



Full wwPDB EM Validation Report ⓘ

Sep 19, 2026 – 01:55 PM EDT

PDB ID : 9Q2L / pdb_00009q2l
EMDB ID : EMD-72167
Title : Rad55-Rad57(E161Q)-SHU-3xRad51 bound to ssDNA with ATP
Authors : Yatskevich, S.; Koo, C.W.; Ciferri, C.
Deposited on : 2025-08-15
Resolution : 3.70 Å(reported)

This is a Full wwPDB EM Validation 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:

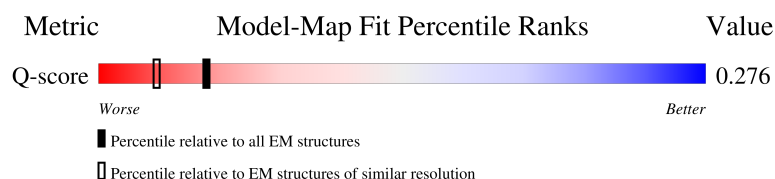
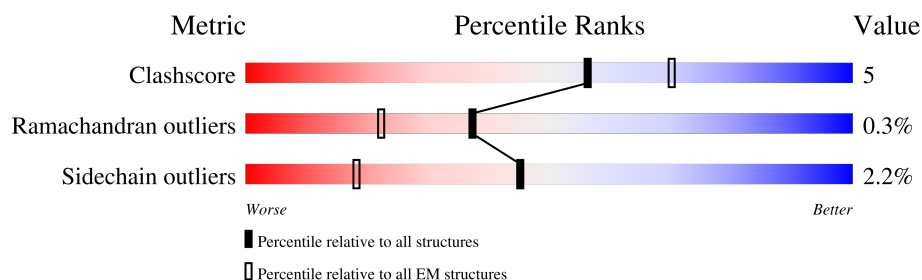
EMDB validation analysis : 0.0.1.dev133
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

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 3.70 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11569 (3.20 - 4.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	631	
2	B	460	
3	C	213	
4	D	281	

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Mol	Chain	Length	Quality of chain
5	E	150	21%81%17%
6	F	262	56%63%15%22%
7	G	418	9%67%10%23%
7	I	418	67%7%27%
7	J	418	48%65%8%27%
8	H	13	77%23%

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 19174 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 Methylated-DNA--protein-cysteine methyltransferase,DNA repair protein RAD55.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	272	2189	1410	373	395	11	0	0

There are 49 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-224	MET	-	expression tag	UNP E5BBQ0
A	-223	SER	-	expression tag	UNP E5BBQ0
A	-222	ALA	-	expression tag	UNP E5BBQ0
A	-221	TRP	-	expression tag	UNP E5BBQ0
A	-220	SER	-	expression tag	UNP E5BBQ0
A	-219	HIS	-	expression tag	UNP E5BBQ0
A	-218	PRO	-	expression tag	UNP E5BBQ0
A	-217	GLN	-	expression tag	UNP E5BBQ0
A	-216	PHE	-	expression tag	UNP E5BBQ0
A	-215	GLU	-	expression tag	UNP E5BBQ0
A	-214	LYS	-	expression tag	UNP E5BBQ0
A	-213	GLY	-	expression tag	UNP E5BBQ0
A	-212	GLY	-	expression tag	UNP E5BBQ0
A	-211	GLY	-	expression tag	UNP E5BBQ0
A	-210	SER	-	expression tag	UNP E5BBQ0
A	-209	GLY	-	expression tag	UNP E5BBQ0
A	-208	GLY	-	expression tag	UNP E5BBQ0
A	-207	GLY	-	expression tag	UNP E5BBQ0
A	-206	SER	-	expression tag	UNP E5BBQ0
A	-205	GLY	-	expression tag	UNP E5BBQ0
A	-204	GLY	-	expression tag	UNP E5BBQ0
A	-203	SER	-	expression tag	UNP E5BBQ0
A	-202	ALA	-	expression tag	UNP E5BBQ0
A	-201	TRP	-	expression tag	UNP E5BBQ0
A	-200	SER	-	expression tag	UNP E5BBQ0
A	-199	HIS	-	expression tag	UNP E5BBQ0
A	-198	PRO	-	expression tag	UNP E5BBQ0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-197	GLN	-	expression tag	UNP E5BBQ0
A	-196	PHE	-	expression tag	UNP E5BBQ0
A	-195	GLU	-	expression tag	UNP E5BBQ0
A	-194	LYS	-	expression tag	UNP E5BBQ0
A	-193	GLY	-	expression tag	UNP E5BBQ0
A	-192	SER	-	expression tag	UNP E5BBQ0
A	-191	MET	-	expression tag	UNP E5BBQ0
A	-162	ARG	GLU	conflict	UNP E5BBQ0
A	-13	PRO	-	linker	UNP E5BBQ0
A	-12	GLY	-	linker	UNP E5BBQ0
A	-11	LEU	-	linker	UNP E5BBQ0
A	-10	GLY	-	linker	UNP E5BBQ0
A	-9	GLY	-	linker	UNP E5BBQ0
A	-8	SER	-	linker	UNP E5BBQ0
A	-7	GLU	-	linker	UNP E5BBQ0
A	-6	ASN	-	linker	UNP E5BBQ0
A	-5	LEU	-	linker	UNP E5BBQ0
A	-4	TYR	-	linker	UNP E5BBQ0
A	-3	PHE	-	linker	UNP E5BBQ0
A	-2	GLN	-	linker	UNP E5BBQ0
A	-1	GLY	-	linker	UNP E5BBQ0
A	0	SER	-	linker	UNP E5BBQ0

- Molecule 2 is a protein called DNA repair protein RAD57.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	410	Total	C	N	O	S	0	0
			3267	2068	556	630	13		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	161	GLN	GLU	engineered mutation	UNP P25301

- Molecule 3 is a protein called Chromosome segregation in meiosis protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	189	Total	C	N	O	S	0	0
			1560	1010	253	291	6		

- Molecule 4 is a protein called Platinum sensitivity protein 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	223	Total 1841	C 1193	N 306	O 334	S 8	0	0

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-38	MET	-	initiating methionine	UNP Q12318
D	-37	SER	-	expression tag	UNP Q12318
D	-36	ALA	-	expression tag	UNP Q12318
D	-35	TRP	-	expression tag	UNP Q12318
D	-34	SER	-	expression tag	UNP Q12318
D	-33	HIS	-	expression tag	UNP Q12318
D	-32	PRO	-	expression tag	UNP Q12318
D	-31	GLN	-	expression tag	UNP Q12318
D	-30	PHE	-	expression tag	UNP Q12318
D	-29	GLU	-	expression tag	UNP Q12318
D	-28	LYS	-	expression tag	UNP Q12318
D	-27	GLY	-	expression tag	UNP Q12318
D	-26	GLY	-	expression tag	UNP Q12318
D	-25	GLY	-	expression tag	UNP Q12318
D	-24	SER	-	expression tag	UNP Q12318
D	-23	GLY	-	expression tag	UNP Q12318
D	-22	GLY	-	expression tag	UNP Q12318
D	-21	GLY	-	expression tag	UNP Q12318
D	-20	SER	-	expression tag	UNP Q12318
D	-19	GLY	-	expression tag	UNP Q12318
D	-18	GLY	-	expression tag	UNP Q12318
D	-17	SER	-	expression tag	UNP Q12318
D	-16	ALA	-	expression tag	UNP Q12318
D	-15	TRP	-	expression tag	UNP Q12318
D	-14	SER	-	expression tag	UNP Q12318
D	-13	HIS	-	expression tag	UNP Q12318
D	-12	PRO	-	expression tag	UNP Q12318
D	-11	GLN	-	expression tag	UNP Q12318
D	-10	PHE	-	expression tag	UNP Q12318
D	-9	GLU	-	expression tag	UNP Q12318
D	-8	LYS	-	expression tag	UNP Q12318
D	-7	SER	-	expression tag	UNP Q12318
D	-6	GLY	-	expression tag	UNP Q12318
D	-5	GLU	-	expression tag	UNP Q12318
D	-4	ASN	-	expression tag	UNP Q12318
D	-3	LEU	-	expression tag	UNP Q12318
D	-2	TYR	-	expression tag	UNP Q12318

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	PHE	-	expression tag	UNP Q12318
D	0	GLN	-	expression tag	UNP Q12318

- Molecule 5 is a protein called Suppressor of HU sensitivity involved in recombination protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	147	Total	C	N	O	S	0	0
			1184	760	207	213	4		

- Molecule 6 is a protein called Suppressor of hydroxyurea sensitivity protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	205	Total	C	N	O	S	0	0
			1703	1112	263	321	7		

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-38	MET	-	initiating methionine	UNP C7GVQ9
F	-37	SER	-	expression tag	UNP C7GVQ9
F	-36	ALA	-	expression tag	UNP C7GVQ9
F	-35	TRP	-	expression tag	UNP C7GVQ9
F	-34	SER	-	expression tag	UNP C7GVQ9
F	-33	HIS	-	expression tag	UNP C7GVQ9
F	-32	PRO	-	expression tag	UNP C7GVQ9
F	-31	GLN	-	expression tag	UNP C7GVQ9
F	-30	PHE	-	expression tag	UNP C7GVQ9
F	-29	GLU	-	expression tag	UNP C7GVQ9
F	-28	LYS	-	expression tag	UNP C7GVQ9
F	-27	GLY	-	expression tag	UNP C7GVQ9
F	-26	GLY	-	expression tag	UNP C7GVQ9
F	-25	GLY	-	expression tag	UNP C7GVQ9
F	-24	SER	-	expression tag	UNP C7GVQ9
F	-23	GLY	-	expression tag	UNP C7GVQ9
F	-22	GLY	-	expression tag	UNP C7GVQ9
F	-21	GLY	-	expression tag	UNP C7GVQ9
F	-20	SER	-	expression tag	UNP C7GVQ9
F	-19	GLY	-	expression tag	UNP C7GVQ9
F	-18	GLY	-	expression tag	UNP C7GVQ9
F	-17	SER	-	expression tag	UNP C7GVQ9
F	-16	ALA	-	expression tag	UNP C7GVQ9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-15	TRP	-	expression tag	UNP C7GVQ9
F	-14	SER	-	expression tag	UNP C7GVQ9
F	-13	HIS	-	expression tag	UNP C7GVQ9
F	-12	PRO	-	expression tag	UNP C7GVQ9
F	-11	GLN	-	expression tag	UNP C7GVQ9
F	-10	PHE	-	expression tag	UNP C7GVQ9
F	-9	GLU	-	expression tag	UNP C7GVQ9
F	-8	LYS	-	expression tag	UNP C7GVQ9
F	-7	SER	-	expression tag	UNP C7GVQ9
F	-6	GLY	-	expression tag	UNP C7GVQ9
F	-5	GLU	-	expression tag	UNP C7GVQ9
F	-4	ASN	-	expression tag	UNP C7GVQ9
F	-3	LEU	-	expression tag	UNP C7GVQ9
F	-2	TYR	-	expression tag	UNP C7GVQ9
F	-1	PHE	-	expression tag	UNP C7GVQ9
F	0	GLN	-	expression tag	UNP C7GVQ9
F	1	GLY	-	expression tag	UNP C7GVQ9

- Molecule 7 is a protein called DNA repair protein RAD51.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	321	Total	C	N	O	S	0	0
			2442	1529	431	464	18		
7	I	307	Total	C	N	O	S	0	0
			2326	1461	412	436	17		
7	J	304	Total	C	N	O	S	0	0
			2308	1449	409	433	17		

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	-17	MET	-	expression tag	UNP P25454
G	-16	HIS	-	expression tag	UNP P25454
G	-15	HIS	-	expression tag	UNP P25454
G	-14	HIS	-	expression tag	UNP P25454
G	-13	HIS	-	expression tag	UNP P25454
G	-12	HIS	-	expression tag	UNP P25454
G	-11	HIS	-	expression tag	UNP P25454
G	-10	HIS	-	expression tag	UNP P25454
G	-9	HIS	-	expression tag	UNP P25454
G	-8	GLY	-	expression tag	UNP P25454
G	-7	GLU	-	expression tag	UNP P25454

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-6	ASN	-	expression tag	UNP P25454
G	-5	LEU	-	expression tag	UNP P25454
G	-4	TYR	-	expression tag	UNP P25454
G	-3	PHE	-	expression tag	UNP P25454
G	-2	GLN	-	expression tag	UNP P25454
G	-1	GLY	-	expression tag	UNP P25454
G	0	SER	-	expression tag	UNP P25454
I	-17	MET	-	expression tag	UNP P25454
I	-16	HIS	-	expression tag	UNP P25454
I	-15	HIS	-	expression tag	UNP P25454
I	-14	HIS	-	expression tag	UNP P25454
I	-13	HIS	-	expression tag	UNP P25454
I	-12	HIS	-	expression tag	UNP P25454
I	-11	HIS	-	expression tag	UNP P25454
I	-10	HIS	-	expression tag	UNP P25454
I	-9	HIS	-	expression tag	UNP P25454
I	-8	GLY	-	expression tag	UNP P25454
I	-7	GLU	-	expression tag	UNP P25454
I	-6	ASN	-	expression tag	UNP P25454
I	-5	LEU	-	expression tag	UNP P25454
I	-4	TYR	-	expression tag	UNP P25454
I	-3	PHE	-	expression tag	UNP P25454
I	-2	GLN	-	expression tag	UNP P25454
I	-1	GLY	-	expression tag	UNP P25454
I	0	SER	-	expression tag	UNP P25454
J	-17	MET	-	expression tag	UNP P25454
J	-16	HIS	-	expression tag	UNP P25454
J	-15	HIS	-	expression tag	UNP P25454
J	-14	HIS	-	expression tag	UNP P25454
J	-13	HIS	-	expression tag	UNP P25454
J	-12	HIS	-	expression tag	UNP P25454
J	-11	HIS	-	expression tag	UNP P25454
J	-10	HIS	-	expression tag	UNP P25454
J	-9	HIS	-	expression tag	UNP P25454
J	-8	GLY	-	expression tag	UNP P25454
J	-7	GLU	-	expression tag	UNP P25454
J	-6	ASN	-	expression tag	UNP P25454
J	-5	LEU	-	expression tag	UNP P25454
J	-4	TYR	-	expression tag	UNP P25454
J	-3	PHE	-	expression tag	UNP P25454
J	-2	GLN	-	expression tag	UNP P25454
J	-1	GLY	-	expression tag	UNP P25454

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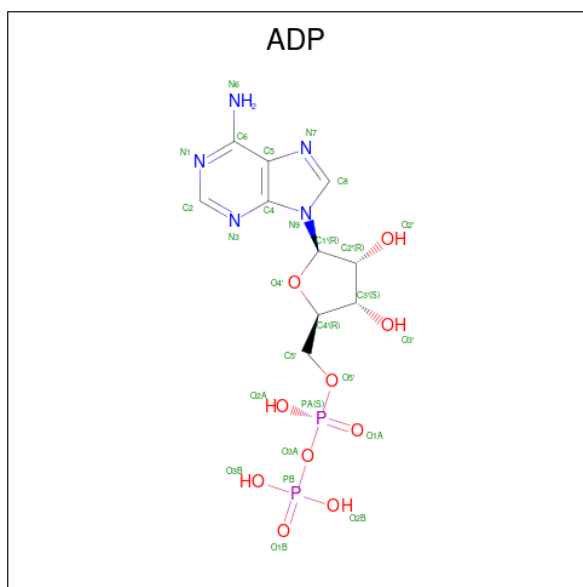
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Chain	Residue	Modelled	Actual	Comment	Reference
J	0	SER	-	expression tag	UNP P25454

- Molecule 8 is a DNA chain called ssDNA (13-mer).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	13	Total	C	N	O	P	0	0
			260	130	26	91	13		

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
10	A	1	Total	Mg	0
			1	1	
10	B	1	Total	Mg	0
			1	1	
10	G	1	Total	Mg	0
			1	1	
10	I	1	Total	Mg	0
			1	1	

- # ATP

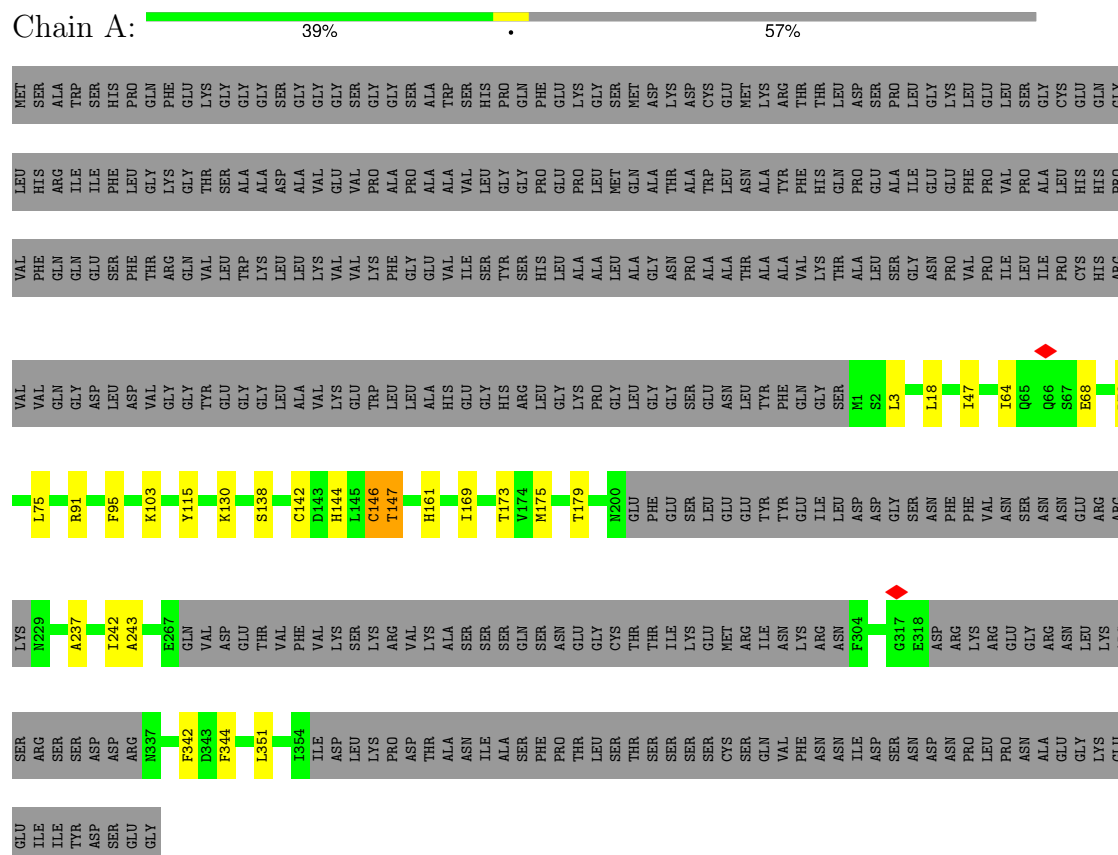
- Molecule 12 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	AltConf
12	F	1	Total Zn 1 1	0

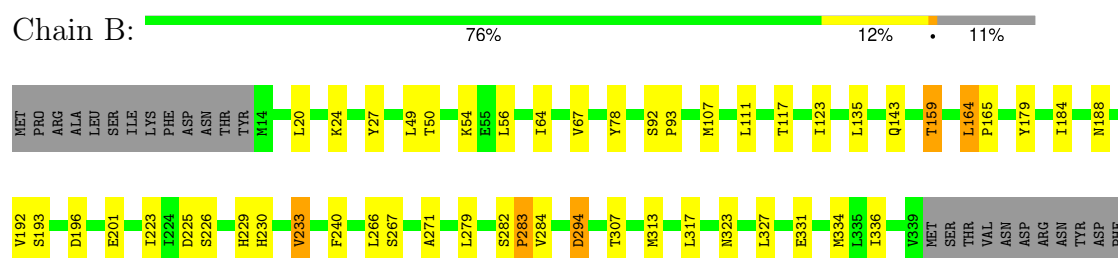
3 Residue-property plots

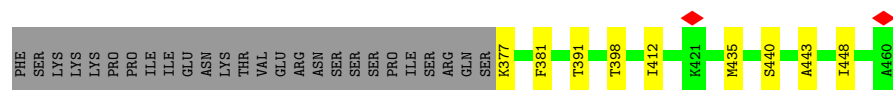
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

- Molecule 1: Methylated-DNA--protein-cysteine methyltransferase,DNA repair protein RAD55

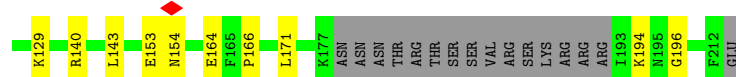
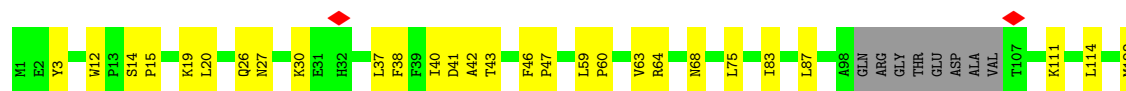


- Molecule 2: DNA repair protein RAD57

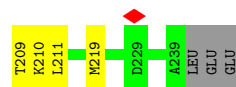
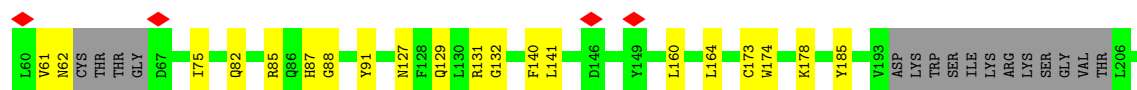
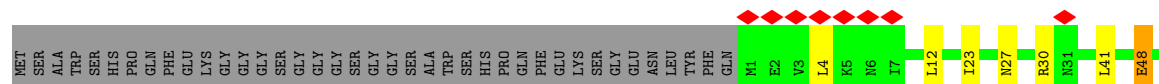




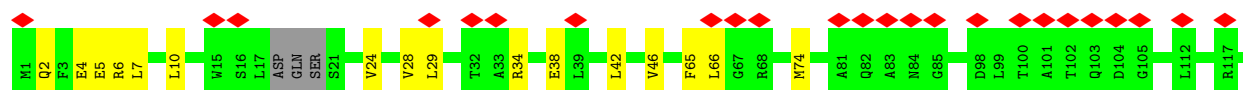
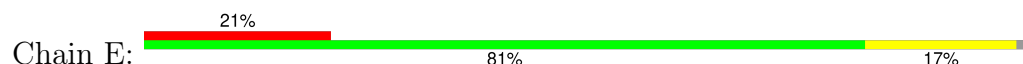
- Molecule 3: Chromosome segregation in meiosis protein 2



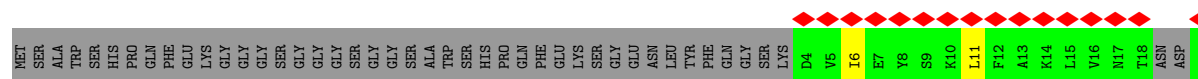
- Molecule 4: Platinum sensitivity protein 3

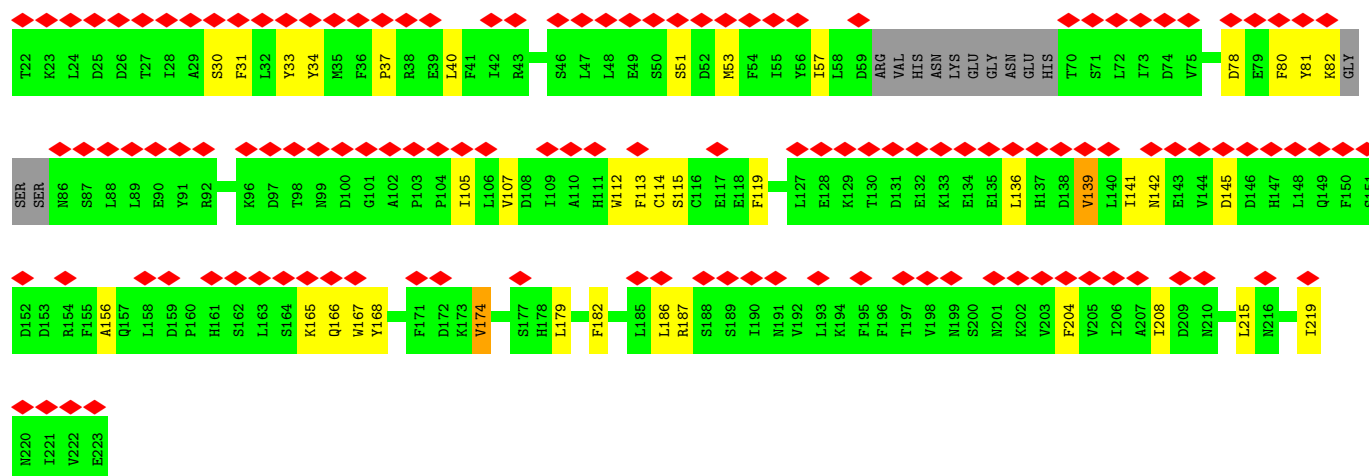


- Molecule 5: Suppressor of HU sensitivity involved in recombination protein 1



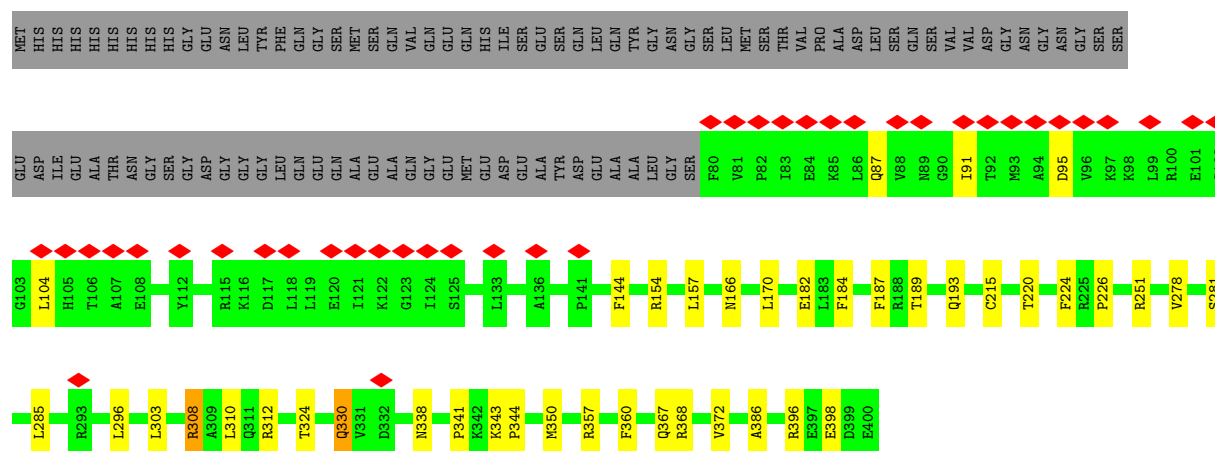
- Molecule 6: Suppressor of hydroxyurea sensitivity protein 2





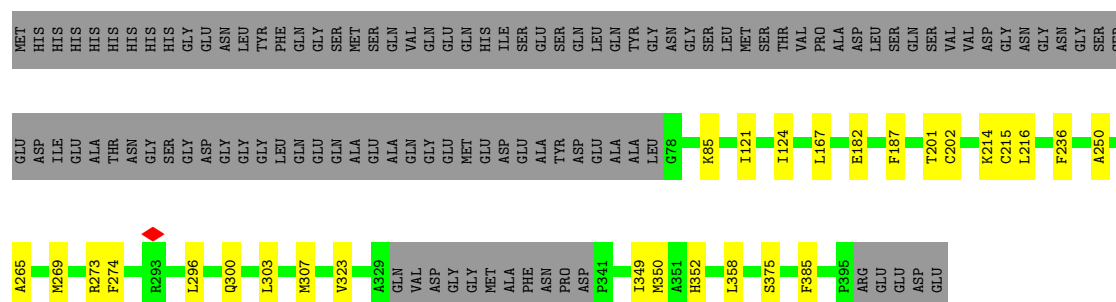
• Molecule 7: DNA repair protein RAD51

Chain G: 9% 67% 10% 23%



• Molecule 7: DNA repair protein RAD51

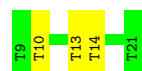
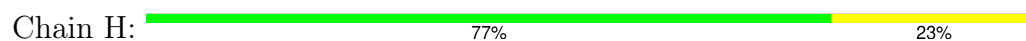
Chain I: 67% 7% 27%



• Molecule 7: DNA repair protein RAD51

Chain J: 48% 65% 8% 27%

- Molecule 8: ssDNA (13-mer)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	23196	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	56	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2600	Depositor
Magnification	Not provided	
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.327	Depositor
Minimum map value	-0.126	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.07	Depositor
Map size (Å)	362.5, 362.5, 362.5	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.90625, 0.90625, 0.90625	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.72	0/2227	0.74	1/2997 (0.0%)
2	B	0.69	0/3325	0.67	0/4494
3	C	0.58	0/1589	0.66	0/2147
4	D	0.49	0/1875	0.66	0/2529
5	E	0.35	0/1210	0.57	0/1645
6	F	0.21	0/1741	0.53	0/2357
7	G	0.60	0/2479	0.70	0/3342
7	I	0.56	1/2360 (0.0%)	0.62	0/3180
7	J	0.29	0/2340	0.61	0/3152
8	H	0.72	0/285	1.12	1/438 (0.2%)
All	All	0.55	1/19431 (0.0%)	0.66	2/26281 (0.0%)

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
5	E	0	1
7	G	0	1
7	J	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	I	375	SER	C-N	-6.17	1.28	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146	CYS	CA-CB-SG	5.14	126.22	114.40
8	H	14	DT	C1'-O4'-C4'	-5.12	102.02	109.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	E	6	ARG	Sidechain
7	G	368	ARG	Sidechain
7	J	312	ARG	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	2189	0	2230	18	0
2	B	3267	0	3264	40	0
3	C	1560	0	1597	30	0
4	D	1841	0	1883	21	0
5	E	1184	0	1202	14	0
6	F	1703	0	1658	21	0
7	G	2442	0	2454	29	0
7	I	2326	0	2364	18	0
7	J	2308	0	2346	16	0
8	H	260	0	157	2	0
9	A	27	0	12	1	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
10	G	1	0	0	0	0
10	I	1	0	0	0	0
11	B	31	0	12	0	0
11	G	31	0	12	0	0
12	F	1	0	0	0	0
All	All	19174	0	19191	189	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 5.

All (189) 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)
4:D:209:THR:HG22	4:D:211:LEU:H	1.49	0.78
7:G:296:LEU:HD21	8:H:13:DT:H2'	1.68	0.76
5:E:10:LEU:HD11	5:E:24:VAL:HG11	1.71	0.71
2:B:164:LEU:HD22	2:B:165:PRO:HD2	1.76	0.68
7:G:308:ARG:HB2	7:G:308:ARG:HH11	1.59	0.68
3:C:15:PRO:HA	3:C:19:LYS:HE2	1.76	0.66
4:D:27:ASN:HA	4:D:30:ARG:HD3	1.77	0.66
3:C:30:LYS:HE3	3:C:68:ASN:HD21	1.60	0.65
7:I:300:GLN:HB3	7:I:349:ILE:HD12	1.81	0.62
7:J:276:LEU:HD11	7:J:278:VAL:HG23	1.82	0.61
3:C:30:LYS:HE3	3:C:68:ASN:ND2	2.15	0.60
3:C:42:ALA:HB1	3:C:75:LEU:HD23	1.83	0.60
6:F:113:PHE:HA	6:F:119:PHE:CZ	2.38	0.59
7:I:303:LEU:HD21	7:I:350:MET:HG2	1.86	0.57
2:B:107:MET:HE3	2:B:435:MET:HE1	1.87	0.56
4:D:132:GLY:HA2	4:D:173:CYS:HB2	1.88	0.56
7:G:303:LEU:HD21	7:G:350:MET:HG2	1.88	0.56
5:E:28:VAL:HG22	5:E:65:PHE:HA	1.87	0.55
2:B:233:VAL:HG22	7:G:308:ARG:HH22	1.72	0.55
7:G:220:THR:HB	7:G:285:LEU:CD2	2.37	0.55
2:B:179:TYR:CG	2:B:184:ILE:HD12	2.42	0.55
2:B:225:ASP:CG	2:B:226:SER:H	2.15	0.55
7:I:216:LEU:HG	7:I:274:PHE:CE2	2.42	0.54
4:D:85:ARG:HA	4:D:91:TYR:OH	2.08	0.54
6:F:112:TRP:CZ3	6:F:179:LEU:HB3	2.43	0.54
7:I:202:CYS:SG	7:I:215:CYS:HB2	2.48	0.54
7:G:251:ARG:HH11	7:G:251:ARG:HG3	1.73	0.53
4:D:61:VAL:HG13	4:D:62:ASN:N	2.23	0.53
6:F:156:ALA:H	6:F:187:ARG:NH2	2.06	0.53
2:B:233:VAL:HG13	7:G:308:ARG:NH2	2.24	0.53
1:A:64:ILE:O	1:A:68:GLU:HA	2.09	0.53
3:C:12:TRP:HZ3	3:C:20:LEU:HD13	1.74	0.53
2:B:20:LEU:HD21	2:B:78:TYR:CE1	2.44	0.53
4:D:185:TYR:HB3	4:D:210:LYS:HD2	1.91	0.53
7:J:196:HIS:HB3	7:J:229:LEU:HD11	1.91	0.53
4:D:23:ILE:HD12	4:D:41:LEU:HB3	1.91	0.52
1:A:144:HIS:O	1:A:147:THR:HG22	2.10	0.52
4:D:12:LEU:HD21	5:E:7:LEU:HD21	1.90	0.52
4:D:61:VAL:HG13	4:D:62:ASN:H	1.74	0.52
4:D:48:GLU:HG3	4:D:178:LYS:NZ	2.24	0.52
3:C:140:ARG:HG2	3:C:140:ARG:HH11	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:131:ARG:CZ	4:D:174:TRP:HE1	2.23	0.51
7:G:330:GLN:HG2	7:G:343:LYS:HD3	1.92	0.51
8:H:10:DT:H3'	7:I:296:LEU:HD21	1.92	0.51
7:J:184:PHE:CD1	7:J:344:PRO:HD3	2.45	0.51
1:A:115:TYR:CD2	2:B:78:TYR:HB3	2.46	0.51
6:F:112:TRP:CE2	6:F:179:LEU:HD13	2.46	0.51
1:A:91:ARG:HG2	1:A:91:ARG:HH11	1.76	0.50
1:A:161:HIS:HA	1:A:243:ALA:HB1	1.92	0.50
2:B:398:THR:HG23	2:B:440:SER:HB2	1.93	0.50
7:G:220:THR:HB	7:G:285:LEU:HD23	1.94	0.50
7:J:214:LYS:HB2	7:J:274:PHE:HA	1.92	0.50
7:G:184:PHE:CG	7:G:344:PRO:HB3	2.48	0.49
2:B:111:LEU:O	2:B:443:ALA:HB2	2.13	0.49
5:E:147:HIS:CE1	5:E:148:ILE:HG12	2.48	0.49
7:J:308:ARG:O	7:J:312:ARG:HG2	2.11	0.49
5:E:2:GLN:O	5:E:5:GLU:HG3	2.12	0.49
3:C:41:ASP:OD1	3:C:42:ALA:N	2.46	0.49
3:C:14:SER:HB3	3:C:166:PRO:HA	1.94	0.49
5:E:133:PRO:HB2	5:E:136:SER:HB3	1.93	0.49
2:B:196:ASP:HA	2:B:230:HIS:CE1	2.47	0.49
7:G:91:ILE:HG23	7:G:95:ASP:HB2	1.94	0.49
2:B:313:MET:HG3	2:B:317:LEU:HD13	1.94	0.49
1:A:138:SER:HB2	1:A:237:ALA:HB3	1.95	0.48
5:E:137:VAL:CG2	5:E:138:PRO:HD3	2.43	0.48
7:G:170:LEU:HD21	7:G:372:VAL:HG23	1.95	0.48
7:J:183:LEU:HB3	7:J:360:PHE:CE1	2.47	0.48
1:A:342:PHE:CE2	1:A:351:LEU:HD23	2.48	0.48
7:J:219:ASP:O	7:J:251:ARG:HD2	2.13	0.48
7:G:87:GLN:HA	7:G:91:ILE:HB	1.96	0.48
7:J:153:ARG:HH11	7:J:153:ARG:HG3	1.79	0.48
3:C:30:LYS:HD2	3:C:64:ARG:HG3	1.96	0.48
4:D:85:ARG:HG2	4:D:91:TYR:CE2	2.48	0.48
7:G:308:ARG:HH11	7:G:308:ARG:CB	2.26	0.48
3:C:60:PRO:HA	3:C:63:VAL:HG22	1.94	0.47
9:A:501:ADP:H2'	2:B:440:SER:O	2.13	0.47
3:C:15:PRO:HD2	3:C:164:GLU:CD	2.39	0.47
1:A:3:LEU:HD12	3:C:38:PHE:CE2	2.49	0.47
2:B:92:SER:HB2	2:B:93:PRO:HD2	1.94	0.47
3:C:46:PHE:CG	3:C:47:PRO:HD2	2.50	0.47
6:F:11:LEU:H	6:F:11:LEU:HD22	1.79	0.47
6:F:30:SER:HA	6:F:33:TYR:CD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:28:VAL:CG2	5:E:65:PHE:HA	2.45	0.47
3:C:171:LEU:CD1	3:C:194:LYS:HE3	2.45	0.47
5:E:29:LEU:HA	5:E:66:LEU:O	2.15	0.46
3:C:3:TYR:CZ	3:C:111:LYS:HB3	2.51	0.46
5:E:137:VAL:HA	5:E:140:ARG:HG2	1.97	0.46
3:C:153:GLU:HG3	3:C:154:ASN:ND2	2.30	0.46
6:F:37:PRO:HB2	6:F:40:LEU:HD13	1.96	0.46
7:J:257:HIS:ND1	7:J:261:LEU:HG	2.31	0.46
3:C:87:LEU:HD11	3:C:114:LEU:HD13	1.97	0.46
1:A:75:LEU:HD12	1:A:103:LYS:O	2.16	0.46
2:B:192:VAL:HG23	7:G:144:PHE:CD1	2.50	0.46
5:E:74:MET:SD	6:F:219:ILE:HA	2.56	0.46
7:G:281:SER:H	7:G:324:THR:HB	1.80	0.46
6:F:80:PHE:C	6:F:82:LYS:H	2.24	0.46
7:I:216:LEU:HD13	7:I:265:ALA:HB1	1.97	0.46
2:B:164:LEU:HD22	2:B:165:PRO:CD	2.45	0.45
7:G:182:GLU:OE1	7:G:357:ARG:HD2	2.16	0.45
7:J:199:ALA:HA	7:J:215:CYS:SG	2.56	0.45
1:A:142:CYS:O	1:A:146:CYS:HB2	2.16	0.45
3:C:41:ASP:CG	3:C:43:THR:H	2.23	0.45
7:G:187:PHE:HA	7:I:352:HIS:CE1	2.51	0.45
3:C:143:LEU:HA	3:C:143:LEU:HD23	1.75	0.45
4:D:87:HIS:CG	4:D:88:GLY:N	2.85	0.45
7:G:338:ASN:ND2	7:G:341:PRO:HA	2.32	0.45
1:A:95:PHE:CE2	2:B:93:PRO:HA	2.52	0.45
3:C:140:ARG:CD	4:D:140:PHE:HB3	2.47	0.45
2:B:117:THR:HG22	2:B:267:SER:HB2	1.99	0.44
3:C:171:LEU:HD11	3:C:194:LYS:HE3	1.99	0.44
1:A:344:PHE:CE1	1:A:351:LEU:HG	2.53	0.44
6:F:105:ILE:HG23	6:F:115:SER:OG	2.17	0.44
3:C:153:GLU:HB3	4:D:82:GLN:HE21	1.83	0.44
4:D:48:GLU:HG3	4:D:178:LYS:HZ3	1.83	0.44
1:A:3:LEU:HD12	3:C:38:PHE:CD2	2.53	0.44
2:B:266:LEU:C	2:B:266:LEU:HD12	2.43	0.44
3:C:40:ILE:HD11	3:C:83:ILE:HG21	2.00	0.44
7:G:224:PHE:CE2	7:G:226:PRO:HG3	2.53	0.44
3:C:26:GLN:C	3:C:27:ASN:HD22	2.25	0.43
7:G:396:ARG:HG2	7:G:398:GLU:H	1.83	0.43
7:I:201:THR:HG22	7:I:236:PHE:CZ	2.52	0.43
2:B:56:LEU:HB2	2:B:67:VAL:HG11	2.00	0.43
7:I:121:ILE:HB	7:I:124:ILE:HG13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:ILE:HG13	1:A:243:ALA:N	2.34	0.43
2:B:24:LYS:CD	2:B:24:LYS:H	2.32	0.43
4:D:75:ILE:HD13	4:D:141:LEU:HD21	1.99	0.43
7:G:310:LEU:HD23	7:G:310:LEU:HA	1.83	0.43
2:B:294:ASP:OD1	2:B:294:ASP:N	2.52	0.43
2:B:331:GLU:HA	2:B:334:MET:SD	2.58	0.43
2:B:240:PHE:CD1	2:B:240:PHE:C	2.97	0.43
7:I:269:MET:HG2	7:I:274:PHE:CD1	2.54	0.43
1:A:72:ASP:HB2	1:A:130:LYS:HE2	2.00	0.42
2:B:49:LEU:HD23	2:B:49:LEU:HA	1.74	0.42
5:E:34:ARG:HG3	5:E:34:ARG:HH11	1.84	0.42
5:E:150:ARG:HG3	5:E:150:ARG:HH11	1.84	0.42
3:C:140:ARG:HD3	4:D:140:PHE:HB3	2.01	0.42
7:I:214:LYS:HE2	7:I:273:ARG:O	2.19	0.42
2:B:50:THR:O	3:C:129:LYS:HE2	2.20	0.42
7:G:154:ARG:HA	7:G:157:LEU:HD23	2.02	0.42
3:C:12:TRP:HH2	3:C:20:LEU:HA	1.85	0.42
6:F:31:PHE:HA	6:F:34:TYR:CE2	2.55	0.42
6:F:142:ASN:OD1	6:F:167:TRP:HA	2.19	0.42
2:B:307:THR:HG23	2:B:381:PHE:HB2	2.01	0.42
7:G:312:ARG:HG2	7:G:312:ARG:HH11	1.85	0.42
7:I:85:LYS:HE2	7:I:85:LYS:HB2	1.87	0.42
6:F:145:ASP:HA	6:F:165:LYS:HA	2.02	0.42
7:G:189:THR:HB	7:G:360:PHE:HB3	2.02	0.42
2:B:159:THR:O	2:B:193:SER:HA	2.20	0.42
2:B:327:LEU:HD12	2:B:327:LEU:HA	1.82	0.42
2:B:54:LYS:HA	2:B:64:ILE:CD1	2.50	0.41
5:E:42:LEU:HB3	5:E:46:VAL:CG1	2.49	0.41
7:I:167:LEU:HD23	7:I:385:PHE:HB3	2.02	0.41
7:J:199:ALA:HB3	7:J:229:LEU:HD21	2.02	0.41
2:B:54:LYS:HA	2:B:64:ILE:HD11	2.01	0.41
3:C:30:LYS:HG3	3:C:64:ARG:NH1	2.35	0.41
7:J:87:GLN:HE22	7:J:93:MET:HB3	1.85	0.41
1:A:130:LYS:HE2	1:A:130:LYS:HB2	1.75	0.41
2:B:123:ILE:O	2:B:271:ALA:HA	2.21	0.41
7:G:367:GLN:HA	7:G:386:ALA:HA	2.02	0.41
7:I:307:MET:HE3	7:I:350:MET:SD	2.60	0.41
2:B:225:ASP:CG	2:B:226:SER:N	2.77	0.41
6:F:51:SER:HB2	6:F:204:PHE:CE2	2.55	0.41
6:F:136:LEU:HA	6:F:139:VAL:HG22	2.02	0.41
7:I:182:GLU:HA	7:I:323:VAL:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:I:250:ALA:HB2	7:J:144:PHE:CD1	2.55	0.41
3:C:37:LEU:HD23	3:C:38:PHE:N	2.36	0.41
6:F:53:MET:HB3	6:F:182:PHE:CZ	2.56	0.41
7:J:159:CYS:HA	7:J:175:VAL:O	2.21	0.41
2:B:229:HIS:O	7:G:308:ARG:NH2	2.53	0.41
6:F:112:TRP:CH2	6:F:179:LEU:HB3	2.55	0.41
2:B:143:GLN:OE1	2:B:188:ASN:HB2	2.20	0.41
2:B:233:VAL:HG21	7:G:308:ARG:HH12	1.85	0.41
4:D:219:MET:HB3	4:D:219:MET:HE2	1.86	0.41
6:F:113:PHE:CG	6:F:114:CYS:N	2.89	0.41
2:B:336:ILE:HD13	2:B:412:ILE:HD13	2.02	0.41
6:F:119:PHE:CE2	6:F:174:VAL:HG12	2.55	0.41
1:A:344:PHE:CD1	1:A:344:PHE:N	2.89	0.41
7:G:166:ASN:N	7:G:166:ASN:HD22	2.18	0.41
7:I:187:PHE:CZ	7:J:351:ALA:HB1	2.56	0.41
7:I:358:LEU:HA	7:I:358:LEU:HD23	1.76	0.41
4:D:160:LEU:O	4:D:164:LEU:HD23	2.21	0.41
6:F:57:ILE:O	6:F:208:ILE:N	2.53	0.41
6:F:166:GLN:HG3	6:F:168:TYR:CE2	2.55	0.41
1:A:175:MET:O	1:A:179:THR:HG23	2.21	0.40
2:B:448:ILE:HD12	2:B:448:ILE:HA	1.96	0.40
2:B:135:LEU:HD13	2:B:223:ILE:HG23	2.03	0.40
7:J:161:THR:HG23	7:J:201:THR:HB	2.03	0.40
4:D:27:ASN:HA	4:D:30:ARG:CD	2.46	0.40
2:B:20:LEU:HD21	2:B:78:TYR:CZ	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	264/631 (42%)	251 (95%)	13 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	406/460 (88%)	382 (94%)	21 (5%)	3 (1%)	18	50
3	C	183/213 (86%)	174 (95%)	7 (4%)	2 (1%)	11	42
4	D	217/281 (77%)	200 (92%)	17 (8%)	0	100	100
5	E	143/150 (95%)	136 (95%)	7 (5%)	0	100	100
6	F	197/262 (75%)	182 (92%)	14 (7%)	1 (0%)	24	56
7	G	319/418 (76%)	299 (94%)	20 (6%)	0	100	100
7	I	303/418 (72%)	290 (96%)	13 (4%)	0	100	100
7	J	300/418 (72%)	287 (96%)	12 (4%)	1 (0%)	36	65
All	All	2332/3251 (72%)	2201 (94%)	124 (5%)	7 (0%)	37	65

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	283	PRO
2	B	284	VAL
3	C	59	LEU
6	F	141	ILE
3	C	196	GLY
7	J	211	GLY
2	B	282	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	246/543 (45%)	241 (98%)	5 (2%)	48	64
2	B	366/415 (88%)	355 (97%)	11 (3%)	36	57
3	C	180/202 (89%)	179 (99%)	1 (1%)	78	79
4	D	210/255 (82%)	206 (98%)	4 (2%)	50	65
5	E	132/135 (98%)	128 (97%)	4 (3%)	36	57
6	F	195/238 (82%)	187 (96%)	8 (4%)	27	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	G	260/336 (77%)	254 (98%)	6 (2%)	44	62
7	I	248/336 (74%)	248 (100%)	0	100	100
7	J	246/336 (73%)	239 (97%)	7 (3%)	38	58
All	All	2083/2796 (74%)	2037 (98%)	46 (2%)	45	63

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	47	ILE
1	A	147	THR
1	A	169	ILE
1	A	173	THR
2	B	27	TYR
2	B	159	THR
2	B	164	LEU
2	B	201	GLU
2	B	233	VAL
2	B	279	LEU
2	B	283	PRO
2	B	294	ASP
2	B	323	ASN
2	B	377	LYS
2	B	391	THR
3	C	122	MET
4	D	4	LEU
4	D	48	GLU
4	D	127	ASN
4	D	129	GLN
5	E	4	GLU
5	E	38	GLU
5	E	120	LEU
5	E	149	ILE
6	F	6	ILE
6	F	78	ASP
6	F	81	TYR
6	F	107	VAL
6	F	139	VAL
6	F	174	VAL
6	F	186	LEU
6	F	215	LEU

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Mol	Chain	Res	Type
7	G	104	LEU
7	G	193	GLN
7	G	215	CYS
7	G	278	VAL
7	G	308	ARG
7	G	330	GLN
7	J	167	LEU
7	J	206	LEU
7	J	208	ILE
7	J	249	TYR
7	J	281	SER
7	J	312	ARG
7	J	324	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	119	ASN
1	A	125	GLN
1	A	144	HIS
2	B	41	ASN
2	B	186	GLN
2	B	230	HIS
2	B	248	ASN
2	B	322	ASN
3	C	27	ASN
3	C	36	GLN
3	C	52	GLN
3	C	68	ASN
4	D	14	ASN
4	D	122	GLN
4	D	233	GLN
5	E	23	ASN
5	E	45	HIS
5	E	118	GLN
7	G	166	ASN
7	G	257	HIS
7	I	87	GLN
7	I	196	HIS
7	I	316	GLN
7	I	326	GLN

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Continued from previous page...

Mol	Chain	Res	Type
7	I	352	HIS
7	J	193	GLN
7	J	246	ASN
7	J	300	GLN
7	J	311	GLN
7	J	316	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 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 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	$\# Z > 2$
9	ADP	A	501	10	28,29,29	1.57	5 (17%)	43,45,45	1.74	12 (27%)
11	ATP	B	501	10	32,33,33	2.39	2 (6%)	48,52,52	0.66	1 (2%)
11	ATP	G	501	10	32,33,33	2.85	2 (6%)	48,52,52	0.84	2 (4%)

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	A	501	10	-	4/16/32/32	0/3/3/3
11	ATP	B	501	10	-	4/22/38/38	0/3/3/3
11	ATP	G	501	10	-	1/22/38/38	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	B	501	ATP	PB-O3B	-12.69	1.45	1.59
11	G	501	ATP	PA-O3A	-11.74	1.46	1.59
11	G	501	ATP	PB-O3B	-10.61	1.48	1.59
9	A	501	ADP	C4-N9	-3.53	1.30	1.37
9	A	501	ADP	C5-C4	3.41	1.45	1.39
9	A	501	ADP	C5-N7	-3.34	1.33	1.39
11	B	501	ATP	PA-O3A	-3.30	1.55	1.59
9	A	501	ADP	C8-N9	-2.27	1.33	1.37
9	A	501	ADP	C2'-C3'	-2.12	1.47	1.53

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	501	ADP	C5-C4-N3	-4.42	120.64	126.72
9	A	501	ADP	N3-C2-N1	-3.29	123.60	128.58
9	A	501	ADP	C2-N3-C4	3.24	119.75	111.83
9	A	501	ADP	C4-C5-N7	-3.15	106.98	110.58
9	A	501	ADP	N3-C4-N9	2.98	132.24	127.17
11	G	501	ATP	O4'-C4'-C3'	-2.63	99.93	105.15
9	A	501	ADP	O3B-PB-O2B	2.57	117.42	107.80
11	G	501	ATP	O5'-PA-O1A	-2.55	98.83	108.94
9	A	501	ADP	O4'-C1'-N9	2.38	112.66	108.09
11	B	501	ATP	O3G-PG-O2G	2.38	116.71	107.80
9	A	501	ADP	C6-C5-N7	2.36	136.63	132.09
9	A	501	ADP	O2A-PA-O1A	2.32	123.23	112.44
9	A	501	ADP	C4-N9-C8	2.27	108.13	105.74
9	A	501	ADP	O2'-C2'-C3'	-2.08	105.15	111.82
9	A	501	ADP	C2-N1-C6	2.02	122.05	118.73

There are no chirality outliers.

All (9) torsion outliers are listed below:

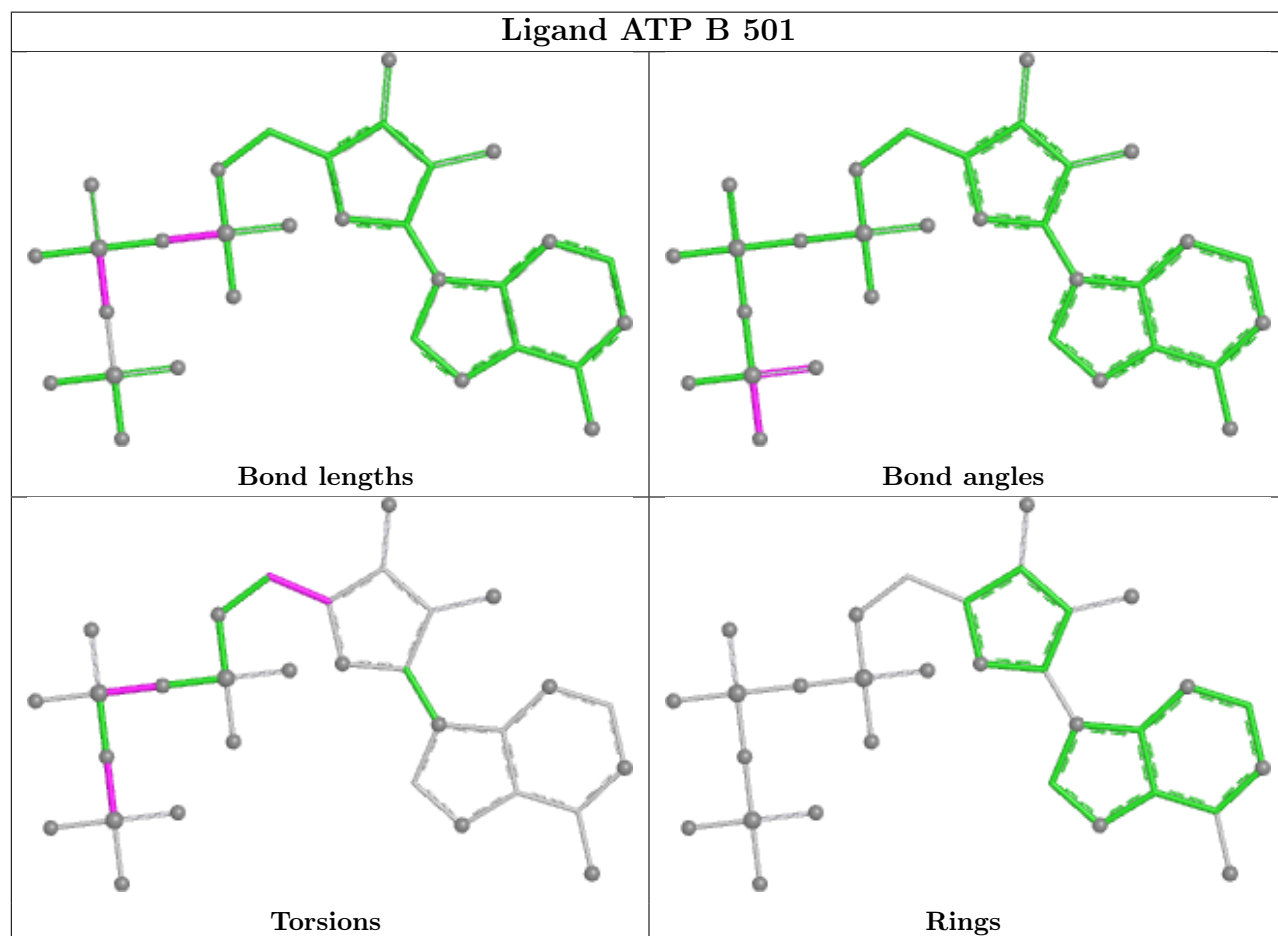
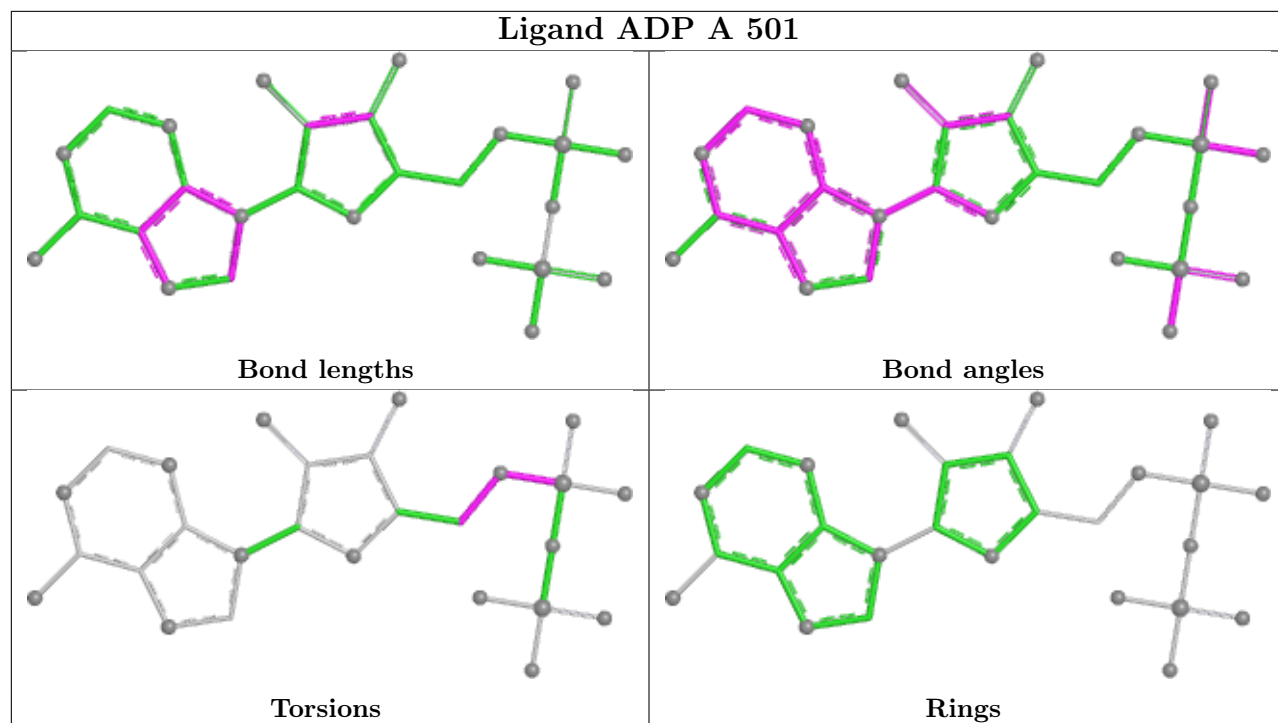
Mol	Chain	Res	Type	Atoms
9	A	501	ADP	C5'-O5'-PA-O1A
9	A	501	ADP	C5'-O5'-PA-O3A
11	B	501	ATP	C3'-C4'-C5'-O5'
11	B	501	ATP	O4'-C4'-C5'-O5'
9	A	501	ADP	C5'-O5'-PA-O2A
11	B	501	ATP	PB-O3B-PG-O1G
11	G	501	ATP	PA-O3A-PB-O1B
11	B	501	ATP	PA-O3A-PB-O2B
9	A	501	ADP	C4'-C5'-O5'-PA

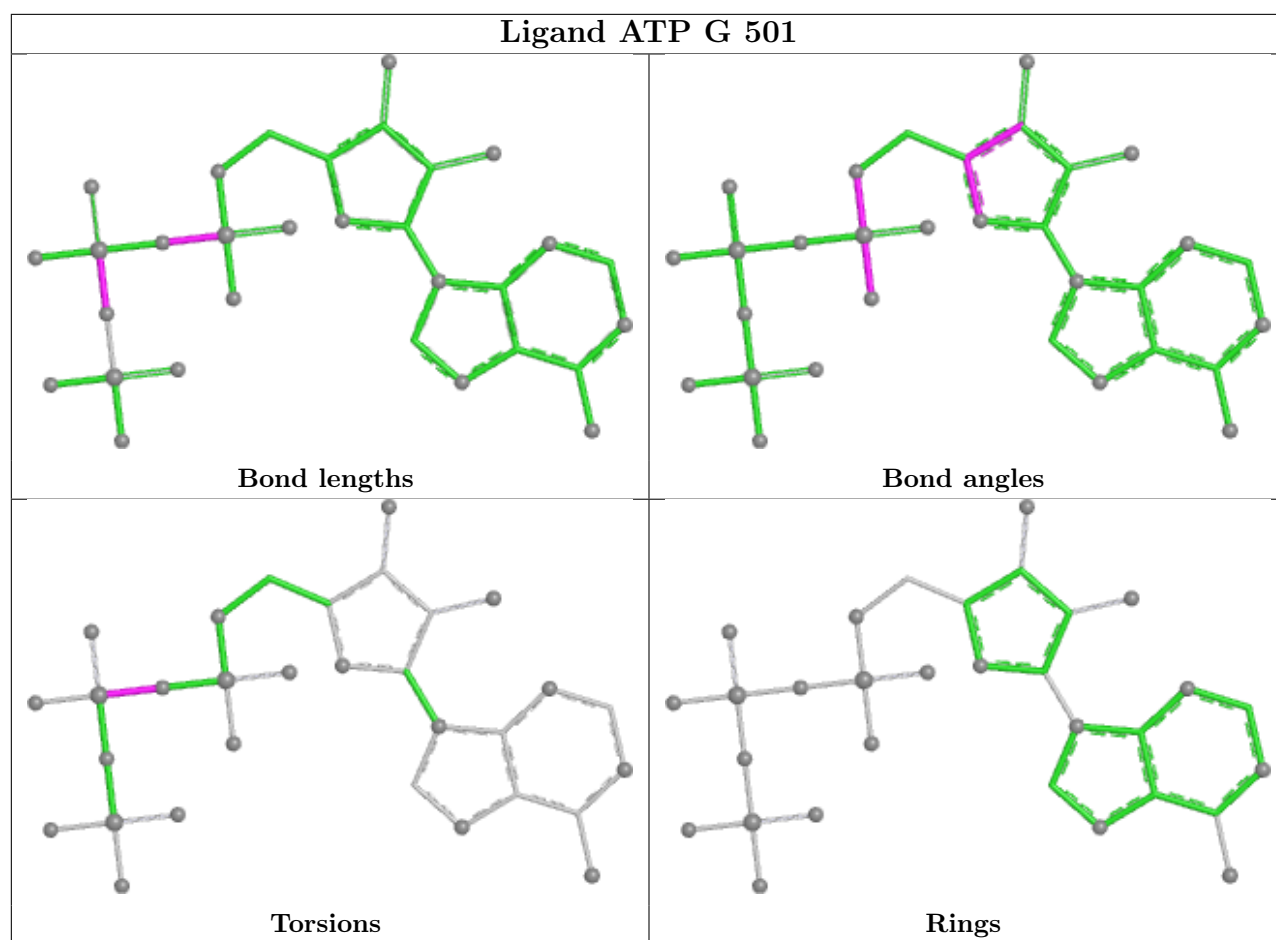
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	501	ADP	1	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.

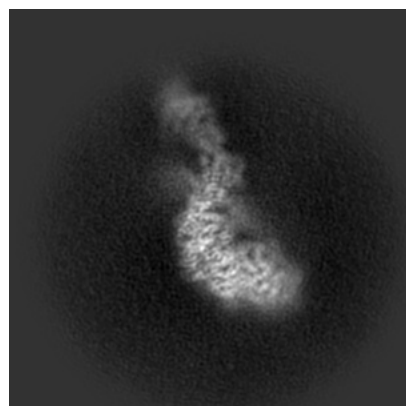
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-72167. These allow visual inspection of the internal detail of the map and identification of artifacts.

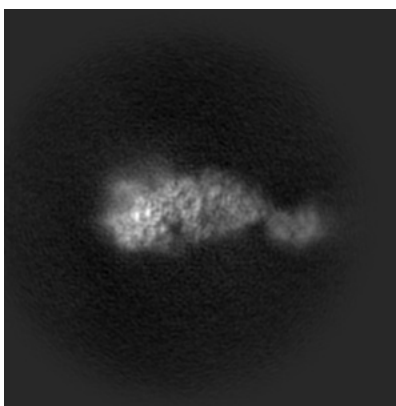
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

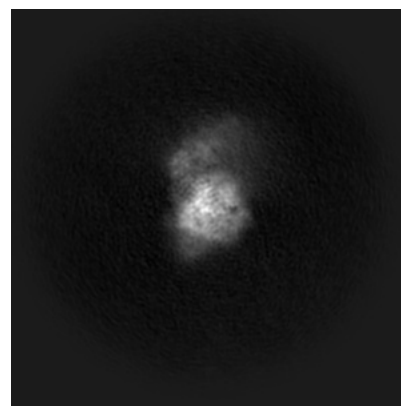
6.1.1 Primary map



X

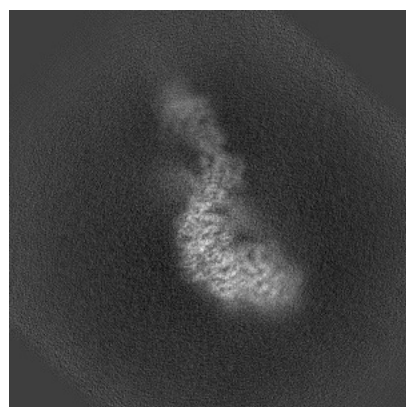


Y

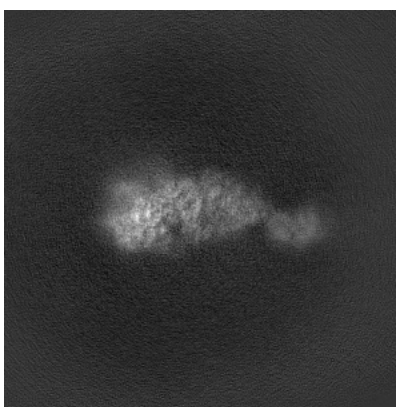


Z

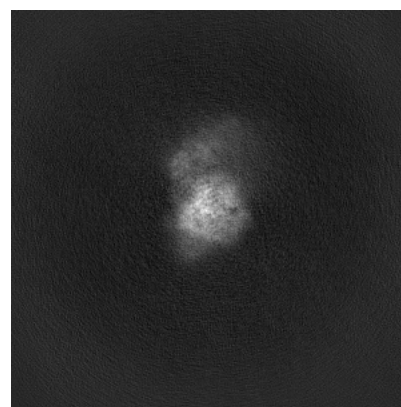
6.1.2 Raw map



X



Y

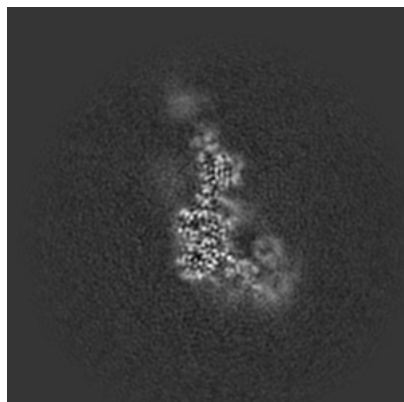


Z

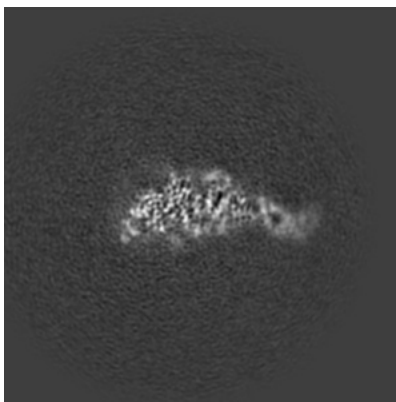
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

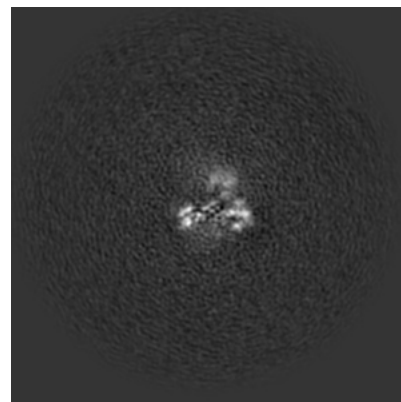
6.2.1 Primary map



X Index: 200

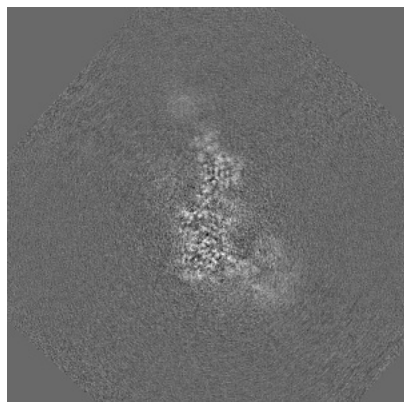


Y Index: 200

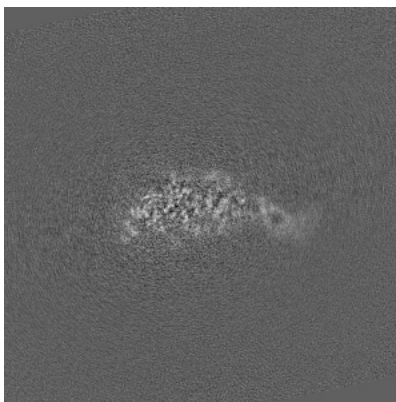


Z Index: 200

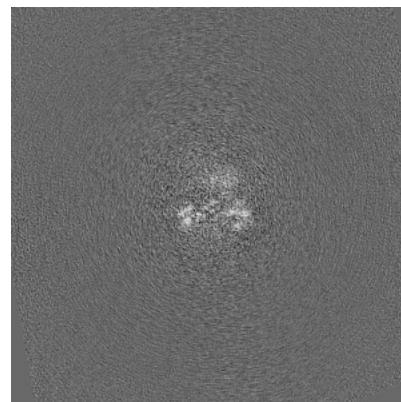
6.2.2 Raw map



X Index: 200



Y Index: 200

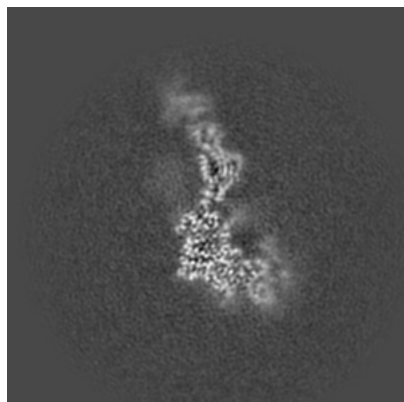


Z Index: 200

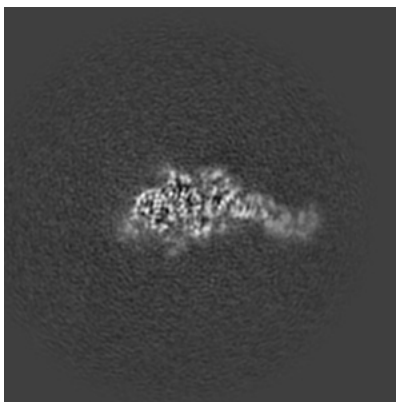
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

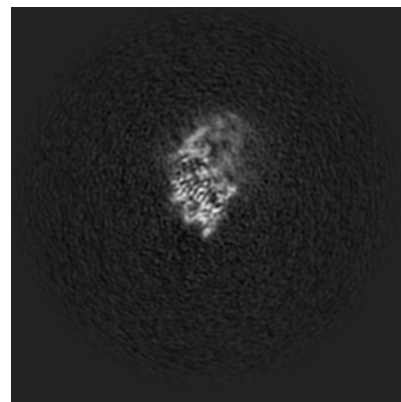
6.3.1 Primary map



X Index: 194

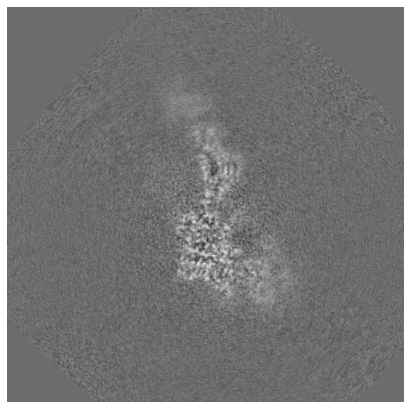


Y Index: 197

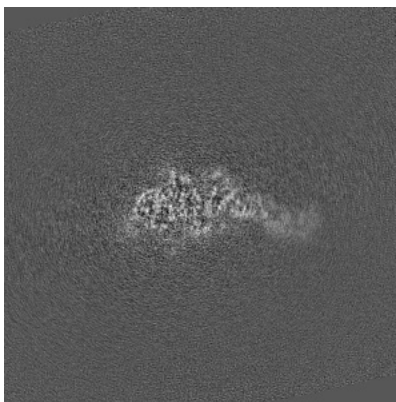


Z Index: 133

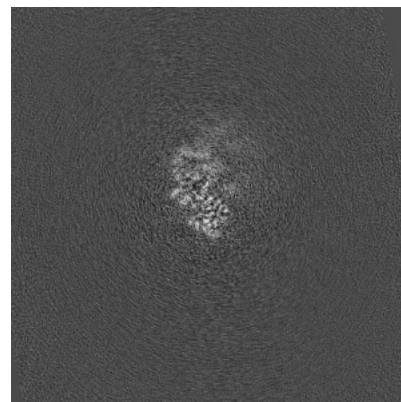
6.3.2 Raw map



X Index: 195



Y Index: 197

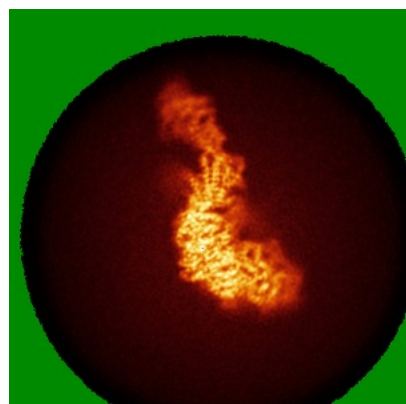


Z Index: 144

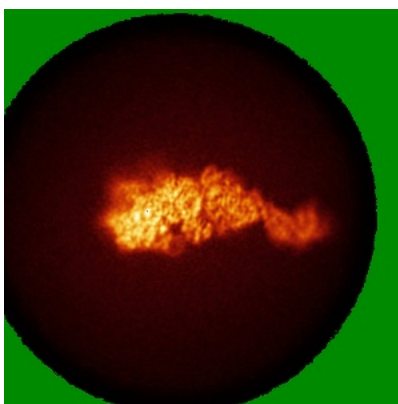
The images above show the largest variance slices of the map in three orthogonal directions.

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

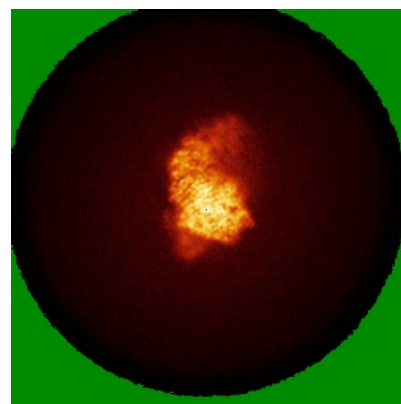
6.4.1 Primary map



X

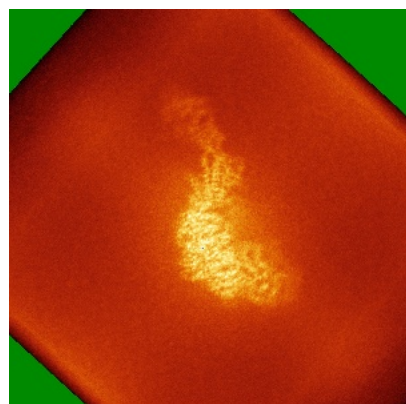


Y

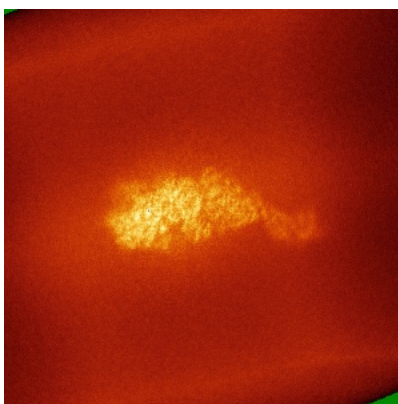


Z

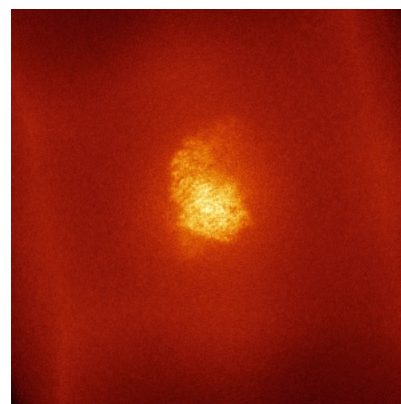
6.4.2 Raw map



X



Y

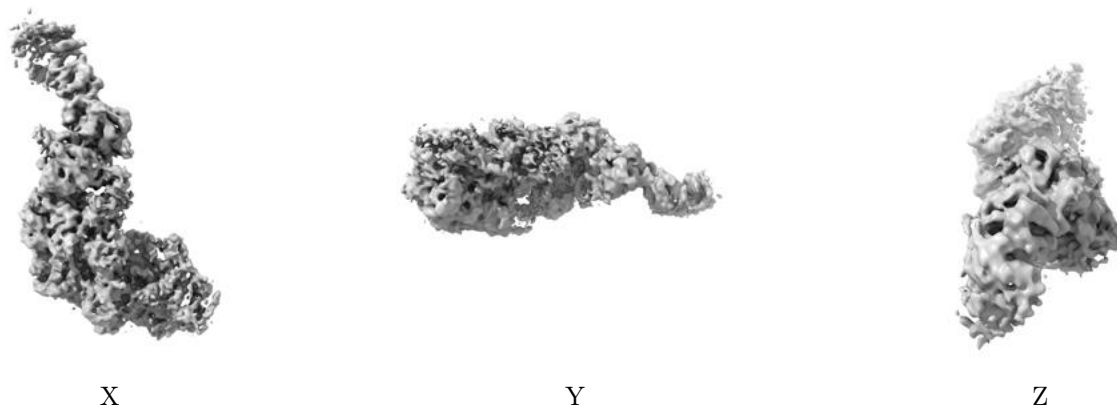


Z

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.

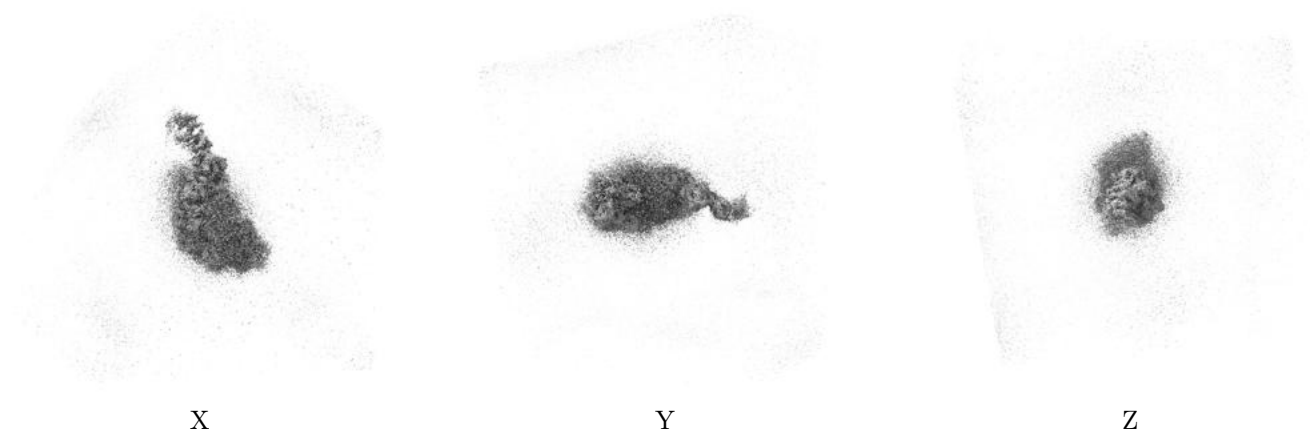
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.07. 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.

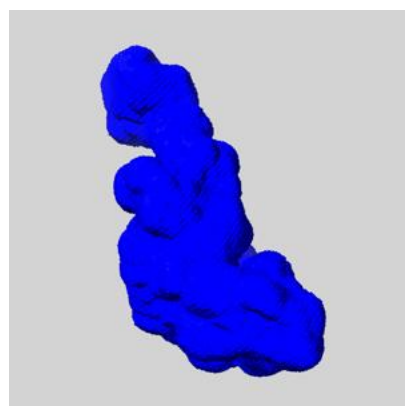
6.6 Mask visualisation [i](#)

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

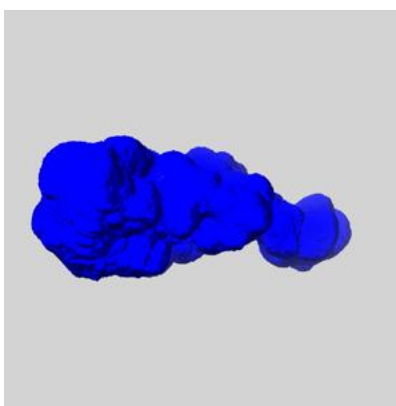
A mask typically either:

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

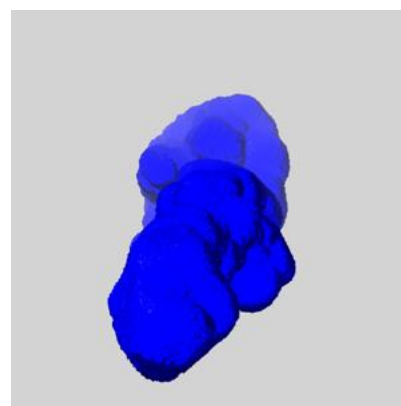
6.6.1 emd_72167_msk_1.map [i](#)



X



Y

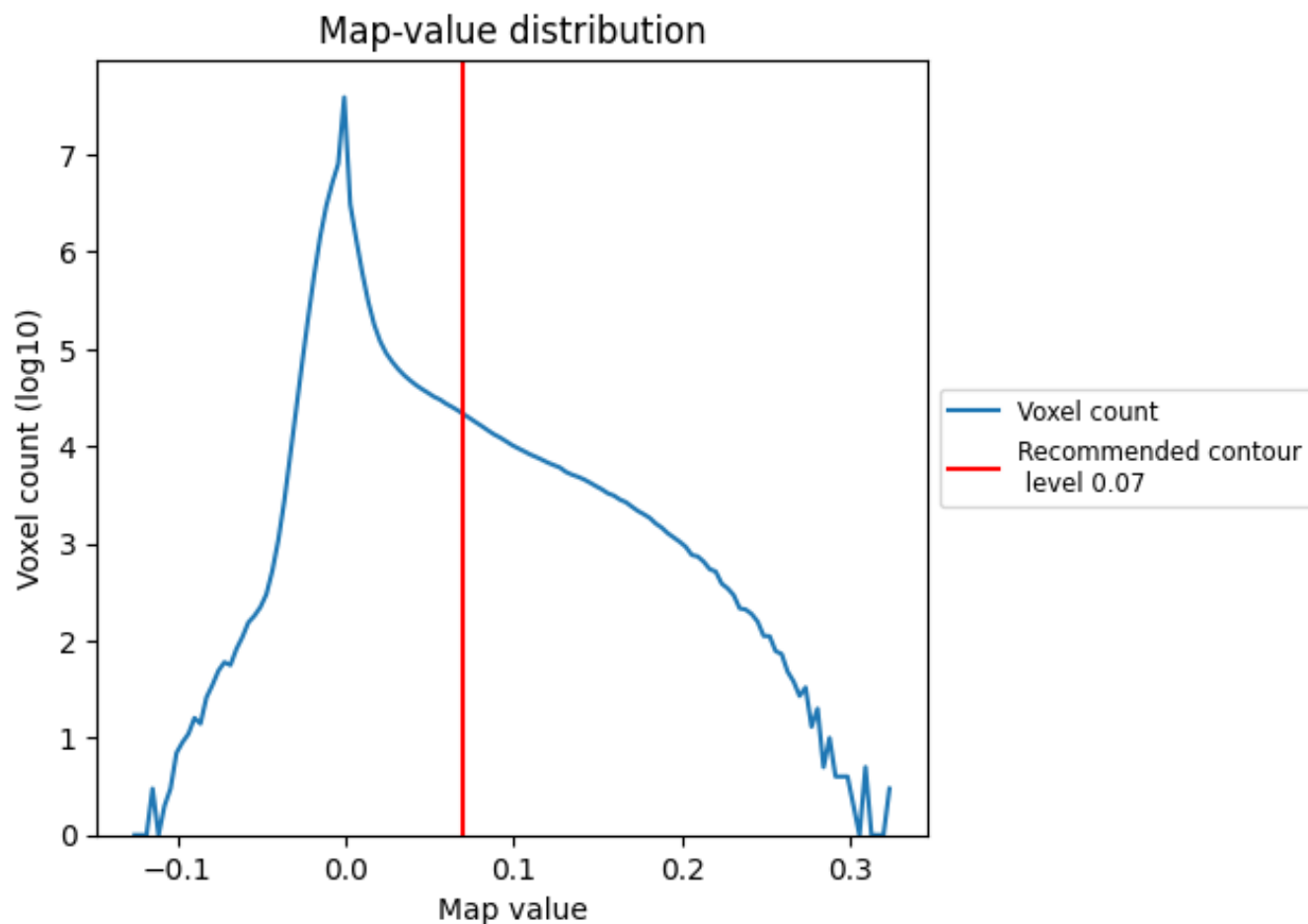


Z

7 Map analysis [i](#)

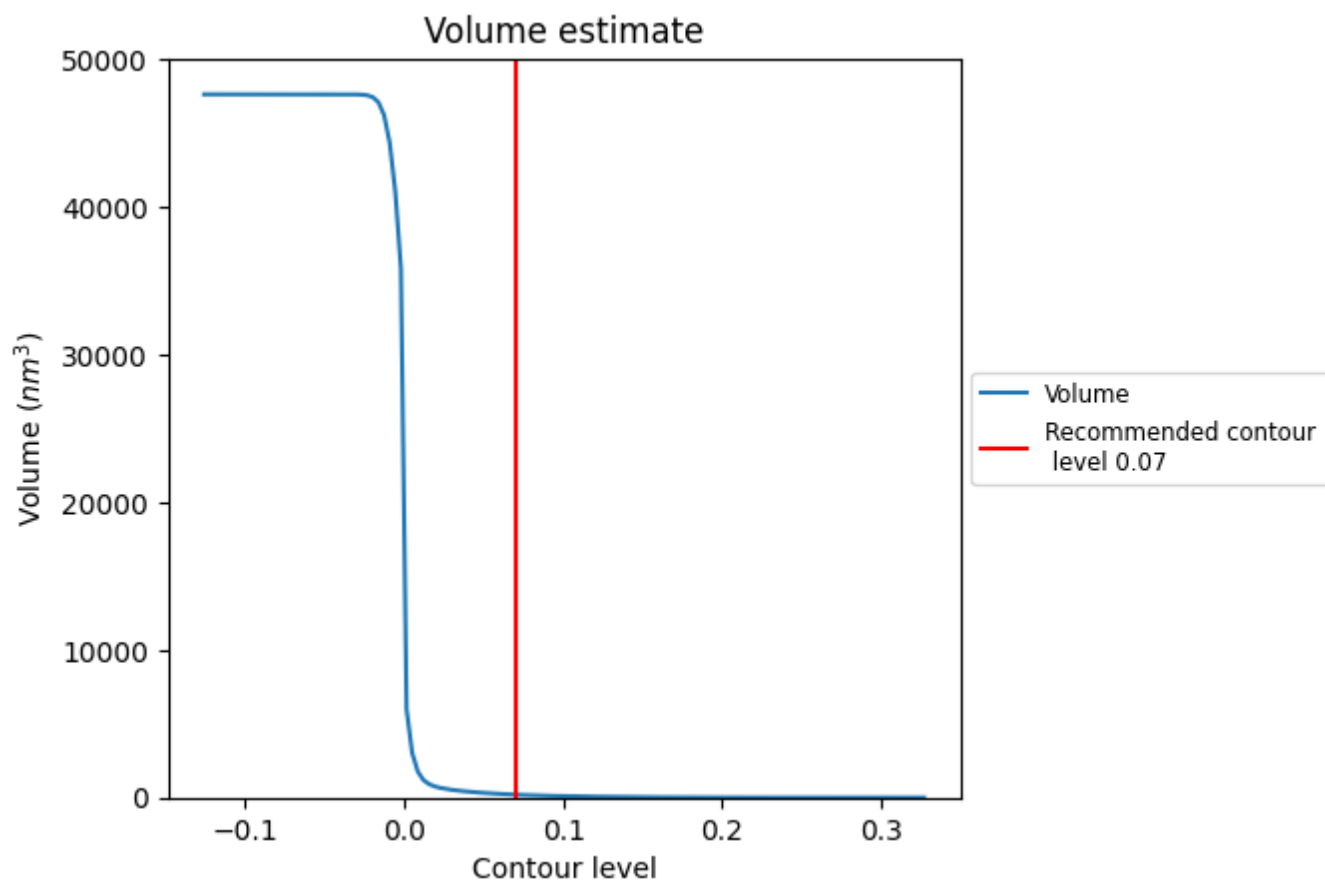
This section contains the results of statistical analysis of the map.

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.

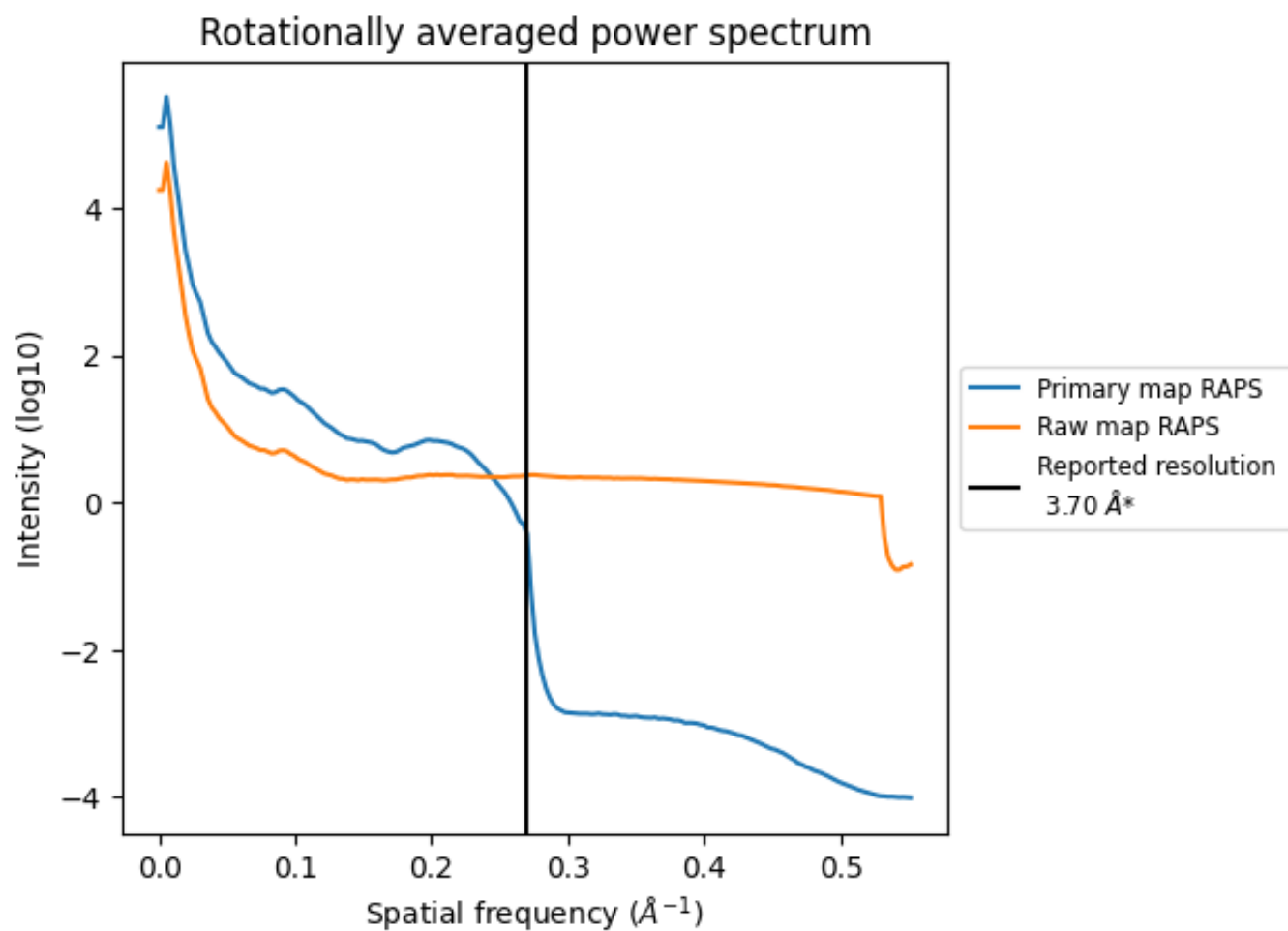
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 197 nm³; this corresponds to an approximate mass of 178 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 ⓘ

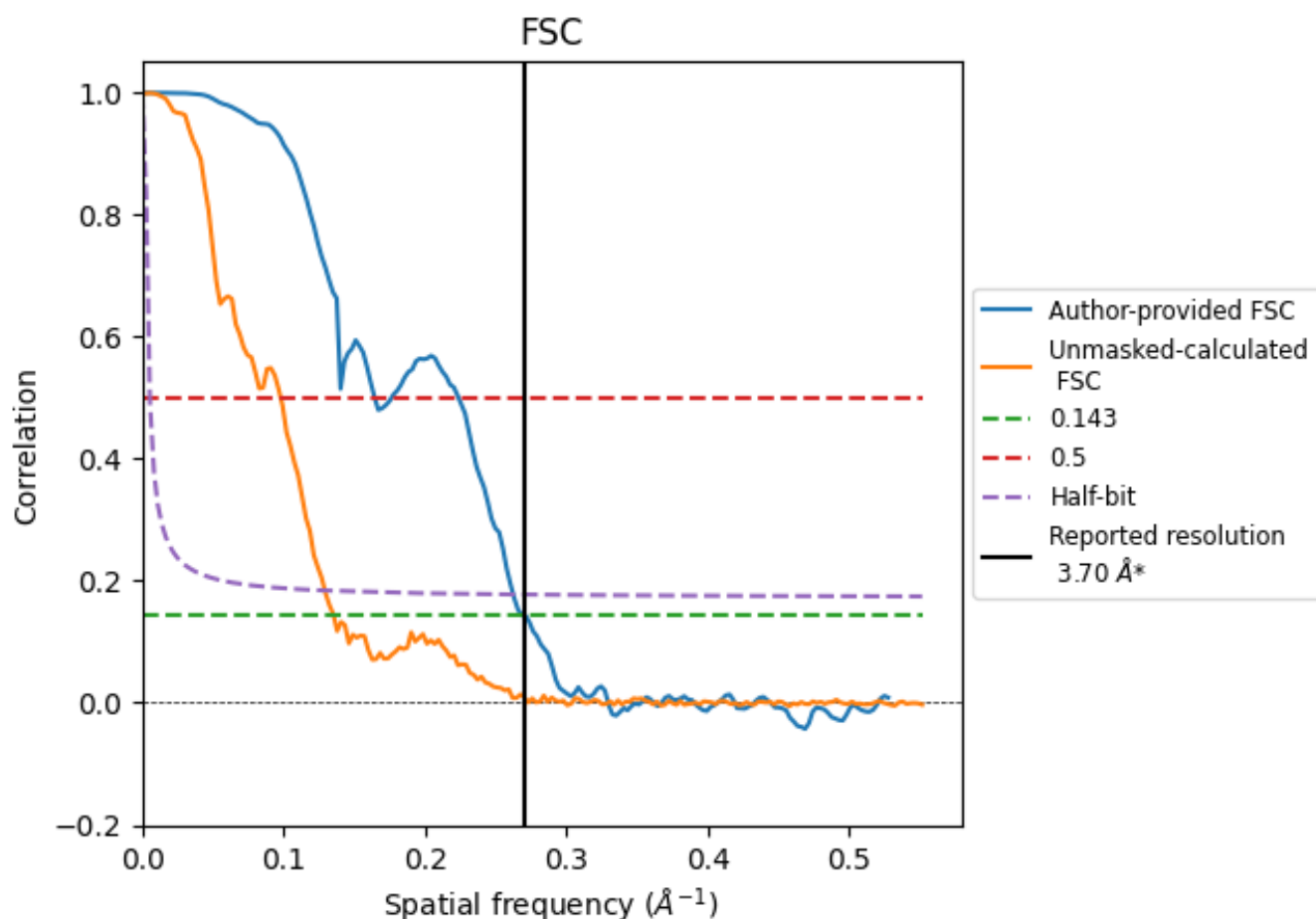


*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

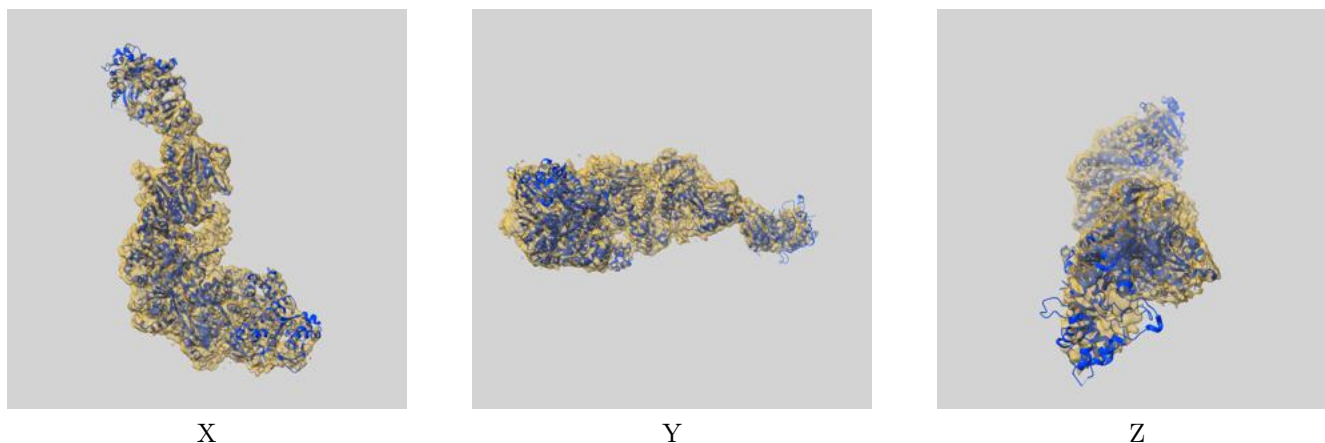
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	3.70	6.08	3.81
Unmasked-calculated*	7.39	10.21	7.74

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 7.39 differs from the reported value 3.7 by more than 10 %

9 Map-model fit [i](#)

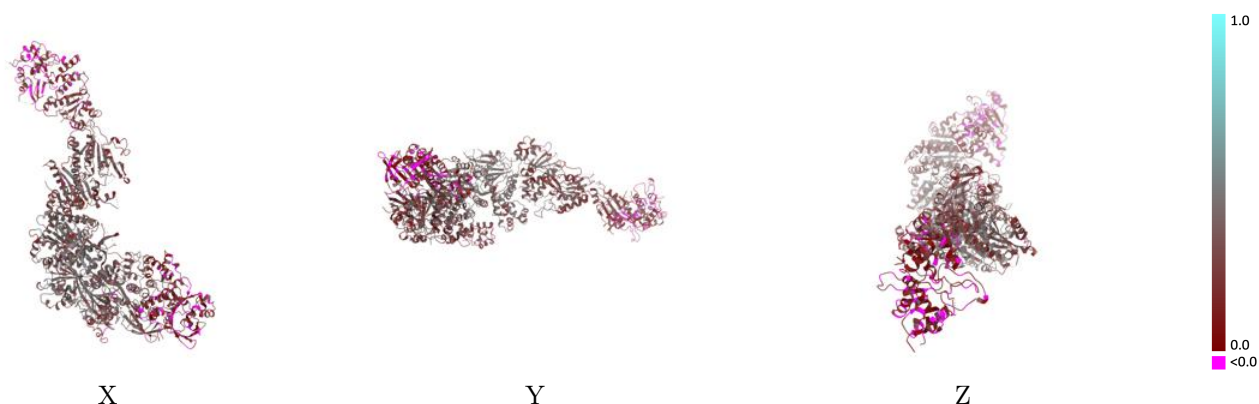
This section contains information regarding the fit between EMDB map EMD-72167 and PDB model 9Q2L. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



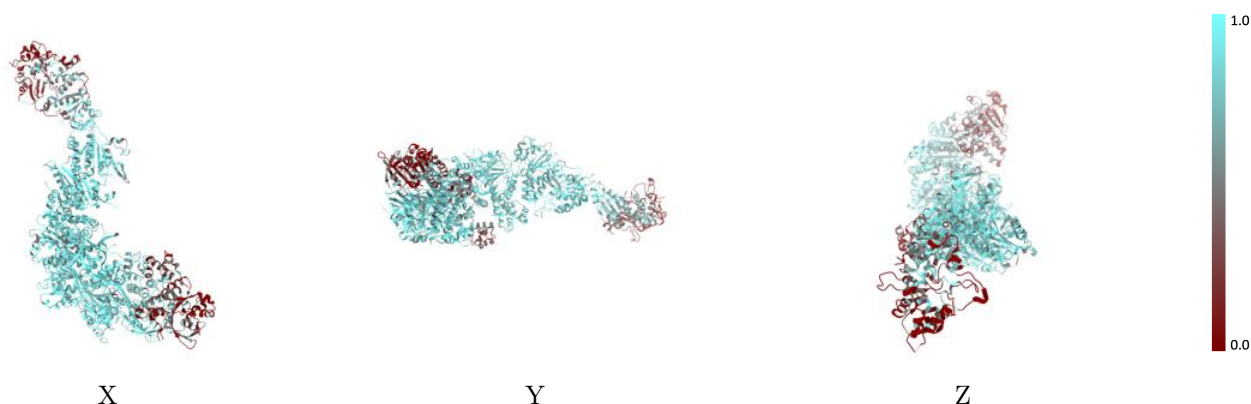
The images above show the 3D surface view of the map at the recommended contour level 0.07 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



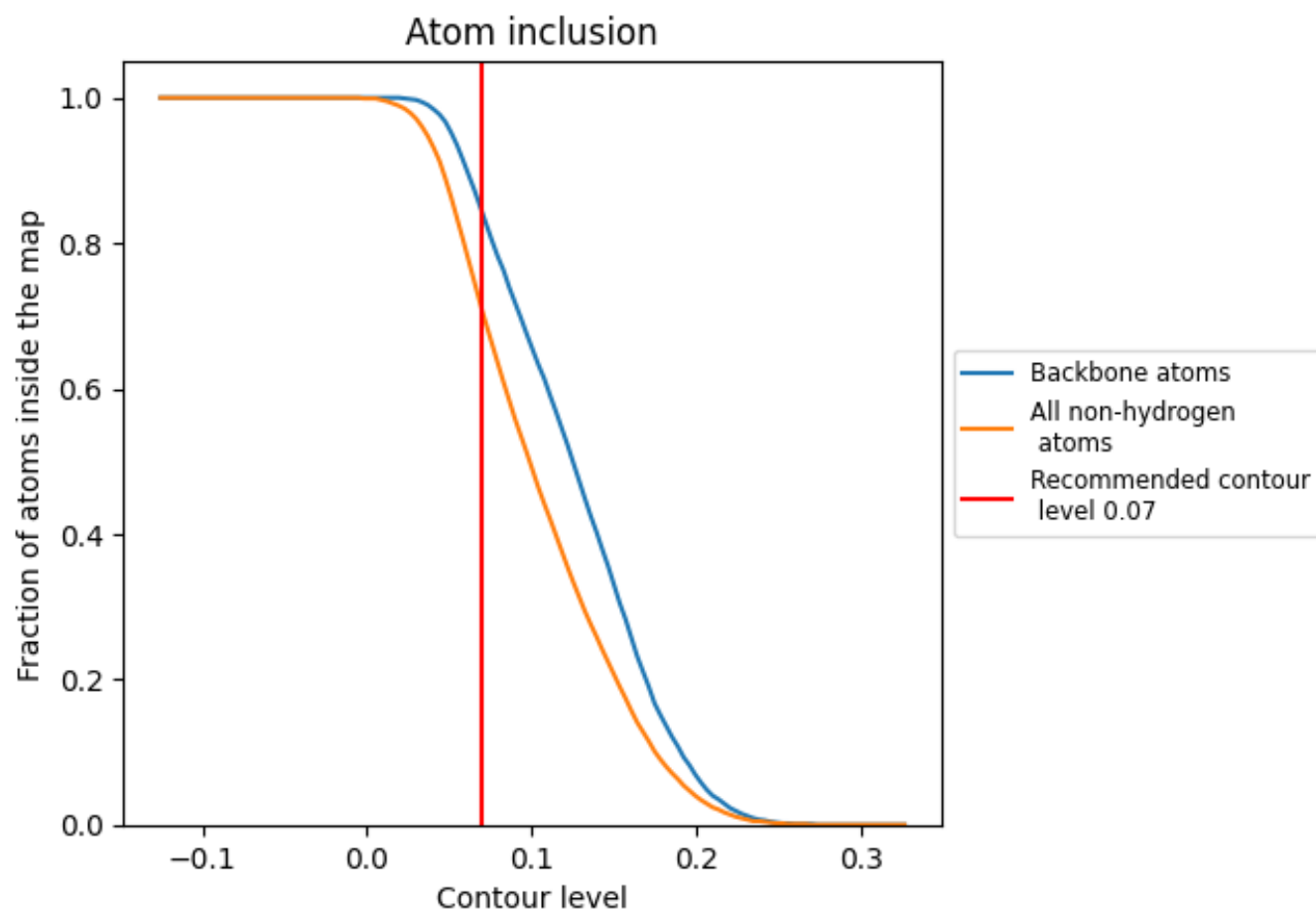
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.07).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 71% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.07) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7070	0.2760
A	0.8870	0.3810
B	0.8860	0.3620
C	0.8170	0.3180
D	0.7890	0.2990
E	0.6000	0.1830
F	0.2460	0.0880
G	0.7780	0.3140
H	0.8730	0.3680
I	0.8430	0.2920
J	0.3040	0.1230

1.0

0.0

<0.0