



wwPDB EM Validation Summary Report ⓘ

Dec 21, 2024 – 04:03 pm GMT

PDB ID : 8RIF
EMDB ID : EMD-19186
Title : Cryo-EM structure of the MCM double hexamer loaded onto dsDNA.
Authors : Miller, T.C.R.; Lim, C.T.; Difley, J.F.X.; Costa, A.
Deposited on : 2023-12-18
Resolution : 2.79 Å (reported)
Based on initial model : 7P30

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.

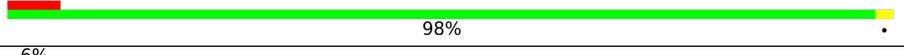
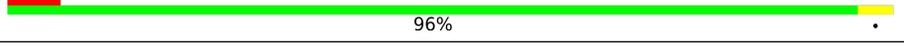
The types of validation reports are described at

<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.dev113
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.40

Continued from previous page...

Mol	Chain	Length	Quality of chain
5	6	1017	 59% 40%
5	E	1017	 58% 40%
6	7	845	 11% 85% 5% 11%
6	F	845	 11% 85% 5% 11%
7	X	53	 6% 98%
8	Y	53	 6% 96%

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 64125 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 replication licensing factor MCM2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	2	624	Total 4938	C 3114	N 878	O 927	S 19	0	0
1	A	624	Total 4938	C 3114	N 878	O 927	S 19	0	0

- Molecule 2 is a protein called DNA replication licensing factor MCM3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	3	614	Total 4811	C 3037	N 858	O 903	S 13	0	0
2	B	614	Total 4811	C 3037	N 858	O 903	S 13	0	0

There are 70 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	-34	MET	-	initiating methionine	UNP P24279
3	-33	LYS	-	expression tag	UNP P24279
3	-32	ARG	-	expression tag	UNP P24279
3	-31	ARG	-	expression tag	UNP P24279
3	-30	TRP	-	expression tag	UNP P24279
3	-29	LYS	-	expression tag	UNP P24279
3	-28	LYS	-	expression tag	UNP P24279
3	-27	ASN	-	expression tag	UNP P24279
3	-26	PHE	-	expression tag	UNP P24279
3	-25	ILE	-	expression tag	UNP P24279
3	-24	ALA	-	expression tag	UNP P24279
3	-23	VAL	-	expression tag	UNP P24279
3	-22	SER	-	expression tag	UNP P24279
3	-21	ALA	-	expression tag	UNP P24279
3	-20	ALA	-	expression tag	UNP P24279
3	-19	ASN	-	expression tag	UNP P24279
3	-18	ARG	-	expression tag	UNP P24279

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
3	-17	PHE	-	expression tag	UNP P24279
3	-16	LYS	-	expression tag	UNP P24279
3	-15	LYS	-	expression tag	UNP P24279
3	-14	ILE	-	expression tag	UNP P24279
3	-13	SER	-	expression tag	UNP P24279
3	-12	SER	-	expression tag	UNP P24279
3	-11	SER	-	expression tag	UNP P24279
3	-10	GLY	-	expression tag	UNP P24279
3	-9	ALA	-	expression tag	UNP P24279
3	-8	LEU	-	expression tag	UNP P24279
3	-7	GLU	-	expression tag	UNP P24279
3	-6	ASN	-	expression tag	UNP P24279
3	-5	LEU	-	expression tag	UNP P24279
3	-4	TYR	-	expression tag	UNP P24279
3	-3	PHE	-	expression tag	UNP P24279
3	-2	GLN	-	expression tag	UNP P24279
3	-1	GLY	-	expression tag	UNP P24279
3	0	GLU	-	expression tag	UNP P24279
B	-34	MET	-	initiating methionine	UNP P24279
B	-33	LYS	-	expression tag	UNP P24279
B	-32	ARG	-	expression tag	UNP P24279
B	-31	ARG	-	expression tag	UNP P24279
B	-30	TRP	-	expression tag	UNP P24279
B	-29	LYS	-	expression tag	UNP P24279
B	-28	LYS	-	expression tag	UNP P24279
B	-27	ASN	-	expression tag	UNP P24279
B	-26	PHE	-	expression tag	UNP P24279
B	-25	ILE	-	expression tag	UNP P24279
B	-24	ALA	-	expression tag	UNP P24279
B	-23	VAL	-	expression tag	UNP P24279
B	-22	SER	-	expression tag	UNP P24279
B	-21	ALA	-	expression tag	UNP P24279
B	-20	ALA	-	expression tag	UNP P24279
B	-19	ASN	-	expression tag	UNP P24279
B	-18	ARG	-	expression tag	UNP P24279
B	-17	PHE	-	expression tag	UNP P24279
B	-16	LYS	-	expression tag	UNP P24279
B	-15	LYS	-	expression tag	UNP P24279
B	-14	ILE	-	expression tag	UNP P24279
B	-13	SER	-	expression tag	UNP P24279
B	-12	SER	-	expression tag	UNP P24279
B	-11	SER	-	expression tag	UNP P24279

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	GLY	-	expression tag	UNP P24279
B	-9	ALA	-	expression tag	UNP P24279
B	-8	LEU	-	expression tag	UNP P24279
B	-7	GLU	-	expression tag	UNP P24279
B	-6	ASN	-	expression tag	UNP P24279
B	-5	LEU	-	expression tag	UNP P24279
B	-4	TYR	-	expression tag	UNP P24279
B	-3	PHE	-	expression tag	UNP P24279
B	-2	GLN	-	expression tag	UNP P24279
B	-1	GLY	-	expression tag	UNP P24279
B	0	GLU	-	expression tag	UNP P24279

- Molecule 3 is a protein called DNA replication licensing factor MCM4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	4	645	Total	C	N	O	S	0	0
			5120	3211	886	993	30		
3	C	645	Total	C	N	O	S	0	0
			5120	3211	886	993	30		

- Molecule 4 is a protein called Minichromosome maintenance protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	5	656	Total	C	N	O	S	0	0
			5136	3219	878	1014	25		
4	D	656	Total	C	N	O	S	0	0
			5136	3219	878	1014	25		

- Molecule 5 is a protein called DNA replication licensing factor MCM6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	6	615	Total	C	N	O	S	0	0
			4874	3077	851	921	25		
5	E	615	Total	C	N	O	S	0	0
			4874	3077	851	921	25		

- Molecule 6 is a protein called DNA replication licensing factor MCM7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	7	755	Total	C	N	O	S	0	0
			5976	3771	1031	1142	32		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	F	755	5976	3771	1031	1142	32	0	0

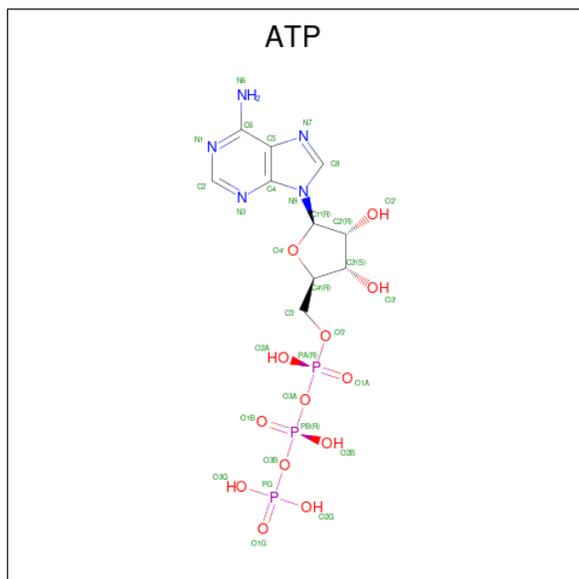
- Molecule 7 is a DNA chain called DNA (53-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
7	X	53	1086	515	199	319	53	0	0

- Molecule 8 is a DNA chain called DNA (53-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
8	Y	53	1087	515	202	317	53	0	0

- Molecule 9 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
9	2	1	31	10	5	13	3	0
9	A	1	31	10	5	13	3	0

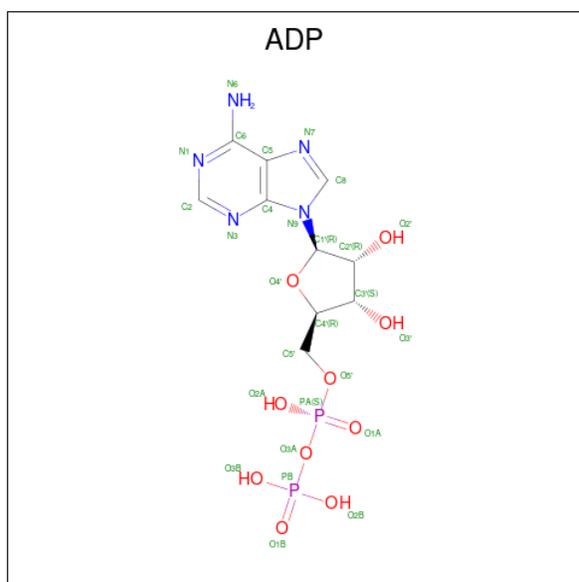
- Molecule 10 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	AltConf
10	2	1	Total Mg 1 1	0
10	3	1	Total Mg 1 1	0
10	5	1	Total Mg 1 1	0
10	7	1	Total Mg 1 1	0
10	A	1	Total Mg 1 1	0
10	B	1	Total Mg 1 1	0
10	D	1	Total Mg 1 1	0
10	F	1	Total Mg 1 1	0

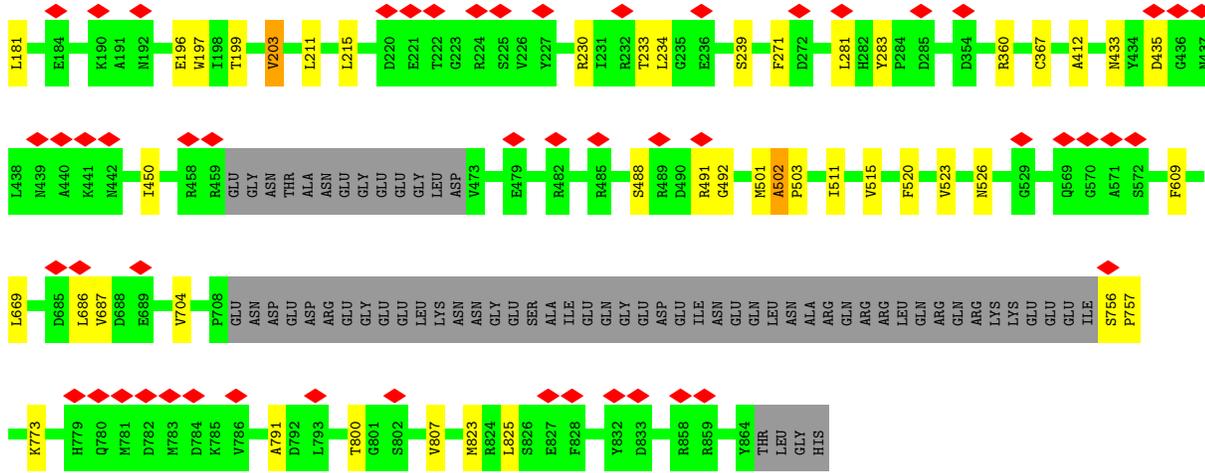
- Molecule 11 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	AltConf
11	2	1	Total Zn 1 1	0
11	4	1	Total Zn 1 1	0
11	5	1	Total Zn 1 1	0
11	6	1	Total Zn 1 1	0
11	7	1	Total Zn 1 1	0
11	A	1	Total Zn 1 1	0
11	C	1	Total Zn 1 1	0
11	D	1	Total Zn 1 1	0
11	E	1	Total Zn 1 1	0
11	F	1	Total Zn 1 1	0

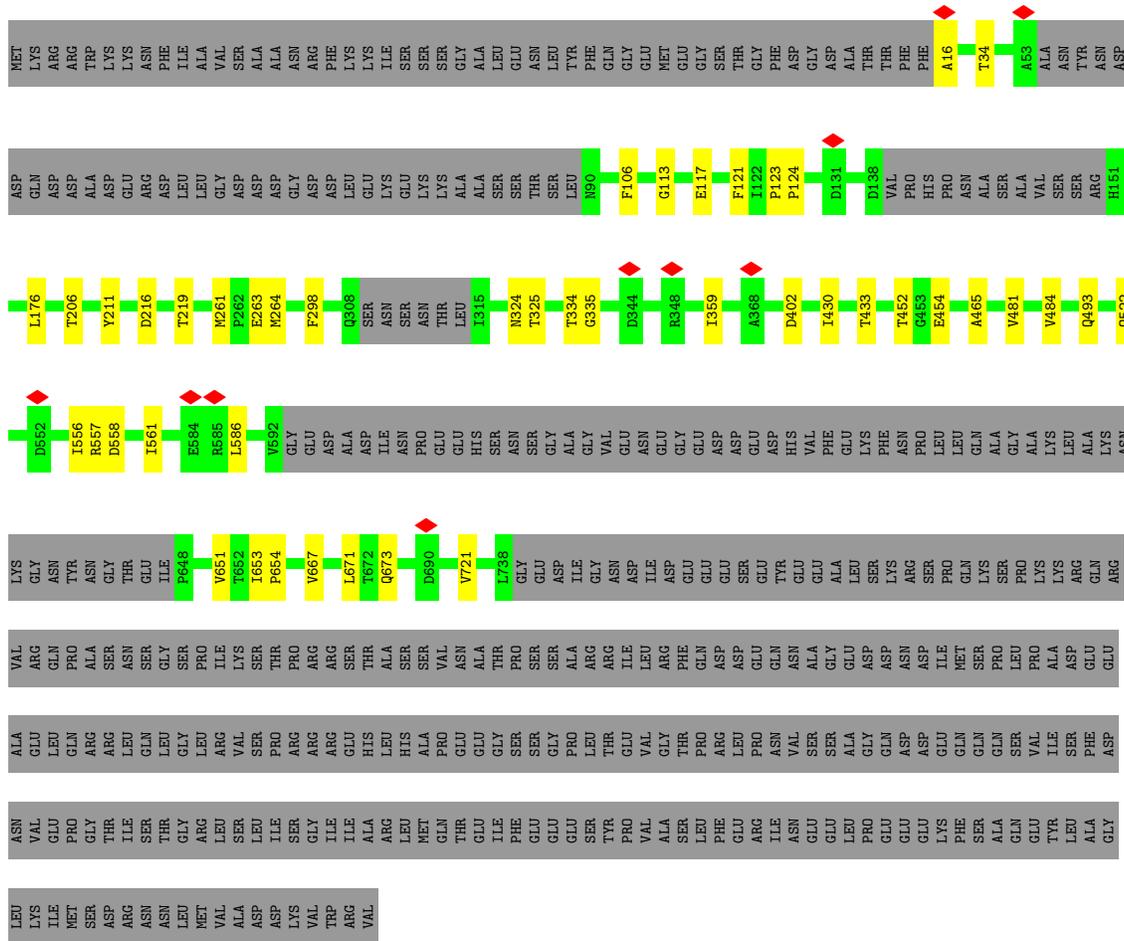
- Molecule 12 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
12	3	1	Total	C	N	O	P	0
			27	10	5	10	2	
12	5	1	Total	C	N	O	P	0
			27	10	5	10	2	
12	7	1	Total	C	N	O	P	0
			27	10	5	10	2	
12	B	1	Total	C	N	O	P	0
			27	10	5	10	2	
12	D	1	Total	C	N	O	P	0
			27	10	5	10	2	
12	F	1	Total	C	N	O	P	0
			27	10	5	10	2	

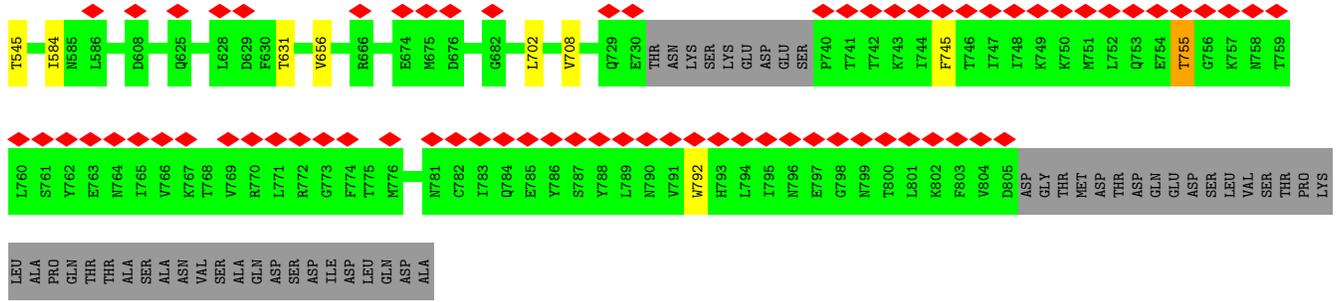


• Molecule 2: DNA replication licensing factor MCM3

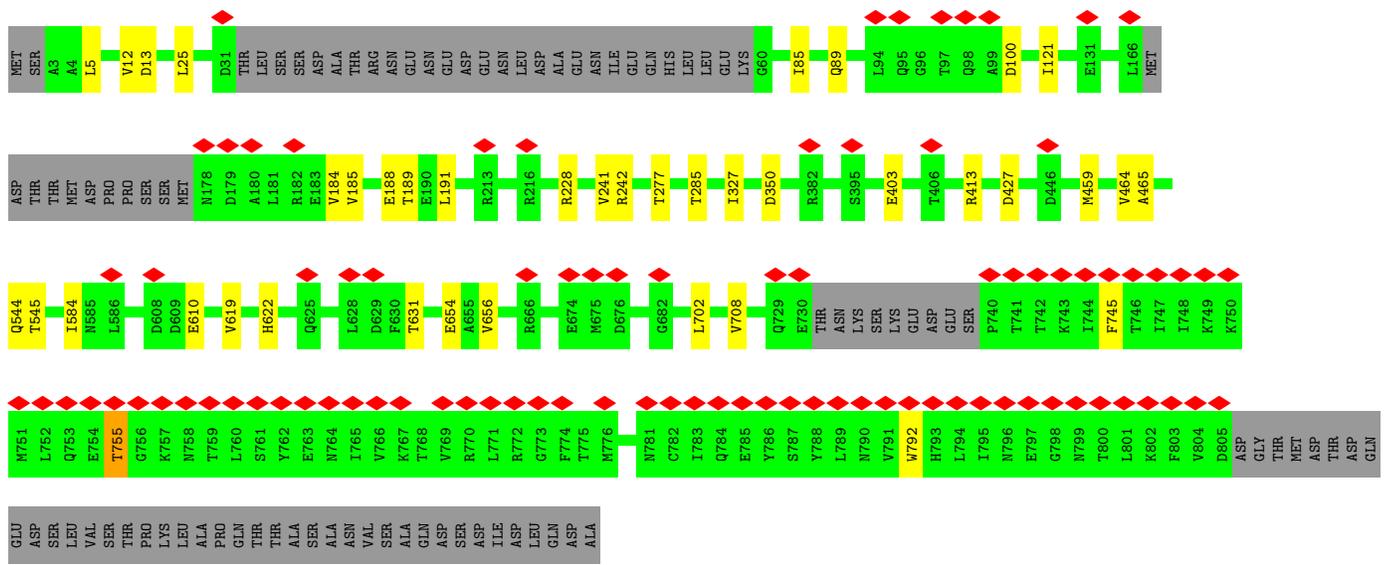
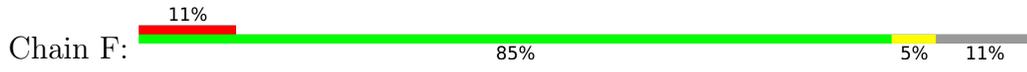


• Molecule 2: DNA replication licensing factor MCM3





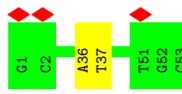
• Molecule 6: DNA replication licensing factor MCM7



• Molecule 7: DNA (53-MER)



• Molecule 8: DNA (53-MER)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	135143	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 ($e^-/\text{\AA}^2$)	51.2	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	4500	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.080	Depositor
Minimum map value	-0.039	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.02	Depositor
Map size (Å)	453.6, 453.6, 453.6	wwPDB
Map dimensions	420, 420, 420	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ATP, MG, ZN, 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	2	0.38	0/5025	0.64	0/6790
1	A	0.38	0/5025	0.64	0/6790
2	3	0.31	0/4894	0.57	0/6635
2	B	0.31	0/4894	0.57	0/6635
3	4	0.36	0/5192	0.60	0/7016
3	C	0.36	0/5192	0.60	0/7016
4	5	0.38	0/5210	0.60	0/7046
4	D	0.38	0/5210	0.60	0/7046
5	6	0.33	0/4952	0.57	0/6679
5	E	0.33	0/4952	0.57	0/6679
6	7	0.36	0/6068	0.61	0/8199
6	F	0.36	0/6068	0.61	0/8199
7	X	0.49	0/1217	0.86	0/1876
8	Y	0.51	0/1219	0.85	0/1879
All	All	0.36	0/65118	0.61	0/88485

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	2	4938	0	4997	25	0
1	A	4938	0	4997	24	0
2	3	4811	0	4893	25	0
2	B	4811	0	4893	25	0
3	4	5120	0	5187	13	0
3	C	5120	0	5187	14	0
4	5	5136	0	5177	25	0
4	D	5136	0	5177	27	0
5	6	4874	0	4918	13	0
5	E	4874	0	4918	14	0
6	7	5976	0	6057	19	0
6	F	5976	0	6057	20	0
7	X	1086	0	596	1	0
8	Y	1087	0	595	1	0
9	2	31	0	12	0	0
9	A	31	0	12	0	0
10	2	1	0	0	0	0
10	3	1	0	0	0	0
10	5	1	0	0	0	0
10	7	1	0	0	0	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
10	D	1	0	0	0	0
10	F	1	0	0	0	0
11	2	1	0	0	0	0
11	4	1	0	0	0	0
11	5	1	0	0	0	0
11	6	1	0	0	0	0
11	7	1	0	0	0	0
11	A	1	0	0	0	0
11	C	1	0	0	0	0
11	D	1	0	0	0	0
11	E	1	0	0	0	0
11	F	1	0	0	0	0
12	3	27	0	12	0	0
12	5	27	0	12	0	0
12	7	27	0	12	0	0
12	B	27	0	12	0	0
12	D	27	0	12	0	0
12	F	27	0	12	0	0
All	All	64125	0	63745	231	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

The worst 5 of 231 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)
1:2:501:MET:O	1:2:502:ALA:C	2.41	0.58
1:A:501:MET:O	1:A:502:ALA:C	2.41	0.58
4:5:410:ILE:HG23	4:5:552:MET:HE1	1.91	0.52
4:D:410:ILE:HG23	4:D:552:MET:HE1	1.93	0.51
3:C:729:LEU:O	3:C:730:GLU:C	2.50	0.50

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	2	618/868 (71%)	583 (94%)	32 (5%)	3 (0%)	25	56
1	A	618/868 (71%)	583 (94%)	32 (5%)	3 (0%)	25	56
2	3	604/1006 (60%)	582 (96%)	22 (4%)	0	100	100
2	B	604/1006 (60%)	582 (96%)	22 (4%)	0	100	100
3	4	637/933 (68%)	604 (95%)	33 (5%)	0	100	100
3	C	637/933 (68%)	604 (95%)	33 (5%)	0	100	100
4	5	648/775 (84%)	603 (93%)	43 (7%)	2 (0%)	37	67
4	D	648/775 (84%)	603 (93%)	43 (7%)	2 (0%)	37	67
5	6	603/1017 (59%)	570 (94%)	33 (6%)	0	100	100
5	E	603/1017 (59%)	570 (94%)	33 (6%)	0	100	100
6	7	747/845 (88%)	718 (96%)	28 (4%)	1 (0%)	48	77
6	F	747/845 (88%)	718 (96%)	28 (4%)	1 (0%)	48	77
All	All	7714/10888 (71%)	7320 (95%)	382 (5%)	12 (0%)	45	73

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	2	435	ASP
6	7	755	THR
1	A	435	ASP
6	F	755	THR
1	2	412	ALA

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	2	548/770 (71%)	540 (98%)	8 (2%)	60 86
1	A	548/770 (71%)	540 (98%)	8 (2%)	60 86
2	3	530/864 (61%)	526 (99%)	4 (1%)	79 93
2	B	530/864 (61%)	526 (99%)	4 (1%)	79 93
3	4	584/848 (69%)	578 (99%)	6 (1%)	73 91
3	C	584/848 (69%)	577 (99%)	7 (1%)	67 89
4	5	589/688 (86%)	579 (98%)	10 (2%)	56 84
4	D	589/688 (86%)	579 (98%)	10 (2%)	56 84
5	6	538/886 (61%)	538 (100%)	0	100 100
5	E	538/886 (61%)	538 (100%)	0	100 100
6	7	671/753 (89%)	664 (99%)	7 (1%)	73 91
6	F	671/753 (89%)	664 (99%)	7 (1%)	73 91
All	All	6920/9618 (72%)	6849 (99%)	71 (1%)	71 91

5 of 71 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	D	365	LYS
4	D	586	GLN
6	F	191	LEU
4	5	613	ARG
4	5	586	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
5	E	274	HIS
6	F	538	HIS
5	E	730	HIS
6	7	538	HIS
4	D	652	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 26 ligands modelled in this entry, 18 are monoatomic - leaving 8 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	$\# Z > 2$
12	ADP	F	1000	10	24,29,29	0.78	0	29,45,45	0.87	1 (3%)
12	ADP	5	1000	10	24,29,29	1.04	2 (8%)	29,45,45	0.84	1 (3%)
12	ADP	3	1000	10	24,29,29	0.77	0	29,45,45	0.91	1 (3%)
12	ADP	7	1000	10	24,29,29	0.78	0	29,45,45	0.87	1 (3%)
12	ADP	B	1000	10	24,29,29	0.77	0	29,45,45	0.91	1 (3%)
12	ADP	D	1000	10	24,29,29	1.04	2 (8%)	29,45,45	0.84	1 (3%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	ATP	A	1000	10	26,33,33	0.80	0	31,52,52	0.89	1 (3%)
9	ATP	2	1000	10	26,33,33	0.80	0	31,52,52	0.89	1 (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
12	ADP	F	1000	10	-	3/12/32/32	0/3/3/3
12	ADP	5	1000	10	-	3/12/32/32	0/3/3/3
12	ADP	3	1000	10	-	2/12/32/32	0/3/3/3
12	ADP	7	1000	10	-	3/12/32/32	0/3/3/3
12	ADP	B	1000	10	-	2/12/32/32	0/3/3/3
12	ADP	D	1000	10	-	3/12/32/32	0/3/3/3
9	ATP	A	1000	10	-	5/18/38/38	0/3/3/3
9	ATP	2	1000	10	-	5/18/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	D	1000	ADP	PB-O2B	-3.15	1.42	1.54
12	5	1000	ADP	PB-O2B	-3.15	1.42	1.54
12	5	1000	ADP	C8-N7	-2.08	1.31	1.34
12	D	1000	ADP	C8-N7	-2.08	1.31	1.34

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	2	1000	ATP	C5-C6-N6	2.85	124.68	120.35
9	A	1000	ATP	C5-C6-N6	2.85	124.68	120.35
12	3	1000	ADP	C5-C6-N6	2.69	124.44	120.35
12	B	1000	ADP	C5-C6-N6	2.69	124.44	120.35
12	7	1000	ADP	C5-C6-N6	2.63	124.35	120.35

There are no chirality outliers.

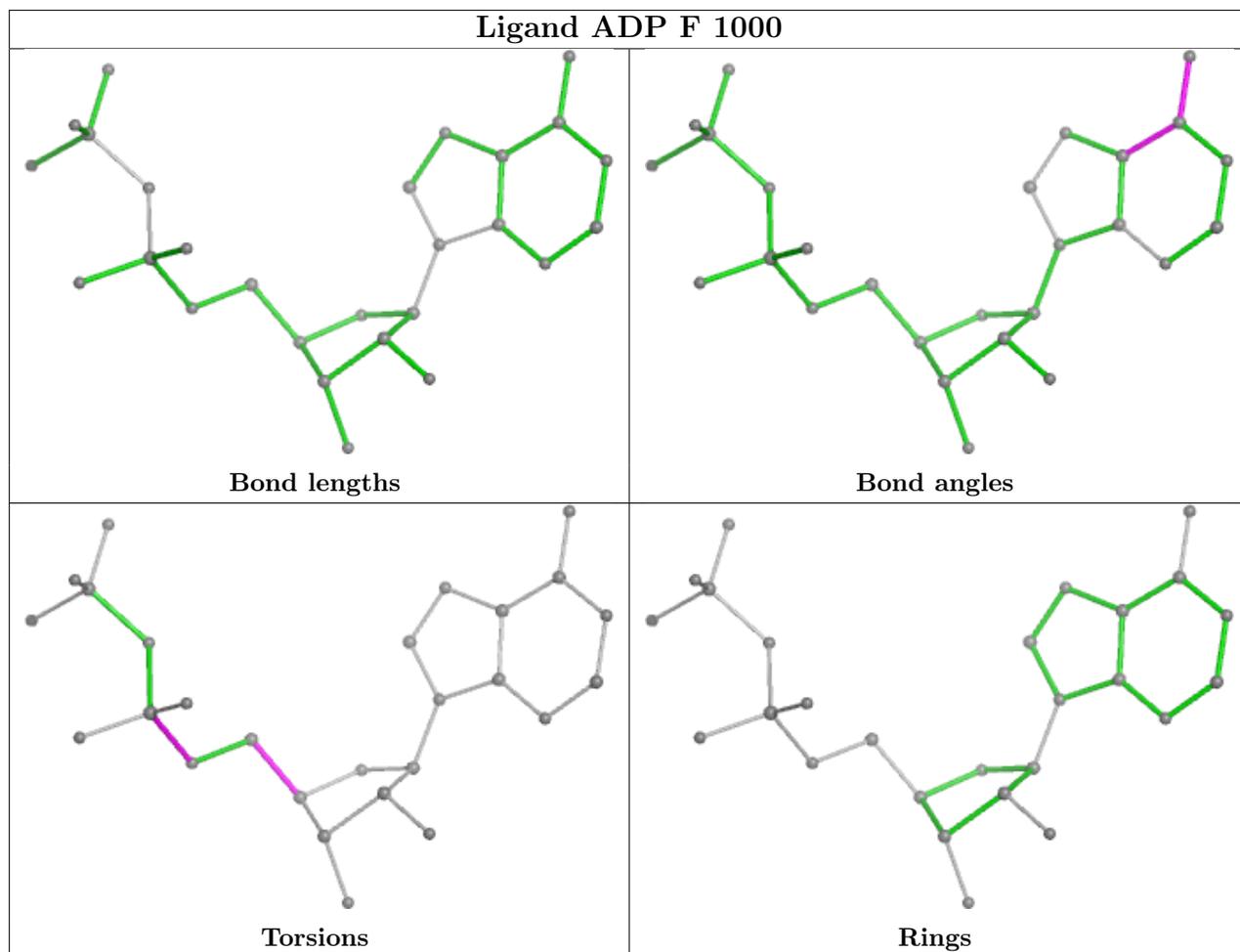
5 of 26 torsion outliers are listed below:

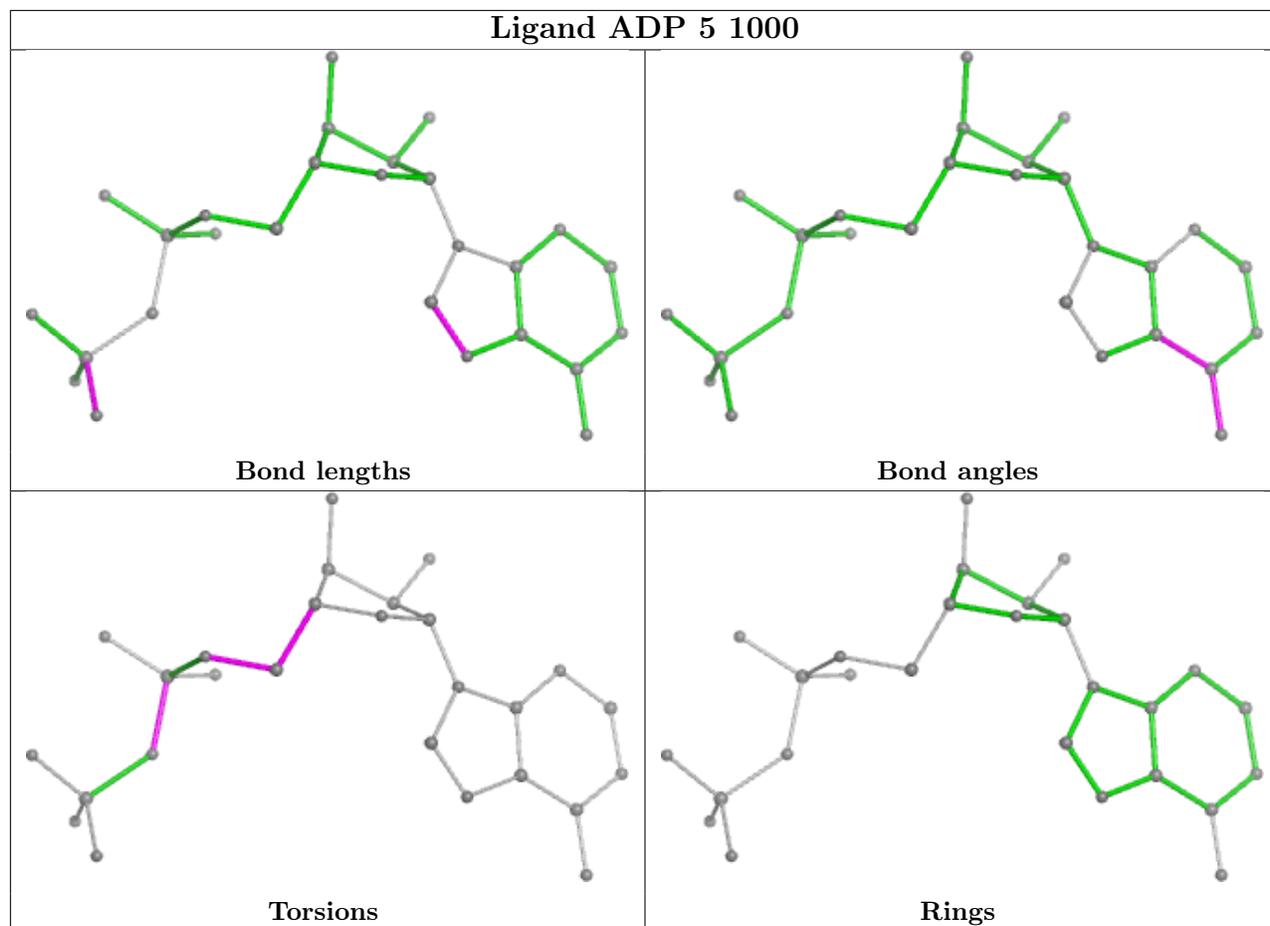
Mol	Chain	Res	Type	Atoms
9	2	1000	ATP	PB-O3B-PG-O3G
9	A	1000	ATP	PB-O3B-PG-O3G
12	7	1000	ADP	C5'-O5'-PA-O1A
12	F	1000	ADP	C5'-O5'-PA-O1A
12	7	1000	ADP	C5'-O5'-PA-O3A

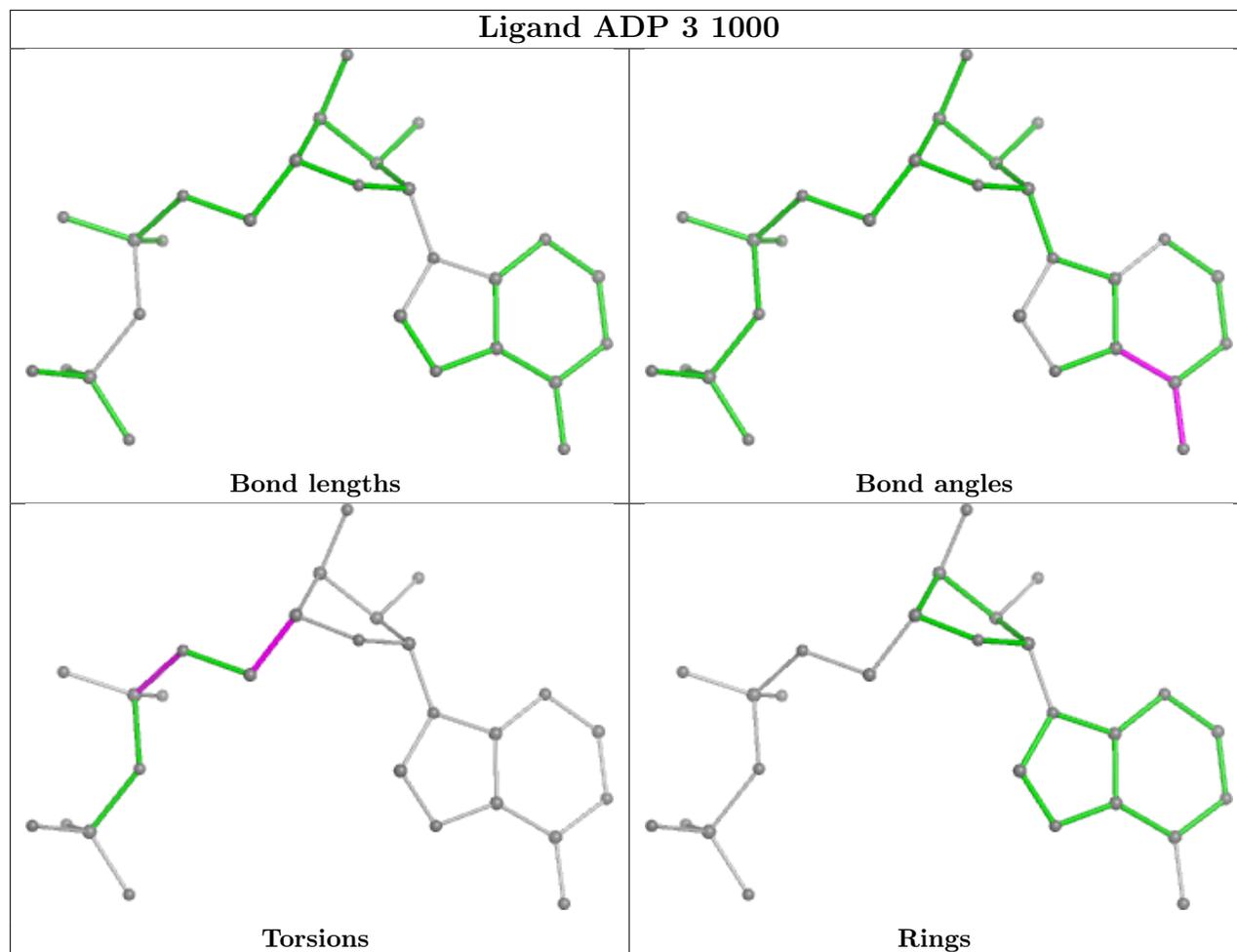
There are no ring outliers.

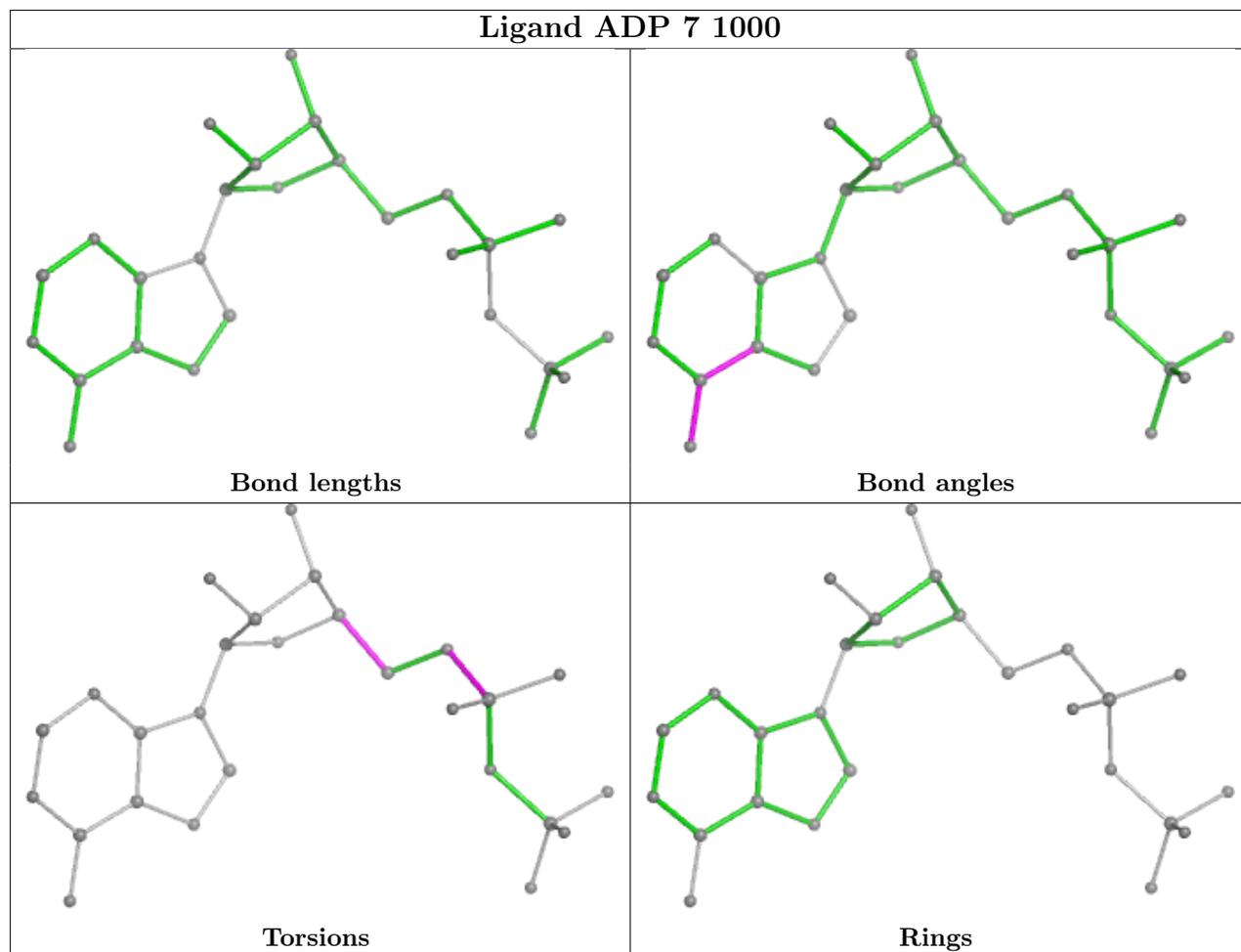
No monomer is involved in short contacts.

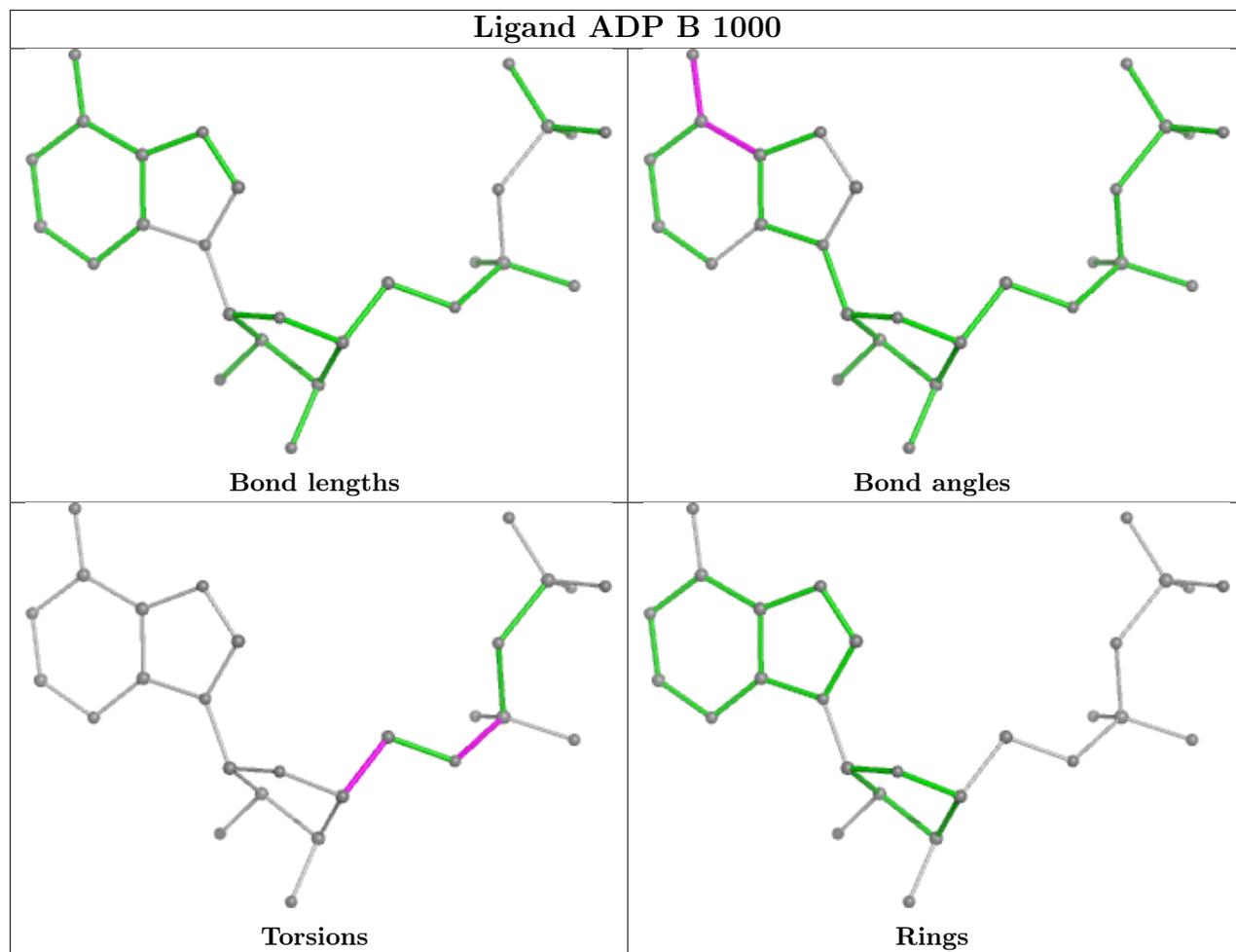
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

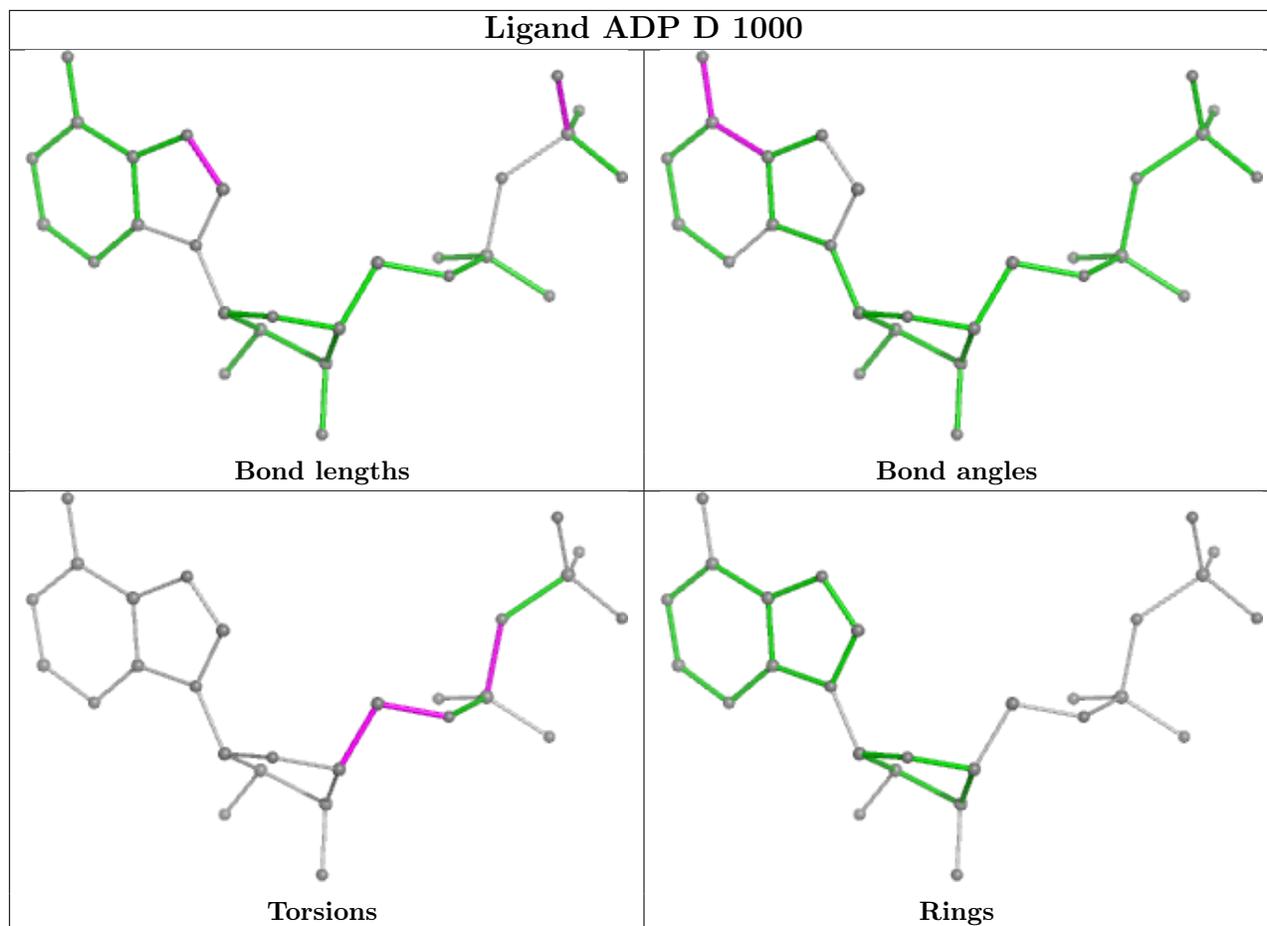


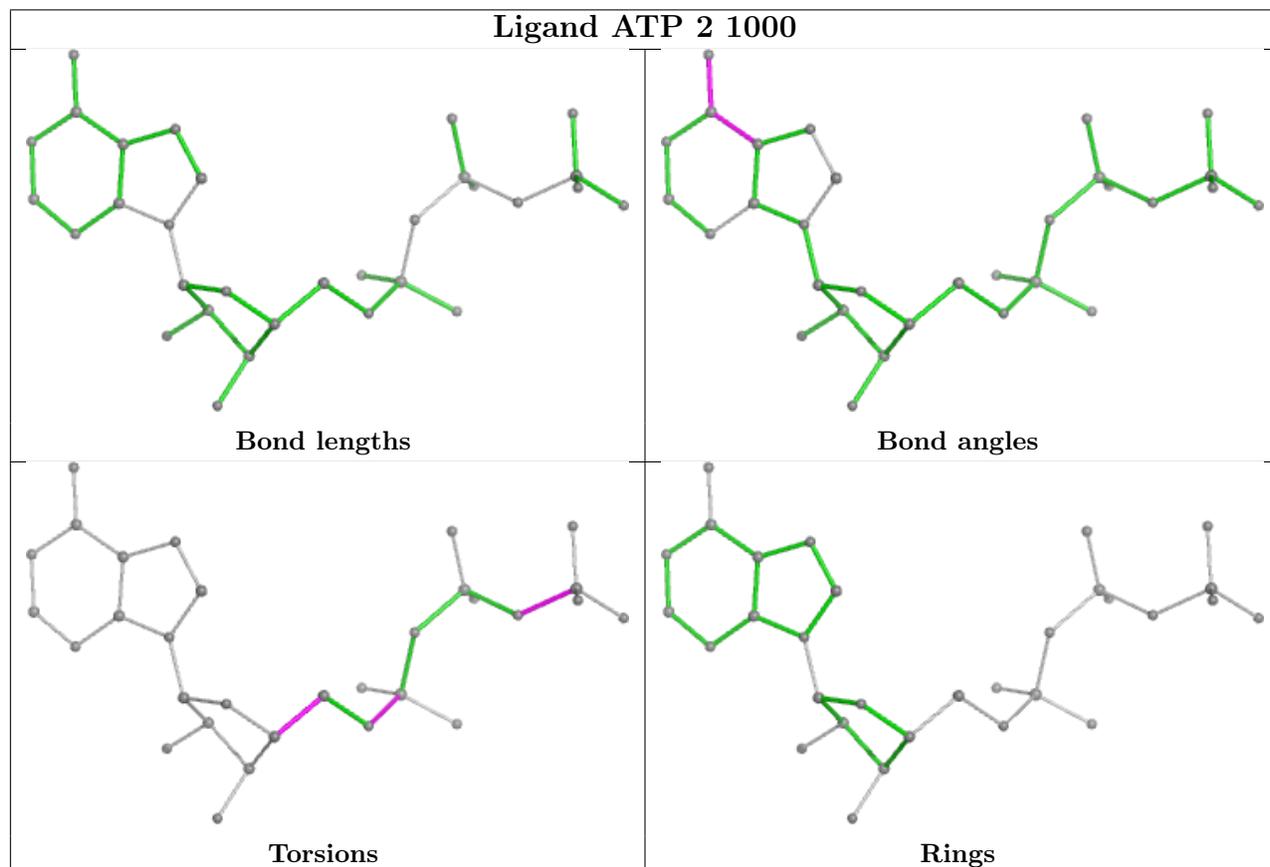
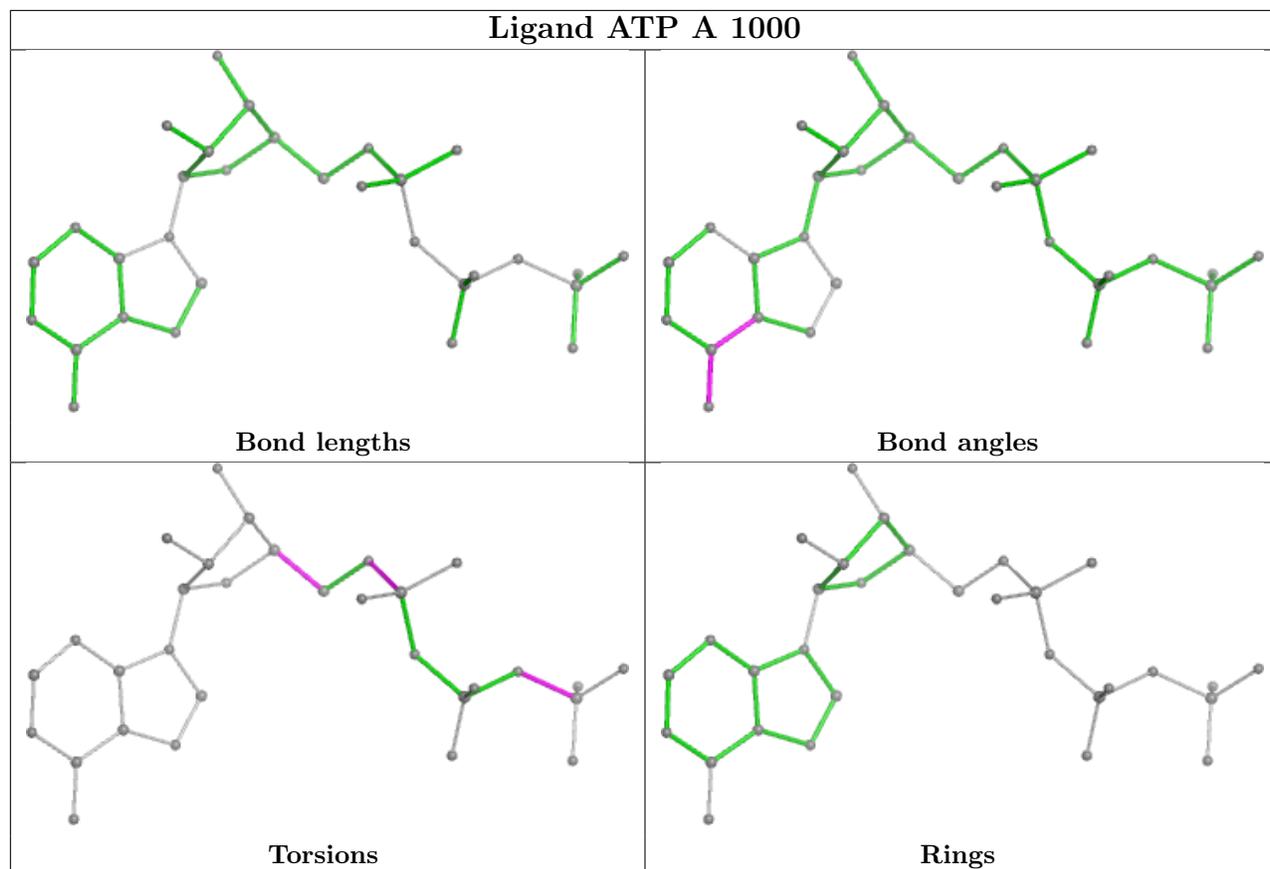












5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

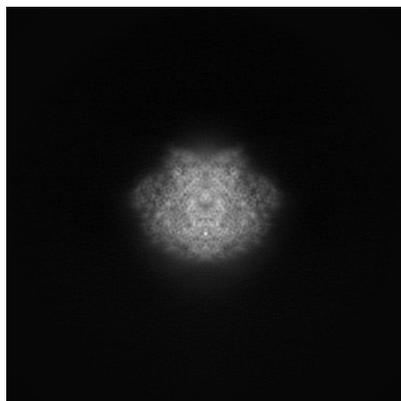
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-19186. These allow visual inspection of the internal detail of the map and identification of artifacts.

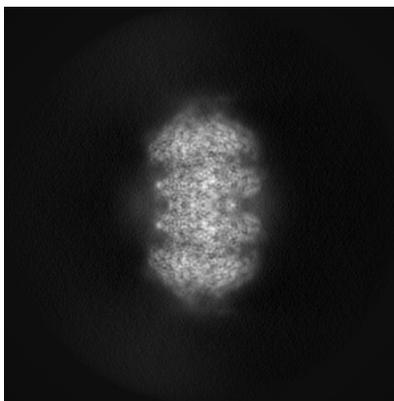
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

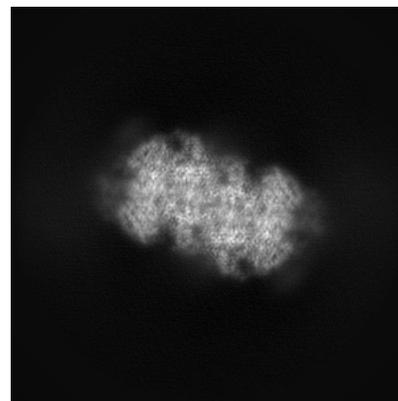
6.1.1 Primary map



X

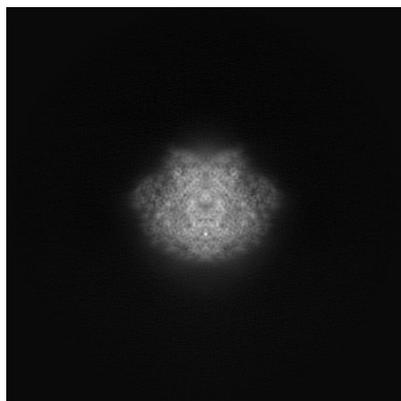


Y

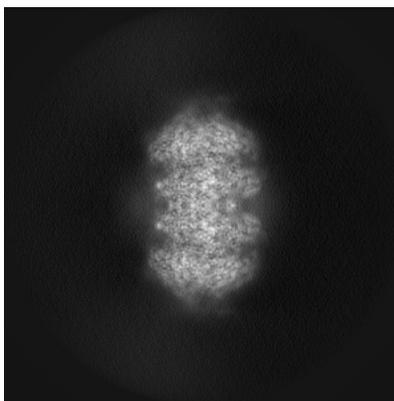


Z

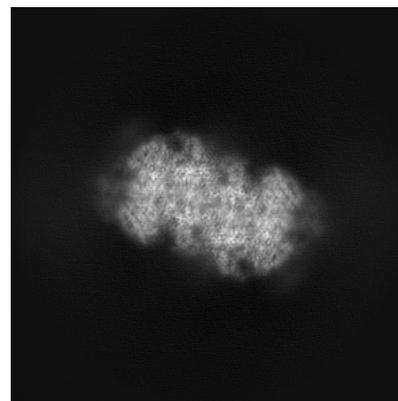
6.1.2 Raw map



X



Y

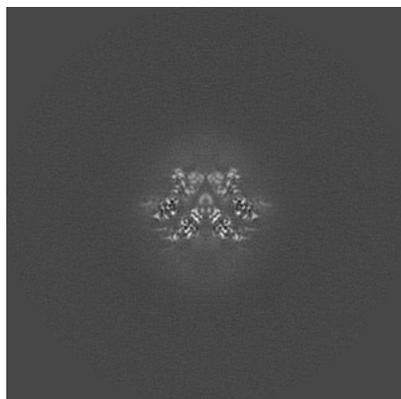


Z

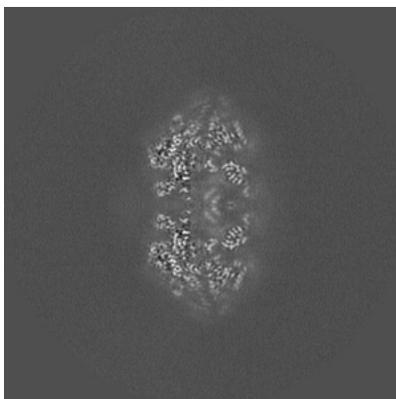
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

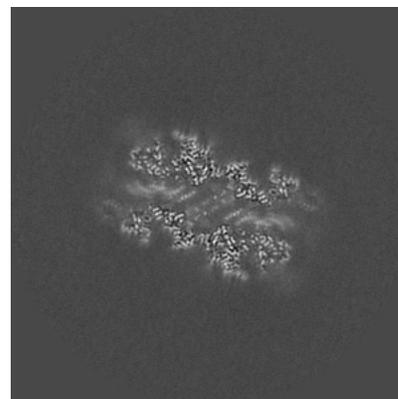
6.2.1 Primary map



X Index: 210

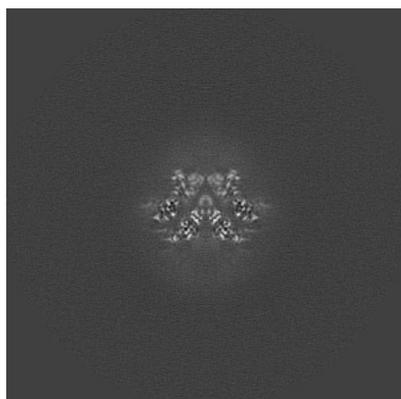


Y Index: 210

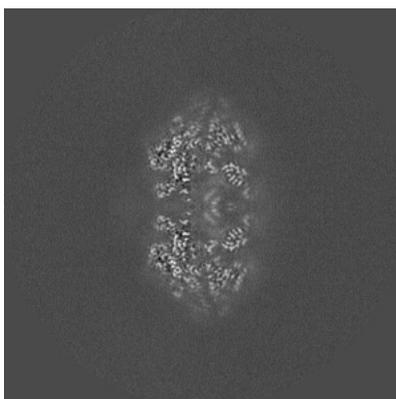


Z Index: 210

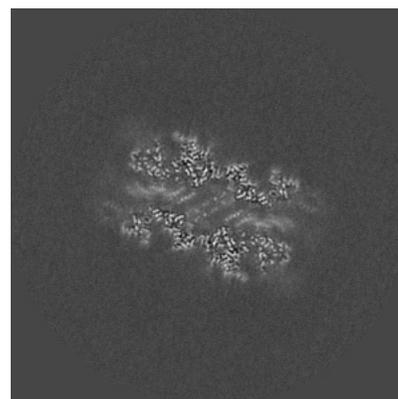
6.2.2 Raw map



X Index: 210



Y Index: 210

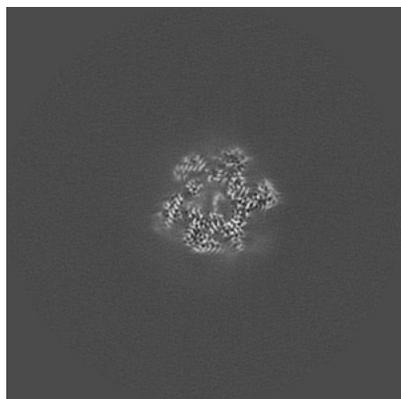


Z Index: 210

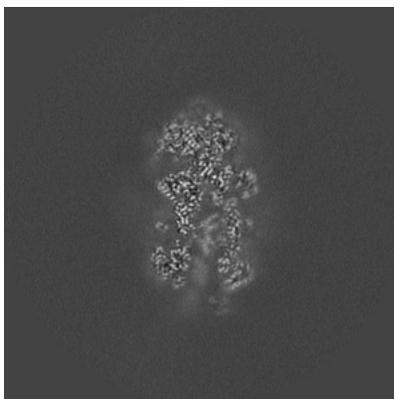
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

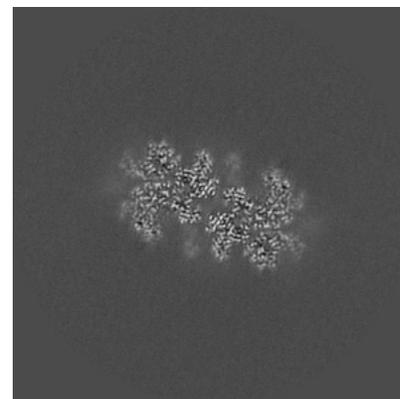
6.3.1 Primary map



X Index: 187

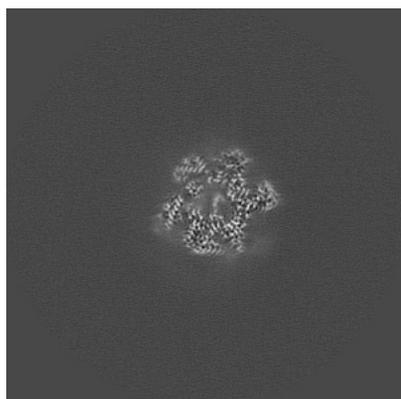


Y Index: 222

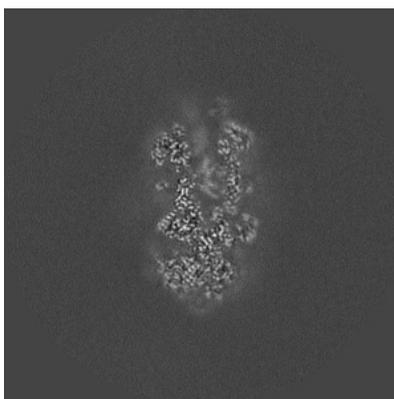


Z Index: 185

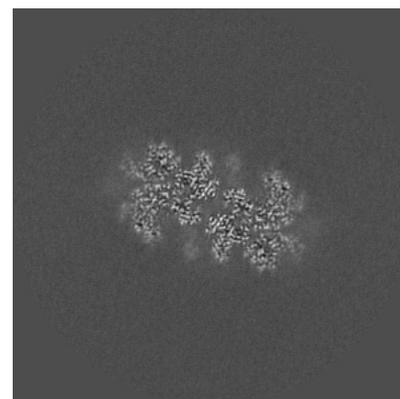
6.3.2 Raw map



X Index: 187



Y Index: 198

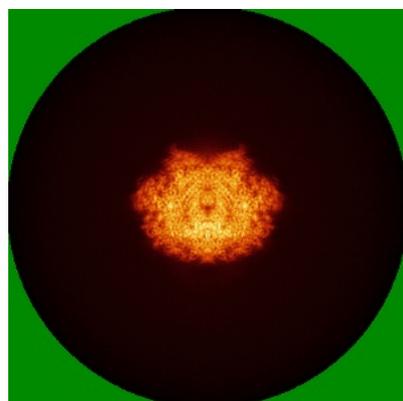


Z Index: 185

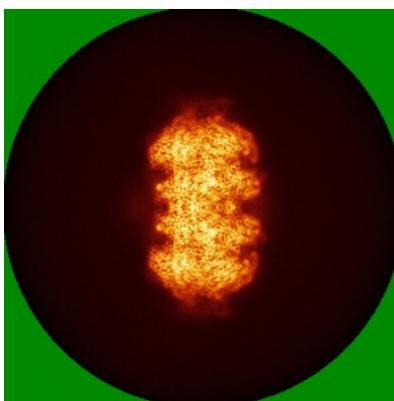
The images above show the largest variance slices of the map in three orthogonal directions.

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

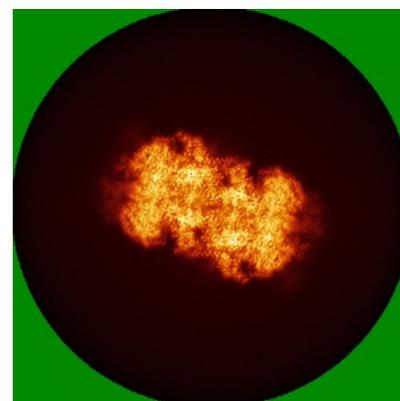
6.4.1 Primary map



X

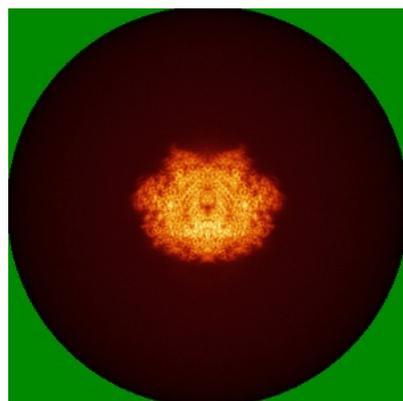


Y

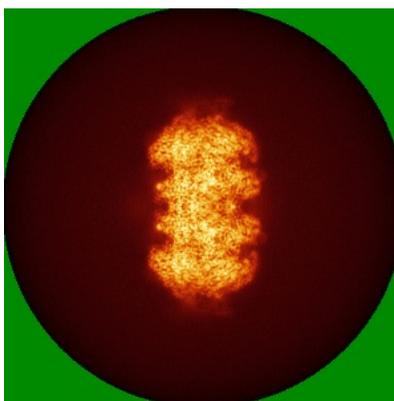


Z

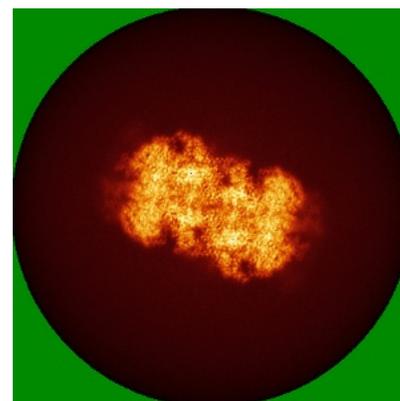
6.4.2 Raw map



X



Y

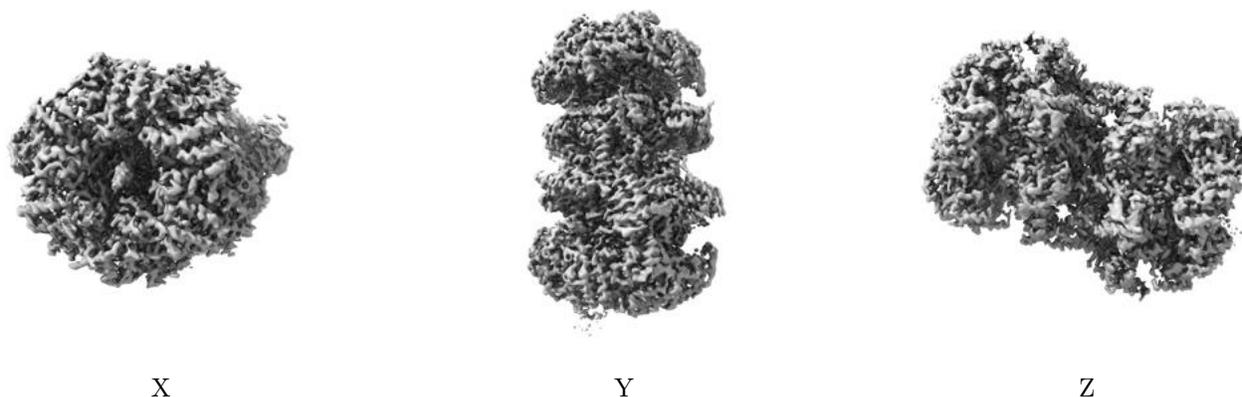


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

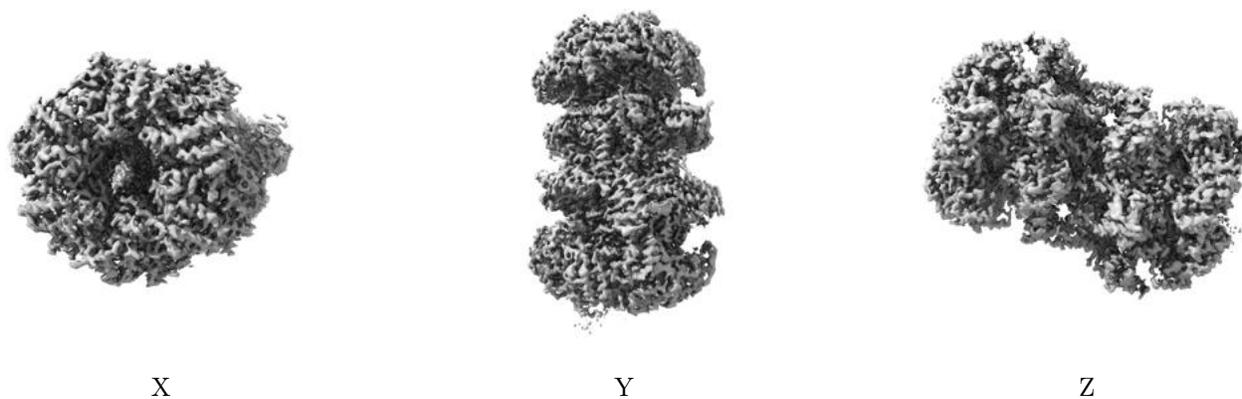
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



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

6.5.2 Raw map



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

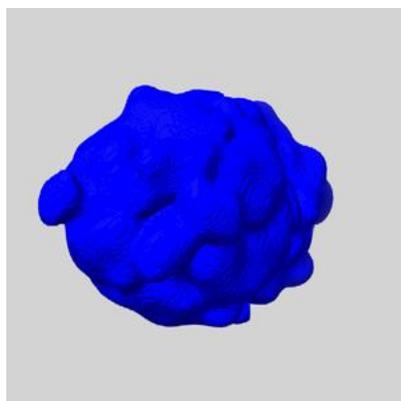
6.6 Mask visualisation [i](#)

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

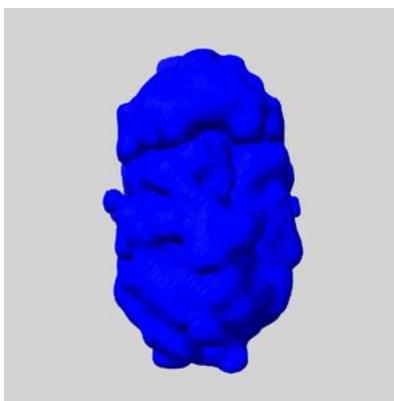
A mask typically either:

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

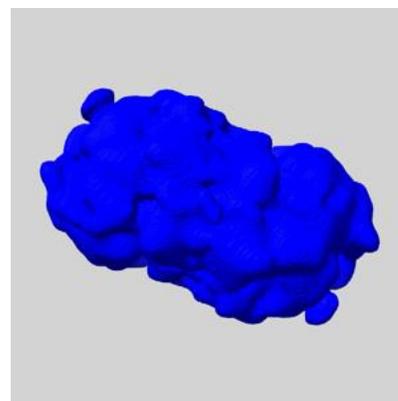
6.6.1 emd_19186_msk_1.map [i](#)



X



Y

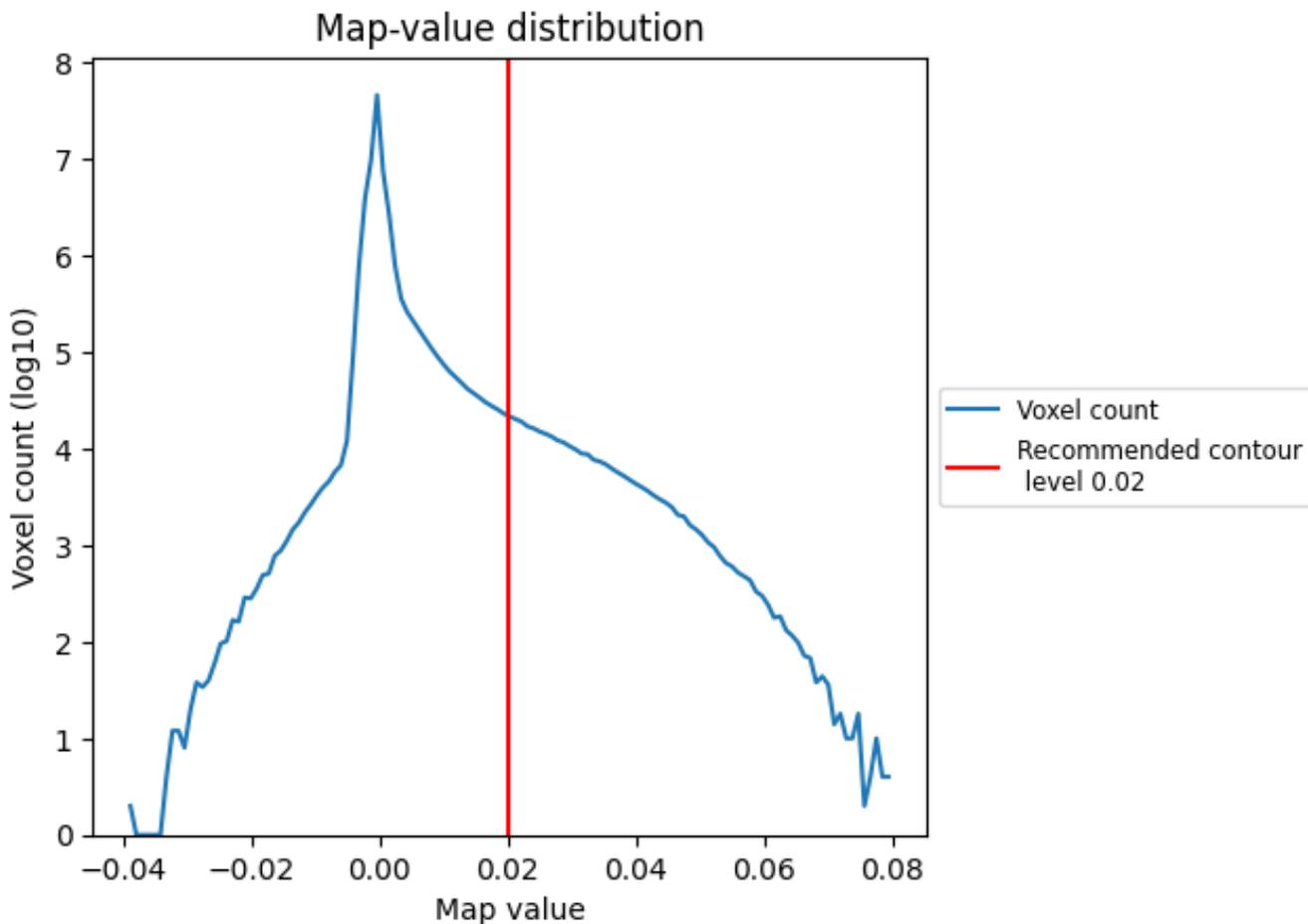


Z

7 Map analysis [i](#)

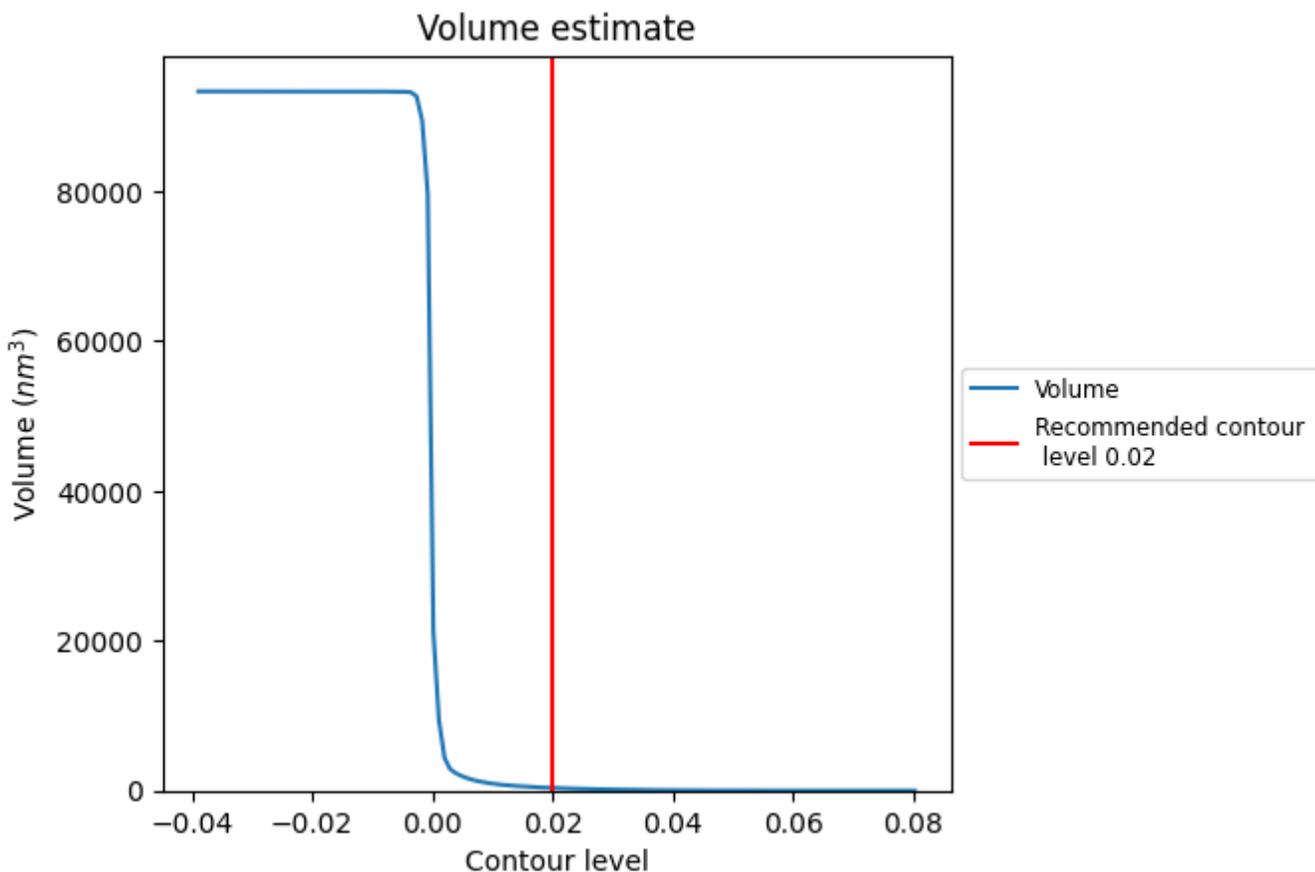
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

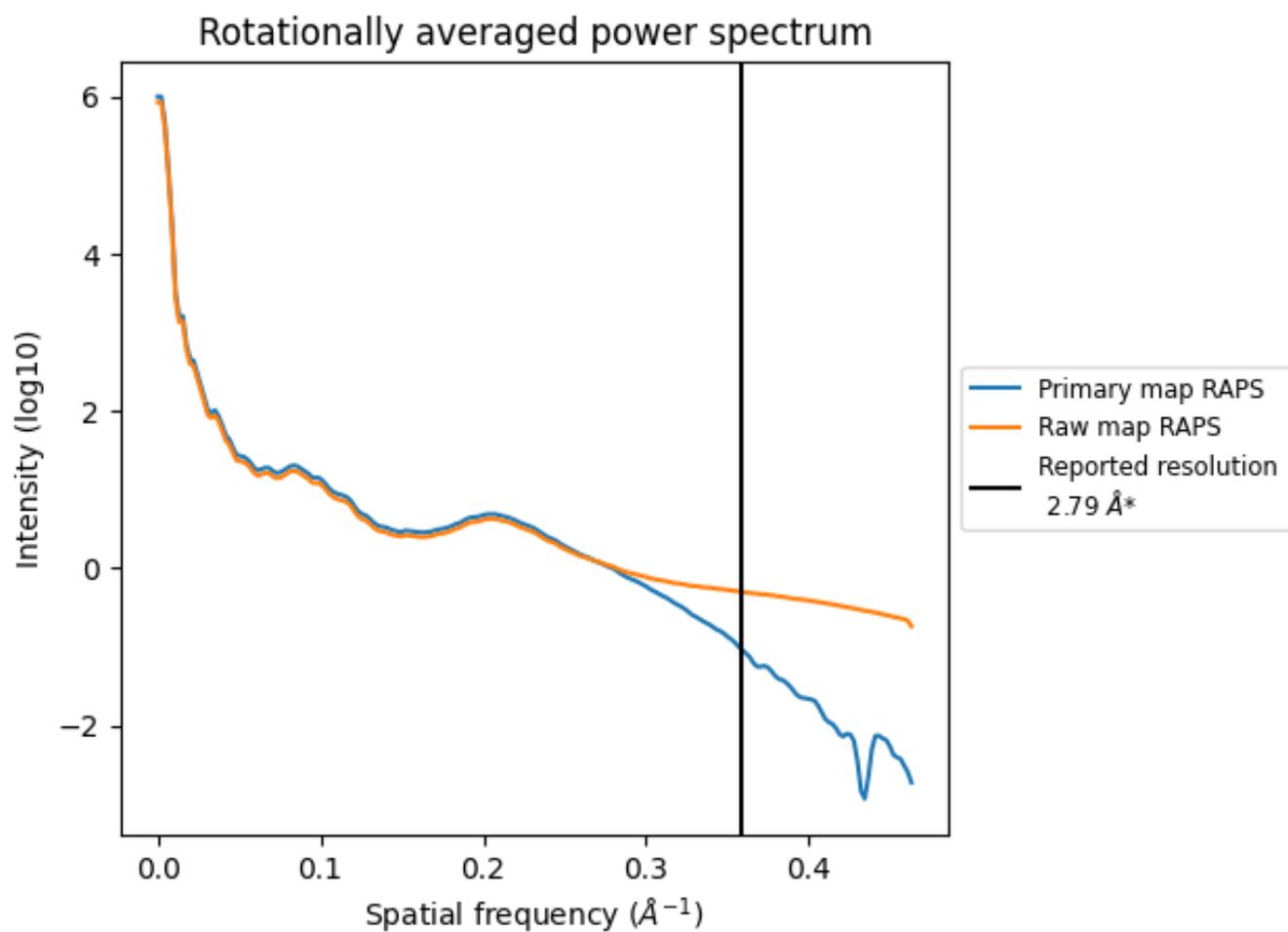
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 363 nm³; this corresponds to an approximate mass of 328 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

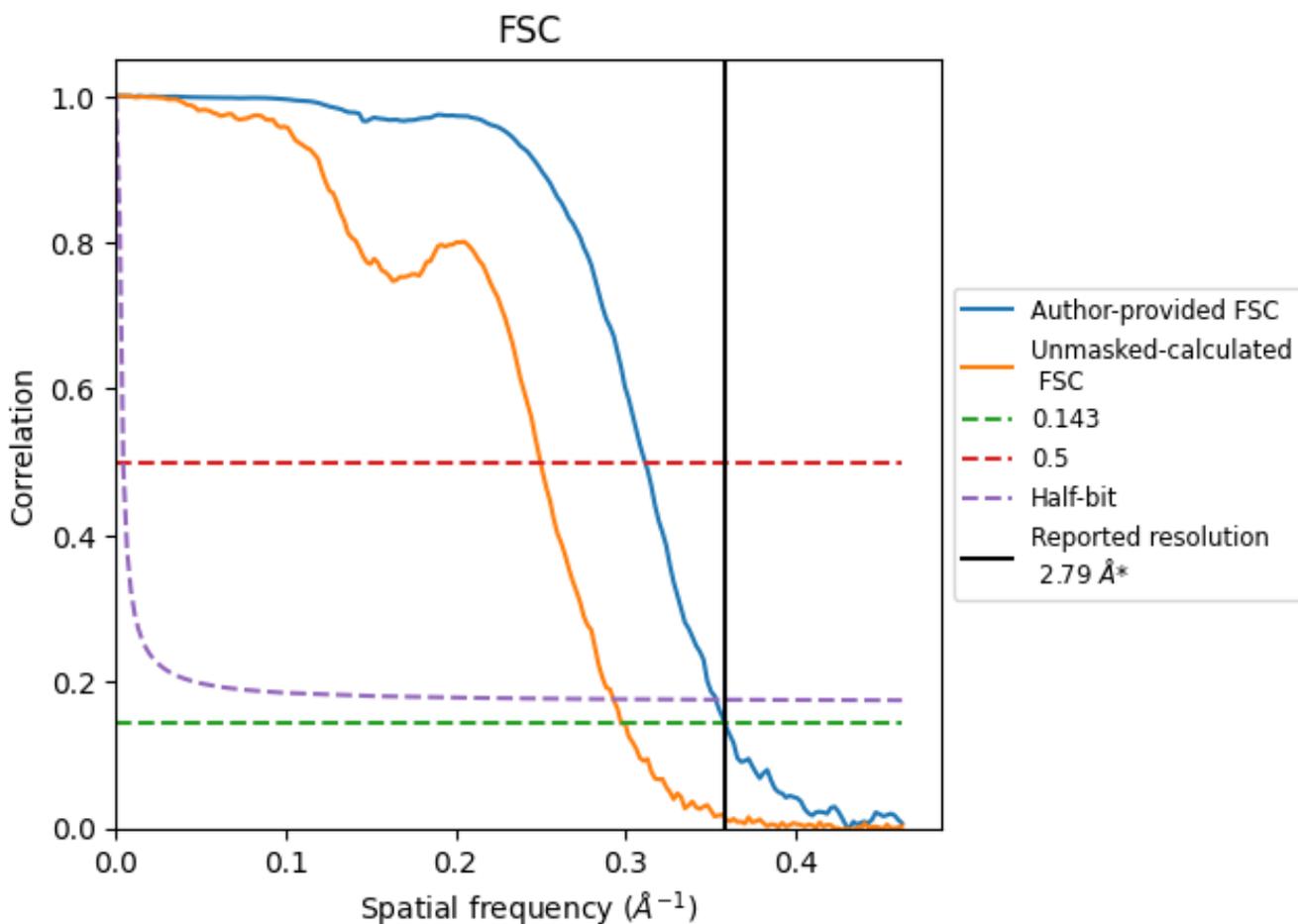


*Reported resolution corresponds to spatial frequency of 0.358 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.358 Å⁻¹

8.2 Resolution estimates [i](#)

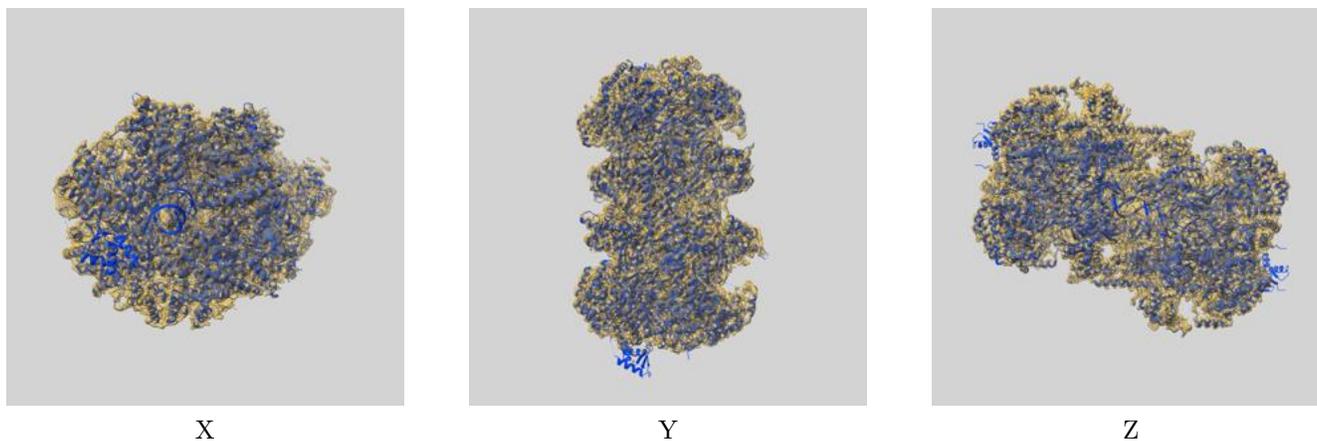
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.79	-	-
Author-provided FSC curve	2.79	3.21	2.83
Unmasked-calculated*	3.36	4.00	3.41

*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 3.36 differs from the reported value 2.79 by more than 10 %

9 Map-model fit [i](#)

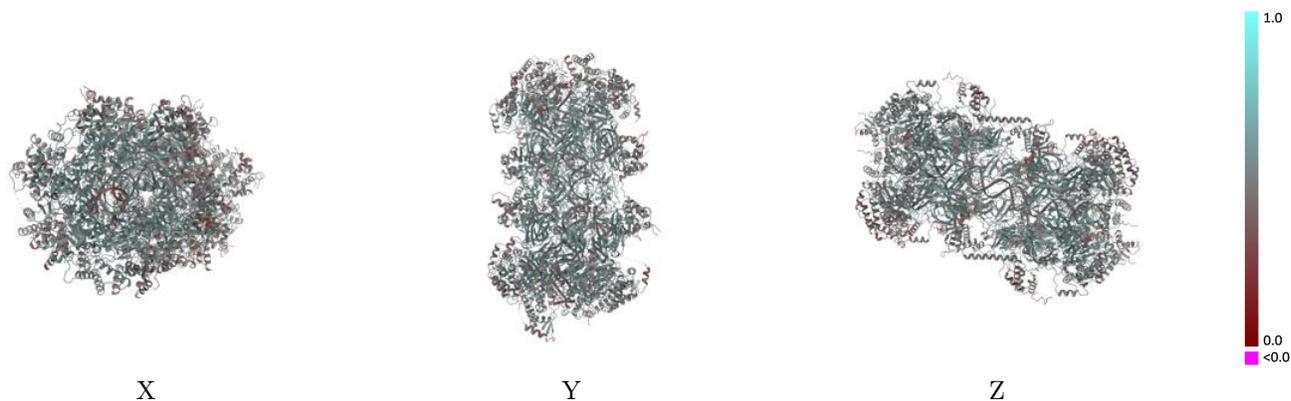
This section contains information regarding the fit between EMDB map EMD-19186 and PDB model 8RIF. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay [i](#)



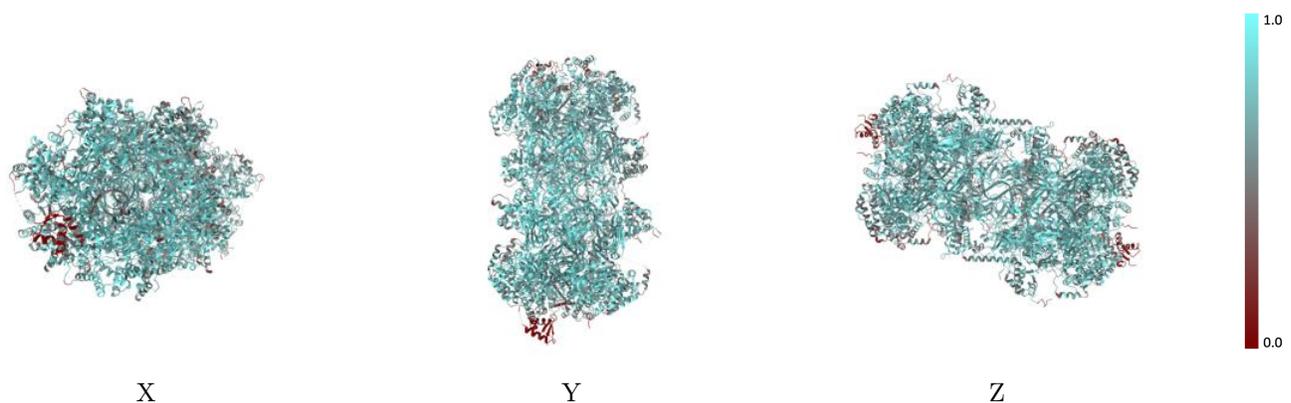
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



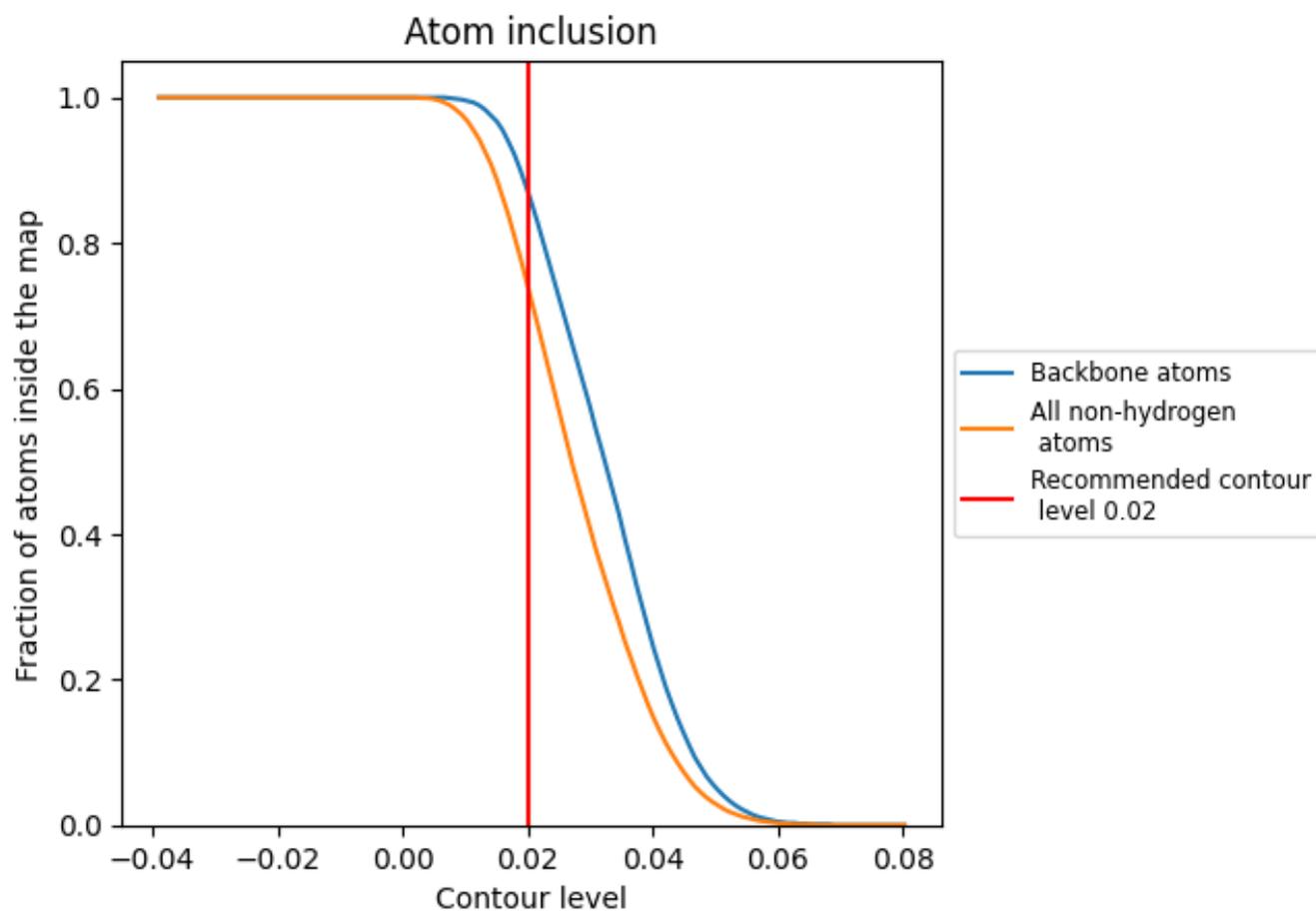
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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, 87% of all backbone atoms, 74% 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	 0.7420	 0.5160
2	 0.7140	 0.4990
3	 0.8070	 0.5360
4	 0.7480	 0.5180
5	 0.7480	 0.5060
6	 0.7400	 0.5220
7	 0.7200	 0.5210
A	 0.7120	 0.5000
B	 0.8070	 0.5380
C	 0.7450	 0.5170
D	 0.7450	 0.5080
E	 0.7380	 0.5230
F	 0.7180	 0.5210
X	 0.6960	 0.4630
Y	 0.6960	 0.4670

