



# wwPDB EM Validation Summary Report ⓘ

Mar 6, 2026 – 12:00 PM UTC

PDB ID : 9DXF / pdb\_00009dx  
EMDB ID : EMD-47286  
Title : attLmm bound serine integrase and RDF complex in the post-rotation state  
Authors : Shin, H.; Rice, P.A.; Olorunniji, F.J.  
Deposited on : 2024-10-11  
Resolution : 4.16 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

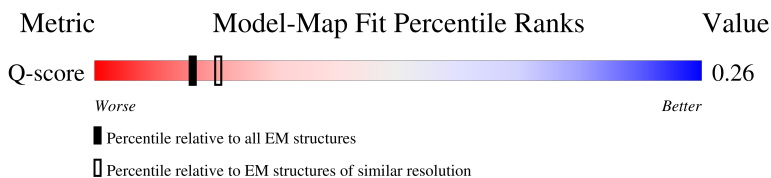
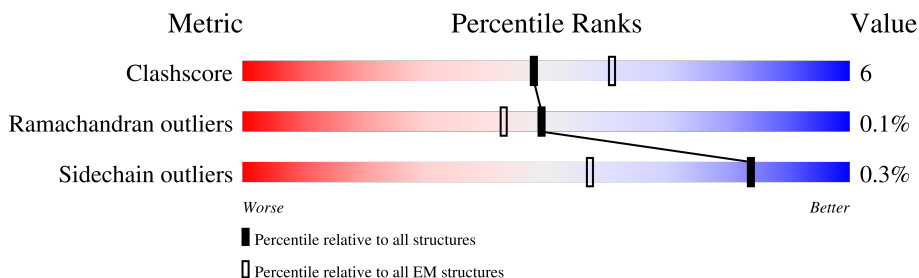
EMDB validation analysis : 0.0.1.dev132  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : 202505.v01 (Using data in the EMD archive up until May 2025)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

# 1 Overall quality at a glance i

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

The reported resolution of this entry is 4.16 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	5480 ( 3.66 - 4.66 )

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	620	<div> <div>7%</div> <div>62%</div> <div>23%</div> <div>14%</div> </div>
1	B	620	<div> <div>5%</div> <div>91%</div> </div>
1	C	620	<div> <div>14%</div> <div>66%</div> <div>19%</div> <div>14%</div> </div>
1	D	620	<div> <div>5%</div> <div>91%</div> </div>

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Mol	Chain	Length	Quality of chain
1	I	620	
1	J	620	
1	K	620	
1	L	620	
2	E	34	
2	M	34	
3	F	33	
3	N	33	
4	G	25	
4	O	25	
5	H	24	
5	P	24	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 46352 atoms, of which 22252 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 Resolvase homolog YokA,SPbeta prophage-derived uncharacterized protein YotN.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	534	Total	C	H	N	O	S	0	0
			8764	2736	4418	752	840	18		
1	B	56	Total	C	H	N	O		0	0
			955	306	474	81	94			
1	C	535	Total	C	H	N	O	S	0	0
			8783	2741	4429	753	841	19		
1	D	57	Total	C	H	N	O		0	0
			970	311	480	82	97			
1	I	535	Total	C	H	N	O	S	0	0
			8783	2741	4429	753	841	19		
1	J	56	Total	C	H	N	O		0	0
			955	306	474	81	94			
1	K	534	Total	C	H	N	O	S	0	0
			8764	2736	4418	752	840	18		
1	L	57	Total	C	H	N	O		0	0
			970	311	480	82	97			

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	546	THR	-	linker	UNP O32006
A	547	SER	-	linker	UNP O32006
A	548	GLY	-	linker	UNP O32006
A	549	SER	-	linker	UNP O32006
A	550	GLY	-	linker	UNP O32006
A	551	GLY	-	linker	UNP O32006
A	552	SER	-	linker	UNP O32006
A	553	GLY	-	linker	UNP O32006
A	554	GLY	-	linker	UNP O32006
A	555	SER	-	linker	UNP O32006
A	556	GLY	-	linker	UNP O32006
A	557	GLY	-	linker	UNP O32006
A	558	SER	-	linker	UNP O32006

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Chain	Residue	Modelled	Actual	Comment	Reference
A	559	GLY	-	linker	UNP O32006
A	560	ARG	-	linker	UNP O32006
A	561	SER	-	linker	UNP O32006
A	562	GLY	-	linker	UNP O32006
A	563	THR	-	linker	UNP O32006
B	-16	THR	-	linker	UNP O32006
B	-15	SER	-	linker	UNP O32006
B	-14	GLY	-	linker	UNP O32006
B	-13	SER	-	linker	UNP O32006
B	-12	GLY	-	linker	UNP O32006
B	-11	GLY	-	linker	UNP O32006
B	-10	SER	-	linker	UNP O32006
B	-9	GLY	-	linker	UNP O32006
B	-8	GLY	-	linker	UNP O32006
B	-7	SER	-	linker	UNP O32006
B	-6	GLY	-	linker	UNP O32006
B	-5	GLY	-	linker	UNP O32006
B	-4	SER	-	linker	UNP O32006
B	-3	GLY	-	linker	UNP O32006
B	-2	ARG	-	linker	UNP O32006
B	-1	SER	-	linker	UNP O32006
B	0	GLY	-	linker	UNP O32006
B	1	THR	-	linker	UNP O32006
C	546	THR	-	linker	UNP O32006
C	547	SER	-	linker	UNP O32006
C	548	GLY	-	linker	UNP O32006
C	549	SER	-	linker	UNP O32006
C	550	GLY	-	linker	UNP O32006
C	551	GLY	-	linker	UNP O32006
C	552	SER	-	linker	UNP O32006
C	553	GLY	-	linker	UNP O32006
C	554	GLY	-	linker	UNP O32006
C	555	SER	-	linker	UNP O32006
C	556	GLY	-	linker	UNP O32006
C	557	GLY	-	linker	UNP O32006
C	558	SER	-	linker	UNP O32006
C	559	GLY	-	linker	UNP O32006
C	560	ARG	-	linker	UNP O32006
C	561	SER	-	linker	UNP O32006
C	562	GLY	-	linker	UNP O32006
C	563	THR	-	linker	UNP O32006
D	-16	THR	-	linker	UNP O32006

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-15	SER	-	linker	UNP O32006
D	-14	GLY	-	linker	UNP O32006
D	-13	SER	-	linker	UNP O32006
D	-12	GLY	-	linker	UNP O32006
D	-11	GLY	-	linker	UNP O32006
D	-10	SER	-	linker	UNP O32006
D	-9	GLY	-	linker	UNP O32006
D	-8	GLY	-	linker	UNP O32006
D	-7	SER	-	linker	UNP O32006
D	-6	GLY	-	linker	UNP O32006
D	-5	GLY	-	linker	UNP O32006
D	-4	SER	-	linker	UNP O32006
D	-3	GLY	-	linker	UNP O32006
D	-2	ARG	-	linker	UNP O32006
D	-1	SER	-	linker	UNP O32006
D	0	GLY	-	linker	UNP O32006
D	1	THR	-	linker	UNP O32006
I	546	THR	-	linker	UNP O32006
I	547	SER	-	linker	UNP O32006
I	548	GLY	-	linker	UNP O32006
I	549	SER	-	linker	UNP O32006
I	550	GLY	-	linker	UNP O32006
I	551	GLY	-	linker	UNP O32006
I	552	SER	-	linker	UNP O32006
I	553	GLY	-	linker	UNP O32006
I	554	GLY	-	linker	UNP O32006
I	555	SER	-	linker	UNP O32006
I	556	GLY	-	linker	UNP O32006
I	557	GLY	-	linker	UNP O32006
I	558	SER	-	linker	UNP O32006
I	559	GLY	-	linker	UNP O32006
I	560	ARG	-	linker	UNP O32006
I	561	SER	-	linker	UNP O32006
I	562	GLY	-	linker	UNP O32006
I	563	THR	-	linker	UNP O32006
J	-16	THR	-	linker	UNP O32006
J	-15	SER	-	linker	UNP O32006
J	-14	GLY	-	linker	UNP O32006
J	-13	SER	-	linker	UNP O32006
J	-12	GLY	-	linker	UNP O32006
J	-11	GLY	-	linker	UNP O32006
J	-10	SER	-	linker	UNP O32006

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Chain	Residue	Modelled	Actual	Comment	Reference
J	-9	GLY	-	linker	UNP O32006
J	-8	GLY	-	linker	UNP O32006
J	-7	SER	-	linker	UNP O32006
J	-6	GLY	-	linker	UNP O32006
J	-5	GLY	-	linker	UNP O32006
J	-4	SER	-	linker	UNP O32006
J	-3	GLY	-	linker	UNP O32006
J	-2	ARG	-	linker	UNP O32006
J	-1	SER	-	linker	UNP O32006
J	0	GLY	-	linker	UNP O32006
J	1	THR	-	linker	UNP O32006
K	546	THR	-	linker	UNP O32006
K	547	SER	-	linker	UNP O32006
K	548	GLY	-	linker	UNP O32006
K	549	SER	-	linker	UNP O32006
K	550	GLY	-	linker	UNP O32006
K	551	GLY	-	linker	UNP O32006
K	552	SER	-	linker	UNP O32006
K	553	GLY	-	linker	UNP O32006
K	554	GLY	-	linker	UNP O32006
K	555	SER	-	linker	UNP O32006
K	556	GLY	-	linker	UNP O32006
K	557	GLY	-	linker	UNP O32006
K	558	SER	-	linker	UNP O32006
K	559	GLY	-	linker	UNP O32006
K	560	ARG	-	linker	UNP O32006
K	561	SER	-	linker	UNP O32006
K	562	GLY	-	linker	UNP O32006
K	563	THR	-	linker	UNP O32006
L	-16	THR	-	linker	UNP O32006
L	-15	SER	-	linker	UNP O32006
L	-14	GLY	-	linker	UNP O32006
L	-13	SER	-	linker	UNP O32006
L	-12	GLY	-	linker	UNP O32006
L	-11	GLY	-	linker	UNP O32006
L	-10	SER	-	linker	UNP O32006
L	-9	GLY	-	linker	UNP O32006
L	-8	GLY	-	linker	UNP O32006
L	-7	SER	-	linker	UNP O32006
L	-6	GLY	-	linker	UNP O32006
L	-5	GLY	-	linker	UNP O32006
L	-4	SER	-	linker	UNP O32006

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Chain	Residue	Modelled	Actual	Comment	Reference
L	-3	GLY	-	linker	UNP O32006
L	-2	ARG	-	linker	UNP O32006
L	-1	SER	-	linker	UNP O32006
L	0	GLY	-	linker	UNP O32006
L	1	THR	-	linker	UNP O32006

- Molecule 2 is a DNA chain called DNA (34-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
2	E	34	Total	C	H	N	O	P	0	0
			1087	337	393	110	213	34		
2	M	34	Total	C	H	N	O	P	0	0
			1087	337	393	110	213	34		

- Molecule 3 is a DNA chain called DNA (33-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
3	F	33	Total	C	H	N	O	P	0	0
			1052	326	373	130	190	33		
3	N	33	Total	C	H	N	O	P	0	0
			1052	326	373	130	190	33		

- Molecule 4 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
4	G	25	Total	C	H	N	O	P	0	0
			796	246	285	90	150	25		
4	O	25	Total	C	H	N	O	P	0	0
			796	246	285	90	150	25		

- Molecule 5 is a DNA chain called DNA (24-MER).

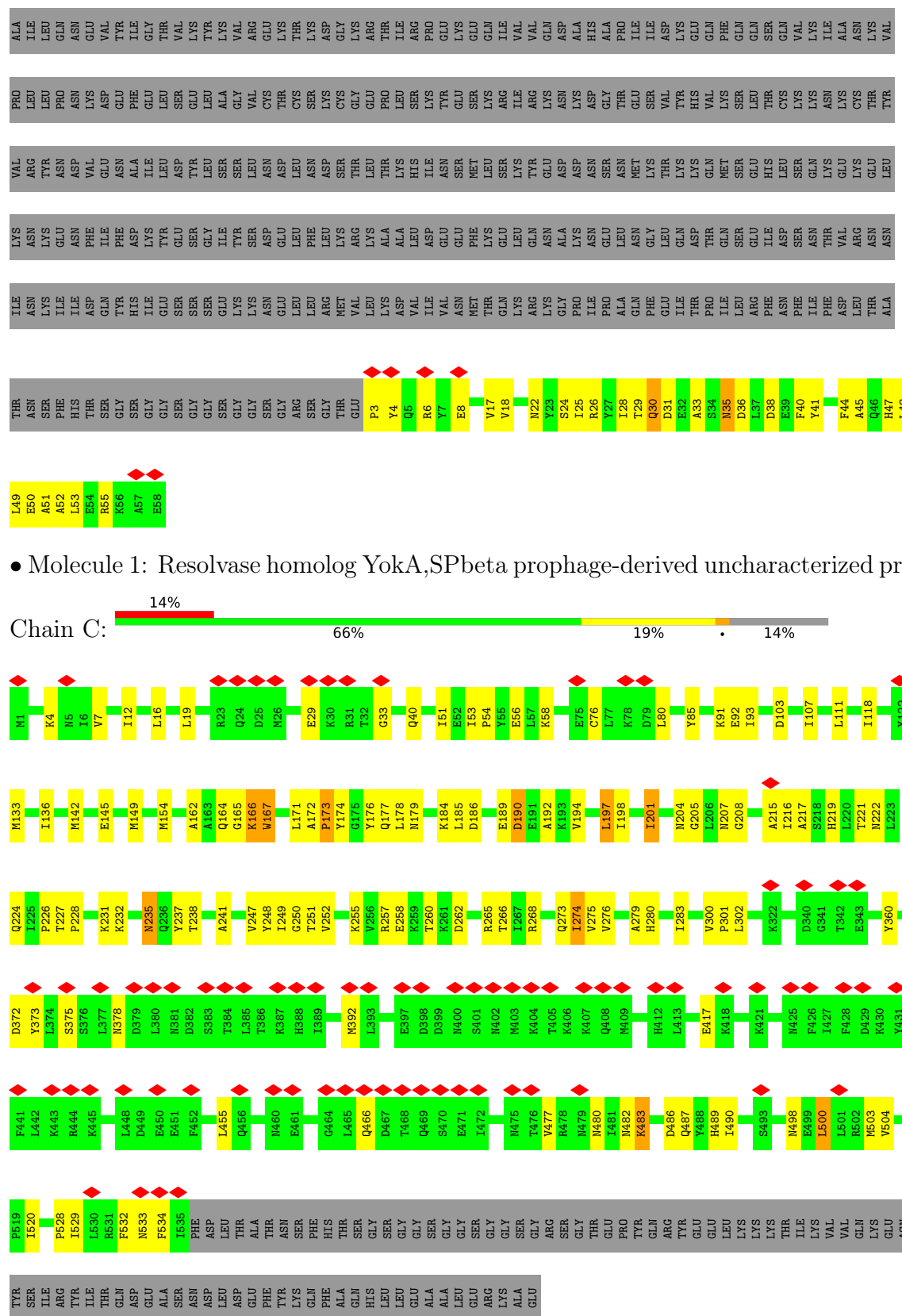
Mol	Chain	Residues	Atoms						AltConf	Trace
5	H	24	Total	C	H	N	O	P	0	0
			767	237	274	87	145	24		
5	P	24	Total	C	H	N	O	P	0	0
			767	237	274	87	145	24		

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).



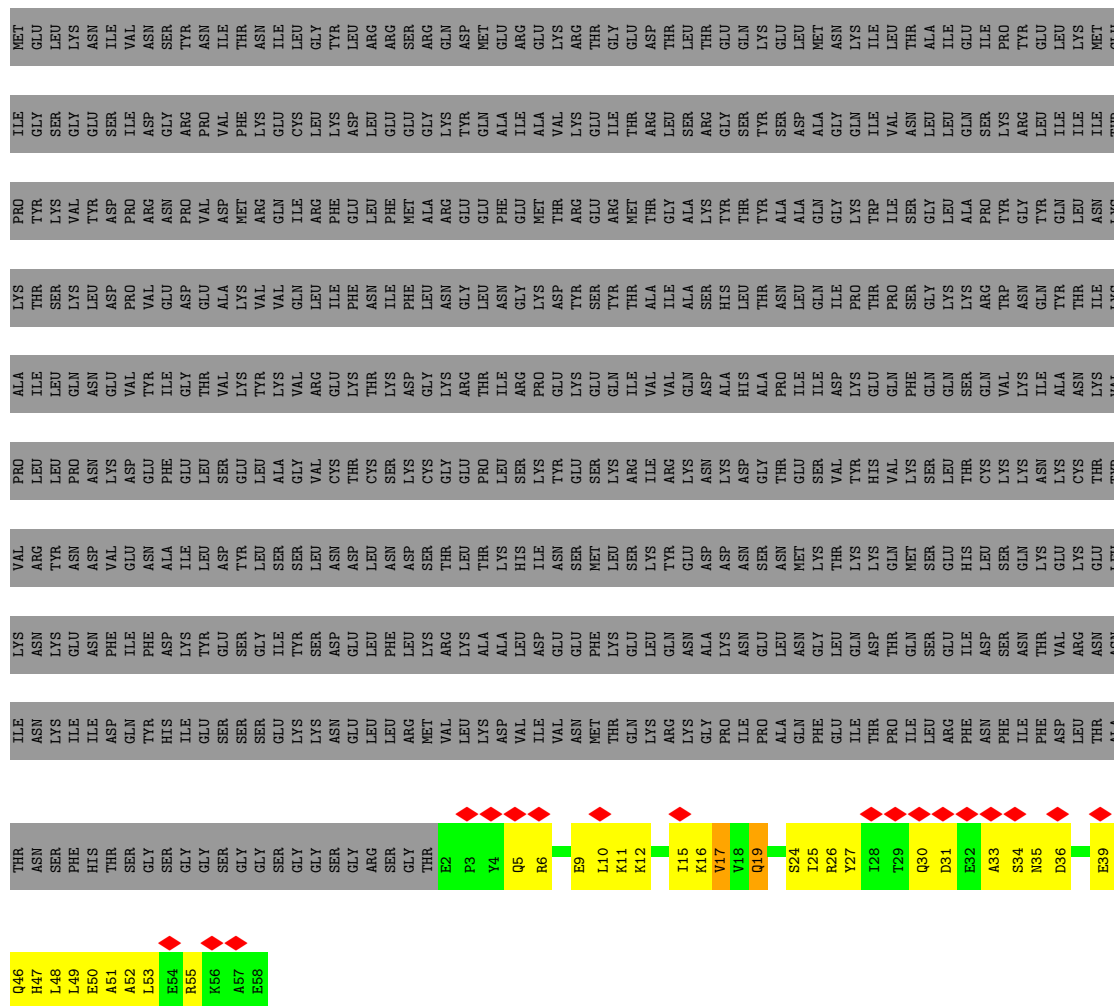
Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total 1	Zn 1	0
6	C	1	Total 1	Zn 1	0
6	I	1	Total 1	Zn 1	0
6	K	1	Total 1	Zn 1	0





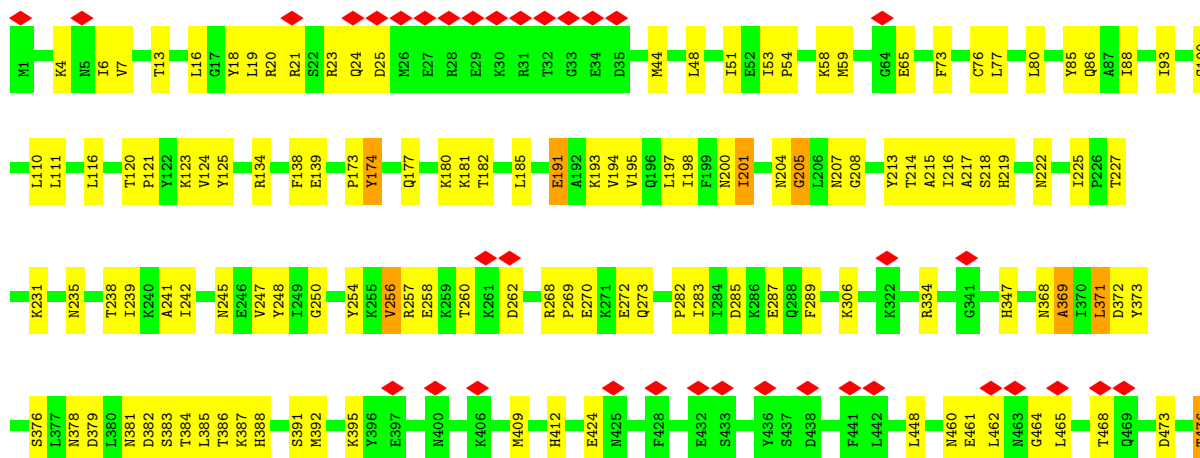
- Molecule 1: Resolvase homolog YokA,SPbeta prophage-derived uncharacterized protein YotN

Chain D:  5% 91%

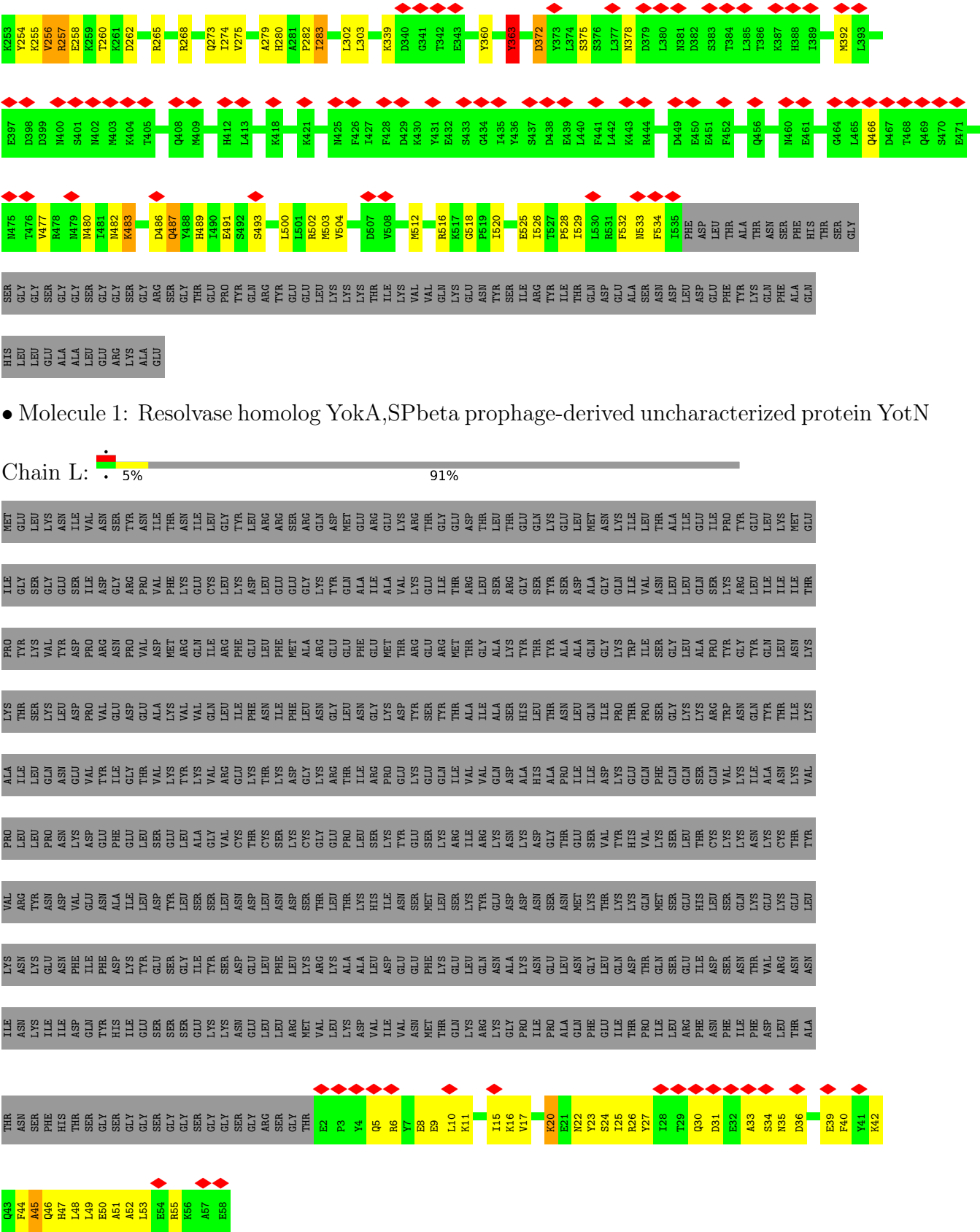



● Molecule 1: Resolvase homolog YokA,SPbeta prophage-derived uncharacterized protein YotN

Chain I:  6% 63% 22% 14%








Chain E:  82% 15%



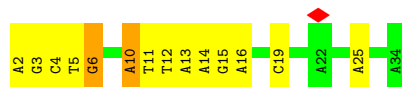
- Molecule 2: DNA (34-MER)

Chain M:  82% 12% 6%



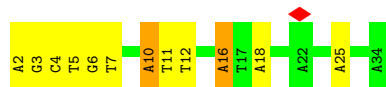
- Molecule 3: DNA (33-MER)

Chain F:  58% 36% 6%



- Molecule 3: DNA (33-MER)

Chain N:  64% 30% 6%



- Molecule 4: DNA (25-MER)

Chain G:  68% 32%



- Molecule 4: DNA (25-MER)

Chain O:  68% 28%




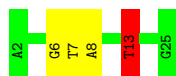
- Molecule 5: DNA (24-MER)

Chain H:  67% 33%



- Molecule 5: DNA (24-MER)

Chain P:  83% 12% .





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	210561	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	65	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.406	Depositor
Minimum map value	-0.006	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.0226	Depositor
Map size ( $\text{\AA}$ )	451.56003, 468.60004, 374.88	wwPDB
Map dimensions	424, 440, 352	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.065, 1.065, 1.065	Depositor

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.47	31/4412 (0.7%)	1.48	85/5927 (1.4%)
1	B	2.44	9/489 (1.8%)	2.55	29/654 (4.4%)
1	C	1.34	23/4420 (0.5%)	1.46	96/5937 (1.6%)
1	D	2.54	13/498 (2.6%)	2.67	57/667 (8.5%)
1	I	1.37	22/4420 (0.5%)	1.44	89/5937 (1.5%)
1	J	2.46	9/489 (1.8%)	2.63	37/654 (5.7%)
1	K	1.36	26/4412 (0.6%)	1.48	94/5927 (1.6%)
1	L	2.56	13/498 (2.6%)	2.69	64/667 (9.6%)
2	E	0.79	0/774	1.24	0/1193
2	M	0.84	0/774	1.23	1/1193 (0.1%)
3	F	1.36	3/764 (0.4%)	1.48	2/1176 (0.2%)
3	N	1.28	2/764 (0.3%)	1.42	2/1176 (0.2%)
4	G	0.84	0/572	1.24	0/880
4	O	0.83	0/572	1.24	0/880
5	H	1.30	0/552	1.45	1/850 (0.1%)
5	P	1.22	0/552	1.42	1/850 (0.1%)
All	All	1.45	151/24962 (0.6%)	1.56	558/34568 (1.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	I	0	3
1	K	0	2
2	E	0	3
2	M	0	3
3	F	0	2
3	N	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
4	G	0	4
4	O	0	3
5	H	0	3
5	P	0	1
All	All	0	27

The worst 5 of 151 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	HIS	CE1-NE2	-9.01	1.23	1.32
1	A	280	HIS	CE1-NE2	-8.97	1.23	1.32
1	B	47	HIS	CE1-NE2	-8.97	1.23	1.32
1	C	280	HIS	CE1-NE2	-8.94	1.23	1.32
1	I	219	HIS	CE1-NE2	-8.92	1.23	1.32

The worst 5 of 558 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	24	SER	CA-C-N	9.27	130.36	121.65
1	J	24	SER	C-N-CA	9.27	130.36	121.65
1	B	24	SER	CA-C-N	8.86	129.98	121.65
1	B	24	SER	C-N-CA	8.86	129.98	121.65
1	K	512	MET	CA-C-N	8.41	131.37	120.44

There are no chirality outliers.

5 of 27 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	176	TYR	Sidechain
2	E	-13	DT	Sidechain
2	E	-16	DT	Sidechain
2	E	-23	DG	Sidechain
3	F	15	DG	Sidechain

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4346	4418	4412	68	0
1	B	481	474	474	9	0
1	C	4354	4429	4424	48	0
1	D	490	480	479	3	0
1	I	4354	4429	4424	73	0
1	J	481	474	474	8	0
1	K	4346	4418	4412	45	0
1	L	490	480	479	3	0
2	E	694	393	388	7	0
2	M	694	393	386	4	0
3	F	679	373	346	10	0
3	N	679	373	348	10	0
4	G	511	285	279	4	0
4	O	511	285	279	6	0
5	H	493	274	257	5	0
5	P	493	274	258	5	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
6	I	1	0	0	0	0
6	K	1	0	0	0	0
All	All	24100	22252	22119	272	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 6.

The worst 5 of 272 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:TYR:O	1:A:248:TYR:OH	1.82	0.97
1:A:80:LEU:HD11	1:A:88:ILE:HD11	1.42	0.97
1:A:120:THR:OG1	1:A:123:LYS:O	1.90	0.90
1:I:473:ASP:OD1	1:I:476:THR:OG1	1.90	0.90
1:I:174:TYR:O	1:I:248:TYR:OH	1.89	0.90

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	532/620 (86%)	494 (93%)	37 (7%)	1 (0%)	43	76
1	B	54/620 (9%)	48 (89%)	6 (11%)	0	100	100
1	C	533/620 (86%)	496 (93%)	37 (7%)	0	100	100
1	D	55/620 (9%)	53 (96%)	2 (4%)	0	100	100
1	I	533/620 (86%)	503 (94%)	29 (5%)	1 (0%)	43	76
1	J	54/620 (9%)	49 (91%)	5 (9%)	0	100	100
1	K	532/620 (86%)	501 (94%)	31 (6%)	0	100	100
1	L	55/620 (9%)	53 (96%)	2 (4%)	0	100	100
All	All	2348/4960 (47%)	2197 (94%)	149 (6%)	2 (0%)	49	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	I	24	GLN

### 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	489/560 (87%)	489 (100%)	0	100	100
1	B	51/560 (9%)	51 (100%)	0	100	100
1	C	490/560 (88%)	488 (100%)	2 (0%)	84	82

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	52/560 (9%)	52 (100%)	0	100	100
1	I	490/560 (88%)	490 (100%)	0	100	100
1	J	51/560 (9%)	51 (100%)	0	100	100
1	K	489/560 (87%)	485 (99%)	4 (1%)	73	77
1	L	52/560 (9%)	52 (100%)	0	100	100
All	All	2164/4480 (48%)	2158 (100%)	6 (0%)	84	83

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	302	LEU
1	K	363	TYR
1	K	493	SER
1	C	302	LEU
1	C	249	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	293	GLN
1	J	35	ASN
1	I	390	ASN
1	K	40	GLN
1	C	207	ASN

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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.

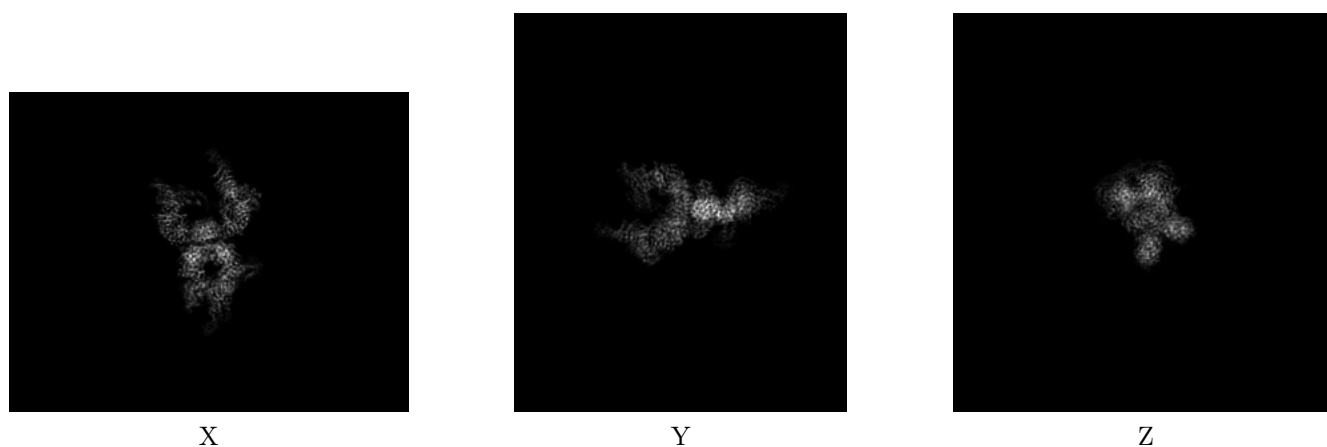
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-47286. 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](#)

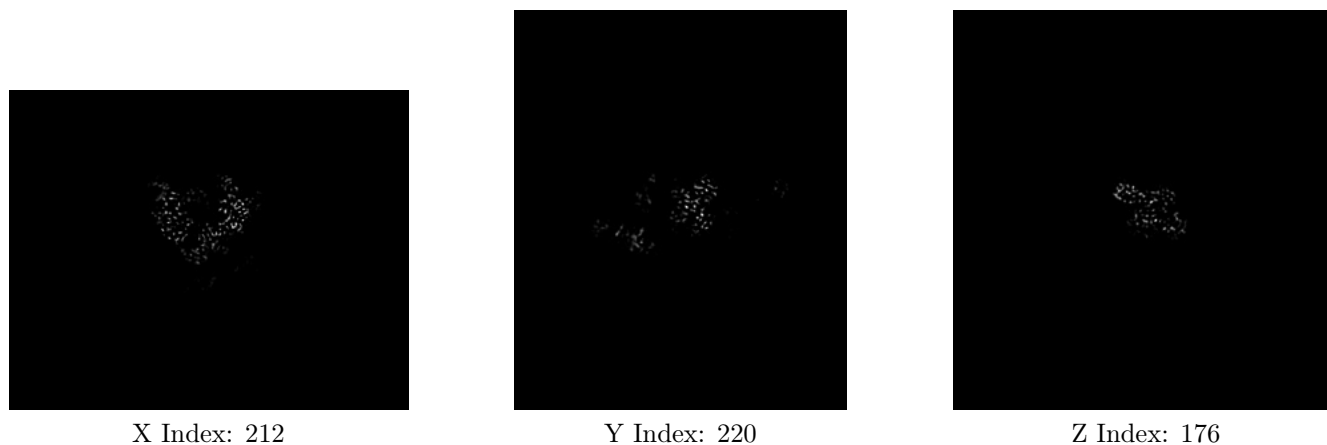
#### 6.1.1 Primary map



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

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

#### 6.2.1 Primary map

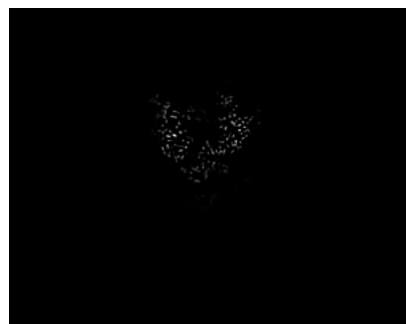




The images above show central slices of the map in three orthogonal directions.

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

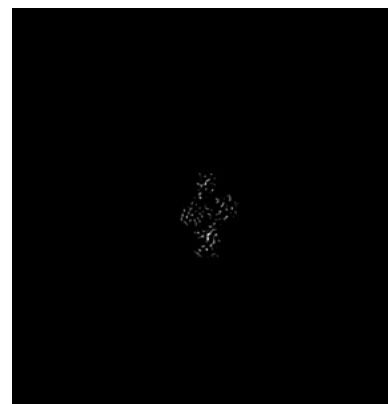
### 6.3.1 Primary map



X Index: 213



Y Index: 238

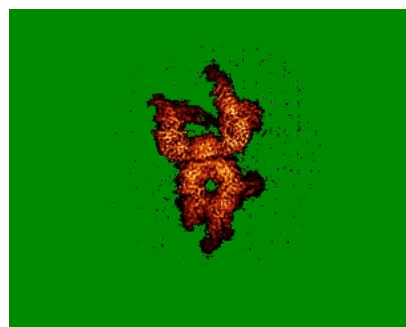


Z Index: 200

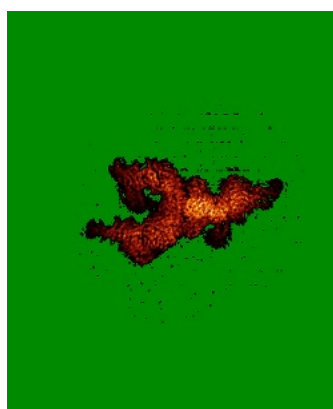
The images above show the largest variance slices of the map in three orthogonal directions.

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

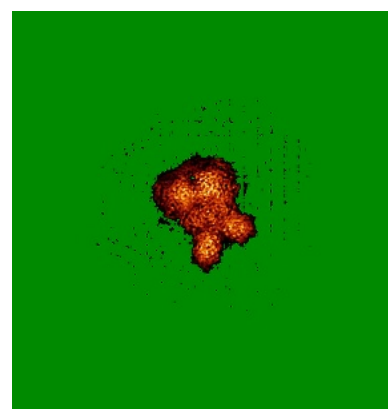
### 6.4.1 Primary map



X



Y

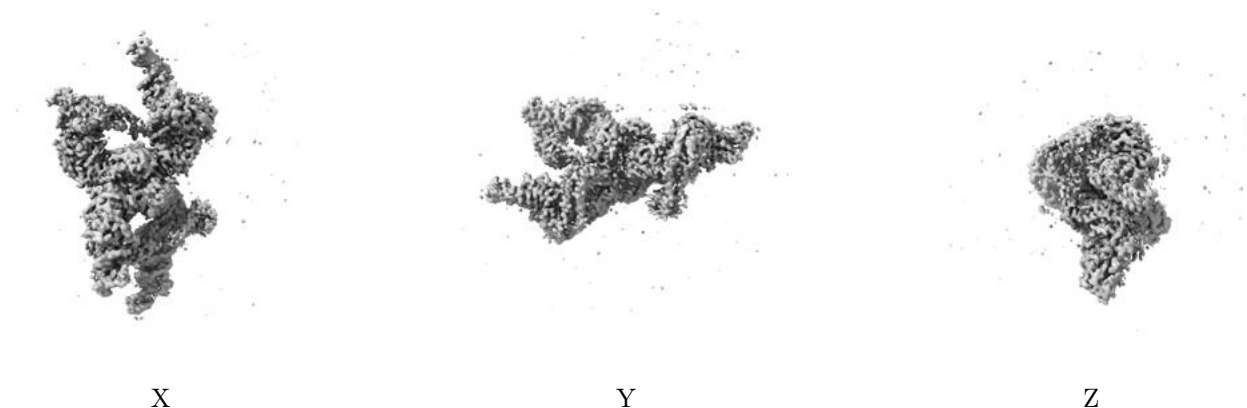


Z

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 0.0226. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

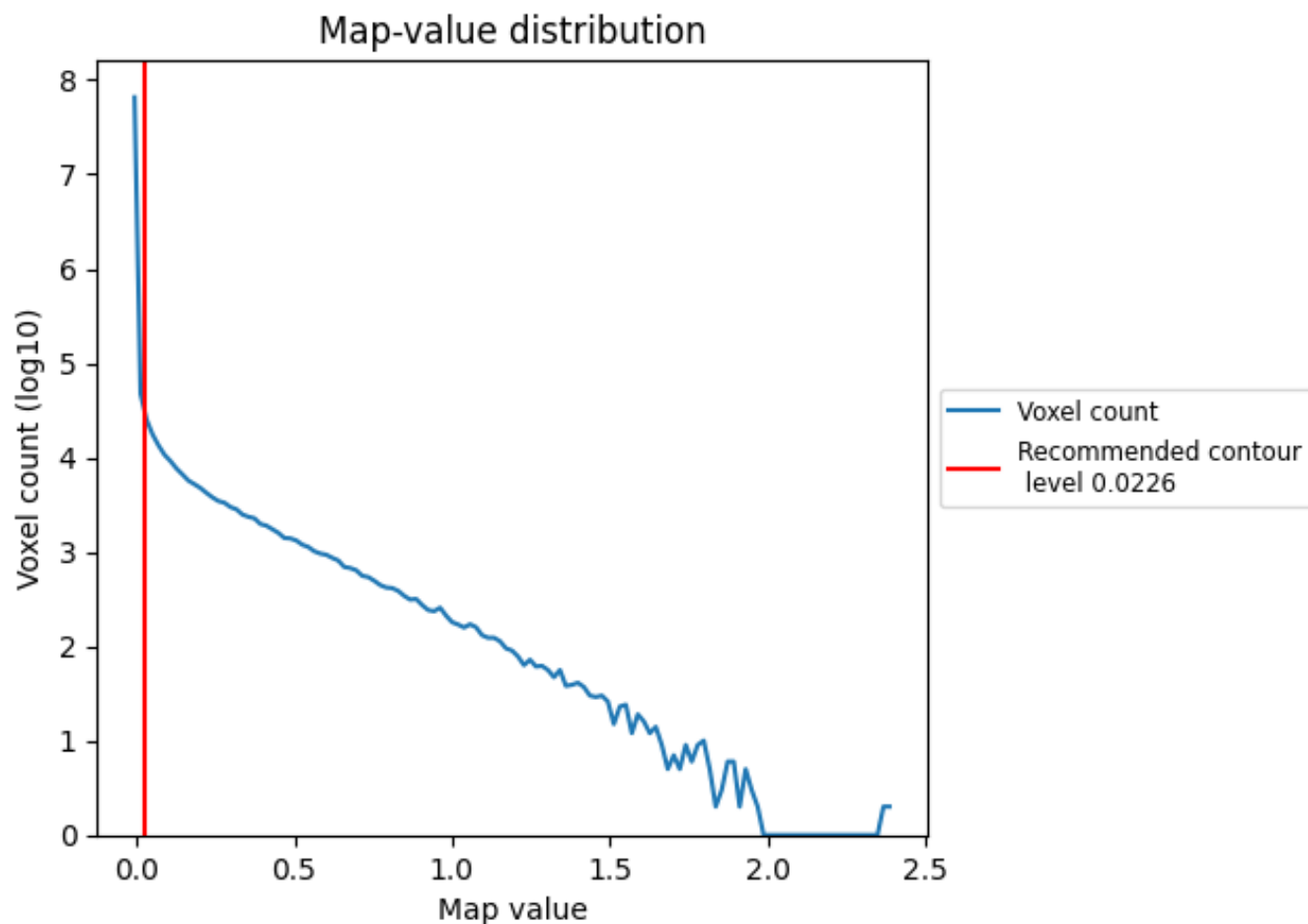
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

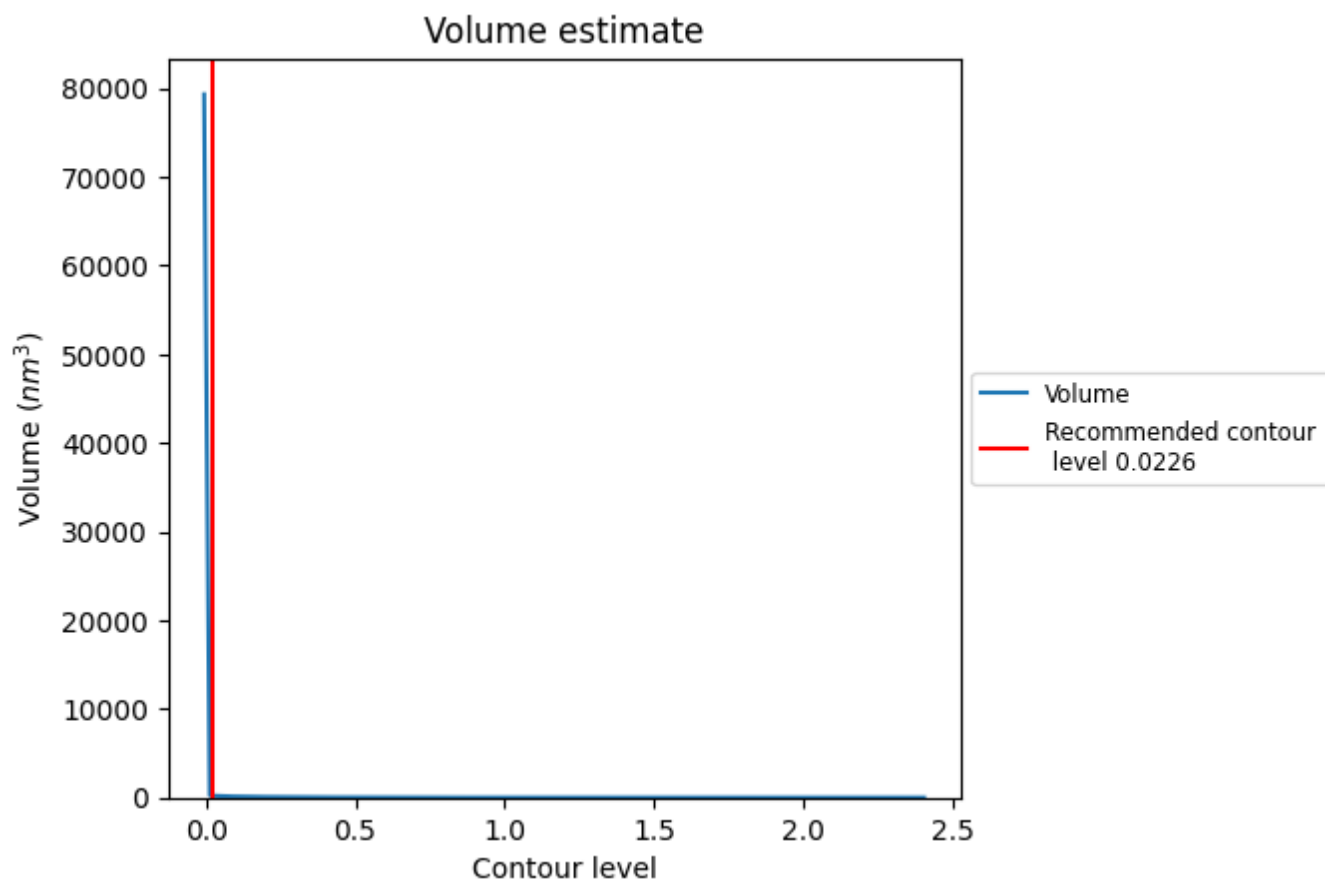
This section contains the results of statistical analysis of the map.

### 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 223 nm<sup>3</sup>; this corresponds to an approximate mass of 201 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

## 7.3 Rotationally averaged power spectrum [i](#)

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

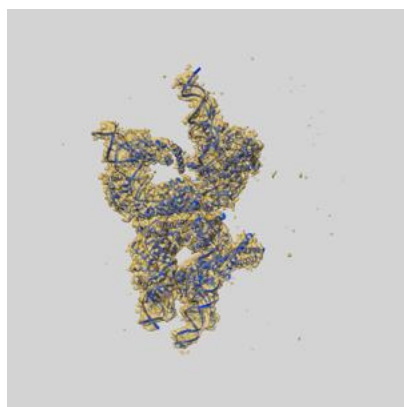
## 8 Fourier-Shell correlation

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

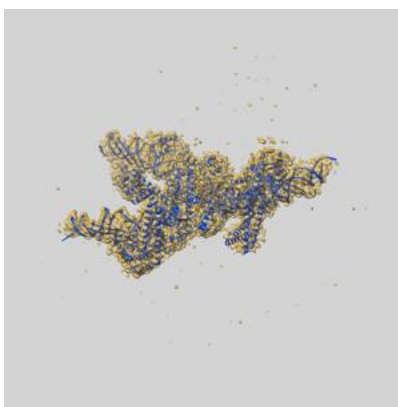
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-47286 and PDB model 9DXF. Per-residue inclusion information can be found in section 3 on page 10.

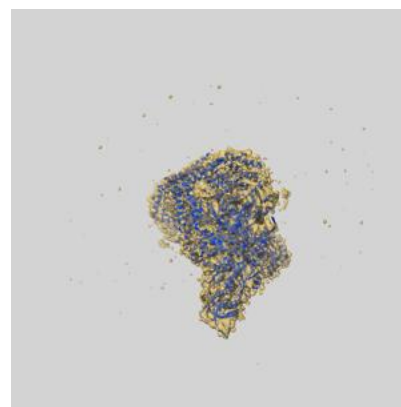
### 9.1 Map-model overlay [i](#)



X



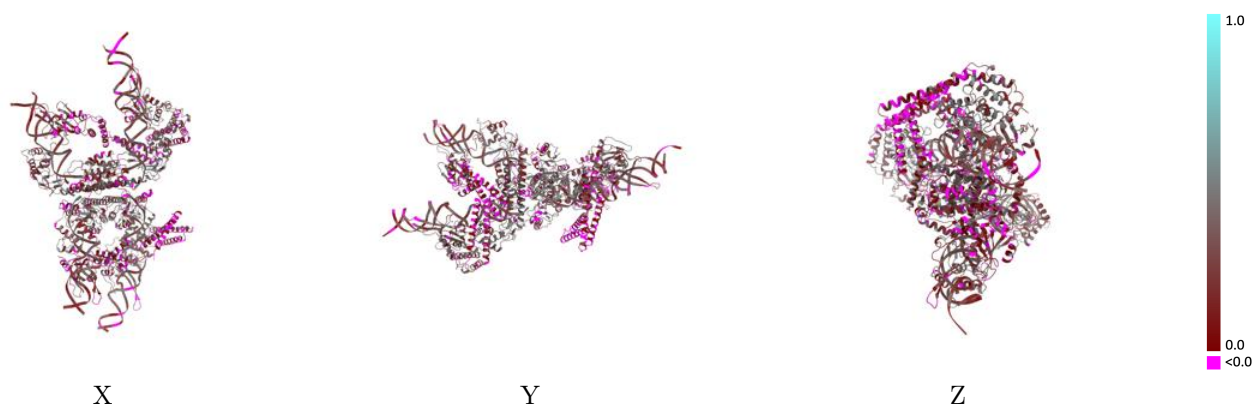
Y



Z

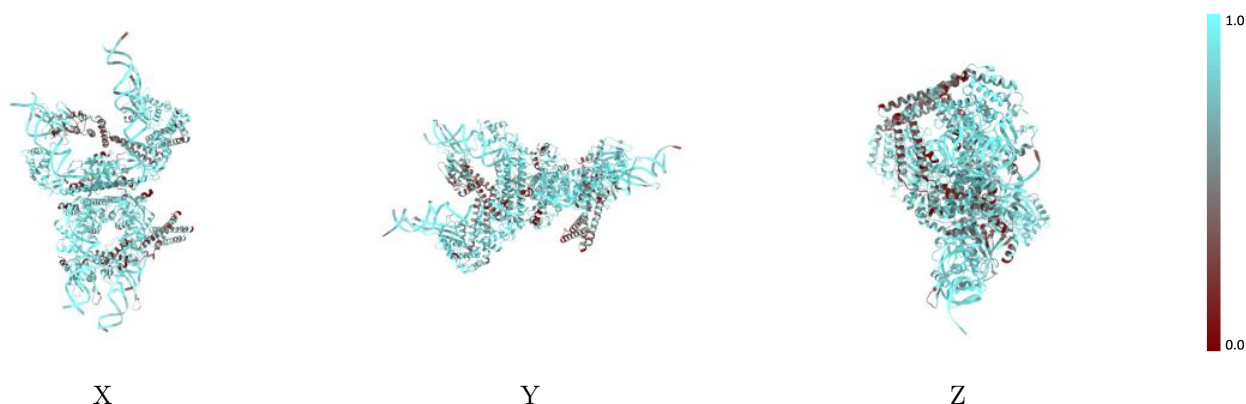
The images above show the 3D surface view of the map at the recommended contour level 0.0226 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



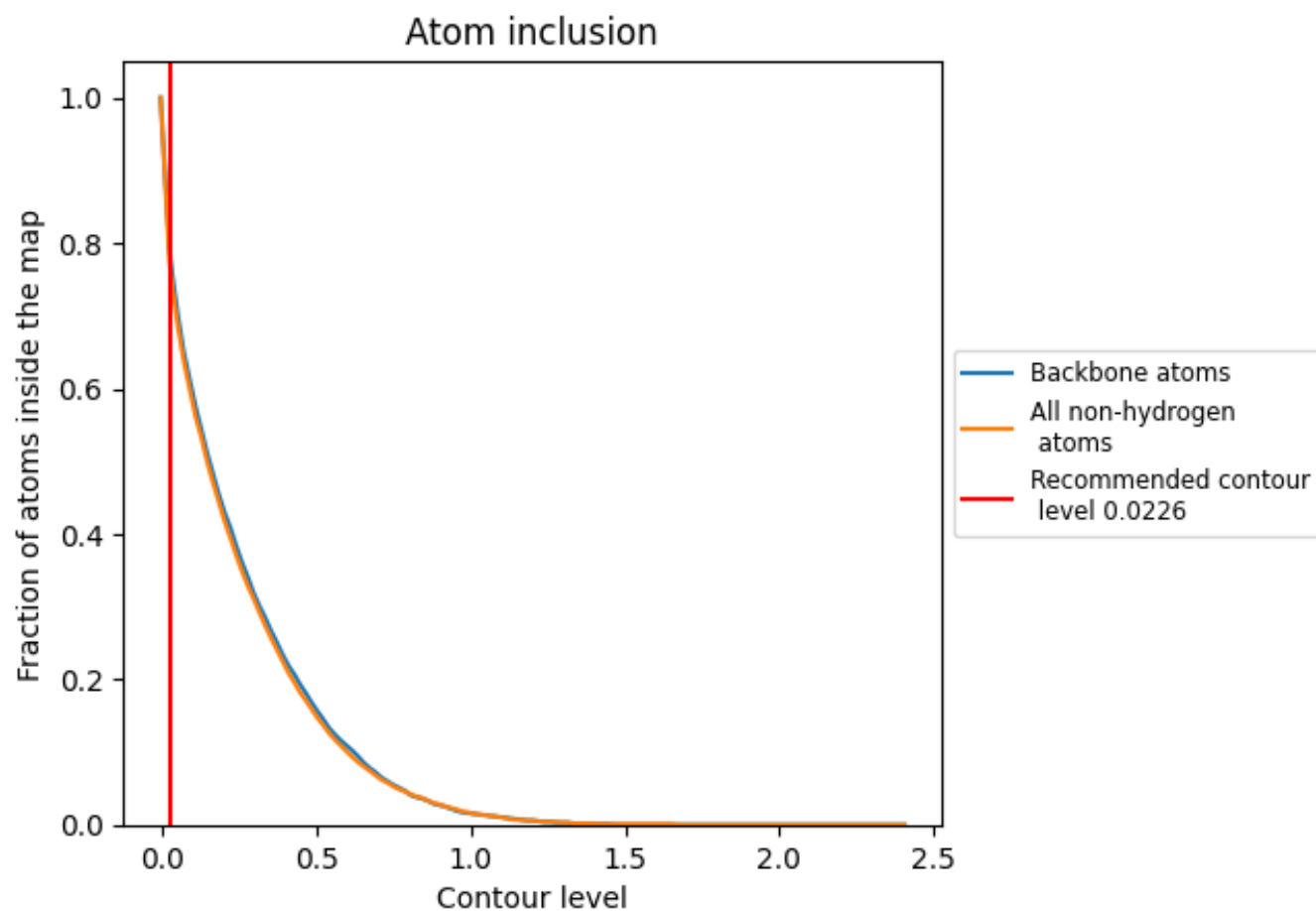
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0226).

## 9.4 Atom inclusion [i](#)





































At the recommended contour level, 79% of all backbone atoms, 77% 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 (0.0226) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7740	 0.2600
A	 0.7990	 0.2910
B	 0.7640	 0.3020
C	 0.6960	 0.2170
D	 0.5200	 0.1340
E	 0.8750	 0.2700
F	 0.8870	 0.2990
G	 0.9330	 0.3040
H	 0.9070	 0.2840
I	 0.8280	 0.2920
J	 0.7810	 0.3020
K	 0.7230	 0.2240
L	 0.5070	 0.1350
M	 0.8950	 0.2750
N	 0.9150	 0.3070
O	 0.9590	 0.3030
P	 0.9370	 0.2910

