



## Full wwPDB EM Validation Report ⓘ

Nov 10, 2024 – 12:21 AM EST

PDB ID : 6D5F  
EMDB ID : EMD-7797  
Title : Cryo-EM reconstruction of membrane-enveloped filamentous virus SFV1 (Sulfolobus filamentous virus 1)  
Authors : Wang, F.; Osinski, T.; Liu, Y.; Krupovic, M.; Prangishvili, D.; Egelman, E.H.  
Deposited on : 2018-04-19  
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](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.dev113  
MolProbity : 4.02b-467  
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) : 2.39

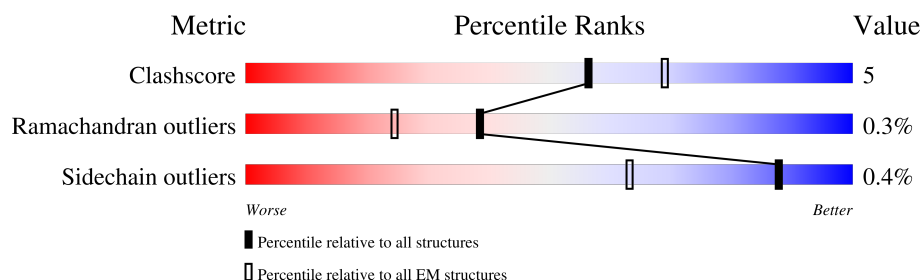
# 1 Overall quality at a glance

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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	199	
1	B	199	
1	C	199	
1	D	199	
1	E	199	
1	F	199	
1	G	199	
1	H	199	

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Mol	Chain	Length	Quality of chain
1	I	199	
1	J	199	
1	K	199	
1	L	199	
1	M	199	
1	N	199	
1	O	199	
1	P	199	
1	Q	199	
1	R	199	
1	S	199	
1	T	199	
1	U	199	
1	V	199	
1	W	199	
1	X	199	
1	Y	199	
1	Z	199	
2	a	137	
2	b	137	
2	c	137	
2	d	137	
2	e	137	
2	f	137	
2	g	137	

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Mol	Chain	Length	Quality of chain
2	h	137	
2	i	137	
2	j	137	
2	k	137	
2	l	137	
2	m	137	
2	n	137	
2	o	137	
2	p	137	
2	q	137	
2	r	137	
2	s	137	
2	t	137	
2	u	137	
2	v	137	
2	w	137	
2	x	137	
2	y	137	
2	z	137	
3	1	336	
3	2	336	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 79946 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 Fimbrial protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	B	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	C	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	D	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	E	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	F	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	G	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	H	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	I	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	J	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	K	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	L	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	M	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	N	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	O	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	P	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	Q	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	R	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	S	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	T	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	U	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	V	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	W	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	X	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	Y	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		
1	Z	194	Total	C	N	O	S	0	0
			1486	949	241	289	7		

- Molecule 2 is a protein called Fimbrial protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	a	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	b	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	c	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	d	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	e	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	f	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	g	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	h	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	i	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	j	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	k	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	l	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	m	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	n	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	o	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	p	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	q	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	r	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	s	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	t	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	u	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	v	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	w	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	x	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	y	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		
2	z	131	Total	C	N	O	S	0	0
			1059	675	183	196	5		

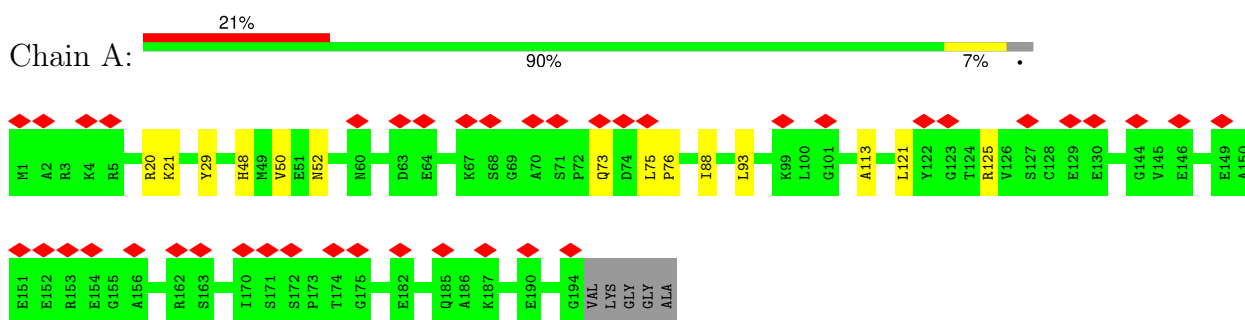
- Molecule 3 is a DNA chain called DNA (336-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	1	336	Total	C	N	O	P	0	0
			6888	3360	1176	2016	336		
3	2	336	Total	C	N	O	P	0	0
			6888	3360	1176	2016	336		

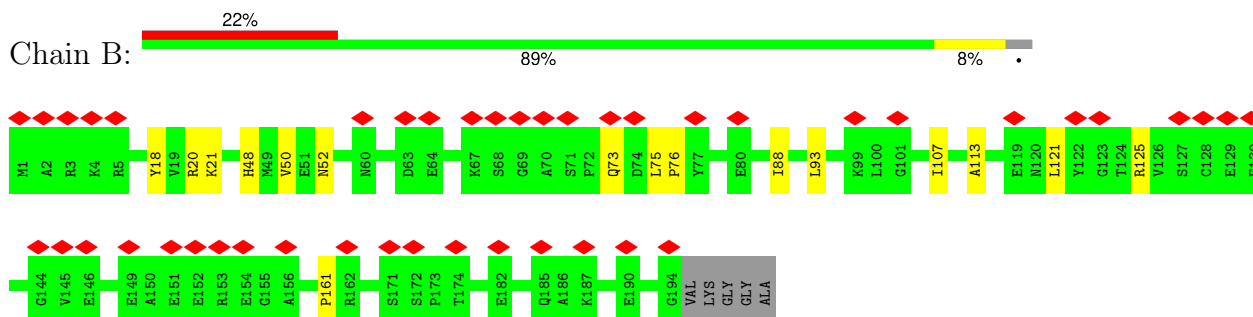
### 3 Residue-property plots [i](#)

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.

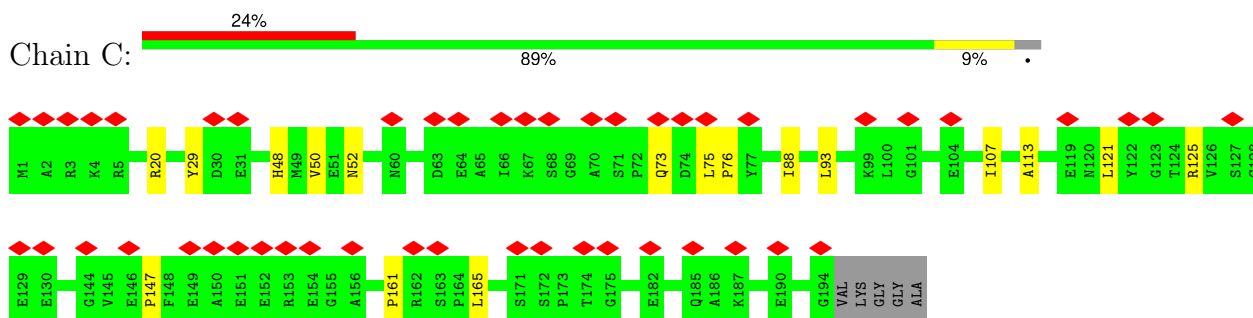
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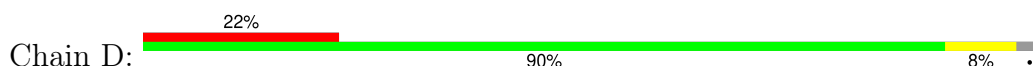
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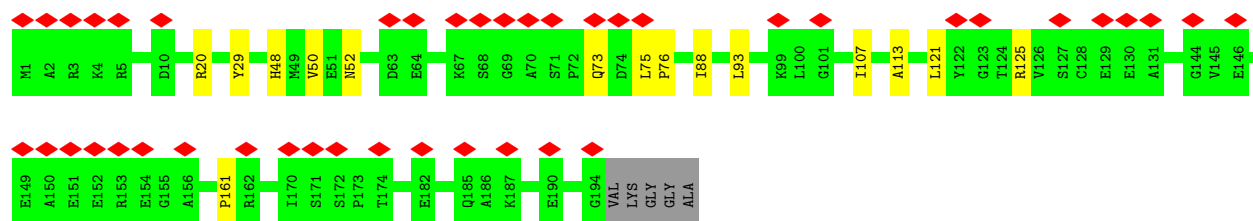
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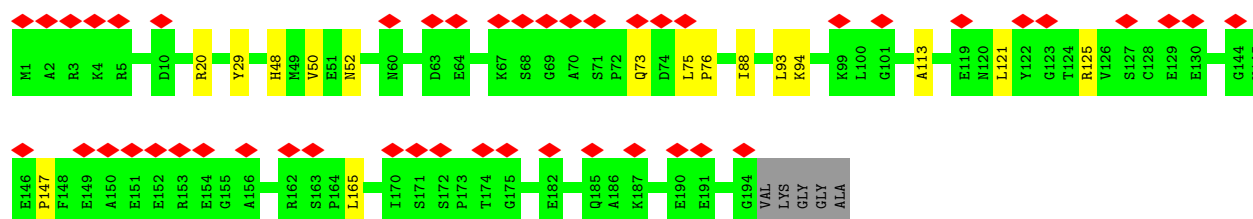
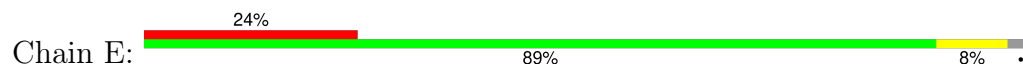
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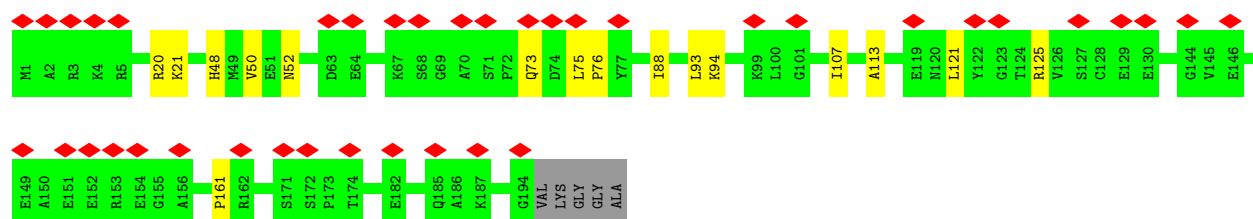
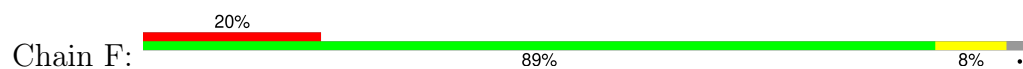




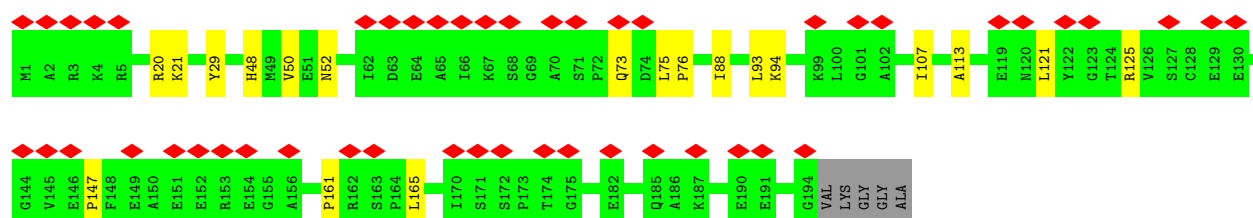
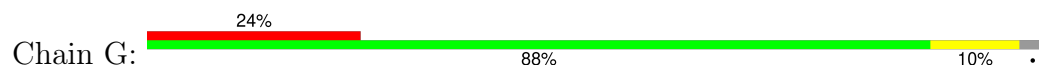
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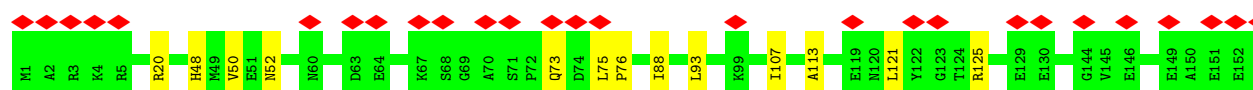
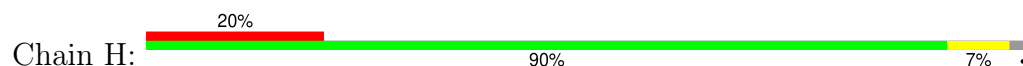
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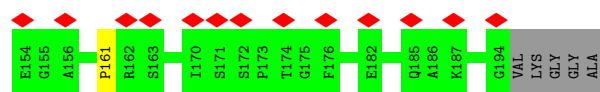


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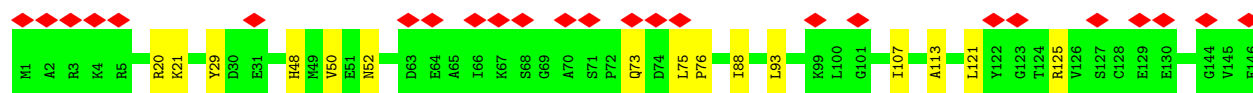
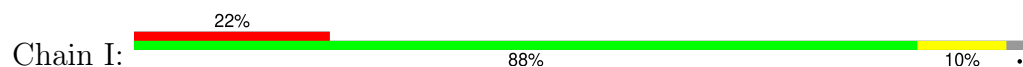


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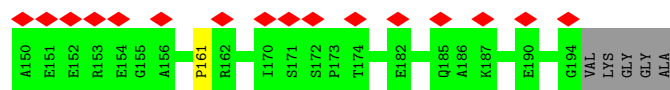
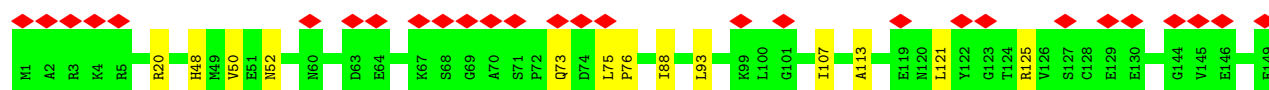
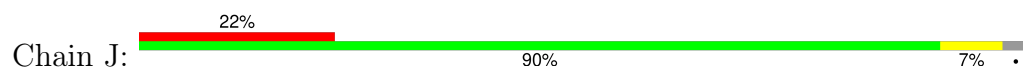




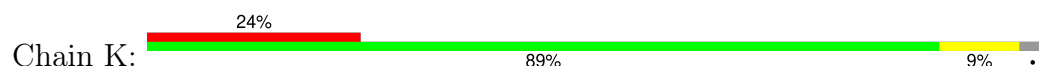
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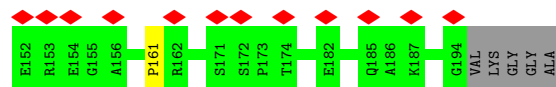
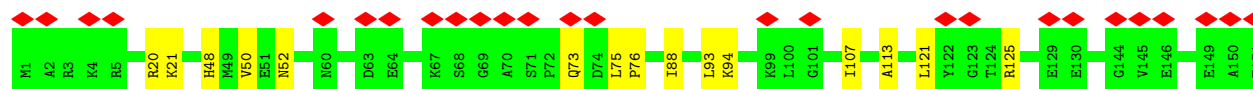
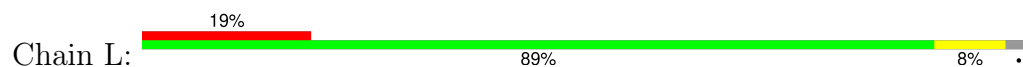
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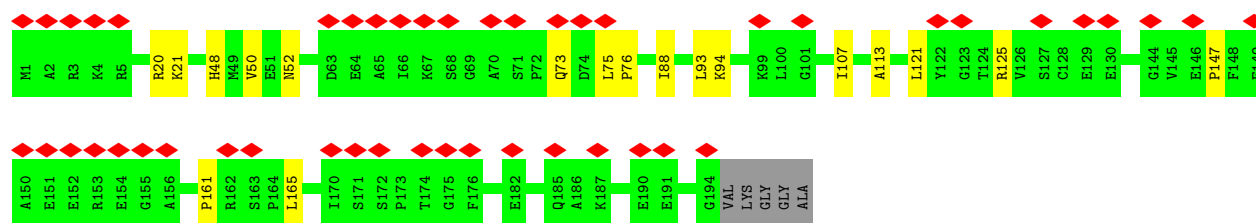
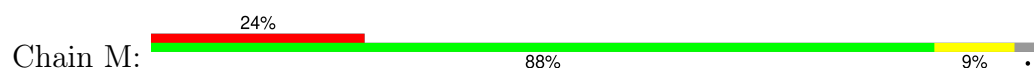
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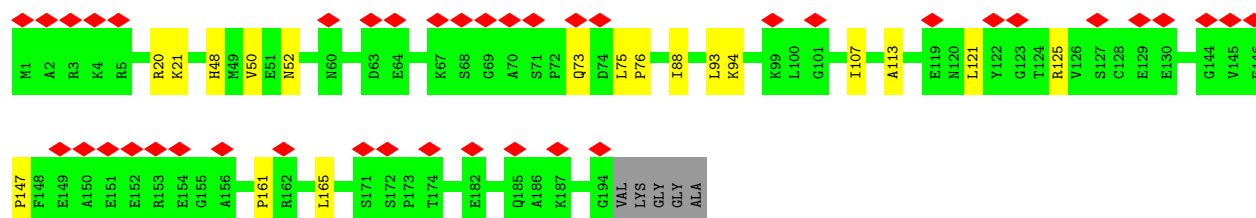
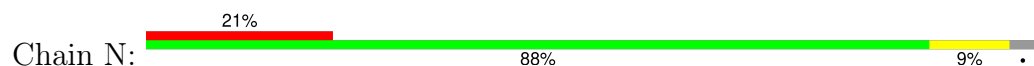
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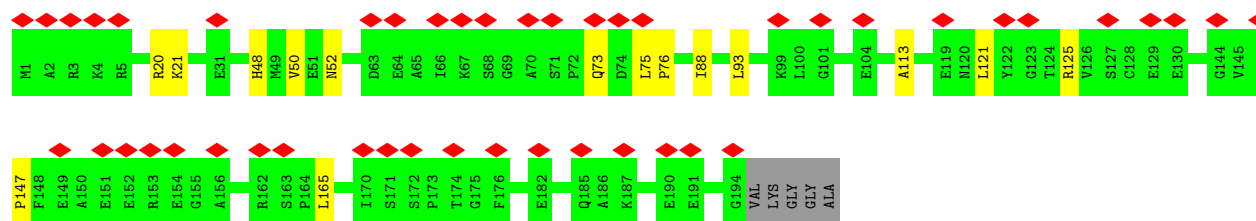
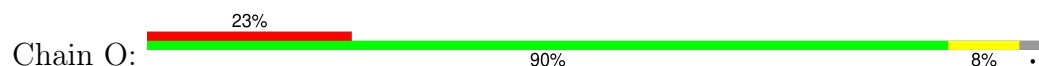
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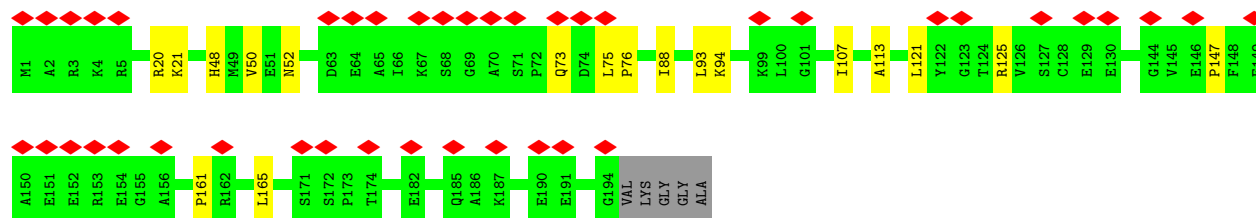
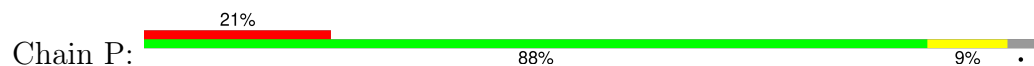
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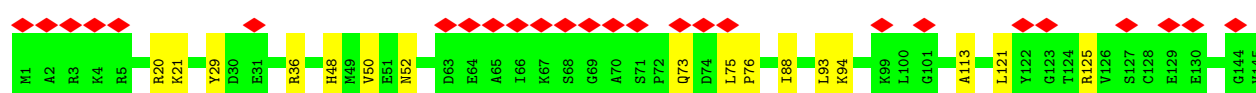
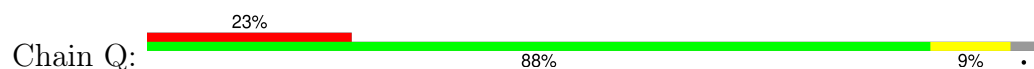
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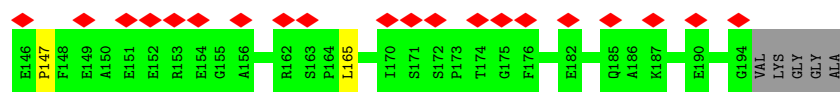


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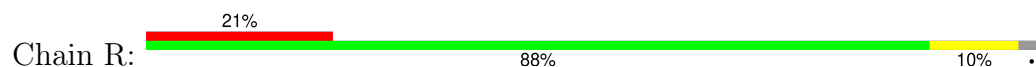


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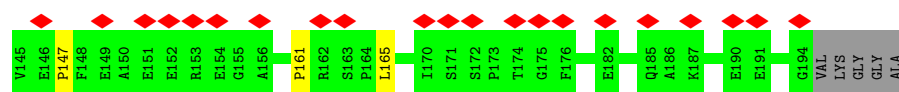
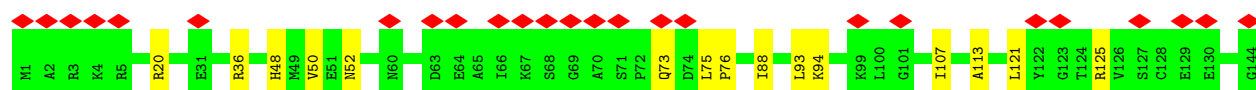
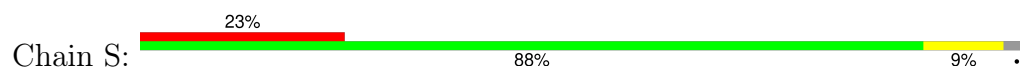




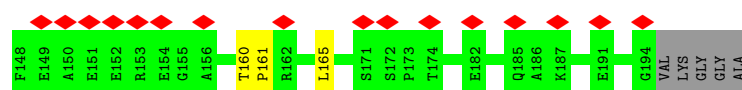
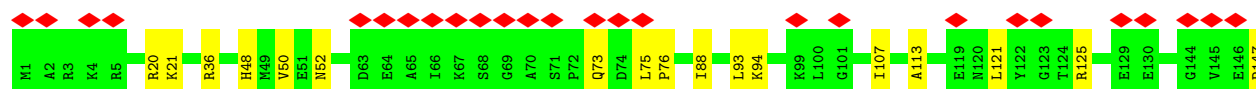
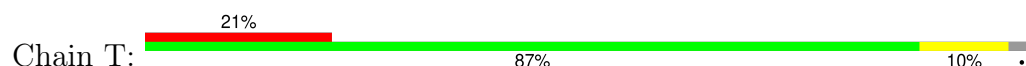
- Molecule 1: Fimbrial protein



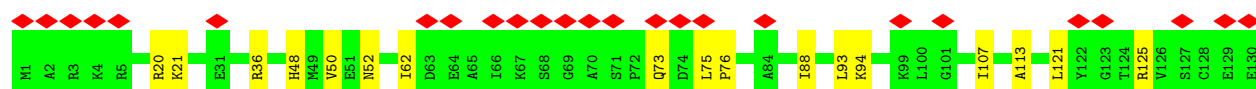
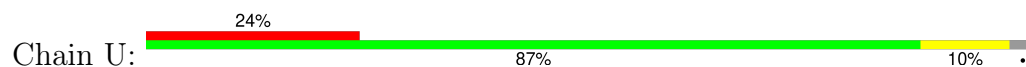
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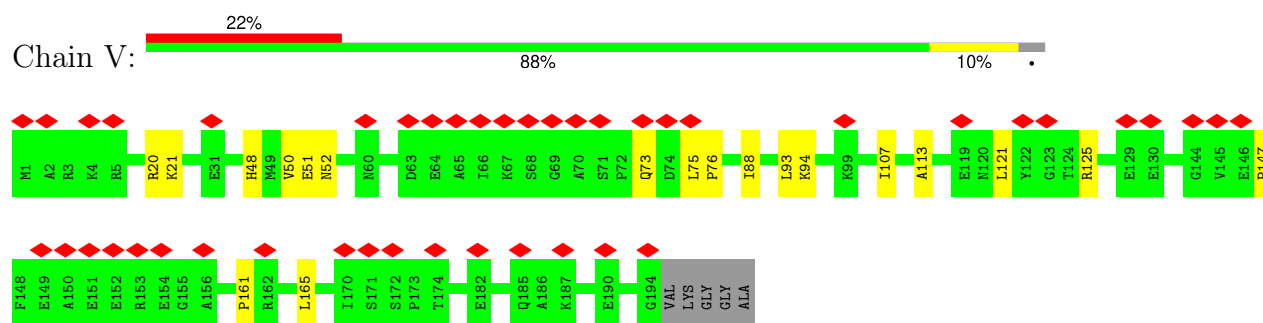
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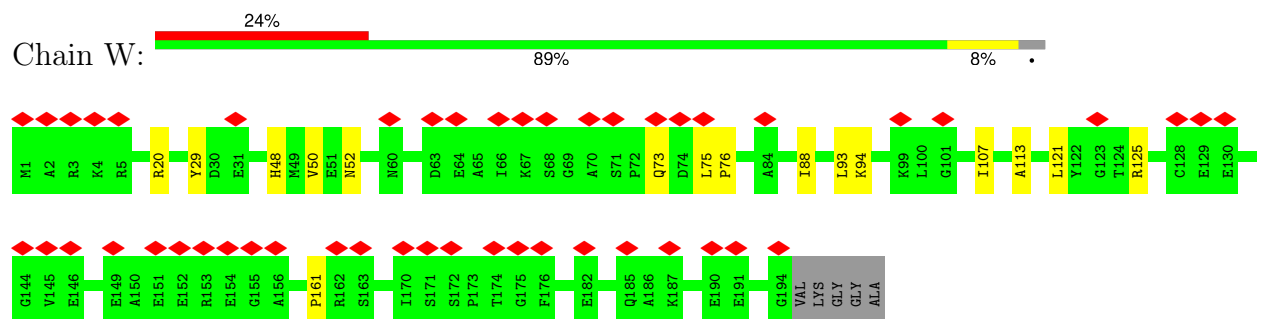
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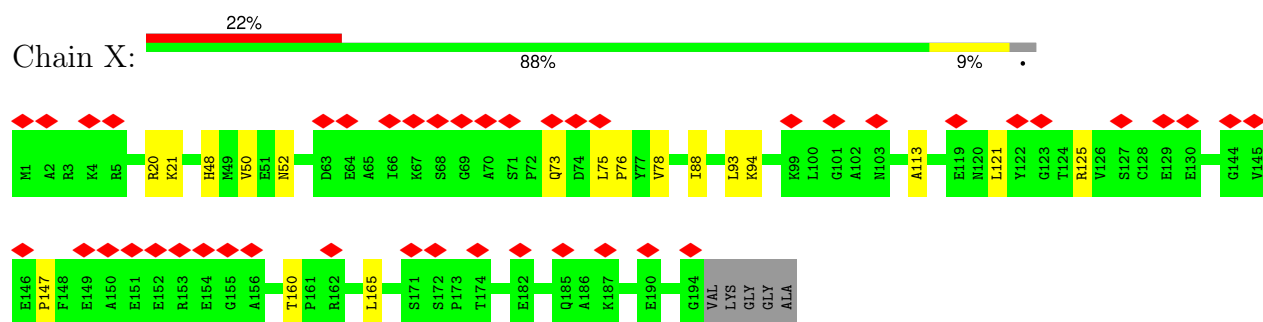
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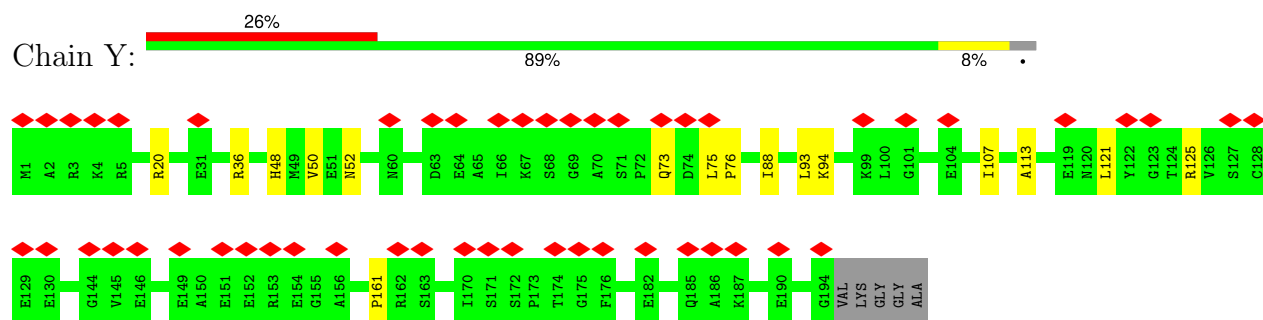
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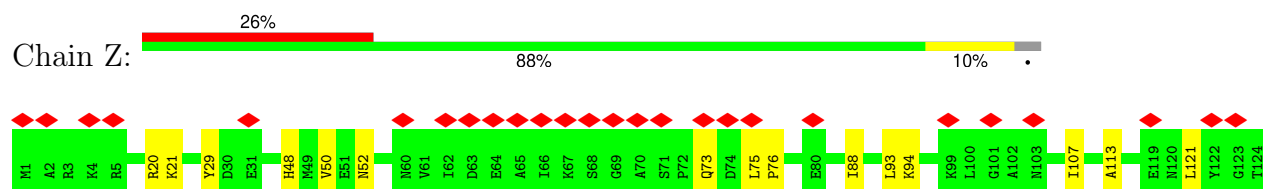
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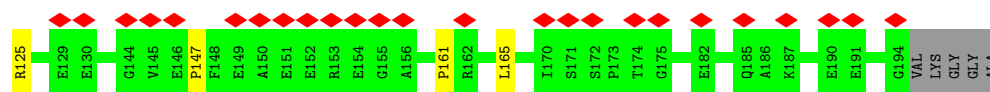


- Molecule 1: Fimbrial protein

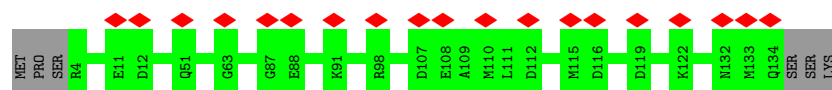


- Molecule 1: Fimbrial protein

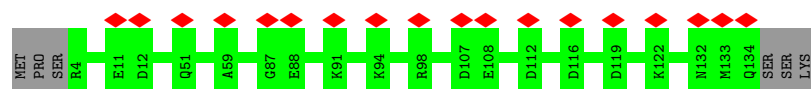




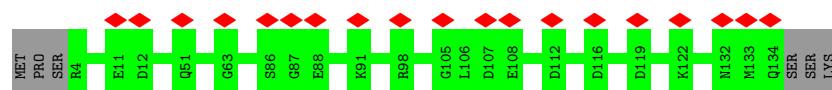
- Molecule 2: Fimbrial protein



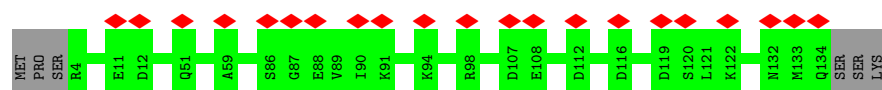
- Molecule 2: Fimbrial protein



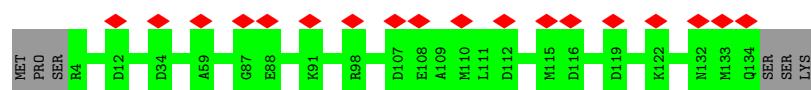
- Molecule 2: Fimbrial protein



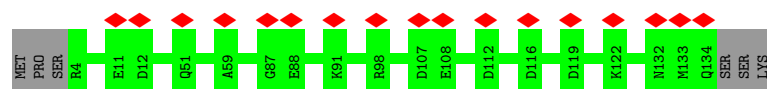
- Molecule 2: Fimbrial protein



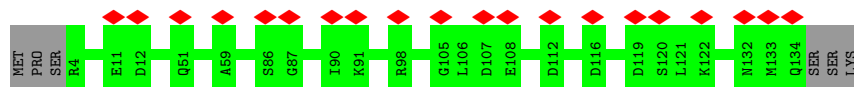
- Molecule 2: Fimbrial protein



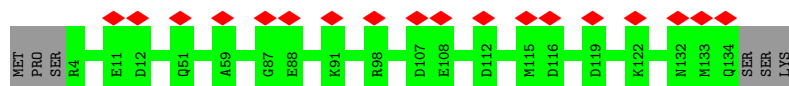
- Molecule 2: Fimbrial protein



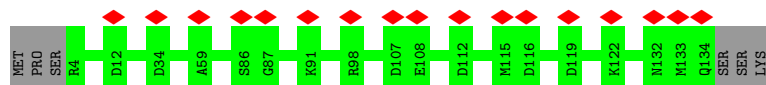
- Molecule 2: Fimbrial protein



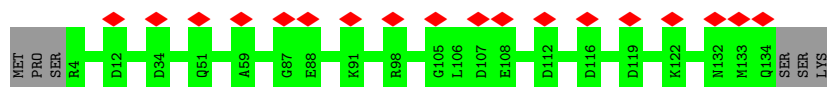
- Molecule 2: Fimbrial protein



- Molecule 2: Fimbrial protein



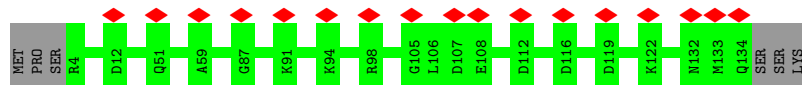
- Molecule 2: Fimbrial protein



- Molecule 2: Fimbrial protein

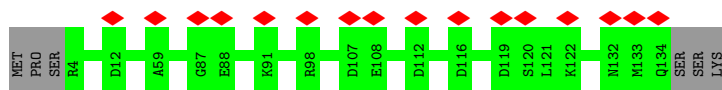


- Molecule 2: Fimbrial protein

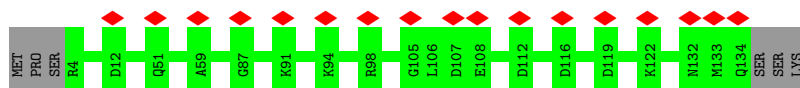


- Molecule 2: Fimbrial protein

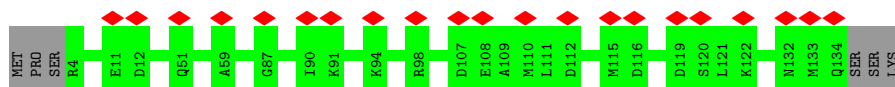




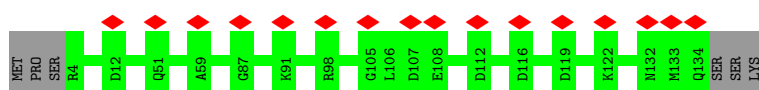
- Molecule 2: Fimbrial protein



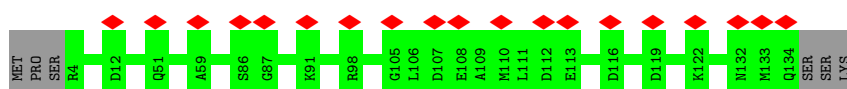
- Molecule 2: Fimbrial protein



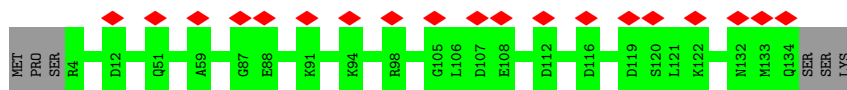
- Molecule 2: Fimbrial protein



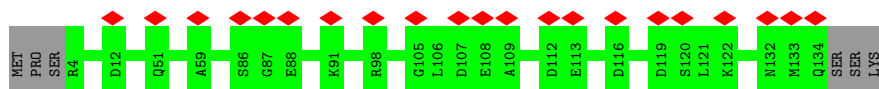
- Molecule 2: Fimbrial protein



- Molecule 2: Fimbrial protein

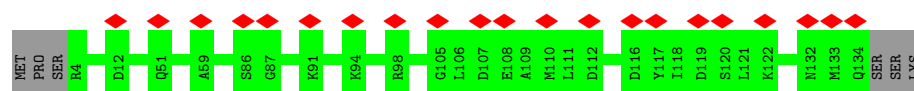


- Molecule 2: Fimbrial protein

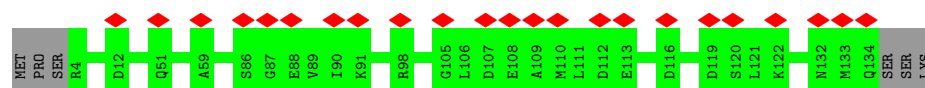


- Molecule 2: Fimbrial protein

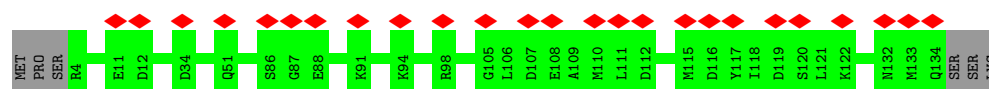




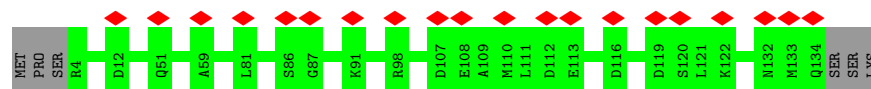
• Molecule 2: Fimbrial protein



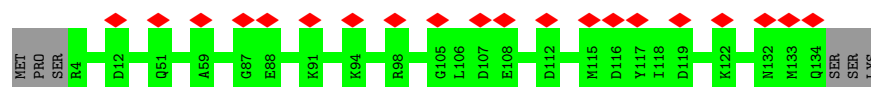
• Molecule 2: Fimbrial protein



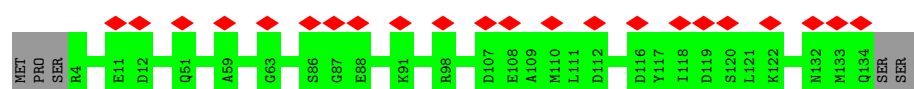
• Molecule 2: Fimbrial protein



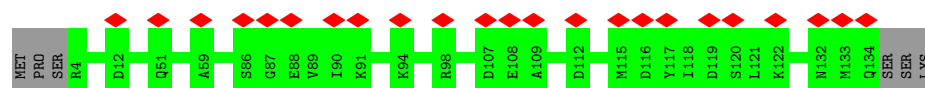
• Molecule 2: Fimbrial protein



• Molecule 2: Fimbrial protein

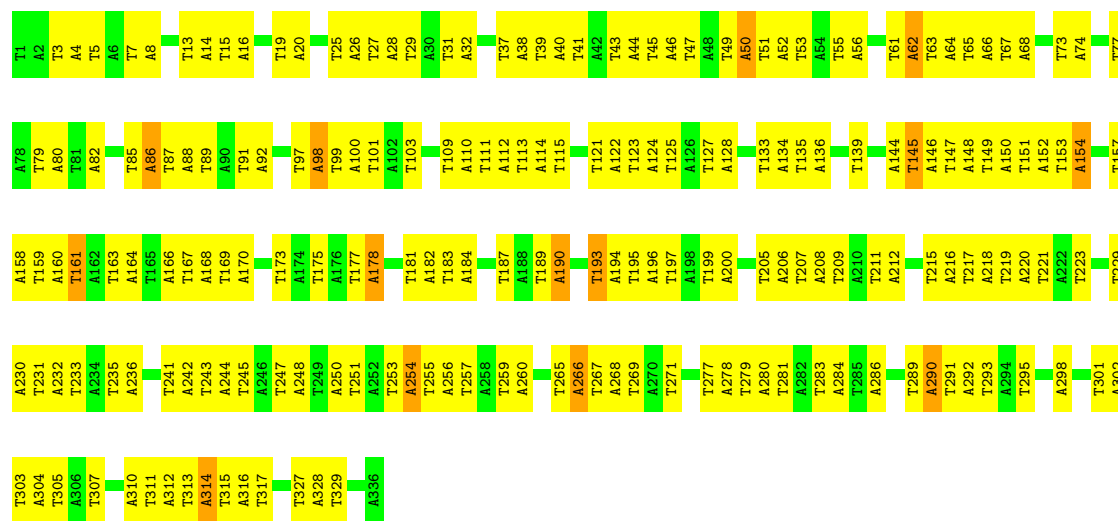


• Molecule 2: Fimbrial protein



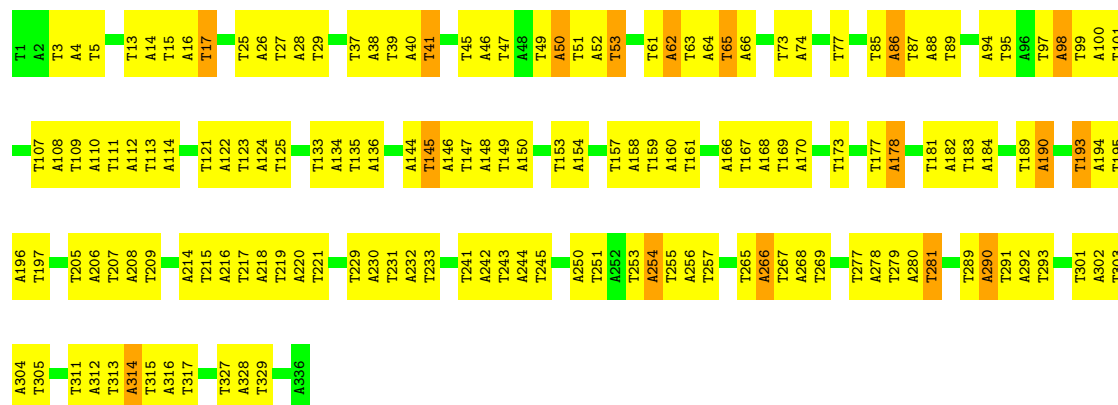
- Molecule 3: DNA (336-MER)

Chain 1:  41% 55%



- Molecule 3: DNA (336-MER)

Chain 2:  53% 42% 5%



## 4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=21.02572°, rise=2.76412 Å, axial sym=C1	Depositor
Number of segments used	87803	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 ( $e^-/\text{Å}^2$ )	48	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.284	Depositor
Minimum map value	-0.175	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.018	Depositor
Recommended contour level	0.07	Depositor
Map size (Å)	418.56, 418.56, 418.56	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.09, 1.09, 1.09	Depositor

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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	0.41	0/1517	0.59	1/2063 (0.0%)
1	B	0.41	0/1517	0.59	1/2063 (0.0%)
1	C	0.41	0/1517	0.59	1/2063 (0.0%)
1	D	0.41	0/1517	0.59	1/2063 (0.0%)
1	E	0.41	0/1517	0.59	1/2063 (0.0%)
1	F	0.41	0/1517	0.59	1/2063 (0.0%)
1	G	0.41	0/1517	0.59	1/2063 (0.0%)
1	H	0.41	0/1517	0.59	1/2063 (0.0%)
1	I	0.41	0/1517	0.59	1/2063 (0.0%)
1	J	0.41	0/1517	0.59	1/2063 (0.0%)
1	K	0.41	0/1517	0.59	1/2063 (0.0%)
1	L	0.41	0/1517	0.59	1/2063 (0.0%)
1	M	0.41	0/1517	0.59	1/2063 (0.0%)
1	N	0.41	0/1517	0.59	1/2063 (0.0%)
1	O	0.41	0/1517	0.59	1/2063 (0.0%)
1	P	0.41	0/1517	0.59	1/2063 (0.0%)
1	Q	0.41	0/1517	0.59	1/2063 (0.0%)
1	R	0.41	0/1517	0.59	1/2063 (0.0%)
1	S	0.41	0/1517	0.59	1/2063 (0.0%)
1	T	0.41	0/1517	0.59	1/2063 (0.0%)
1	U	0.41	0/1517	0.59	1/2063 (0.0%)
1	V	0.41	0/1517	0.59	1/2063 (0.0%)
1	W	0.41	0/1517	0.59	1/2063 (0.0%)
1	X	0.41	0/1517	0.59	1/2063 (0.0%)
1	Y	0.41	0/1517	0.59	1/2063 (0.0%)
1	Z	0.41	0/1517	0.59	1/2063 (0.0%)
2	a	0.46	0/1076	0.62	0/1453
2	b	0.46	0/1076	0.62	0/1453
2	c	0.46	0/1076	0.62	0/1453
2	d	0.46	0/1076	0.62	0/1453
2	e	0.46	0/1076	0.62	0/1453
2	f	0.46	0/1076	0.62	0/1453
2	g	0.46	0/1076	0.62	0/1453
2	h	0.46	0/1076	0.62	0/1453

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
2	i	0.46	0/1076	0.62	0/1453
2	j	0.46	0/1076	0.62	0/1453
2	k	0.46	0/1076	0.62	0/1453
2	l	0.46	0/1076	0.62	0/1453
2	m	0.46	0/1076	0.62	0/1453
2	n	0.46	0/1076	0.62	0/1453
2	o	0.46	0/1076	0.62	0/1453
2	p	0.46	0/1076	0.62	0/1453
2	q	0.46	0/1076	0.62	0/1453
2	r	0.46	0/1076	0.62	0/1453
2	s	0.46	0/1076	0.62	0/1453
2	t	0.46	0/1076	0.62	0/1453
2	u	0.46	0/1076	0.62	0/1453
2	v	0.46	0/1076	0.62	0/1453
2	w	0.46	0/1076	0.62	0/1453
2	x	0.46	0/1076	0.62	0/1453
2	y	0.46	0/1076	0.62	0/1453
2	z	0.46	0/1076	0.62	0/1453
3	1	1.30	8/7727 (0.1%)	1.15	33/11924 (0.3%)
3	2	1.30	7/7727 (0.1%)	1.15	34/11924 (0.3%)
All	All	0.68	15/82872 (0.0%)	0.75	93/115264 (0.1%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	1	66	DA	C3'-O3'	-6.97	1.34	1.44
3	2	66	DA	C3'-O3'	-6.93	1.34	1.44
3	1	144	DA	N9-C4	-6.17	1.34	1.37
3	2	144	DA	N9-C4	-6.04	1.34	1.37
3	2	178	DA	C3'-O3'	-5.87	1.36	1.44
3	1	178	DA	C3'-O3'	-5.86	1.36	1.44
3	1	290	DA	N9-C4	-5.36	1.34	1.37
3	2	190	DA	C3'-O3'	-5.31	1.37	1.44
3	1	190	DA	C3'-O3'	-5.29	1.37	1.44
3	2	290	DA	N9-C4	-5.28	1.34	1.37
3	1	166	DA	C3'-O3'	-5.09	1.37	1.44
3	2	166	DA	C3'-O3'	-5.07	1.37	1.44
3	1	114	DA	C3'-O3'	-5.06	1.37	1.44
3	2	114	DA	C3'-O3'	-5.05	1.37	1.44
3	1	154	DA	C3'-O3'	-5.00	1.37	1.44

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	2	314	DA	O4'-C4'-C3'	-6.35	101.96	104.50
3	1	314	DA	O4'-C4'-C3'	-6.26	101.99	104.50
3	1	197	DT	O4'-C4'-C3'	-6.11	102.06	104.50
3	2	197	DT	O4'-C4'-C3'	-6.04	102.08	104.50
3	1	101	DT	O4'-C4'-C3'	-6.02	102.09	104.50
3	2	257	DT	O4'-C4'-C3'	-5.95	102.12	104.50
3	2	98	DA	O4'-C4'-C3'	-5.93	102.13	104.50
3	1	98	DA	O4'-C4'-C3'	-5.89	102.14	104.50
3	1	89	DT	O4'-C4'-C3'	-5.83	102.17	104.50
3	2	101	DT	O4'-C4'-C3'	-5.83	102.17	104.50
3	2	173	DT	O4'-C4'-C3'	-5.82	102.17	104.50
3	1	257	DT	O4'-C4'-C3'	-5.80	102.18	104.50
3	2	53	DT	O4'-C4'-C3'	-5.70	102.22	104.50
3	1	62	DA	O4'-C4'-C3'	-5.67	102.23	104.50
3	1	266	DA	O4'-C4'-C3'	-5.67	102.23	104.50
3	1	53	DT	O4'-C4'-C3'	-5.67	102.23	104.50
3	1	269	DT	O4'-C4'-C3'	-5.67	102.23	104.50
3	1	145	DT	O4'-C4'-C3'	-5.66	102.24	104.50
3	2	269	DT	O4'-C4'-C3'	-5.65	102.24	104.50
3	2	89	DT	O4'-C4'-C3'	-5.64	102.24	104.50
3	1	173	DT	O4'-C4'-C3'	-5.63	102.25	104.50
3	1	41	DT	O4'-C4'-C3'	-5.62	102.25	104.50
3	2	145	DT	O4'-C4'-C3'	-5.62	102.25	104.50
3	1	281	DT	O4'-C4'-C3'	-5.61	102.26	104.50
3	2	41	DT	O4'-C4'-C3'	-5.61	102.26	104.50
3	2	281	DT	O4'-C4'-C3'	-5.60	102.26	104.50
3	2	86	DA	O4'-C4'-C3'	-5.59	102.26	104.50
3	2	62	DA	O4'-C4'-C3'	-5.58	102.27	104.50
3	2	266	DA	O4'-C4'-C3'	-5.56	102.28	104.50
3	1	29	DT	O4'-C4'-C3'	-5.55	102.28	104.50
3	2	233	DT	O4'-C4'-C3'	-5.54	102.28	104.50
3	1	50	DA	O4'-C4'-C3'	-5.52	102.29	104.50
3	1	233	DT	O4'-C4'-C3'	-5.50	102.30	104.50
3	2	293	DT	O4'-C4'-C3'	-5.50	102.30	104.50
3	1	86	DA	O4'-C4'-C3'	-5.49	102.31	104.50
3	1	125	DT	O4'-C4'-C3'	-5.45	102.32	104.50
3	2	317	DT	O4'-C4'-C3'	-5.43	102.33	104.50
3	2	125	DT	O4'-C4'-C3'	-5.42	102.33	104.50
3	1	293	DT	O4'-C4'-C3'	-5.42	102.33	104.50
3	1	305	DT	O4'-C4'-C3'	-5.42	102.33	104.50
3	2	29	DT	O4'-C4'-C3'	-5.42	102.33	104.50
3	2	50	DA	O4'-C4'-C3'	-5.41	102.33	104.50
3	1	77	DT	O4'-C4'-C3'	-5.38	102.35	104.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	2	305	DT	O4'-C4'-C3'	-5.38	102.35	104.50
3	1	209	DT	O4'-C4'-C3'	-5.35	102.36	104.50
3	2	209	DT	O4'-C4'-C3'	-5.35	102.36	104.50
1	F	121	LEU	CA-CB-CG	5.34	127.59	115.30
1	Z	121	LEU	CA-CB-CG	5.34	127.58	115.30
3	1	245	DT	O4'-C4'-C3'	-5.34	102.36	104.50
3	2	245	DT	O4'-C4'-C3'	-5.34	102.36	104.50
1	J	121	LEU	CA-CB-CG	5.33	127.57	115.30
1	X	121	LEU	CA-CB-CG	5.33	127.57	115.30
1	G	121	LEU	CA-CB-CG	5.33	127.56	115.30
1	L	121	LEU	CA-CB-CG	5.33	127.56	115.30
1	S	121	LEU	CA-CB-CG	5.33	127.56	115.30
1	B	121	LEU	CA-CB-CG	5.33	127.56	115.30
1	T	121	LEU	CA-CB-CG	5.33	127.56	115.30
1	Y	121	LEU	CA-CB-CG	5.33	127.56	115.30
1	E	121	LEU	CA-CB-CG	5.33	127.55	115.30
1	P	121	LEU	CA-CB-CG	5.33	127.55	115.30
1	R	121	LEU	CA-CB-CG	5.33	127.55	115.30
1	D	121	LEU	CA-CB-CG	5.32	127.55	115.30
1	I	121	LEU	CA-CB-CG	5.32	127.55	115.30
1	N	121	LEU	CA-CB-CG	5.32	127.54	115.30
1	K	121	LEU	CA-CB-CG	5.32	127.54	115.30
1	M	121	LEU	CA-CB-CG	5.32	127.54	115.30
1	V	121	LEU	CA-CB-CG	5.32	127.54	115.30
1	O	121	LEU	CA-CB-CG	5.32	127.53	115.30
1	W	121	LEU	CA-CB-CG	5.32	127.53	115.30
1	C	121	LEU	CA-CB-CG	5.31	127.52	115.30
1	Q	121	LEU	CA-CB-CG	5.31	127.52	115.30
1	U	121	LEU	CA-CB-CG	5.31	127.52	115.30
3	1	317	DT	O4'-C4'-C3'	-5.31	102.38	104.50
3	1	329	DT	O4'-C4'-C3'	-5.31	102.38	104.50
1	A	121	LEU	CA-CB-CG	5.30	127.49	115.30
1	H	121	LEU	CA-CB-CG	5.30	127.49	115.30
3	1	65	DT	O4'-C4'-C3'	-5.30	102.38	104.50
3	2	77	DT	O4'-C4'-C3'	-5.30	102.38	104.50
3	2	254	DA	O4'-C4'-C3'	-5.26	102.39	104.50
3	1	221	DT	O4'-C4'-C3'	-5.26	102.39	104.50
3	2	65	DT	O4'-C4'-C3'	-5.25	102.40	104.50
3	1	113	DT	O4'-C4'-C3'	-5.23	102.41	104.50
3	1	254	DA	O4'-C4'-C3'	-5.22	102.41	104.50
3	1	5	DT	O4'-C4'-C3'	-5.21	102.41	104.50
3	2	329	DT	O4'-C4'-C3'	-5.21	102.42	104.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	1	161	DT	O4'-C4'-C3'	-5.20	102.42	104.50
3	2	221	DT	O4'-C4'-C3'	-5.18	102.43	104.50
3	2	161	DT	O4'-C4'-C3'	-5.17	102.43	104.50
3	2	113	DT	O4'-C4'-C3'	-5.16	102.44	104.50
3	2	5	DT	O4'-C4'-C3'	-5.13	102.45	104.50
3	1	193	DT	O4'-C4'-C3'	-5.09	102.46	104.50
3	2	193	DT	O4'-C4'-C3'	-5.09	102.46	104.50
3	2	17	DT	O4'-C4'-C3'	-5.00	102.50	104.50

There are no chirality outliers.

There are no planarity outliers.

## 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	1486	0	1450	7	0
1	B	1486	0	1450	8	0
1	C	1486	0	1450	8	0
1	D	1486	0	1450	7	0
1	E	1486	0	1450	9	0
1	F	1486	0	1450	8	0
1	G	1486	0	1450	11	0
1	H	1486	0	1450	6	0
1	I	1486	0	1450	10	0
1	J	1486	0	1450	6	0
1	K	1486	0	1450	9	0
1	L	1486	0	1450	9	0
1	M	1486	0	1450	9	0
1	N	1486	0	1450	10	0
1	O	1486	0	1450	7	0
1	P	1486	0	1450	10	0
1	Q	1486	0	1450	11	0
1	R	1486	0	1450	11	0
1	S	1486	0	1450	9	0
1	T	1486	0	1450	12	0
1	U	1486	0	1450	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	V	1486	0	1450	12	0
1	W	1486	0	1450	8	0
1	X	1486	0	1450	10	0
1	Y	1486	0	1450	8	0
1	Z	1486	0	1450	11	0
2	a	1059	0	1076	0	0
2	b	1059	0	1076	0	0
2	c	1059	0	1076	0	0
2	d	1059	0	1076	0	0
2	e	1059	0	1076	0	0
2	f	1059	0	1076	0	0
2	g	1059	0	1076	0	0
2	h	1059	0	1076	0	0
2	i	1059	0	1076	0	0
2	j	1059	0	1076	0	0
2	k	1059	0	1076	0	0
2	l	1059	0	1076	0	0
2	m	1059	0	1076	0	0
2	n	1059	0	1076	0	0
2	o	1059	0	1076	0	0
2	p	1059	0	1076	0	0
2	q	1059	0	1076	0	0
2	r	1059	0	1076	0	0
2	s	1059	0	1076	0	0
2	t	1059	0	1076	0	0
2	u	1059	0	1076	0	0
2	v	1059	0	1076	0	0
2	w	1059	0	1076	0	0
2	x	1059	0	1076	0	0
2	y	1059	0	1076	0	0
2	z	1059	0	1076	0	0
3	1	6888	0	3865	122	0
3	2	6888	0	3865	93	0
All	All	79946	0	73406	365	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 (365) 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:P:20:ARG:NH2	3:1:67:DT:OP1	2.30	0.65
1:S:20:ARG:NH2	3:1:271:DT:OP1	2.31	0.64
1:H:20:ARG:NH2	3:1:115:DT:OP1	2.31	0.64
1:I:20:ARG:NH2	3:1:211:DT:OP1	2.30	0.63
1:U:20:ARG:NH2	3:1:283:DT:OP1	2.31	0.63
1:G:94:LYS:HB2	3:2:123:DT:H5''	1.79	0.63
1:L:20:ARG:NH2	3:1:91:DT:OP1	2.33	0.62
1:N:20:ARG:NH2	3:1:79:DT:OP1	2.33	0.62
1:Y:20:ARG:NH2	3:1:307:DT:OP1	2.34	0.61
1:T:20:ARG:NH2	3:1:43:DT:OP1	2.34	0.61
1:W:20:ARG:NH2	3:1:295:DT:OP1	2.34	0.61
1:E:20:ARG:NH2	3:1:187:DT:OP1	2.32	0.61
1:Z:94:LYS:HB2	3:2:315:DT:H5''	1.81	0.61
1:G:20:ARG:NH2	3:1:199:DT:OP1	2.35	0.59
1:V:94:LYS:HB2	3:2:291:DT:H5''	1.84	0.59
1:J:20:ARG:NH2	3:1:103:DT:OP1	2.36	0.59
1:Q:20:ARG:NH2	3:1:259:DT:OP1	2.34	0.59
1:K:20:ARG:NH2	3:1:223:DT:OP1	2.35	0.59
1:O:20:ARG:NH2	3:1:247:DT:OP1	2.36	0.59
1:X:20:ARG:NH2	3:1:19:DT:OP1	2.36	0.59
1:C:20:ARG:NH2	3:1:175:DT:OP1	2.37	0.58
1:Z:20:ARG:NH2	3:1:7:DT:OP1	2.35	0.58
1:A:20:ARG:NH2	3:1:163:DT:OP1	2.36	0.57
1:D:20:ARG:NH2	3:1:139:DT:OP1	2.36	0.57
1:I:73:GLN:O	1:I:125:ARG:NH2	2.38	0.57
1:R:20:ARG:NH2	3:1:55:DT:OP1	2.37	0.57
1:Y:73:GLN:O	1:Y:125:ARG:NH2	2.38	0.57
1:Z:73:GLN:O	1:Z:125:ARG:NH2	2.38	0.57
1:W:73:GLN:O	1:W:125:ARG:NH2	2.38	0.57
3:2:217:DT:H2'	3:2:218:DA:C8	2.39	0.57
1:B:20:ARG:NH2	3:1:151:DT:OP1	2.37	0.57
1:A:73:GLN:O	1:A:125:ARG:NH2	2.38	0.56
1:H:73:GLN:O	1:H:125:ARG:NH2	2.38	0.56
1:K:94:LYS:HB2	3:2:99:DT:H5''	1.86	0.56
1:S:73:GLN:O	1:S:125:ARG:NH2	2.38	0.56
1:U:73:GLN:O	1:U:125:ARG:NH2	2.38	0.56
1:X:73:GLN:O	1:X:125:ARG:NH2	2.38	0.56
1:F:73:GLN:O	1:F:125:ARG:NH2	2.38	0.56
1:G:73:GLN:O	1:G:125:ARG:NH2	2.38	0.56
1:C:73:GLN:O	1:C:125:ARG:NH2	2.38	0.56
1:J:73:GLN:O	1:J:125:ARG:NH2	2.38	0.56
1:K:73:GLN:O	1:K:125:ARG:NH2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:73:GLN:O	1:L:125:ARG:NH2	2.38	0.56
1:O:73:GLN:O	1:O:125:ARG:NH2	2.38	0.56
1:Q:73:GLN:O	1:Q:125:ARG:NH2	2.38	0.56
3:1:181:DT:H2'	3:1:182:DA:C8	2.41	0.56
3:1:217:DT:H2'	3:1:218:DA:C8	2.40	0.56
1:D:73:GLN:O	1:D:125:ARG:NH2	2.38	0.56
1:N:73:GLN:O	1:N:125:ARG:NH2	2.38	0.56
1:P:73:GLN:O	1:P:125:ARG:NH2	2.38	0.56
3:2:181:DT:H2'	3:2:182:DA:C8	2.41	0.56
1:M:73:GLN:O	1:M:125:ARG:NH2	2.38	0.56
1:R:73:GLN:O	1:R:125:ARG:NH2	2.38	0.56
1:B:73:GLN:O	1:B:125:ARG:NH2	2.38	0.56
1:T:73:GLN:O	1:T:125:ARG:NH2	2.38	0.56
1:V:73:GLN:O	1:V:125:ARG:NH2	2.38	0.56
1:X:21:LYS:NZ	3:1:20:DA:OP2	2.39	0.56
1:X:94:LYS:HB2	3:2:303:DT:H5''	1.87	0.56
1:E:73:GLN:O	1:E:125:ARG:NH2	2.38	0.55
3:2:157:DT:H2'	3:2:158:DA:C8	2.41	0.55
1:T:94:LYS:HB2	3:2:279:DT:H5''	1.89	0.55
3:2:45:DT:O4	3:2:46:DA:N6	2.40	0.55
1:M:21:LYS:NZ	3:1:236:DA:OP2	2.39	0.55
1:M:20:ARG:NH2	3:1:235:DT:OP1	2.39	0.55
1:S:94:LYS:HB2	3:2:51:DT:H5''	1.90	0.55
3:1:45:DT:O4	3:1:46:DA:N6	2.40	0.54
3:1:291:DT:H2'	3:1:292:DA:C8	2.42	0.54
3:2:291:DT:H2'	3:2:292:DA:C8	2.42	0.54
3:1:157:DT:H2'	3:1:158:DA:C8	2.41	0.54
1:F:94:LYS:HB2	3:2:195:DT:H5''	1.90	0.54
1:X:160:THR:OG1	1:Z:73:GLN:NE2	2.40	0.54
3:2:215:DT:H72	3:2:216:DA:H62	1.73	0.53
1:G:21:LYS:NZ	3:1:200:DA:OP2	2.41	0.53
3:2:315:DT:H2'	3:2:316:DA:C8	2.44	0.53
1:P:94:LYS:HB2	3:2:255:DT:H5''	1.91	0.53
3:1:215:DT:H72	3:1:216:DA:H62	1.73	0.53
1:G:73:GLN:NE2	1:I:160:THR:OG1	2.42	0.53
1:R:21:LYS:NZ	3:1:56:DA:OP2	2.41	0.53
1:W:94:LYS:HB2	3:2:27:DT:H5''	1.91	0.53
3:1:135:DT:H2'	3:1:136:DA:C8	2.44	0.53
3:1:315:DT:H2'	3:1:316:DA:C8	2.44	0.53
3:1:167:DT:H72	3:1:168:DA:H62	1.74	0.52
1:V:51:GLU:OE1	3:1:47:DT:H5''	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2:167:DT:H72	3:2:168:DA:H62	1.74	0.52
1:F:20:ARG:NH2	3:1:127:DT:OP1	2.43	0.52
1:L:94:LYS:HB2	3:2:231:DT:H5''	1.92	0.52
3:2:135:DT:H2'	3:2:136:DA:C8	2.44	0.52
1:Q:94:LYS:HB2	3:2:63:DT:H5''	1.92	0.52
1:V:20:ARG:NH2	3:1:31:DT:OP1	2.43	0.52
3:2:123:DT:H2'	3:2:124:DA:C8	2.46	0.51
1:U:94:LYS:HB2	3:2:39:DT:H5''	1.93	0.51
3:1:123:DT:H2'	3:1:124:DA:C8	2.46	0.51
3:1:255:DT:H2'	3:1:256:DA:C8	2.46	0.51
1:E:94:LYS:HB2	3:2:135:DT:H5''	1.94	0.50
3:2:255:DT:H2'	3:2:256:DA:C8	2.46	0.50
3:1:327:DT:H2'	3:1:328:DA:C8	2.47	0.50
1:Q:21:LYS:NZ	3:1:260:DA:OP2	2.41	0.50
3:2:73:DT:H2'	3:2:74:DA:C8	2.46	0.50
3:1:73:DT:H2'	3:1:74:DA:C8	2.46	0.50
3:1:145:DT:H2'	3:1:146:DA:C8	2.47	0.50
3:2:327:DT:H2'	3:2:328:DA:C8	2.47	0.50
3:1:99:DT:H2'	3:1:100:DA:C8	2.47	0.50
3:2:27:DT:H2'	3:2:28:DA:C8	2.47	0.50
3:2:169:DT:H2'	3:2:170:DA:C8	2.47	0.50
1:M:94:LYS:HB2	3:2:87:DT:H5''	1.93	0.50
3:2:231:DT:H2'	3:2:232:DA:C8	2.47	0.50
1:R:94:LYS:HB2	3:2:267:DT:H5''	1.94	0.50
3:2:145:DT:H2'	3:2:146:DA:C8	2.47	0.50
3:2:243:DT:H2'	3:2:244:DA:C8	2.47	0.50
3:1:195:DT:H2'	3:1:196:DA:C8	2.47	0.50
3:2:195:DT:H2'	3:2:196:DA:C8	2.47	0.50
3:1:243:DT:H2'	3:1:244:DA:C8	2.47	0.49
3:1:231:DT:H2'	3:1:232:DA:C8	2.47	0.49
3:1:27:DT:H2'	3:1:28:DA:C8	2.47	0.49
3:2:303:DT:H2'	3:2:304:DA:C8	2.47	0.49
3:2:253:DT:H2'	3:2:254:DA:C8	2.48	0.49
3:2:267:DT:H2'	3:2:268:DA:C8	2.47	0.49
3:1:303:DT:H2'	3:1:304:DA:C8	2.47	0.49
3:2:99:DT:H2'	3:2:100:DA:C8	2.47	0.49
3:1:169:DT:H2'	3:1:170:DA:C8	2.47	0.49
3:2:39:DT:H2'	3:2:40:DA:C8	2.48	0.49
3:1:253:DT:H2'	3:1:254:DA:C8	2.48	0.49
3:1:87:DT:H2'	3:1:88:DA:C8	2.48	0.49
3:2:147:DT:H2'	3:2:148:DA:C8	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:1:39:DT:H2'	3:1:40:DA:C8	2.48	0.49
3:1:147:DT:H2'	3:1:148:DA:C8	2.48	0.49
3:1:267:DT:H2'	3:1:268:DA:C8	2.47	0.49
3:1:279:DT:H2'	3:1:280:DA:C8	2.48	0.49
1:C:29:TYR:OH	3:1:187:DT:O2	2.17	0.48
3:2:279:DT:H2'	3:2:280:DA:C8	2.48	0.48
1:I:29:TYR:OH	3:1:223:DT:O2	2.18	0.48
1:N:21:LYS:NZ	3:1:80:DA:OP2	2.43	0.48
3:1:51:DT:H2'	3:1:52:DA:C8	2.49	0.48
3:1:205:DT:H2'	3:1:206:DA:C8	2.49	0.48
3:2:46:DA:H2''	3:2:47:DT:H5'	1.96	0.48
3:2:63:DT:H2'	3:2:64:DA:C8	2.49	0.48
3:2:87:DT:H2'	3:2:88:DA:C8	2.48	0.48
1:A:29:TYR:OH	3:1:175:DT:O2	2.13	0.48
1:I:21:LYS:NZ	3:1:212:DA:OP2	2.40	0.48
3:1:15:DT:H2'	3:1:16:DA:C8	2.49	0.47
3:2:49:DT:H2'	3:2:50:DA:C8	2.49	0.47
3:2:51:DT:H2'	3:2:52:DA:C8	2.49	0.47
1:A:21:LYS:NZ	3:1:164:DA:OP2	2.43	0.47
3:1:63:DT:H2'	3:1:64:DA:C8	2.49	0.47
3:2:205:DT:H2'	3:2:206:DA:C8	2.49	0.47
3:2:289:DT:H2'	3:2:290:DA:C8	2.50	0.47
3:2:193:DT:H2'	3:2:194:DA:C8	2.50	0.47
1:P:21:LYS:NZ	3:1:68:DA:OP2	2.43	0.47
3:2:265:DT:H2'	3:2:266:DA:C8	2.50	0.47
3:2:97:DT:H2'	3:2:98:DA:C8	2.50	0.47
3:1:46:DA:H2''	3:1:47:DT:H5'	1.96	0.47
3:1:97:DT:H2'	3:1:98:DA:C8	2.50	0.47
3:1:313:DT:H2'	3:1:314:DA:C8	2.49	0.47
3:2:313:DT:H2'	3:2:314:DA:C8	2.49	0.47
3:1:49:DT:H2'	3:1:50:DA:C8	2.49	0.47
3:2:15:DT:H2'	3:2:16:DA:C8	2.50	0.47
3:2:229:DT:H2'	3:2:230:DA:C8	2.50	0.47
3:1:289:DT:H2'	3:1:290:DA:C8	2.50	0.47
3:1:229:DT:H2'	3:1:230:DA:C8	2.50	0.46
1:N:94:LYS:HB2	3:2:243:DT:H5''	1.98	0.46
1:O:21:LYS:NZ	3:1:248:DA:OP2	2.41	0.46
1:T:160:THR:OG1	1:V:73:GLN:NE2	2.48	0.46
1:E:29:TYR:OH	3:1:199:DT:O2	2.19	0.46
1:T:21:LYS:NZ	3:1:44:DA:OP2	2.47	0.46
3:2:25:DT:H2'	3:2:26:DA:C8	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:1:193:DT:H2'	3:1:194:DA:C8	2.50	0.46
3:1:265:DT:H2'	3:1:266:DA:C8	2.50	0.46
1:Y:94:LYS:HB2	3:2:15:DT:H5''	1.98	0.46
3:1:25:DT:H2'	3:1:26:DA:C8	2.50	0.46
3:2:107:DT:H72	3:2:108:DA:N6	2.31	0.46
1:V:21:LYS:NZ	3:1:32:DA:OP2	2.46	0.46
3:1:277:DT:H2'	3:1:278:DA:C8	2.51	0.46
3:1:121:DT:H2'	3:1:122:DA:C8	2.50	0.45
3:2:3:DT:H2'	3:2:4:DA:C8	2.51	0.45
3:2:121:DT:H2'	3:2:122:DA:C8	2.50	0.45
3:2:85:DT:H2'	3:2:86:DA:C8	2.52	0.45
1:Z:21:LYS:NZ	3:1:8:DA:OP2	2.44	0.45
3:1:3:DT:H2'	3:1:4:DA:C8	2.51	0.45
3:1:13:DT:H2'	3:1:14:DA:C8	2.52	0.45
3:1:133:DT:H2'	3:1:134:DA:C8	2.52	0.45
1:R:29:TYR:OH	3:1:67:DT:O2	2.23	0.45
3:2:13:DT:H2'	3:2:14:DA:C8	2.52	0.45
3:2:37:DT:H2'	3:2:38:DA:C8	2.52	0.45
1:Q:29:TYR:OH	3:1:271:DT:O2	2.23	0.44
1:U:21:LYS:NZ	3:1:284:DA:OP2	2.45	0.44
3:1:183:DT:H2'	3:1:184:DA:C8	2.53	0.44
3:2:61:DT:H2'	3:2:62:DA:C8	2.53	0.44
3:2:277:DT:H2'	3:2:278:DA:C8	2.51	0.44
3:2:133:DT:H2'	3:2:134:DA:C8	2.52	0.44
3:1:219:DT:H2'	3:1:220:DA:C8	2.53	0.44
3:2:183:DT:H2'	3:2:184:DA:C8	2.53	0.44
3:2:219:DT:H2'	3:2:220:DA:C8	2.53	0.44
3:1:37:DT:H2'	3:1:38:DA:C8	2.52	0.44
1:B:48:HIS:CE1	1:B:52:ASN:HD21	2.37	0.43
1:D:50:VAL:HG11	1:D:93:LEU:HD22	2.01	0.43
1:Y:48:HIS:CE1	1:Y:52:ASN:HD21	2.36	0.43
1:Y:50:VAL:HG11	1:Y:93:LEU:HD22	2.01	0.43
3:2:111:DT:H2'	3:2:112:DA:C8	2.53	0.43
1:D:48:HIS:CE1	1:D:52:ASN:HD21	2.37	0.43
1:J:50:VAL:HG11	1:J:93:LEU:HD22	2.01	0.43
1:N:88:ILE:HD12	1:N:113:ALA:HB2	2.00	0.43
1:P:48:HIS:CE1	1:P:52:ASN:HD21	2.37	0.43
3:2:159:DT:H2'	3:2:160:DA:C8	2.54	0.43
1:G:48:HIS:CE1	1:G:52:ASN:HD21	2.37	0.43
1:J:48:HIS:CE1	1:J:52:ASN:HD21	2.37	0.43
1:N:48:HIS:CE1	1:N:52:ASN:HD21	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:48:HIS:CE1	1:R:52:ASN:HD21	2.37	0.43
1:S:50:VAL:HG11	1:S:93:LEU:HD22	2.01	0.43
1:U:48:HIS:CE1	1:U:52:ASN:HD21	2.37	0.43
3:1:311:DT:H72	3:1:312:DA:N6	2.33	0.43
3:2:189:DT:H2'	3:2:190:DA:C8	2.54	0.43
3:2:301:DT:H2'	3:2:302:DA:C8	2.53	0.43
1:C:50:VAL:HG11	1:C:93:LEU:HD22	2.01	0.43
1:E:48:HIS:CE1	1:E:52:ASN:HD21	2.37	0.43
1:I:48:HIS:CE1	1:I:52:ASN:HD21	2.37	0.43
1:L:48:HIS:CE1	1:L:52:ASN:HD21	2.37	0.43
1:N:50:VAL:HG11	1:N:93:LEU:HD22	2.01	0.43
1:P:50:VAL:HG11	1:P:93:LEU:HD22	2.01	0.43
1:U:88:ILE:HD12	1:U:113:ALA:HB2	2.00	0.43
1:W:48:HIS:CE1	1:W:52:ASN:HD21	2.37	0.43
3:1:61:DT:H2'	3:1:62:DA:C8	2.53	0.43
3:1:85:DT:H2'	3:1:86:DA:C8	2.52	0.43
1:B:50:VAL:HG11	1:B:93:LEU:HD22	2.01	0.43
1:N:93:LEU:HD23	1:N:93:LEU:HA	4.59	0.43
1:O:88:ILE:HD12	1:O:113:ALA:HB2	2.00	0.43
1:T:48:HIS:CE1	1:T:52:ASN:HD21	2.37	0.43
1:T:50:VAL:HG11	1:T:93:LEU:HD22	2.01	0.43
1:T:88:ILE:HD12	1:T:113:ALA:HB2	2.00	0.43
1:W:50:VAL:HG11	1:W:93:LEU:HD22	2.01	0.43
1:W:88:ILE:HD12	1:W:113:ALA:HB2	2.00	0.43
1:Z:29:TYR:OH	3:1:19:DT:O2	2.17	0.43
1:H:50:VAL:HG11	1:H:93:LEU:HD22	2.01	0.43
1:Q:88:ILE:HD12	1:Q:113:ALA:HB2	2.00	0.43
1:S:48:HIS:CE1	1:S:52:ASN:HD21	2.37	0.43
3:1:189:DT:H2'	3:1:190:DA:C8	2.54	0.43
1:T:36:ARG:HG3	3:2:281:DT:H2''	2.01	0.43
3:1:111:DT:H2'	3:1:112:DA:C8	2.53	0.43
3:2:177:DT:H2'	3:2:178:DA:C8	2.54	0.43
1:H:48:HIS:CE1	1:H:52:ASN:HD21	2.37	0.43
1:H:88:ILE:HD12	1:H:113:ALA:HB2	2.01	0.43
1:K:48:HIS:CE1	1:K:52:ASN:HD21	2.37	0.43
3:2:250:DA:N7	3:2:251:DT:H72	2.34	0.43
1:C:48:HIS:CE1	1:C:52:ASN:HD21	2.37	0.43
1:I:50:VAL:HG11	1:I:93:LEU:HD22	2.01	0.43
1:M:50:VAL:HG11	1:M:93:LEU:HD22	2.01	0.43
1:O:48:HIS:CE1	1:O:52:ASN:HD21	2.37	0.43
1:Q:36:ARG:HG3	3:2:65:DT:H2''	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:48:HIS:CE1	1:Q:52:ASN:HD21	2.37	0.43
1:R:93:LEU:HD23	1:R:93:LEU:HA	4.59	0.43
1:V:48:HIS:CE1	1:V:52:ASN:HD21	2.37	0.43
1:Z:50:VAL:HG11	1:Z:93:LEU:HD22	2.01	0.43
3:1:250:DA:N7	3:1:251:DT:H72	2.34	0.43
1:A:48:HIS:CE1	1:A:52:ASN:HD21	2.37	0.42
1:F:48:HIS:CE1	1:F:52:ASN:HD21	2.37	0.42
1:L:88:ILE:HD12	1:L:113:ALA:HB2	2.00	0.42
1:U:93:LEU:HD23	1:U:93:LEU:HA	4.59	0.42
3:1:301:DT:H2'	3:1:302:DA:C8	2.53	0.42
1:C:88:ILE:HD12	1:C:113:ALA:HB2	2.01	0.42
1:E:88:ILE:HD12	1:E:113:ALA:HB2	2.00	0.42
1:G:50:VAL:HG11	1:G:93:LEU:HD22	2.01	0.42
1:J:88:ILE:HD12	1:J:113:ALA:HB2	2.00	0.42
1:P:88:ILE:HD12	1:P:113:ALA:HB2	2.01	0.42
1:R:88:ILE:HD12	1:R:113:ALA:HB2	2.00	0.42
1:U:36:ARG:HG3	3:2:41:DT:H2''	2.01	0.42
1:U:50:VAL:HG11	1:U:93:LEU:HD22	2.01	0.42
1:A:50:VAL:HG11	1:A:93:LEU:HD22	2.01	0.42
1:A:88:ILE:HD12	1:A:113:ALA:HB2	2.00	0.42
1:O:50:VAL:HG11	1:O:93:LEU:HD22	2.01	0.42
1:Q:50:VAL:HG11	1:Q:93:LEU:HD22	2.01	0.42
1:Y:36:ARG:HG3	3:2:17:DT:H2''	2.02	0.42
3:1:177:DT:H2'	3:1:178:DA:C8	2.54	0.42
3:2:311:DT:H72	3:2:312:DA:N6	2.34	0.42
1:M:48:HIS:CE1	1:M:52:ASN:HD21	2.36	0.42
1:V:50:VAL:HG11	1:V:93:LEU:HD22	2.01	0.42
1:V:88:ILE:HD12	1:V:113:ALA:HB2	2.01	0.42
1:W:29:TYR:OH	3:1:307:DT:O2	2.23	0.42
3:1:159:DT:H2'	3:1:160:DA:C8	2.54	0.42
1:F:50:VAL:HG11	1:F:93:LEU:HD22	2.01	0.42
1:Q:93:LEU:HD23	1:Q:93:LEU:HA	4.59	0.42
1:B:18:TYR:OH	3:1:161:DT:O2	2.31	0.42
1:E:50:VAL:HG11	1:E:93:LEU:HD22	2.01	0.42
1:F:88:ILE:HD12	1:F:113:ALA:HB2	2.01	0.42
1:G:88:ILE:HD12	1:G:113:ALA:HB2	2.00	0.42
1:B:88:ILE:HD12	1:B:113:ALA:HB2	2.00	0.42
1:L:50:VAL:HG11	1:L:93:LEU:HD22	2.01	0.42
1:X:48:HIS:CE1	1:X:52:ASN:HD21	2.37	0.42
1:X:50:VAL:HG11	1:X:93:LEU:HD22	2.01	0.42
1:S:88:ILE:HD12	1:S:113:ALA:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:48:HIS:CE1	1:Z:52:ASN:HD21	2.37	0.42
1:Z:88:ILE:HD12	1:Z:113:ALA:HB2	2.01	0.42
3:1:149:DT:H2'	3:1:150:DA:C8	2.55	0.42
1:I:88:ILE:HD12	1:I:113:ALA:HB2	2.01	0.42
1:R:50:VAL:HG11	1:R:93:LEU:HD22	2.01	0.42
3:2:153:DT:H2'	3:2:154:DA:C8	2.55	0.42
1:B:21:LYS:NZ	3:1:152:DA:OP2	2.41	0.42
1:F:21:LYS:NZ	3:1:128:DA:OP2	2.43	0.42
1:K:88:ILE:HD12	1:K:113:ALA:HB2	2.01	0.42
1:S:36:ARG:HG3	3:2:53:DT:H2''	2.02	0.42
1:L:21:LYS:NZ	3:1:92:DA:OP2	2.48	0.41
1:L:93:LEU:HD23	1:L:93:LEU:HA	4.59	0.41
1:M:88:ILE:HD12	1:M:113:ALA:HB2	2.00	0.41
1:Y:88:ILE:HD12	1:Y:113:ALA:HB2	2.00	0.41
1:D:88:ILE:HD12	1:D:113:ALA:HB2	2.00	0.41
1:K:93:LEU:HD23	1:K:93:LEU:HA	4.59	0.41
3:1:153:DT:H2'	3:1:154:DA:C8	2.55	0.41
3:2:149:DT:H2'	3:2:150:DA:C8	2.55	0.41
3:2:241:DT:H2'	3:2:242:DA:C8	2.54	0.41
1:D:29:TYR:OH	3:1:151:DT:O2	2.31	0.41
1:J:107:ILE:HG22	1:J:161:PRO:HG3	2.02	0.41
1:X:88:ILE:HD12	1:X:113:ALA:HB2	2.00	0.41
1:Z:107:ILE:HG22	1:Z:161:PRO:HG3	2.02	0.41
1:E:94:LYS:HD3	3:2:135:DT:H5''	2.02	0.41
1:Y:107:ILE:HG22	1:Y:161:PRO:HG3	2.02	0.41
3:1:241:DT:H2'	3:1:242:DA:C8	2.54	0.41
1:F:107:ILE:HG22	1:F:161:PRO:HG3	2.02	0.41
1:H:107:ILE:HG22	1:H:161:PRO:HG3	2.02	0.41
1:K:50:VAL:HG11	1:K:93:LEU:HD22	2.01	0.41
1:X:147:PRO:HG3	1:X:165:LEU:HD13	2.03	0.41
3:1:123:DT:C2	3:2:214:DA:H2	2.39	0.41
1:L:107:ILE:HG22	1:L:161:PRO:HG3	2.02	0.41
1:M:107:ILE:HG22	1:M:161:PRO:HG3	2.02	0.41
1:V:107:ILE:HD11	1:X:78:VAL:HG22	2.03	0.41
1:V:147:PRO:HG3	1:V:165:LEU:HD13	2.03	0.41
1:W:107:ILE:HG22	1:W:161:PRO:HG3	2.02	0.41
1:Z:147:PRO:HG3	1:Z:165:LEU:HD13	2.03	0.41
3:1:109:DT:H2'	3:1:110:DA:C8	2.56	0.41
1:G:107:ILE:HG22	1:G:161:PRO:HG3	2.02	0.41
1:N:107:ILE:HG22	1:N:161:PRO:HG3	2.02	0.41
1:V:107:ILE:HG22	1:V:161:PRO:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2:109:DT:H2'	3:2:110:DA:C8	2.56	0.41
3:2:207:DT:H2'	3:2:208:DA:C8	2.56	0.41
1:D:107:ILE:HG22	1:D:161:PRO:HG3	2.02	0.41
1:G:147:PRO:HG3	1:G:165:LEU:HD13	2.03	0.41
1:I:107:ILE:HG22	1:I:161:PRO:HG3	2.02	0.41
1:K:147:PRO:HG3	1:K:165:LEU:HD13	2.03	0.41
1:O:147:PRO:HG3	1:O:165:LEU:HD13	2.03	0.41
1:P:107:ILE:HG22	1:P:161:PRO:HG3	2.02	0.41
1:R:147:PRO:HG3	1:R:165:LEU:HD13	2.03	0.41
1:T:93:LEU:HD23	1:T:93:LEU:HA	4.59	0.41
1:T:147:PRO:HG3	1:T:165:LEU:HD13	2.03	0.41
1:U:107:ILE:HG22	1:U:161:PRO:HG3	2.02	0.41
1:I:147:PRO:HG3	1:I:165:LEU:HD13	2.03	0.41
1:M:147:PRO:HG3	1:M:165:LEU:HD13	2.03	0.41
1:P:147:PRO:HG3	1:P:165:LEU:HD13	2.03	0.41
3:1:207:DT:H2'	3:1:208:DA:C8	2.56	0.41
3:1:310:DA:H2	3:2:27:DT:C2	2.39	0.41
1:B:107:ILE:HG22	1:B:161:PRO:HG3	2.02	0.40
1:E:147:PRO:HG3	1:E:165:LEU:HD13	2.03	0.40
3:2:94:DA:N7	3:2:95:DT:H72	2.37	0.40
1:C:107:ILE:HG22	1:C:161:PRO:HG3	2.02	0.40
1:K:107:ILE:HG22	1:K:161:PRO:HG3	2.02	0.40
1:Q:147:PRO:HG3	1:Q:165:LEU:HD13	2.03	0.40
3:1:286:DA:H2	3:2:51:DT:C2	2.40	0.40
1:C:147:PRO:HG3	1:C:165:LEU:HD13	2.03	0.40
1:N:147:PRO:HG3	1:N:165:LEU:HD13	2.03	0.40
1:P:93:LEU:HD23	1:P:93:LEU:HA	4.59	0.40
1:T:107:ILE:HG22	1:T:161:PRO:HG3	2.02	0.40
1:U:62:ILE:HD13	1:U:62:ILE:HA	1.96	0.40
1:G:29:TYR:OH	3:1:211:DT:O2	2.20	0.40
1:S:107:ILE:HG22	1:S:161:PRO:HG3	2.02	0.40
1:S:147:PRO:HG3	1:S:165:LEU:HD13	2.03	0.40
3:1:298:DA:H2	3:2:39:DT:C2	2.39	0.40
1:R:107:ILE:HG22	1:R:161:PRO:HG3	2.02	0.40
1:U:147:PRO:HG3	1:U:165:LEU:HD13	2.03	0.40
3:1:82:DA:H2	3:2:255:DT:C2	2.40	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	192/199 (96%)	172 (90%)	19 (10%)	1 (0%)	25	57
1	B	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	C	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	D	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	E	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	F	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	G	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	H	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	I	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	J	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	K	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	L	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	M	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	N	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	O	192/199 (96%)	172 (90%)	19 (10%)	1 (0%)	25	57
1	P	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	Q	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	R	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	S	192/199 (96%)	172 (90%)	19 (10%)	1 (0%)	25	57
1	T	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	U	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	V	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	W	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
1	X	192/199 (96%)	172 (90%)	19 (10%)	1 (0%)	25	57
1	Y	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Z	192/199 (96%)	171 (89%)	20 (10%)	1 (0%)	25	57
2	a	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	b	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	c	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	d	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	e	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	f	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	g	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	h	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	i	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	j	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	k	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	l	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	m	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	n	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	o	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	p	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	q	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	r	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	s	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	t	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	u	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	v	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	w	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	x	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	y	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
2	z	129/137 (94%)	122 (95%)	7 (5%)	0	100	100
All	All	8346/8736 (96%)	7622 (91%)	698 (8%)	26 (0%)	38	67

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	PRO

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Mol	Chain	Res	Type
1	B	76	PRO
1	C	76	PRO
1	D	76	PRO
1	E	76	PRO
1	F	76	PRO
1	G	76	PRO
1	H	76	PRO
1	I	76	PRO
1	J	76	PRO
1	K	76	PRO
1	L	76	PRO
1	M	76	PRO
1	N	76	PRO
1	O	76	PRO
1	P	76	PRO
1	Q	76	PRO
1	R	76	PRO
1	S	76	PRO
1	T	76	PRO
1	U	76	PRO
1	V	76	PRO
1	W	76	PRO
1	X	76	PRO
1	Y	76	PRO
1	Z	76	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	B	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	C	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	D	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	E	158/168 (94%)	157 (99%)	1 (1%)	84	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	G	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	H	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	I	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	J	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	K	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	L	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	M	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	N	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	O	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	P	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	Q	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	R	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	S	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	T	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	U	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	V	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	W	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	X	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	Y	158/168 (94%)	157 (99%)	1 (1%)	84	90
1	Z	158/168 (94%)	157 (99%)	1 (1%)	84	90
2	a	114/121 (94%)	114 (100%)	0	100	100
2	b	114/121 (94%)	114 (100%)	0	100	100
2	c	114/121 (94%)	114 (100%)	0	100	100
2	d	114/121 (94%)	114 (100%)	0	100	100
2	e	114/121 (94%)	114 (100%)	0	100	100
2	f	114/121 (94%)	114 (100%)	0	100	100
2	g	114/121 (94%)	114 (100%)	0	100	100
2	h	114/121 (94%)	114 (100%)	0	100	100
2	i	114/121 (94%)	114 (100%)	0	100	100
2	j	114/121 (94%)	114 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	k	114/121 (94%)	114 (100%)	0	100	100
2	l	114/121 (94%)	114 (100%)	0	100	100
2	m	114/121 (94%)	114 (100%)	0	100	100
2	n	114/121 (94%)	114 (100%)	0	100	100
2	o	114/121 (94%)	114 (100%)	0	100	100
2	p	114/121 (94%)	114 (100%)	0	100	100
2	q	114/121 (94%)	114 (100%)	0	100	100
2	r	114/121 (94%)	114 (100%)	0	100	100
2	s	114/121 (94%)	114 (100%)	0	100	100
2	t	114/121 (94%)	114 (100%)	0	100	100
2	u	114/121 (94%)	114 (100%)	0	100	100
2	v	114/121 (94%)	114 (100%)	0	100	100
2	w	114/121 (94%)	114 (100%)	0	100	100
2	x	114/121 (94%)	114 (100%)	0	100	100
2	y	114/121 (94%)	114 (100%)	0	100	100
2	z	114/121 (94%)	114 (100%)	0	100	100
All	All	7072/7514 (94%)	7046 (100%)	26 (0%)	88	93

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	LEU
1	B	75	LEU
1	C	75	LEU
1	D	75	LEU
1	E	75	LEU
1	F	75	LEU
1	G	75	LEU
1	H	75	LEU
1	I	75	LEU
1	J	75	LEU
1	K	75	LEU
1	L	75	LEU
1	M	75	LEU
1	N	75	LEU
1	O	75	LEU

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Mol	Chain	Res	Type
1	P	75	LEU
1	Q	75	LEU
1	R	75	LEU
1	S	75	LEU
1	T	75	LEU
1	U	75	LEU
1	V	75	LEU
1	W	75	LEU
1	X	75	LEU
1	Y	75	LEU
1	Z	75	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (103) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	120	ASN
1	B	52	ASN
1	B	120	ASN
1	C	52	ASN
1	C	120	ASN
1	D	52	ASN
1	D	120	ASN
1	E	52	ASN
1	E	120	ASN
1	F	52	ASN
1	F	120	ASN
1	G	52	ASN
1	G	120	ASN
1	H	52	ASN
1	H	120	ASN
1	I	52	ASN
1	I	120	ASN
1	J	52	ASN
1	J	120	ASN
1	K	52	ASN
1	K	120	ASN
1	L	52	ASN
1	L	120	ASN
1	M	52	ASN
1	M	120	ASN
1	N	52	ASN

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Mol	Chain	Res	Type
1	N	120	ASN
1	O	52	ASN
1	O	120	ASN
1	P	52	ASN
1	P	120	ASN
1	Q	52	ASN
1	Q	120	ASN
1	R	52	ASN
1	R	120	ASN
1	S	52	ASN
1	S	120	ASN
1	T	52	ASN
1	T	120	ASN
1	U	52	ASN
1	U	120	ASN
1	V	52	ASN
1	V	120	ASN
1	W	52	ASN
1	W	120	ASN
1	X	52	ASN
1	X	120	ASN
1	Y	52	ASN
1	Z	52	ASN
1	Z	120	ASN
2	a	41	ASN
2	a	44	ASN
2	b	41	ASN
2	b	44	ASN
2	c	41	ASN
2	c	44	ASN
2	d	41	ASN
2	d	44	ASN
2	e	41	ASN
2	e	44	ASN
2	f	41	ASN
2	f	44	ASN
2	g	41	ASN
2	g	44	ASN
2	h	41	ASN
2	h	44	ASN
2	i	41	ASN
2	i	44	ASN

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Mol	Chain	Res	Type
2	j	41	ASN
2	j	44	ASN
2	k	41	ASN
2	k	44	ASN
2	l	41	ASN
2	l	44	ASN
2	m	41	ASN
2	m	44	ASN
2	n	41	ASN
2	n	44	ASN
2	o	41	ASN
2	o	44	ASN
2	p	41	ASN
2	p	44	ASN
2	q	41	ASN
2	q	44	ASN
2	r	41	ASN
2	r	44	ASN
2	s	41	ASN
2	s	44	ASN
2	t	41	ASN
2	t	44	ASN
2	u	41	ASN
2	u	44	ASN
2	v	41	ASN
2	v	44	ASN
2	w	41	ASN
2	w	44	ASN
2	x	41	ASN
2	x	44	ASN
2	y	41	ASN
2	y	44	ASN
2	z	41	ASN
2	z	44	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 [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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.

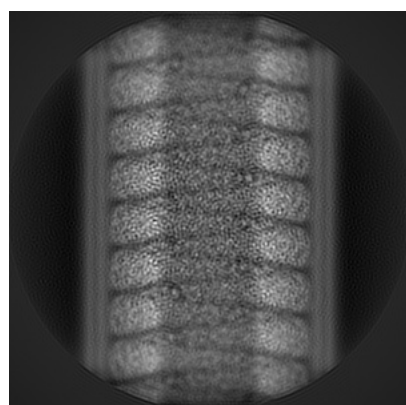
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-7797. These allow visual inspection of the internal detail of the map and identification of artifacts.

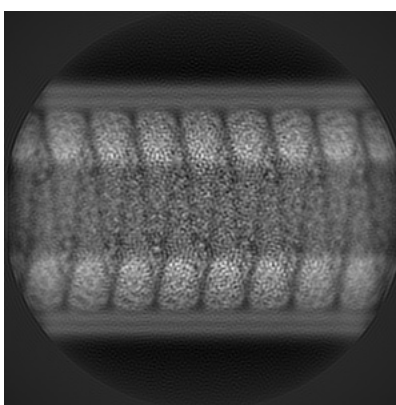
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

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

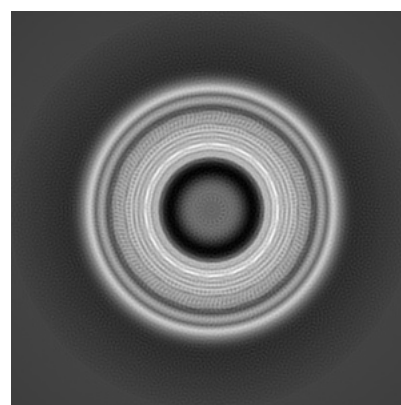
#### 6.1.1 Primary map



X



Y

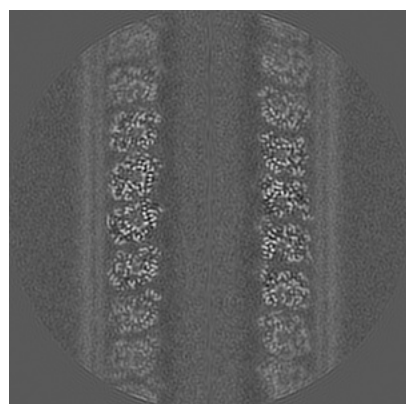


Z

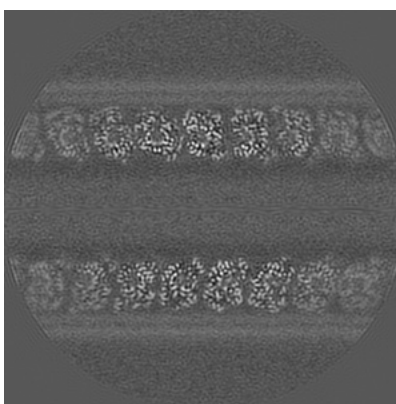
The images above show the map projected in three orthogonal directions.

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

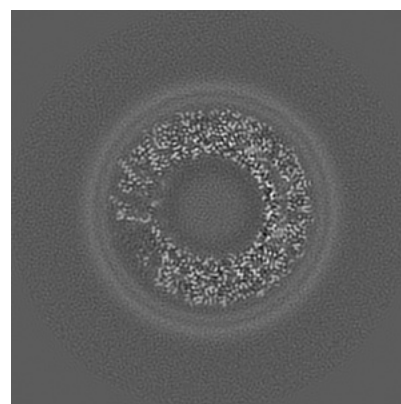
#### 6.2.1 Primary map



X Index: 192



Y Index: 192

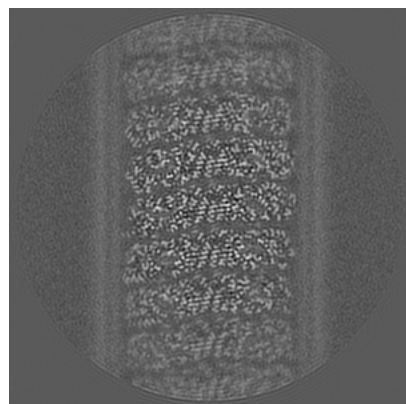


Z Index: 192

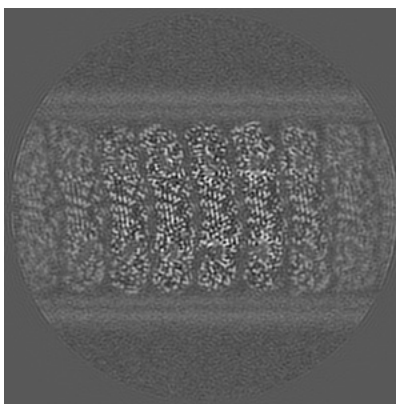
The images above show central slices of the map in three orthogonal directions.

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

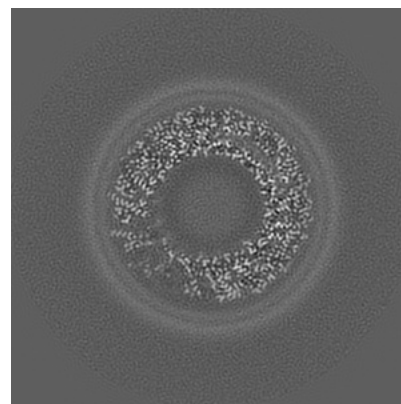
### 6.3.1 Primary map



X Index: 246



Y Index: 246

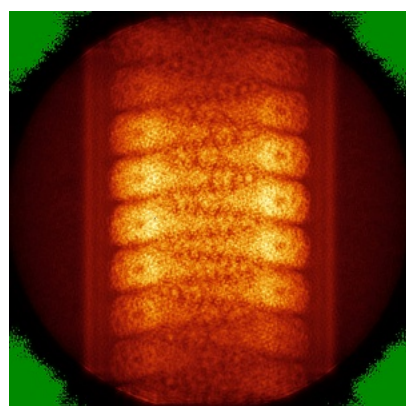


Z Index: 196

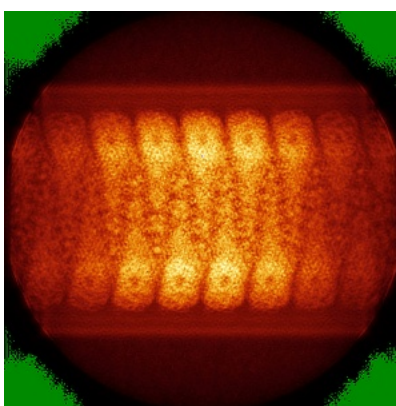
The images above show the largest variance slices of the map in three orthogonal directions.

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

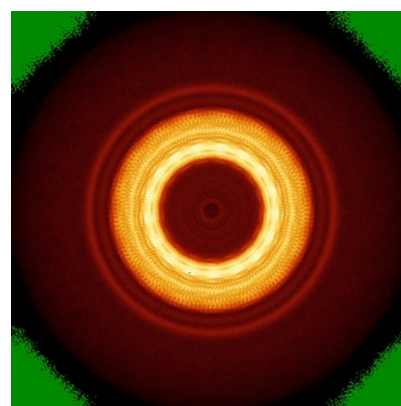
### 6.4.1 Primary map



X



Y



Z

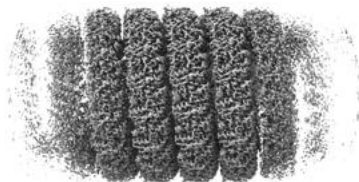
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

## 6.5 Orthogonal surface views [i](#)

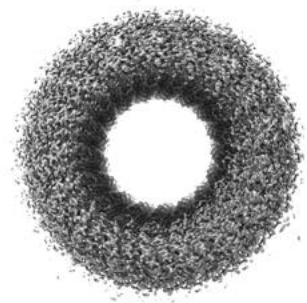
### 6.5.1 Primary map



X



Y



Z

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.

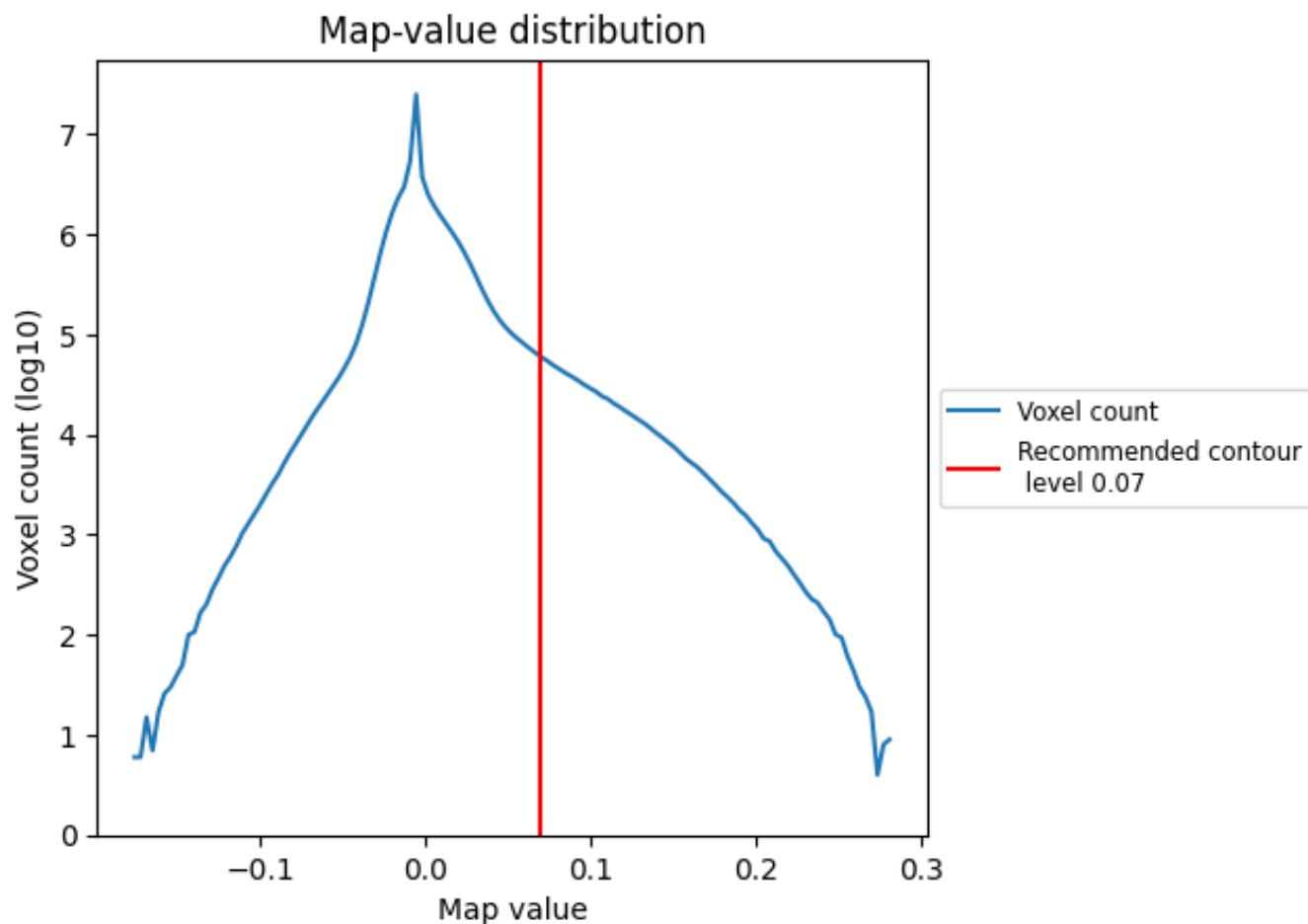
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

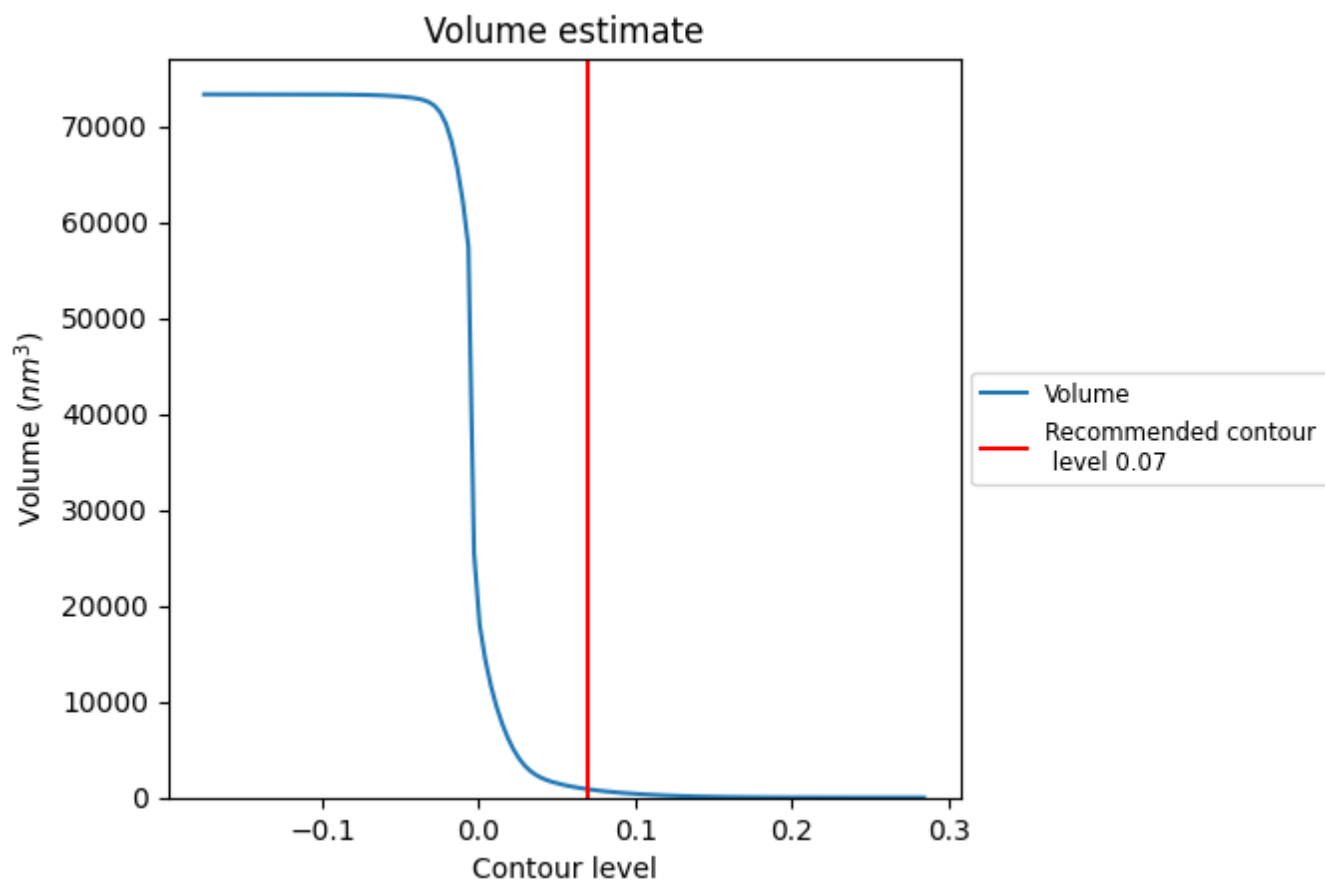
This section contains the results of statistical analysis of the map.

### 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](#)

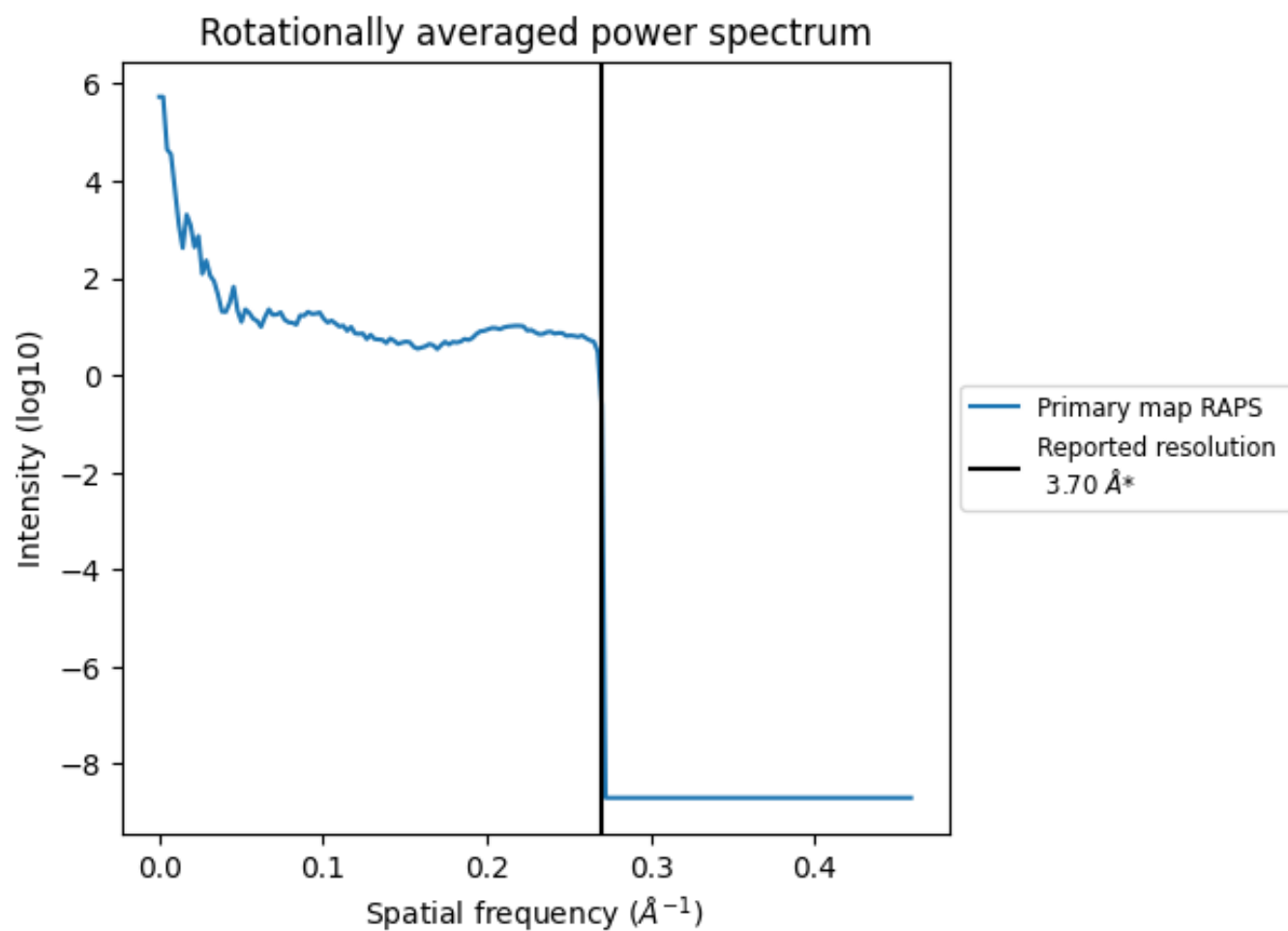


The volume at the recommended contour level is 864 nm<sup>3</sup>; this corresponds to an approximate mass of 780 kDa.

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 Å<sup>-1</sup>

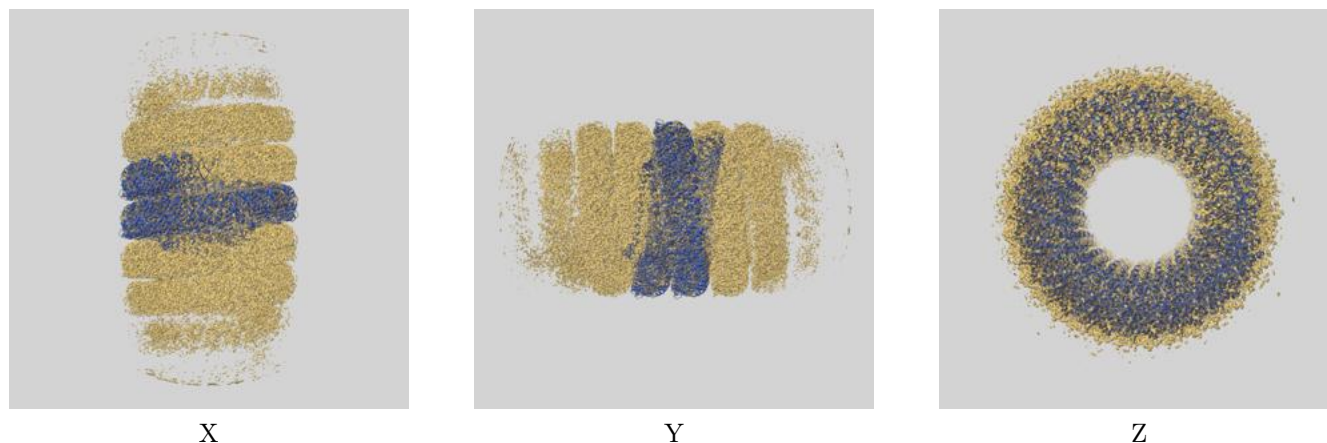
## 8 Fourier-Shell correlation ⓘ

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

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

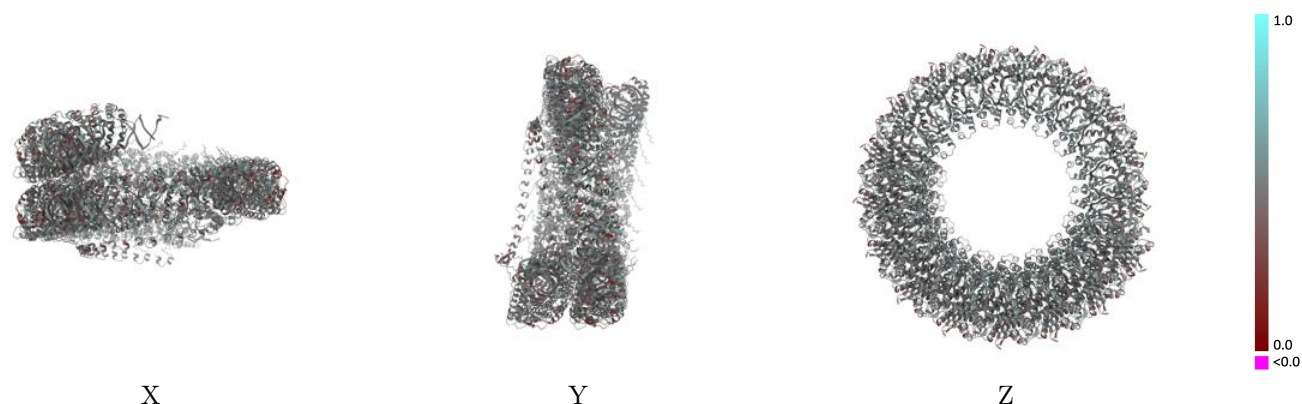
This section contains information regarding the fit between EMDB map EMD-7797 and PDB model 6D5F. Per-residue inclusion information can be found in section [3](#) on page [8](#).

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



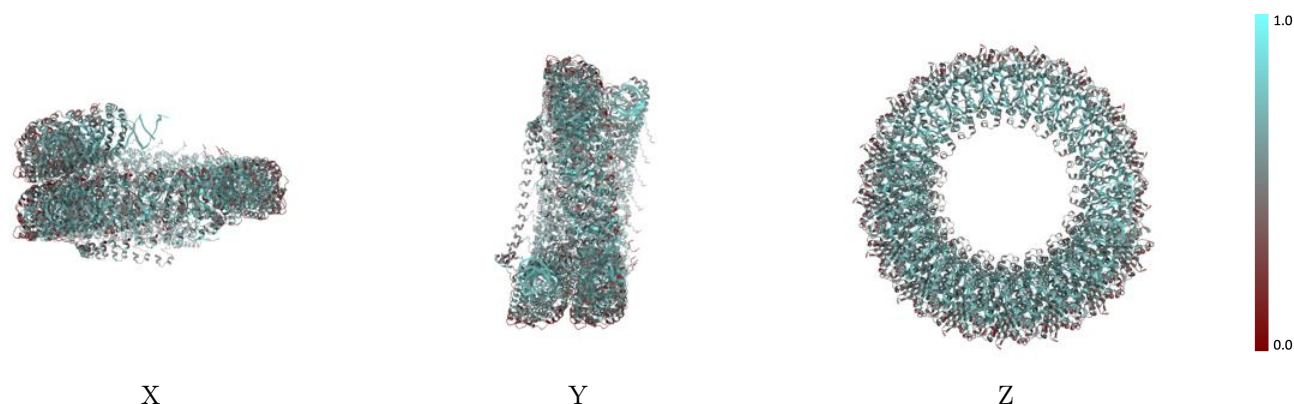
The images above show the 3D surface view of the map at the recommended contour level 0.07 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



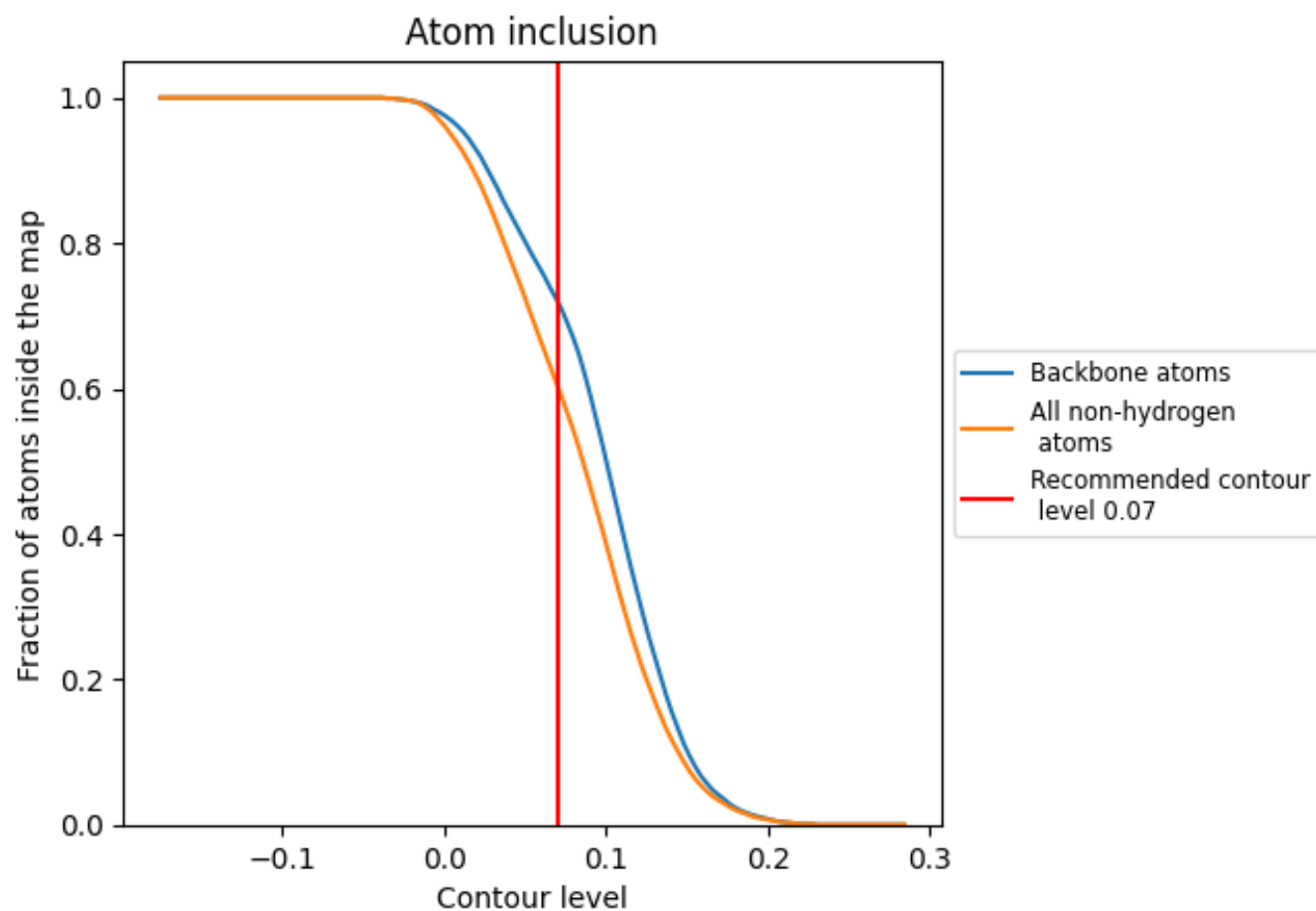
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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, 72% of all backbone atoms, 60% 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	 0.6030	 0.4790
1	 0.7620	 0.4600
2	 0.7570	 0.4750
A	 0.5630	 0.4740
B	 0.5590	 0.4730
C	 0.5500	 0.4730
D	 0.5570	 0.4710
E	 0.5570	 0.4740
F	 0.5620	 0.4740
G	 0.5520	 0.4750
H	 0.5560	 0.4710
I	 0.5540	 0.4740
J	 0.5550	 0.4720
K	 0.5510	 0.4740
L	 0.5580	 0.4710
M	 0.5480	 0.4740
N	 0.5570	 0.4710
O	 0.5480	 0.4730
P	 0.5480	 0.4680
Q	 0.5460	 0.4720
R	 0.5540	 0.4700
S	 0.5420	 0.4720
T	 0.5520	 0.4690
U	 0.5400	 0.4710
V	 0.5450	 0.4690
W	 0.5410	 0.4730
X	 0.5410	 0.4680
Y	 0.5380	 0.4730
Z	 0.5330	 0.4660
a	 0.6100	 0.4950
b	 0.6000	 0.4970
c	 0.6030	 0.4970
d	 0.6020	 0.4950
e	 0.6080	 0.4990
f	 0.6070	 0.4950



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Chain	Atom inclusion	Q-score
g	 0.5960	 0.4970
h	 0.6070	 0.4960
i	 0.6040	 0.4950
j	 0.6060	 0.4950
k	 0.6090	 0.4960
l	 0.6040	 0.4970
m	 0.6040	 0.4940
n	 0.6060	 0.4950
o	 0.6030	 0.4960
p	 0.6080	 0.4960
q	 0.5950	 0.4970
r	 0.6020	 0.4970
s	 0.5960	 0.4890
t	 0.5930	 0.4980
u	 0.5890	 0.4900
v	 0.5870	 0.4950
w	 0.5890	 0.4890
x	 0.5820	 0.4940
y	 0.5820	 0.4880
z	 0.5840	 0.4940