bacterial transcription intermediate complex mediated by activator PspF  
Authors : Ye, F.Z.; Zhang, X.D.  
Deposited on : 2022-01-20  
Resolution : 3.50 Å(reported)  
Based on initial model : 5NSS

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>.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

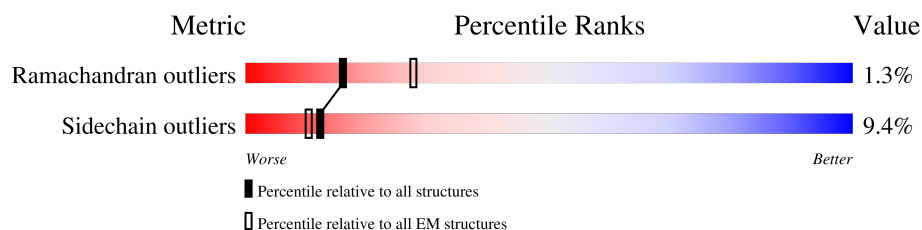
EMDB validation analysis : 0.0.1.dev92  
Mogul : 1.8.4, CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.37.1

# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:  
*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.50 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)
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	329	 91% 6%
1	B	329	 65% 32%
2	C	1342	 97%
3	D	1407	 92% 5%
4	E	91	 81% 19%
5	N	63	 41% 16% 43%
6	T	63	 38% 19% 43%
7	a	295	 50% 33% 12%
7	b	295	 58% 25% 13%

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
7	c	295	
7	d	295	
7	e	295	
7	f	295	
8	M	477	

## 2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 41491 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	309	Total	C	N	O	S	0	0
			2322	1453	407	455	7		
1	B	223	Total	C	N	O	S	0	0
			1676	1045	294	332	5		

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	1341	Total	C	N	O	S	0	0
			10125	6360	1761	1964	40		

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	1334	Total	C	N	O	S	0	0
			9634	6052	1730	1814	38		

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	74	Total	C	N	O	S	0	0
			546	335	109	101	1		

- Molecule 5 is a DNA chain called Non-template promoter DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	N	36	Total	C	N	O	P	0	0
			738	349	137	216	36		

- Molecule 6 is a DNA chain called Template promoter DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	T	36	Total	C	N	O	P	0	0
			738	349	137	216	36		

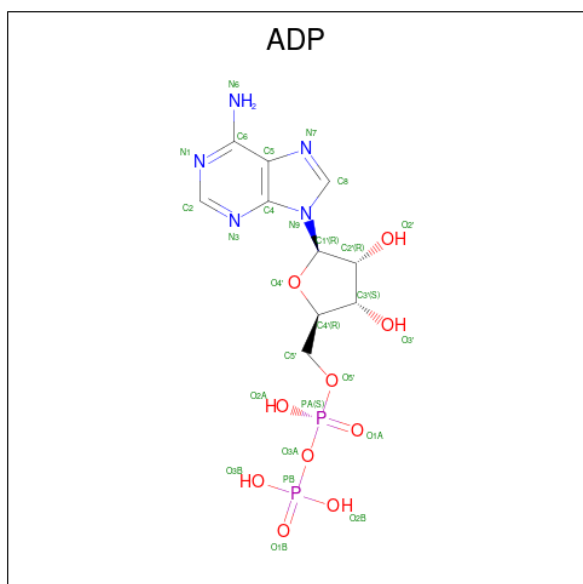
- Molecule 7 is a protein called Transcription activator PspF.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	a	259	Total	C	N	O	S	0	0
			2052	1301	361	380	10		
7	b	258	Total	C	N	O	S	0	0
			2052	1302	360	380	10		
7	c	256	Total	C	N	O	S	0	0
			2000	1270	349	372	9		
7	d	256	Total	C	N	O	S	0	0
			2041	1295	361	376	9		
7	f	259	Total	C	N	O	S	0	0
			2053	1304	358	381	10		
7	e	259	Total	C	N	O	S	0	0
			2062	1308	364	381	9		

- Molecule 8 is a protein called RNA polymerase sigma-54 factor.

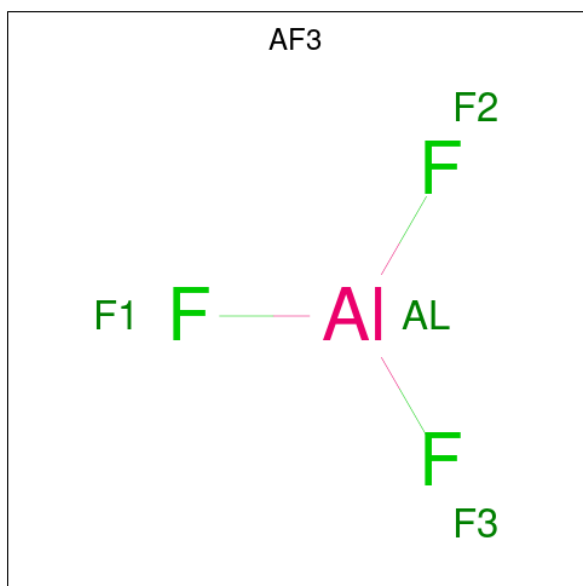
Mol	Chain	Residues	Atoms					AltConf	Trace
8	M	417	Total	C	N	O	S	0	0
			3301	2071	572	646	12		

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
9	a	1	Total	C	N	O	P	0
			27	10	5	10	2	
9	b	1	Total	C	N	O	P	0
			27	10	5	10	2	
9	c	1	Total	C	N	O	P	0
			27	10	5	10	2	
9	f	1	Total	C	N	O	P	0
			27	10	5	10	2	
9	e	1	Total	C	N	O	P	0
			27	10	5	10	2	

- Molecule 10 is ALUMINUM FLUORIDE (three-letter code: AF3) (formula:  $\text{AlF}_3$ ) (labeled as "Ligand of Interest" by depositor).




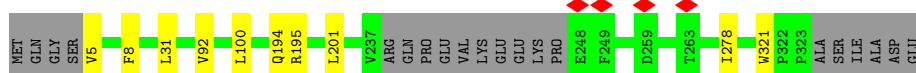
Mol	Chain	Residues	Atoms			AltConf
10	a	1	Total	Al	F	0
			4	1	3	
10	b	1	Total	Al	F	0
			4	1	3	
10	f	1	Total	Al	F	0
			4	1	3	
10	e	1	Total	Al	F	0
			4	1	3	

### 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.

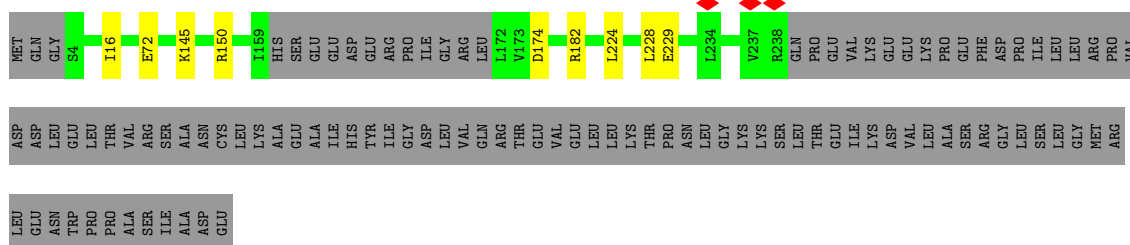
- Molecule 1: DNA-directed RNA polymerase subunit alpha

Chain A: 



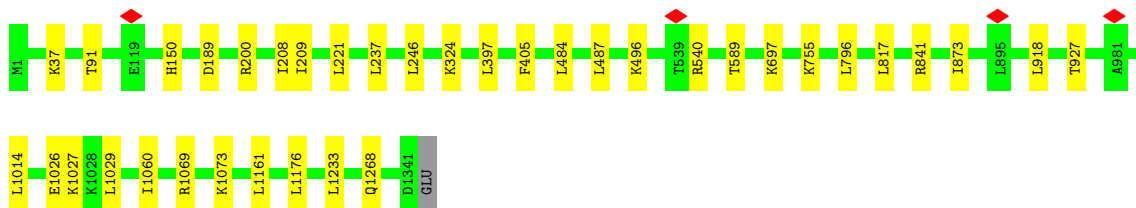
- Molecule 1: DNA-directed RNA polymerase subunit alpha

Chain B: 




- Molecule 2: DNA-directed RNA polymerase subunit beta

Chain C: 

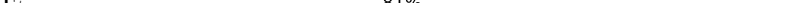


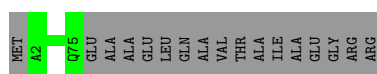
- Molecule 3: DNA-directed RNA polymerase subunit beta'

Chain D: 



- Molecule 4: DNA-directed RNA polymerase subunit omega

Chain E:  81% 19%



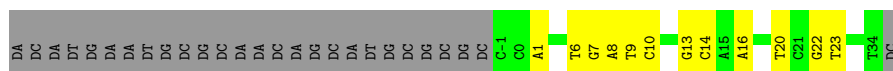
- Molecule 5: Non-template promoter DNA

Chain N:  41% 16% 43%



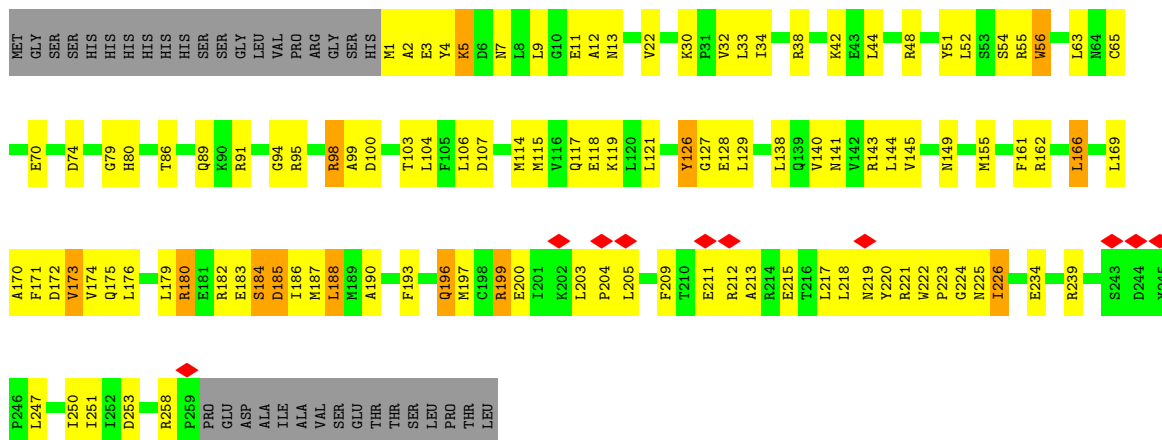
- Molecule 6: Template promoter DNA

Chain T:  38% 19% 43%



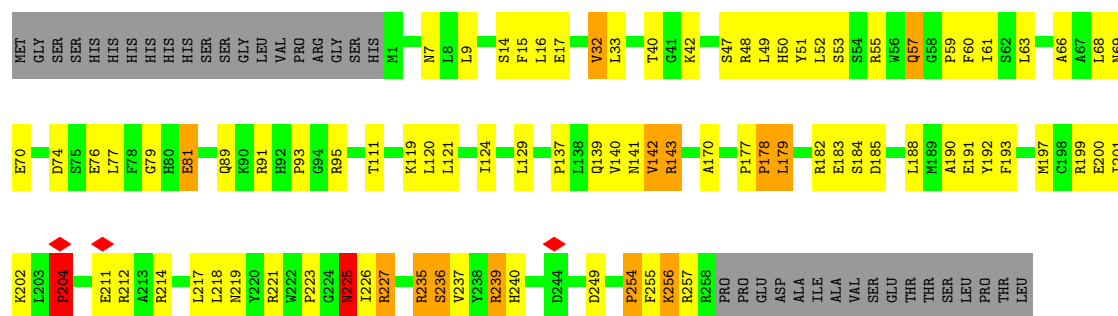
- Molecule 7: Transcription activator PspF

Chain a:  50% 33% 0 12%



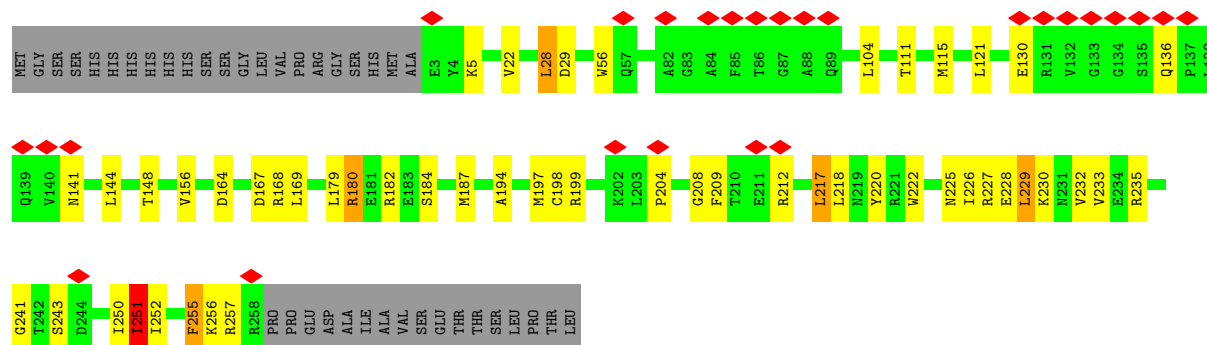
- Molecule 7: Transcription activator PspF

Chain b: 



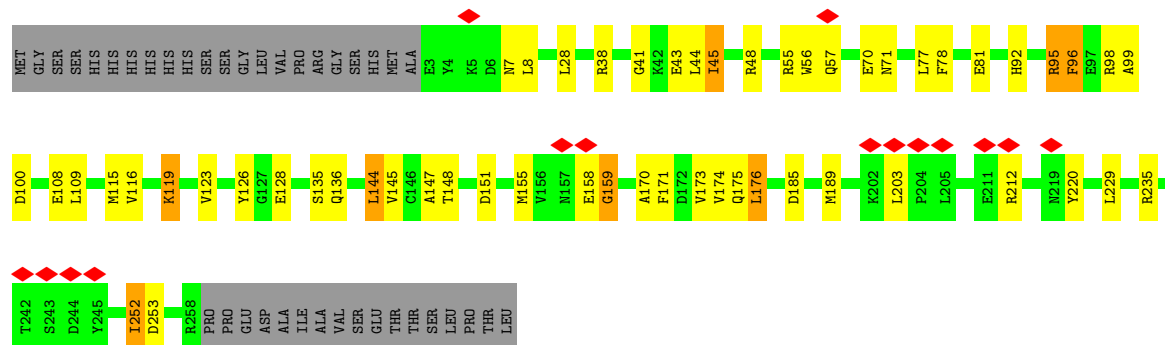
• Molecule 7: Transcription activator PspF

Chain c: 



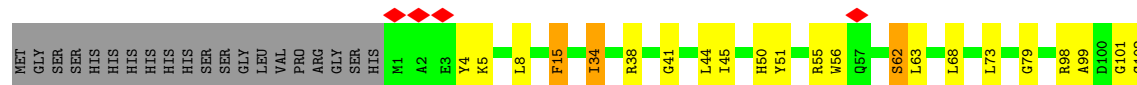
• Molecule 7: Transcription activator PspF

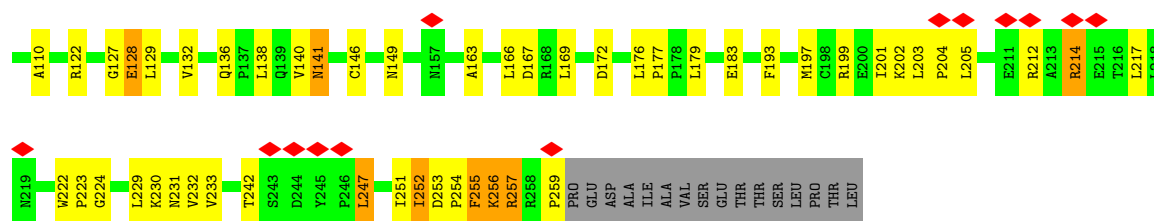
Chain d: 



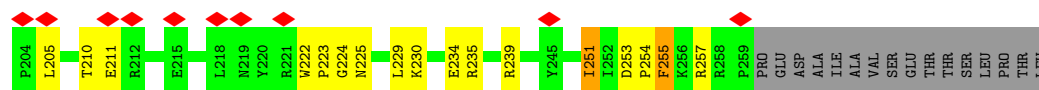
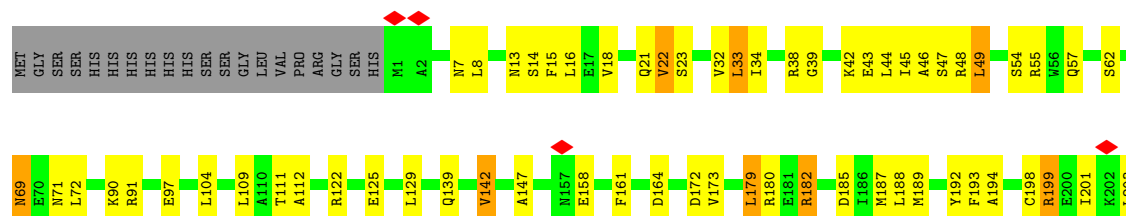
• Molecule 7: Transcription activator PspF

Chain f: 

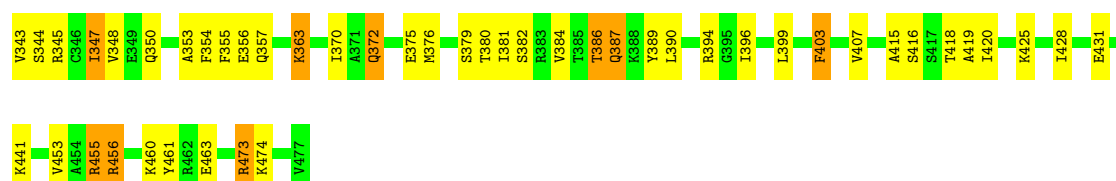
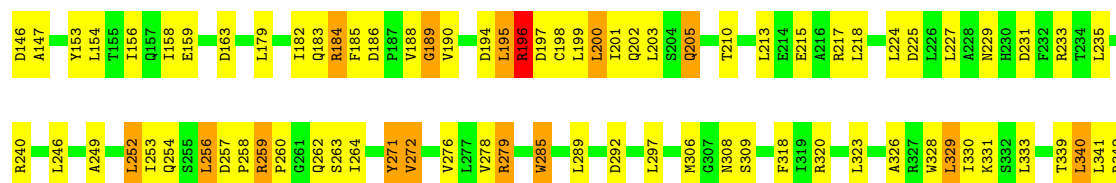
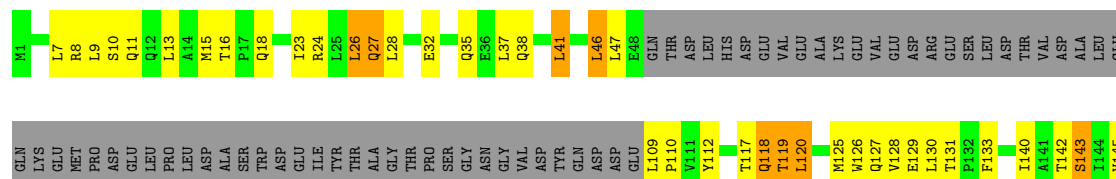




• Molecule 7: Transcription activator PspF



• Molecule 8: RNA polymerase sigma-54 factor



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	33285	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 ( $e^-/\text{\AA}^2$ )	50	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	2600	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.302	Depositor
Minimum map value	-0.152	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.012	Depositor
Map size (Å)	308.0, 308.0, 308.0	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	Depositor

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AF3, ADP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.23	0/2351	0.44	0/3202
1	B	0.22	0/1693	0.43	0/2300
2	C	0.23	0/10283	0.40	0/13940
3	D	0.23	0/9768	0.41	1/13270 (0.0%)
4	E	0.21	0/547	0.36	0/740
5	N	0.90	5/827 (0.6%)	1.32	12/1274 (0.9%)
6	T	1.28	9/827 (1.1%)	1.43	18/1274 (1.4%)
7	a	1.89	61/2093 (2.9%)	2.03	83/2833 (2.9%)
7	b	1.87	69/2093 (3.3%)	2.04	78/2832 (2.8%)
7	c	1.21	27/2038 (1.3%)	1.49	37/2762 (1.3%)
7	d	1.17	14/2082 (0.7%)	1.54	42/2817 (1.5%)
7	e	1.52	44/2104 (2.1%)	1.73	66/2848 (2.3%)
7	f	1.74	52/2095 (2.5%)	1.74	70/2837 (2.5%)
8	M	1.80	84/3348 (2.5%)	1.95	145/4533 (3.2%)
All	All	1.04	365/42149 (0.9%)	1.19	552/57462 (1.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
7	a	0	1
7	b	0	1
7	d	0	1
All	All	0	3

The worst 5 of 365 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	a	98	ARG	CZ-NH2	28.28	1.69	1.33

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	b	143	ARG	CZ-NH1	24.18	1.64	1.33
7	f	128	GLU	CD-OE1	21.25	1.49	1.25
7	a	48	ARG	CZ-NH1	20.99	1.60	1.33
7	f	193	PHE	CG-CD2	19.56	1.68	1.38

The worst 5 of 552 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	a	48	ARG	NE-CZ-NH1	38.39	139.50	120.30
7	b	143	ARG	NE-CZ-NH2	-38.03	101.29	120.30
7	a	98	ARG	NE-CZ-NH1	-34.60	103.00	120.30
7	d	100	ASP	CB-CG-OD2	-30.82	90.56	118.30
7	a	48	ARG	NE-CZ-NH2	-22.94	108.83	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
7	a	162	ARG	Mainchain
7	b	219	ASN	Mainchain
7	d	92	HIS	Mainchain

## 5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

## 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	305/329 (93%)	282 (92%)	22 (7%)	1 (0%)	41	75
1	B	219/329 (67%)	203 (93%)	16 (7%)	0	100	100

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	1339/1342 (100%)	1264 (94%)	75 (6%)	0	100	100
3	D	1322/1407 (94%)	1226 (93%)	93 (7%)	3 (0%)	47	81
4	E	72/91 (79%)	70 (97%)	2 (3%)	0	100	100
7	a	257/295 (87%)	200 (78%)	43 (17%)	14 (5%)	2	17
7	b	256/295 (87%)	216 (84%)	27 (10%)	13 (5%)	2	19
7	c	254/295 (86%)	221 (87%)	31 (12%)	2 (1%)	19	58
7	d	254/295 (86%)	226 (89%)	23 (9%)	5 (2%)	7	39
7	e	257/295 (87%)	217 (84%)	30 (12%)	10 (4%)	3	25
7	f	257/295 (87%)	225 (88%)	22 (9%)	10 (4%)	3	25
8	M	413/477 (87%)	360 (87%)	42 (10%)	11 (3%)	5	33
All	All	5205/5745 (91%)	4710 (90%)	426 (8%)	69 (1%)	16	48

5 of 69 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	a	12	ALA
7	a	56	TRP
7	a	121	LEU
7	b	81	GLU
7	b	93	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	A	249/286 (87%)	240 (96%)	9 (4%)	35	66
1	B	180/286 (63%)	171 (95%)	9 (5%)	24	58
2	C	1047/1157 (90%)	1010 (96%)	37 (4%)	36	67
3	D	913/1168 (78%)	872 (96%)	41 (4%)	27	61
4	E	53/75 (71%)	53 (100%)	0	100	100
7	a	219/252 (87%)	156 (71%)	63 (29%)	0	2

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	b	219/252 (87%)	181 (83%)	38 (17%)	2	11
7	c	208/252 (82%)	181 (87%)	27 (13%)	4	21
7	d	218/252 (86%)	191 (88%)	27 (12%)	4	23
7	e	220/252 (87%)	192 (87%)	28 (13%)	4	22
7	f	219/252 (87%)	192 (88%)	27 (12%)	4	23
8	M	367/423 (87%)	286 (78%)	81 (22%)	1	5
All	All	4112/4907 (84%)	3725 (91%)	387 (9%)	12	35

5 of 387 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
7	d	57	GLN
7	e	49	LEU
7	d	98	ARG
7	f	63	LEU
7	e	179	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 49 such sidechains are listed below:

Mol	Chain	Res	Type
7	f	80	HIS
7	e	139	GLN
7	f	117	GLN
7	e	7	ASN
8	M	3	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

9 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
10	AF3	e	602	-	0,3,3	-	-	-		
9	ADP	f	601	-	24,29,29	3.84	12 (50%)	29,45,45	4.30	24 (82%)
9	ADP	b	601	-	24,29,29	0.96	1 (4%)	29,45,45	1.49	4 (13%)
10	AF3	b	602	-	0,3,3	-	-	-		
9	ADP	a	601	-	24,29,29	0.95	1 (4%)	29,45,45	1.49	4 (13%)
9	ADP	c	400	-	24,29,29	0.96	1 (4%)	29,45,45	1.47	4 (13%)
9	ADP	e	601	-	24,29,29	2.14	7 (29%)	29,45,45	3.31	19 (65%)
10	AF3	f	602	-	0,3,3	-	-	-		
10	AF3	a	602	-	0,3,3	-	-	-		

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
9	ADP	f	601	-	-	5/12/32/32	0/3/3/3
9	ADP	b	601	-	-	5/12/32/32	0/3/3/3
9	ADP	a	601	-	-	3/12/32/32	0/3/3/3
9	ADP	c	400	-	-	2/12/32/32	0/3/3/3
9	ADP	e	601	-	-	1/12/32/32	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	f	601	ADP	O4'-C1'	9.06	1.53	1.41
9	f	601	ADP	C2'-C1'	-9.05	1.40	1.53
9	f	601	ADP	O2'-C2'	7.49	1.60	1.43
9	f	601	ADP	PA-O1A	5.67	1.71	1.50

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	e	601	ADP	PA-O1A	-4.86	1.33	1.50

The worst 5 of 55 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	f	601	ADP	PA-O3A-PB	-8.19	104.71	132.83
9	e	601	ADP	PA-O3A-PB	-6.94	109.00	132.83
9	f	601	ADP	O4'-C4'-C5'	6.52	130.81	109.37
9	f	601	ADP	C5'-C4'-C3'	-6.47	90.95	115.18
9	f	601	ADP	O2A-PA-O1A	6.24	143.08	112.24

There are no chirality outliers.

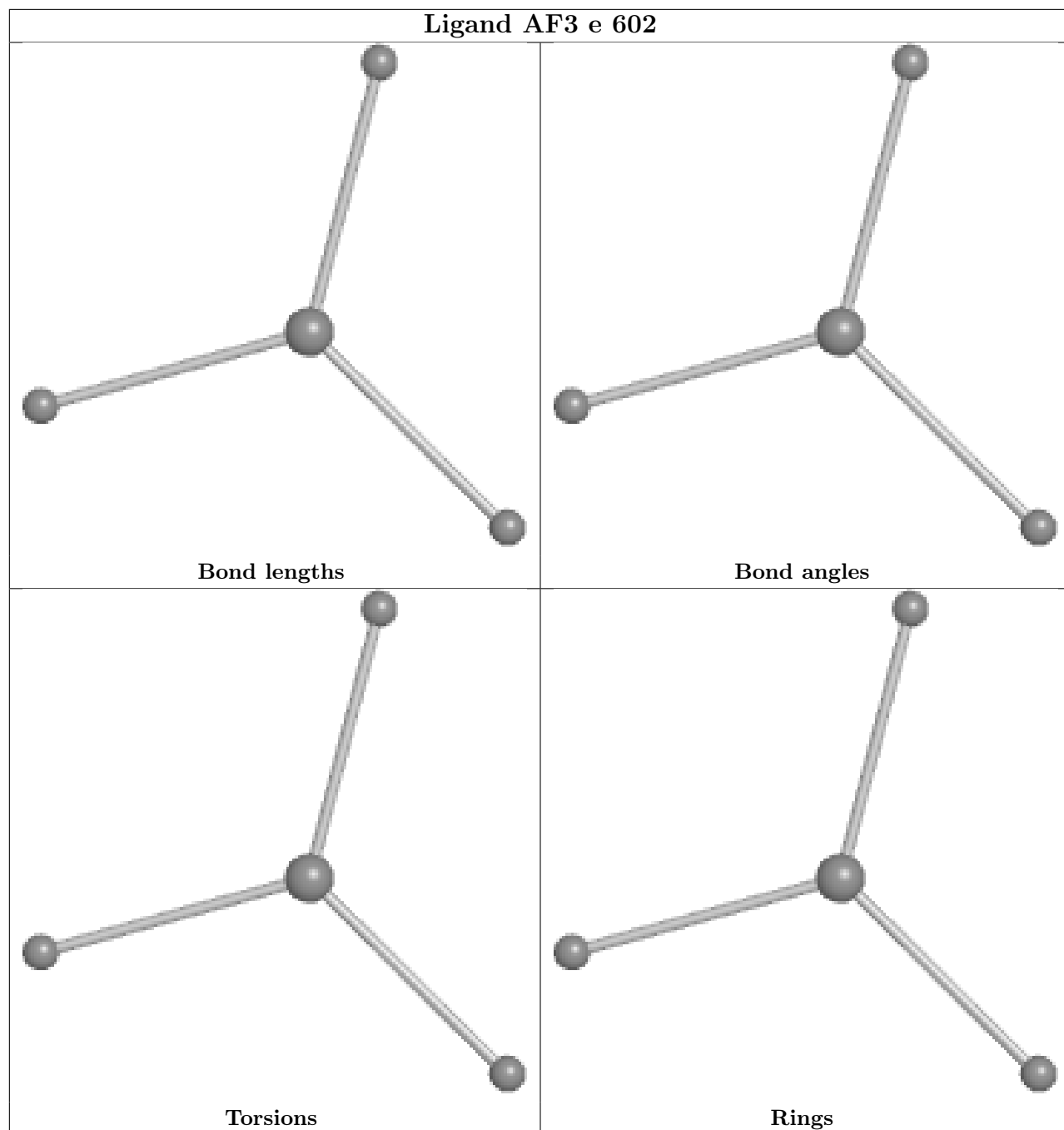
5 of 16 torsion outliers are listed below:

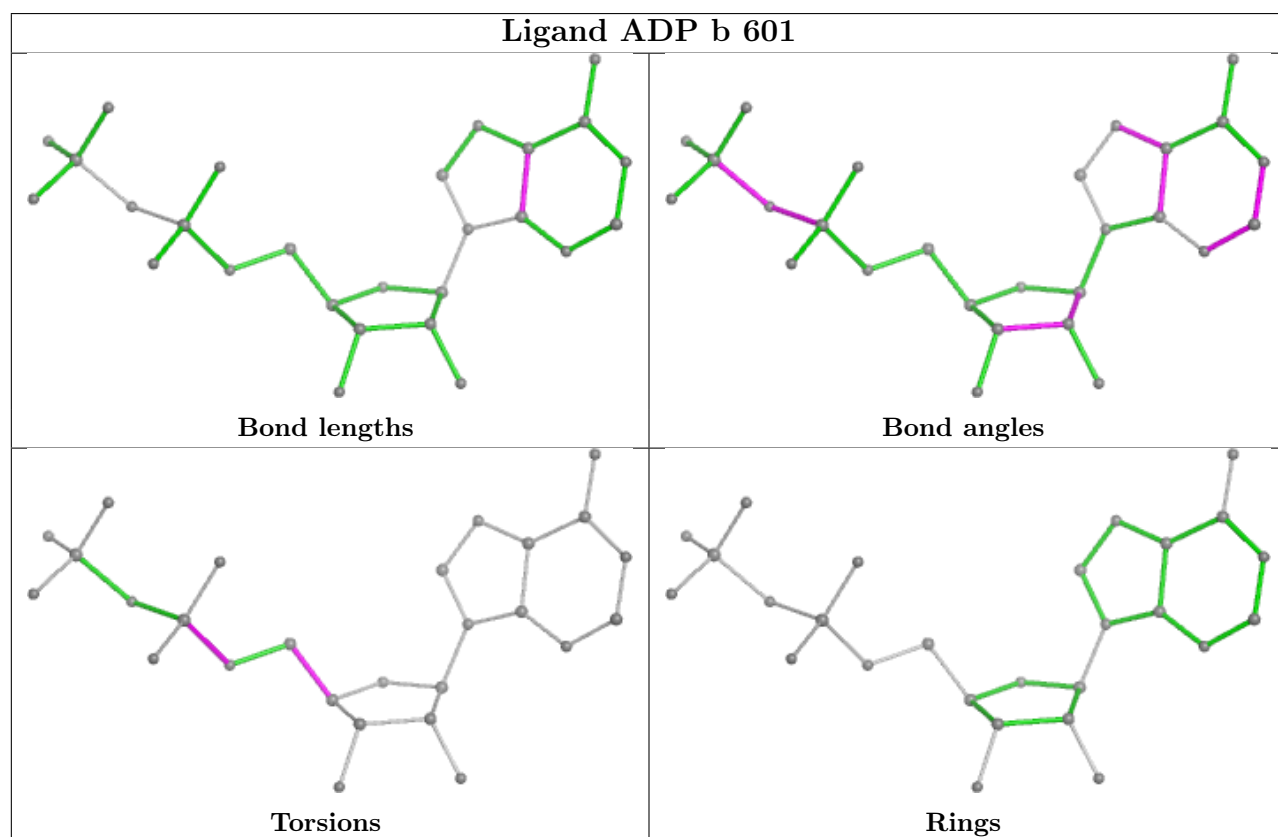
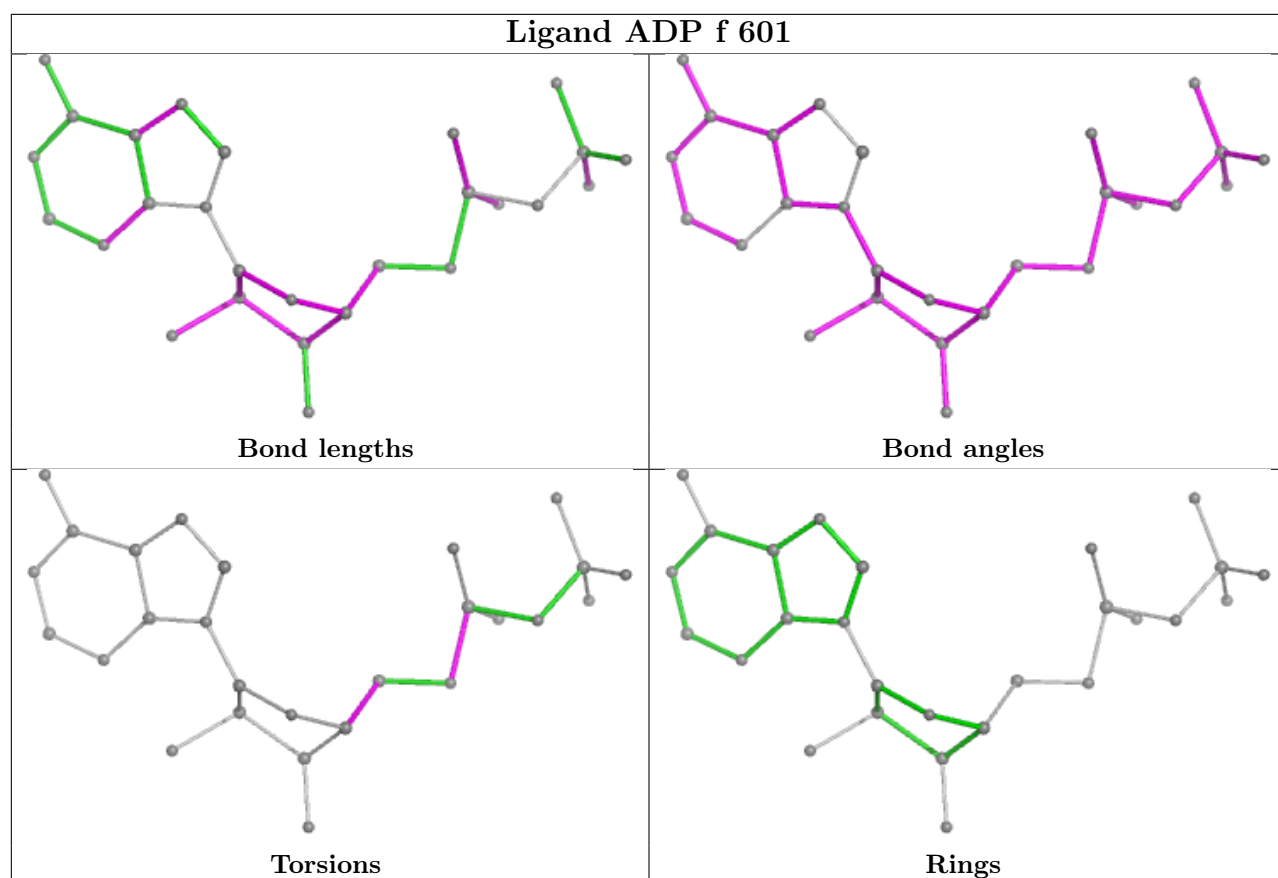
Mol	Chain	Res	Type	Atoms
9	b	601	ADP	C5'-O5'-PA-O2A
9	c	400	ADP	C5'-O5'-PA-O1A
9	f	601	ADP	C5'-O5'-PA-O2A
9	f	601	ADP	C3'-C4'-C5'-O5'
9	f	601	ADP	O4'-C4'-C5'-O5'

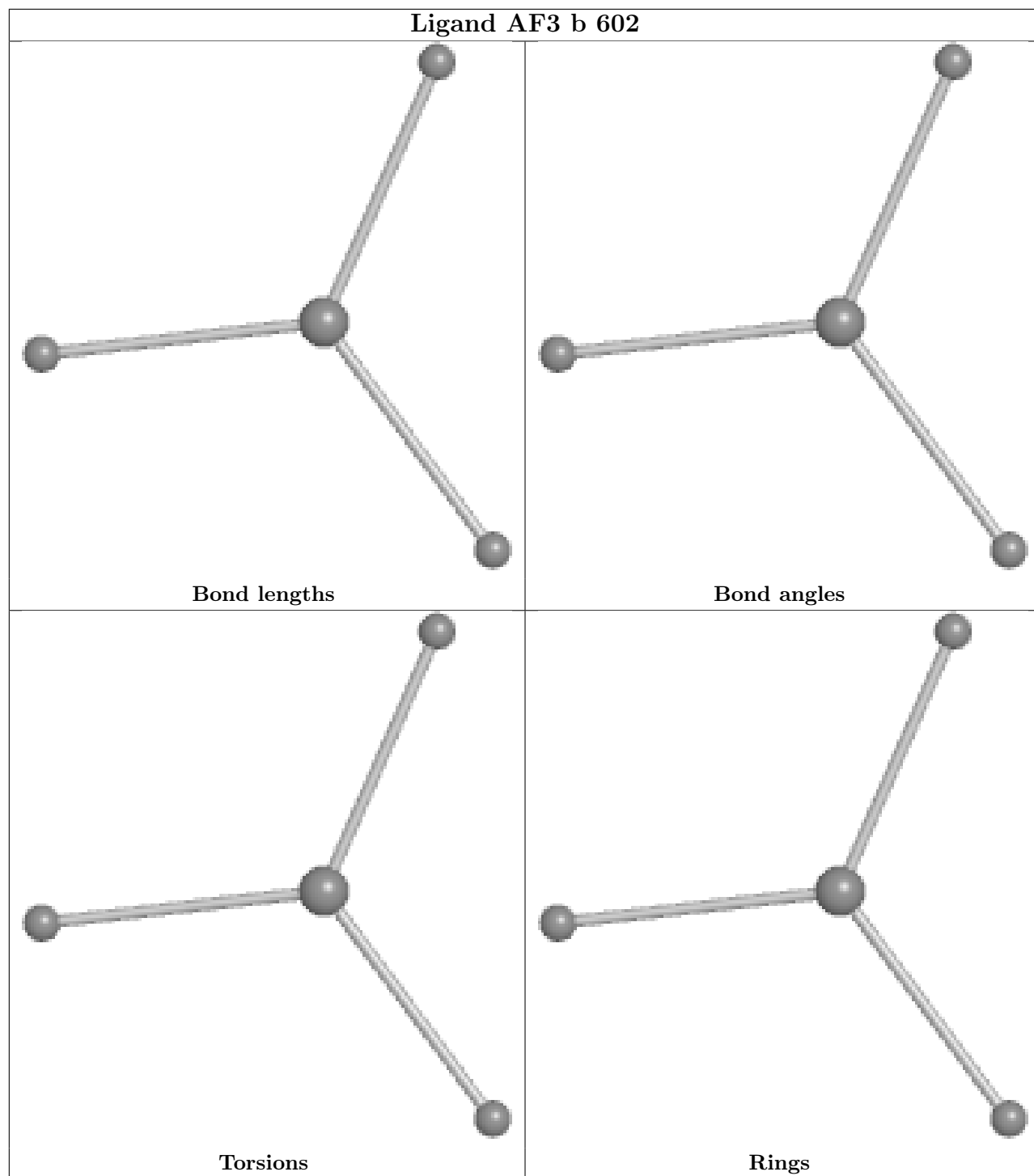
There are no ring outliers.

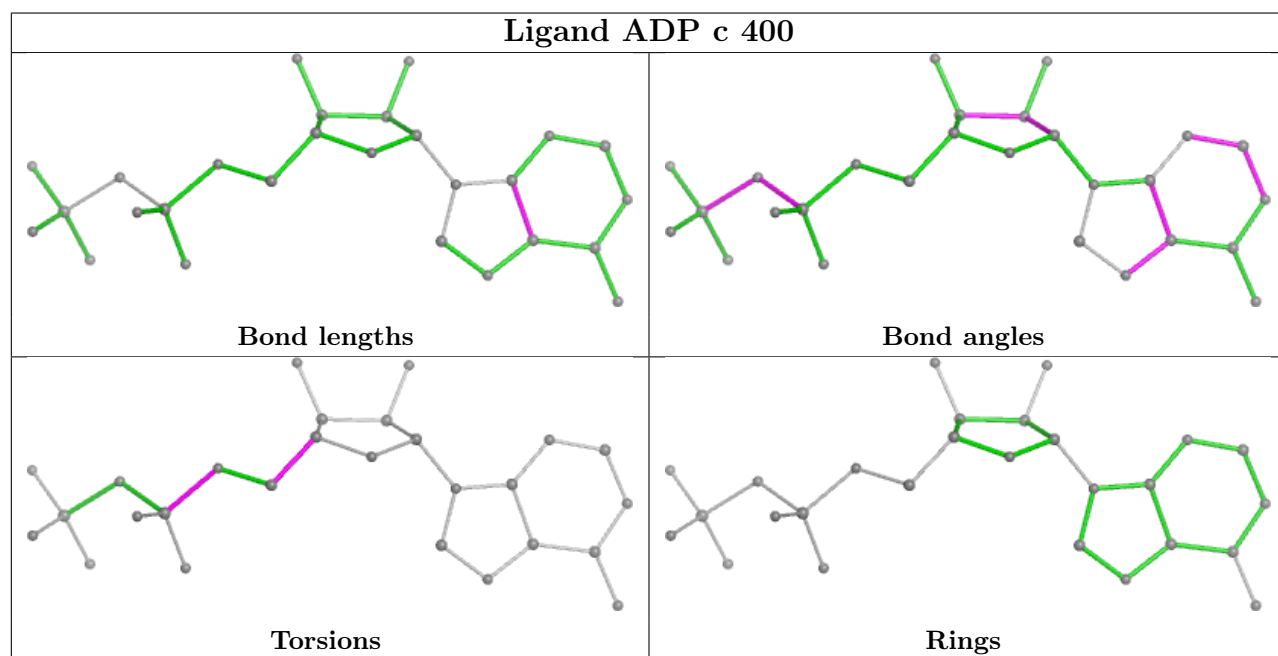
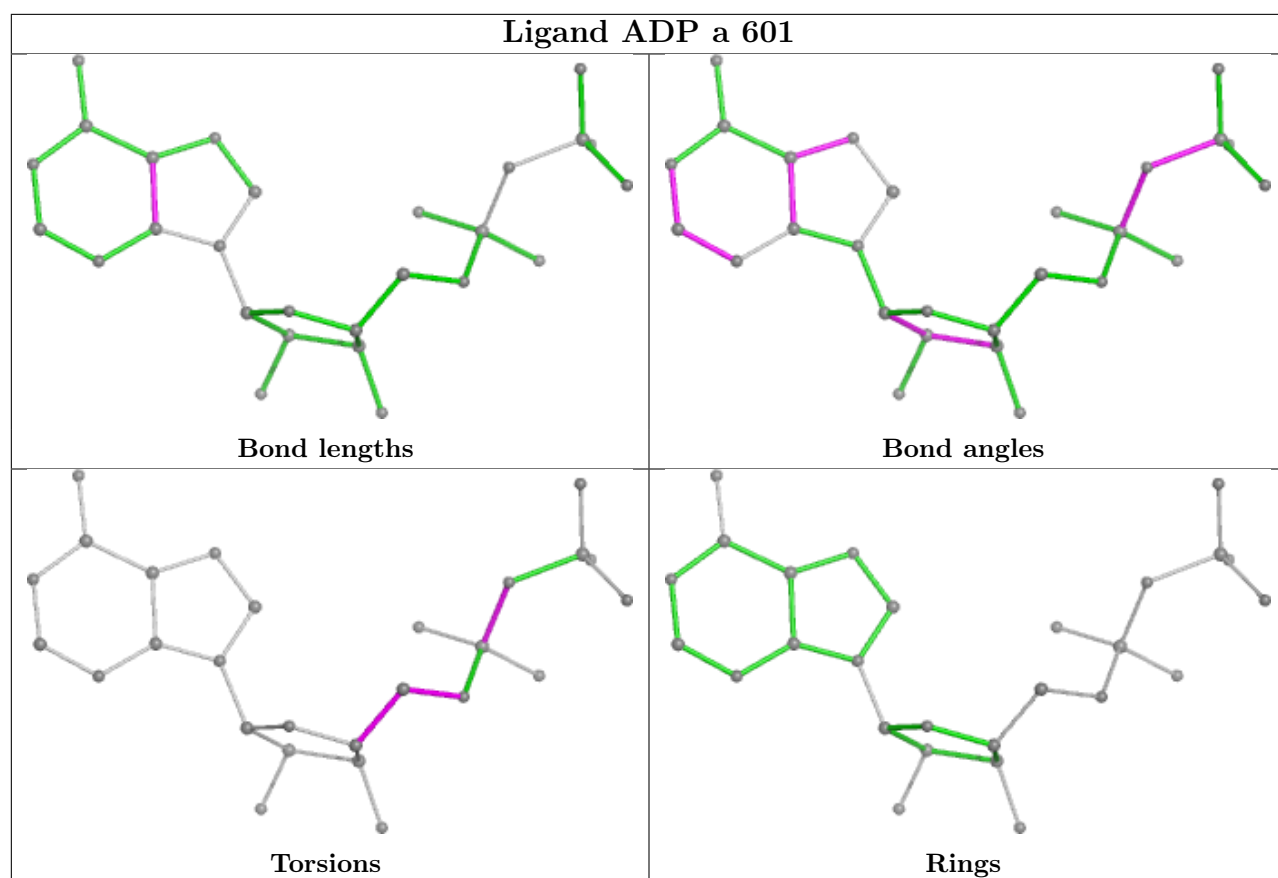
No monomer is involved in short contacts.

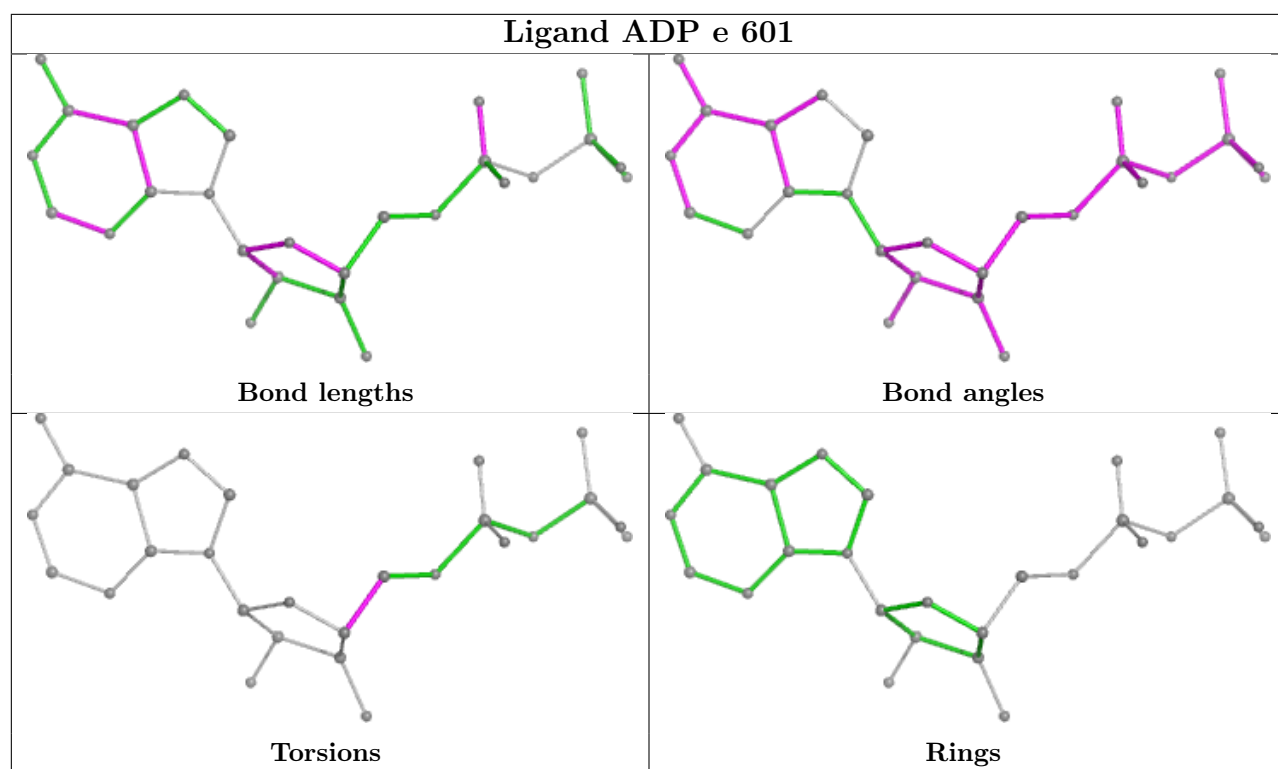
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

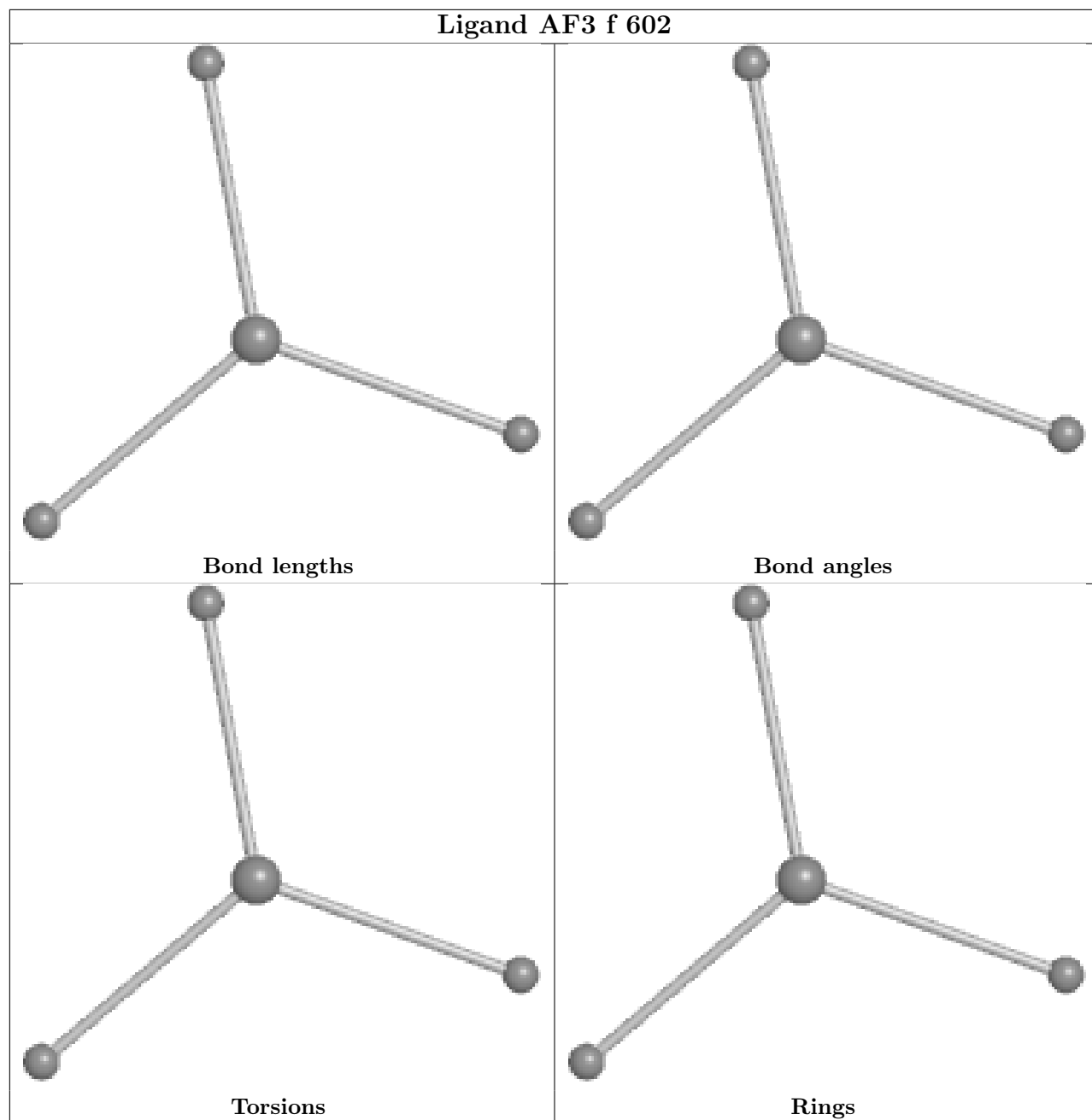


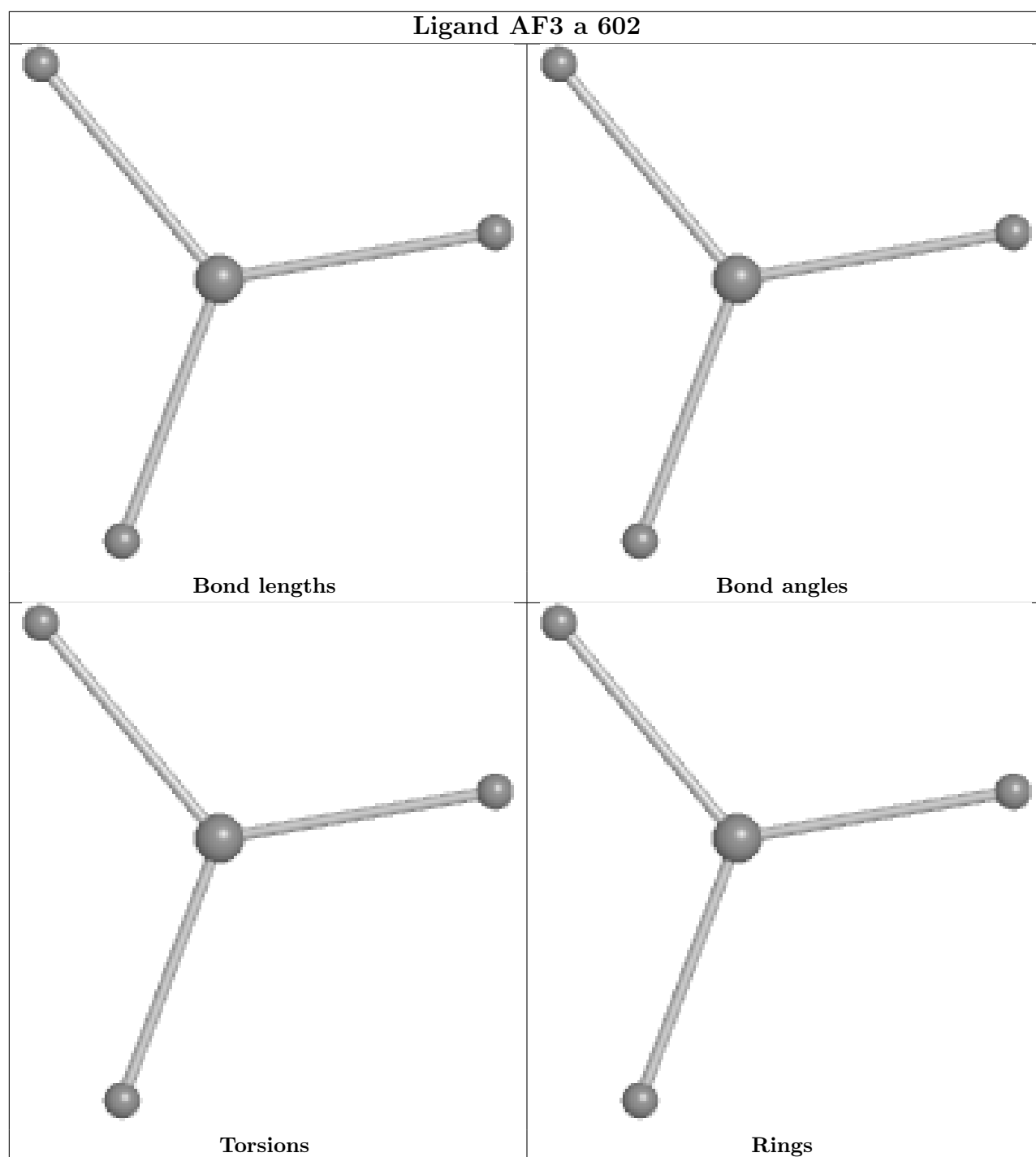












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

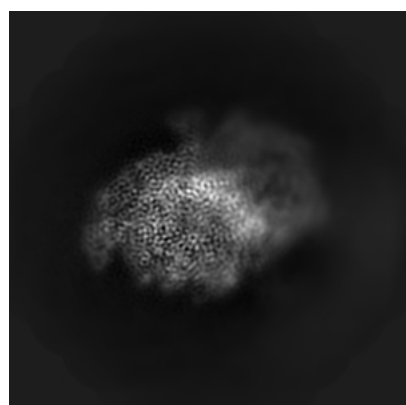
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-14171. These allow visual inspection of the internal detail of the map and identification of artifacts.

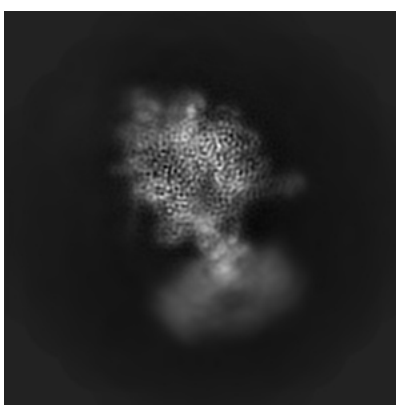
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

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

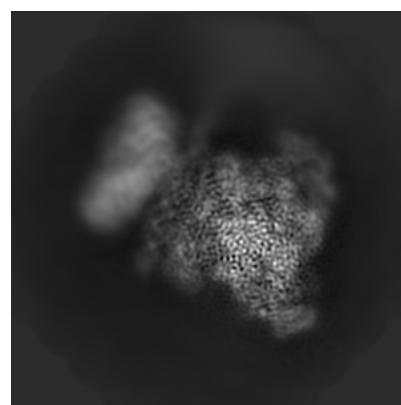
#### 6.1.1 Primary map



X



Y

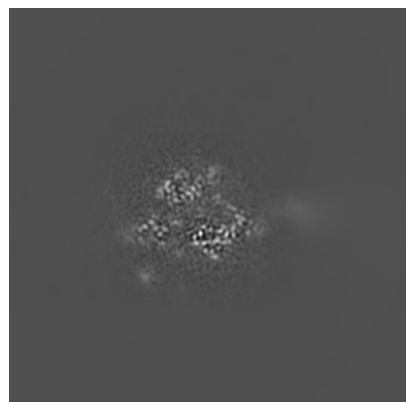


Z

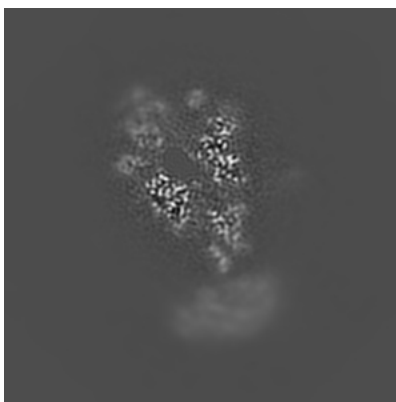
The images above show the map projected in three orthogonal directions.

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

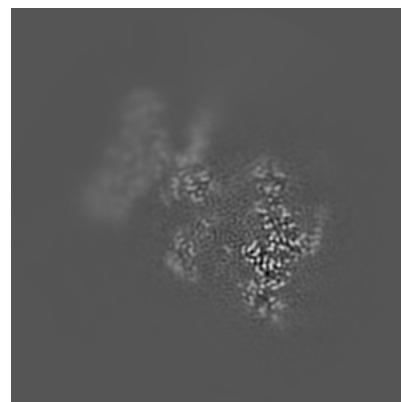
#### 6.2.1 Primary map



X Index: 140



Y Index: 140

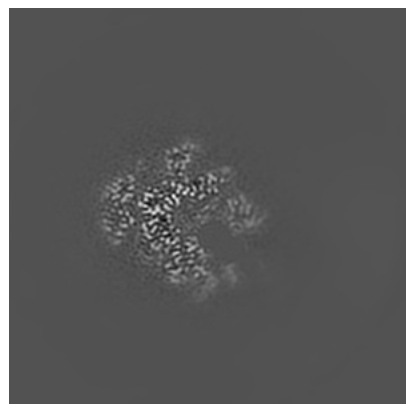


Z Index: 140

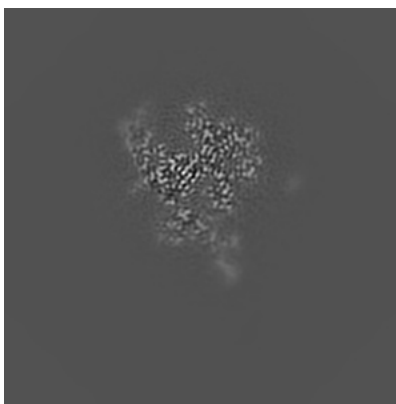
The images above show central slices of the map in three orthogonal directions.

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

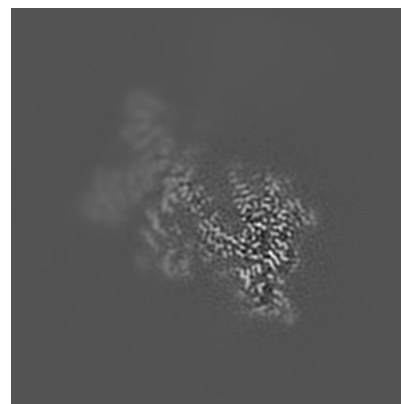
### 6.3.1 Primary map



X Index: 173



Y Index: 111

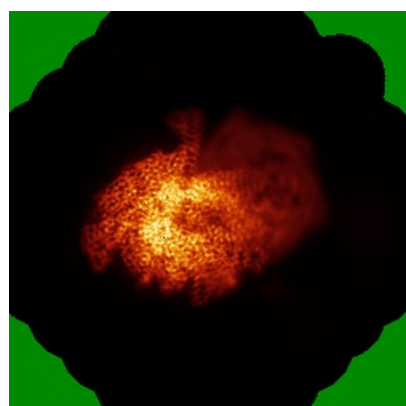


Z Index: 149

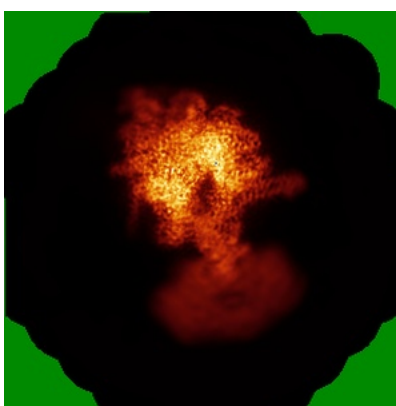
The images above show the largest variance slices of the map in three orthogonal directions.

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

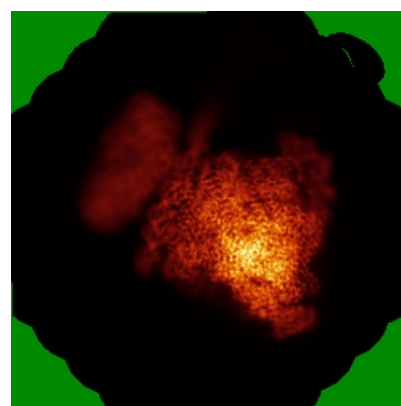
### 6.4.1 Primary map



X



Y

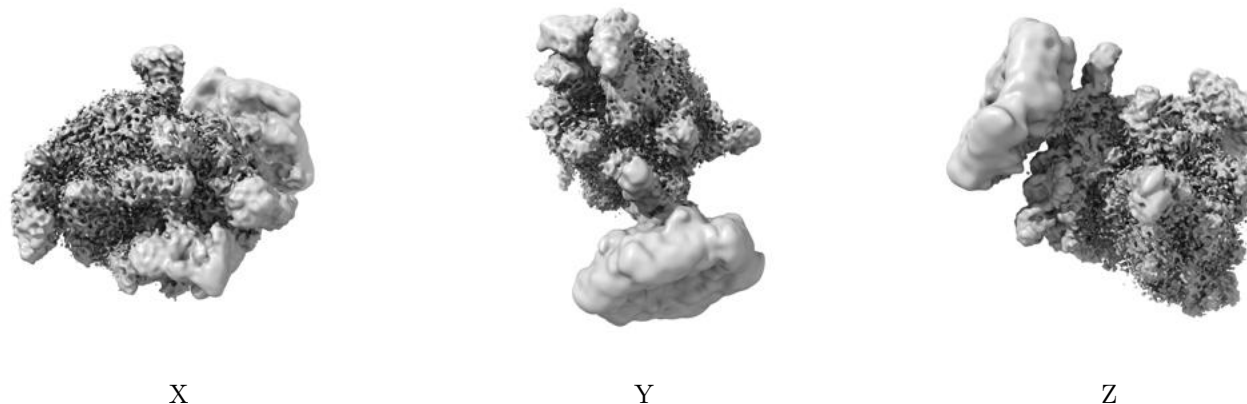


Z

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 0.012. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

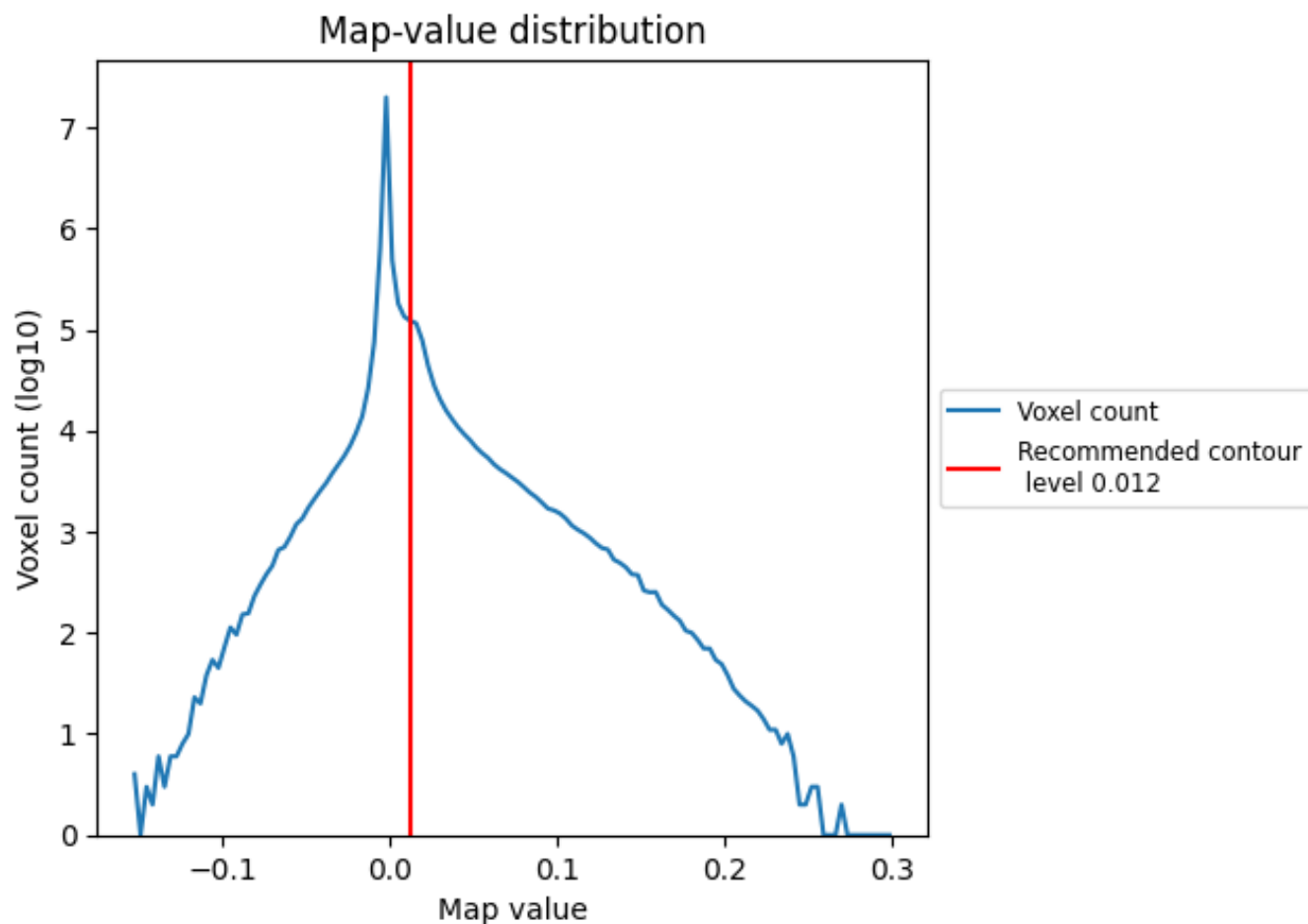
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

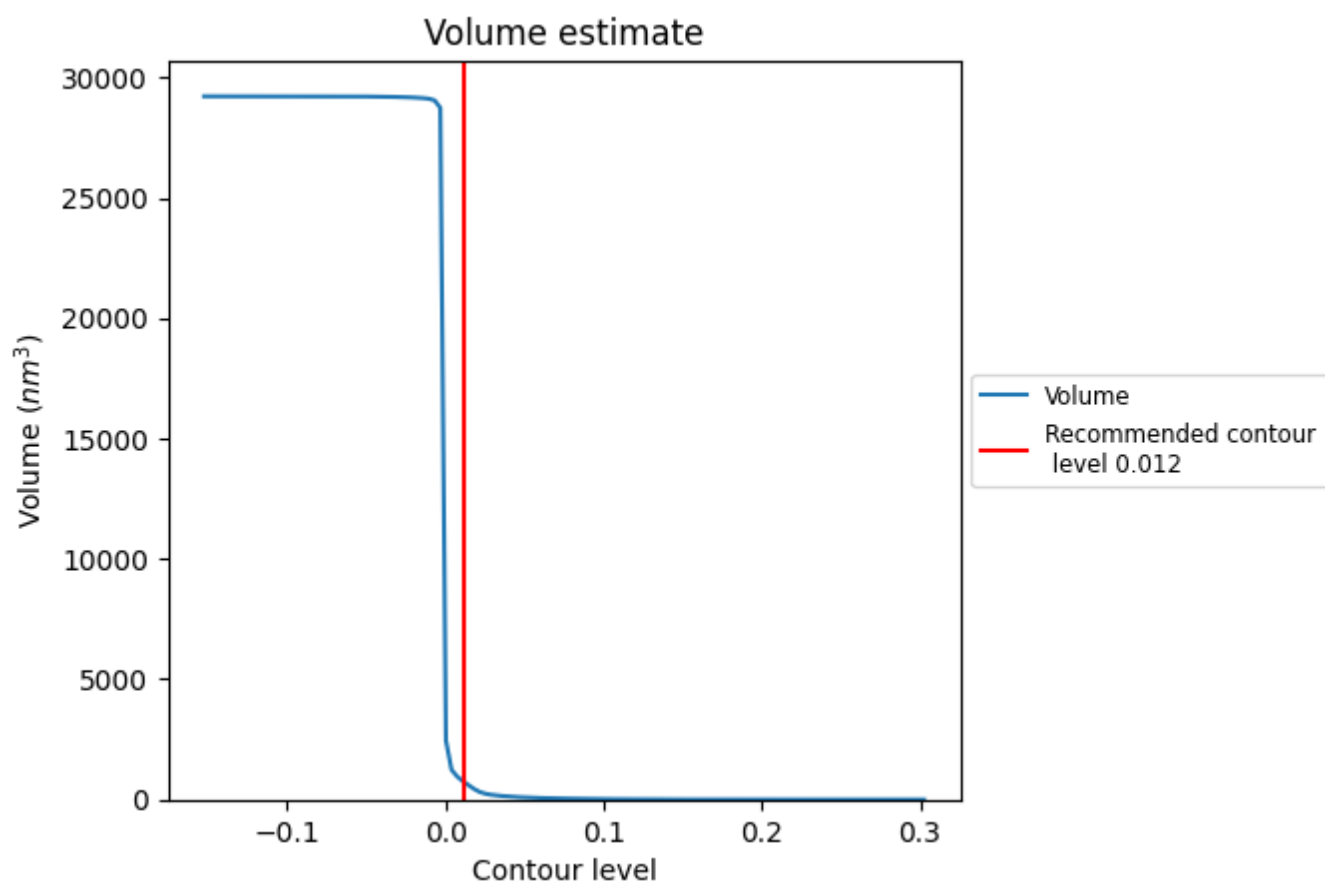
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

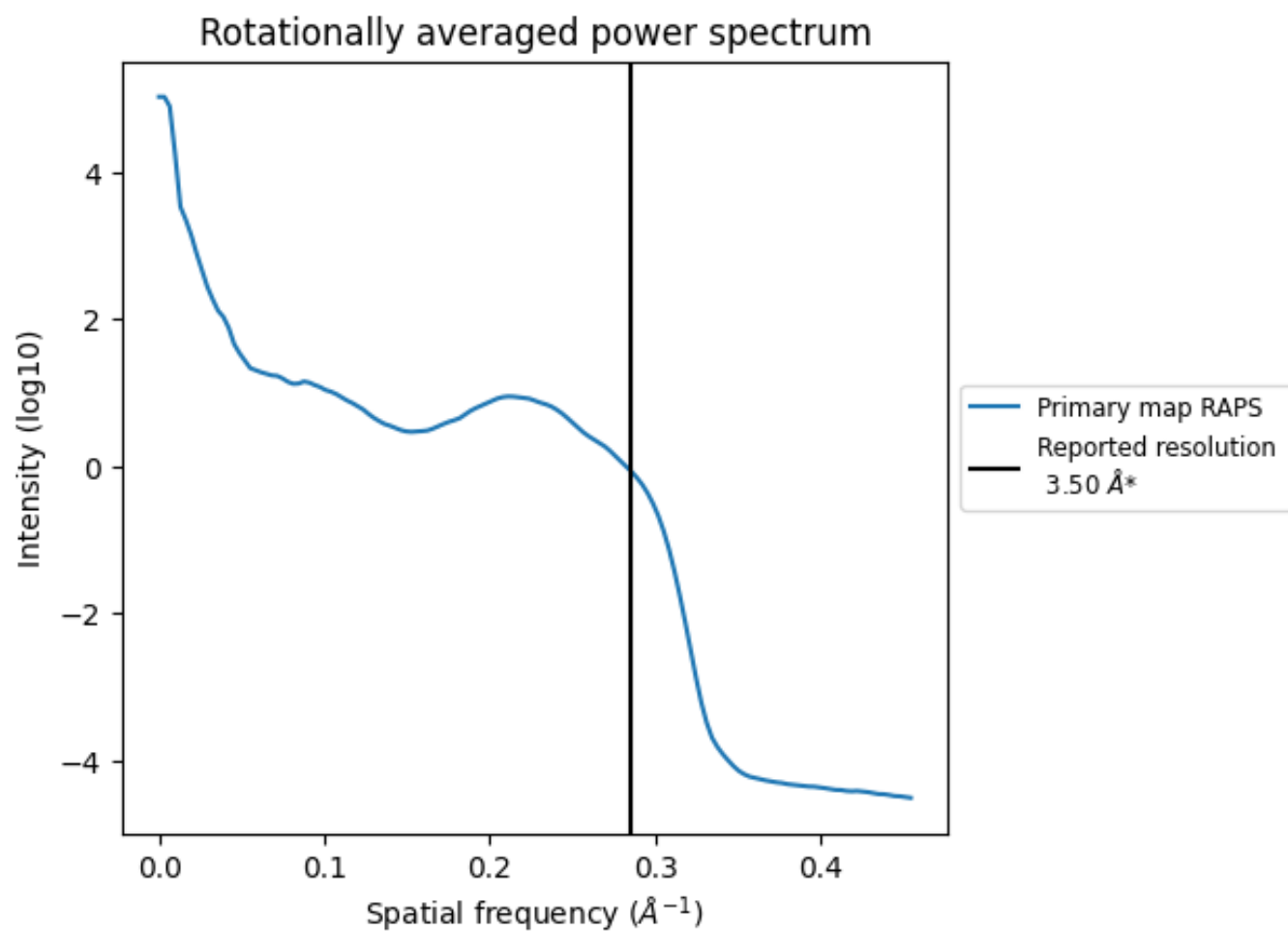
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 723  $\text{nm}^3$ ; this corresponds to an approximate mass of 653 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 0.286 Å<sup>-1</sup>

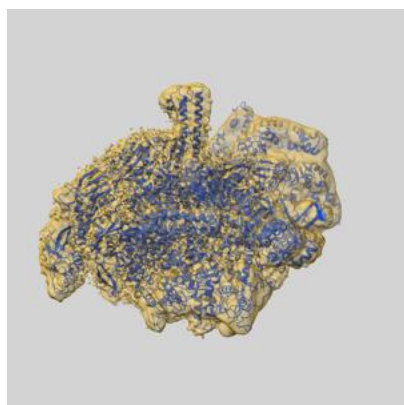
## 8 Fourier-Shell correlation

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

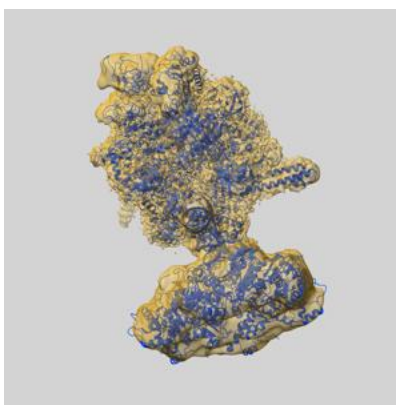
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-14171 and PDB model 7QV9. Per-residue inclusion information can be found in [section 3](#) on [page 7](#).

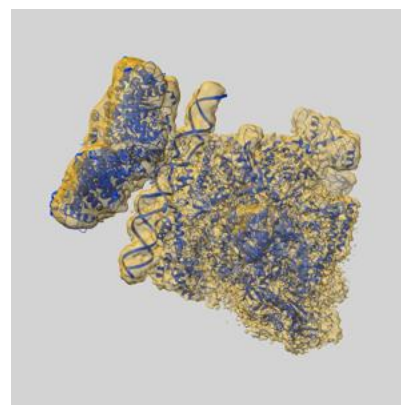
### 9.1 Map-model overlay [i](#)



X



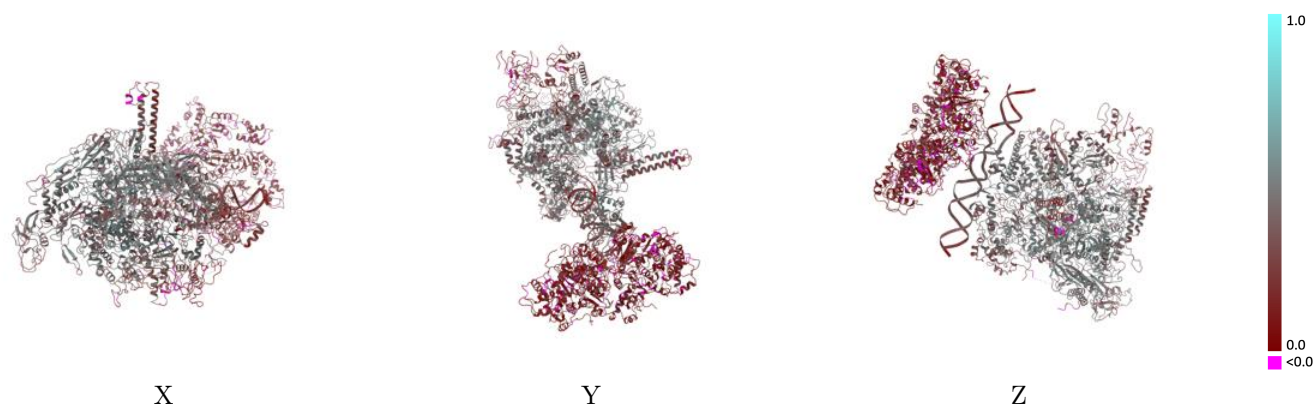
Y



Z

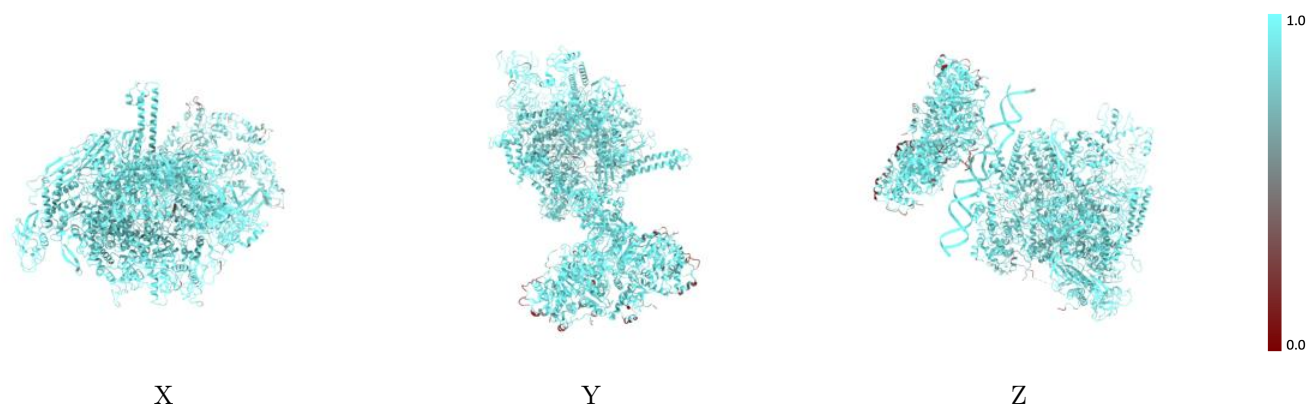
The images above show the 3D surface view of the map at the recommended contour level 0.012 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



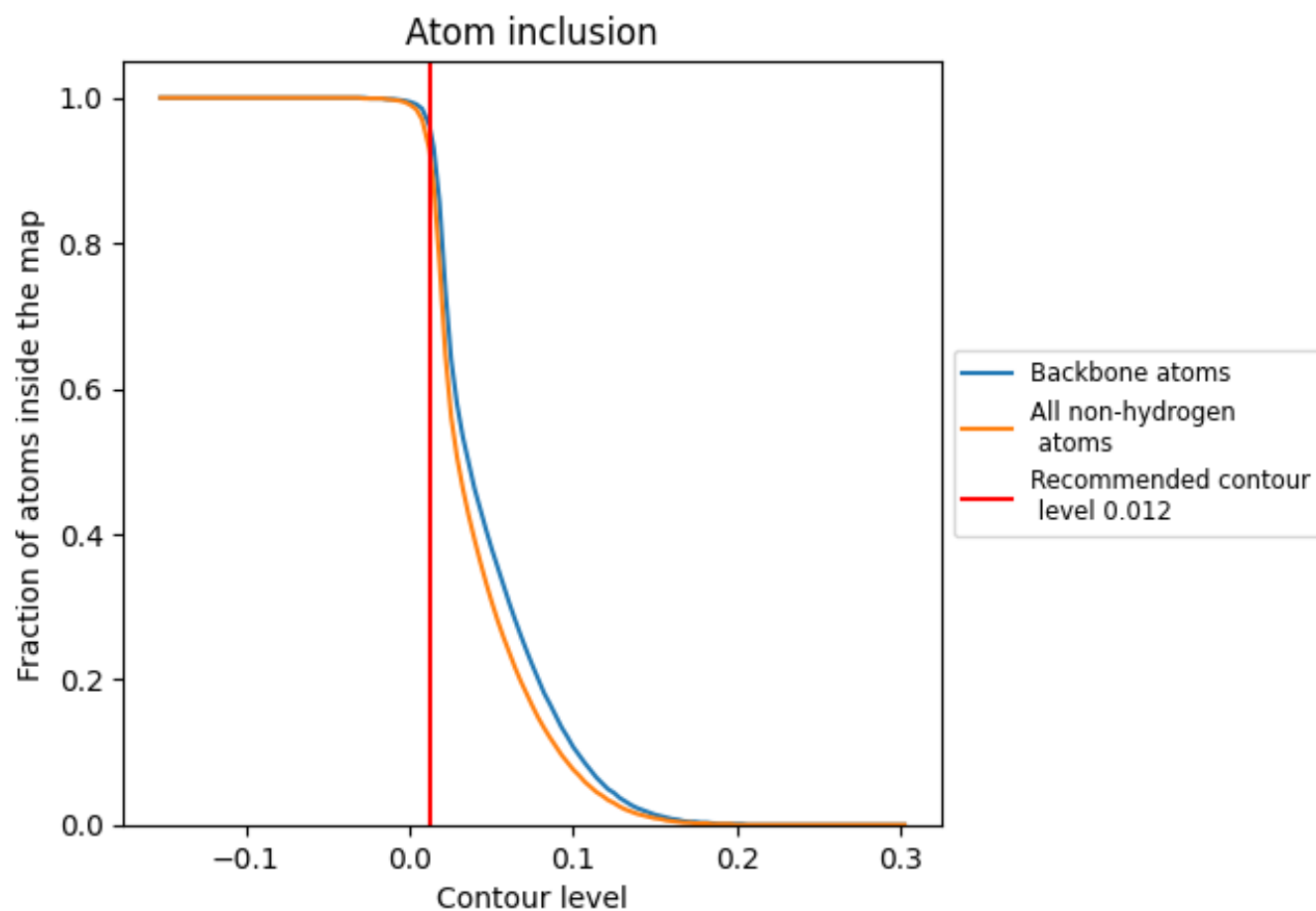
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

## 9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.012).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 93% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.012) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.9290</div>	<div><div></div>0.3230</div>
A	<div><div></div>0.8950</div>	<div><div></div>0.3980</div>
B	<div><div></div>0.9260</div>	<div><div></div>0.3830</div>
C	<div><div></div>0.9410</div>	<div><div></div>0.4200</div>
D	<div><div></div>0.9290</div>	<div><div></div>0.3910</div>
E	<div><div></div>0.9490</div>	<div><div></div>0.4320</div>
M	<div><div></div>0.9590</div>	<div><div></div>0.4280</div>
N	<div><div></div>0.9860</div>	<div><div></div>0.3110</div>
T	<div><div></div>0.9770</div>	<div><div></div>0.3090</div>
a	<div><div></div>0.9170</div>	<div><div></div>0.1500</div>
b	<div><div></div>0.9390</div>	<div><div></div>0.1620</div>
c	<div><div></div>0.8730</div>	<div><div></div>0.1240</div>
d	<div><div></div>0.9080</div>	<div><div></div>0.1270</div>
e	<div><div></div>0.9290</div>	<div><div></div>0.1350</div>
f	<div><div></div>0.9020</div>	<div><div></div>0.1310</div>

1.0

0.0

<0.0