



## Full wwPDB EM Validation Report ⓘ

Jun 25, 2025 – 03:57 am BST

PDB ID : 7AMY / pdb\_00007amy  
EMDB ID : EMD-11827  
Title : Nonameric cytoplasmic domain of FlhA from *Vibrio parahaemolyticus*  
Authors : Kuhlen, L.; Johnson, S.; Lea, S.  
Deposited on : 2020-10-09  
Resolution : 3.75 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.dev118  
MolProbity : 4-5-2 with Phenix2.0rc1  
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) : 2.44

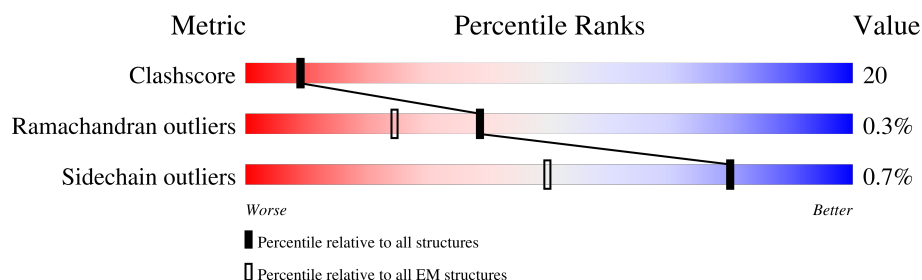
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.75 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	702	<div> <div>24%</div> <div>30% 20% 50%</div> </div>
1	B	702	<div> <div>23%</div> <div>30% 21% 50%</div> </div>
1	C	702	<div> <div>23%</div> <div>30% 20% 50%</div> </div>
1	D	702	<div> <div>23%</div> <div>30% 20% 50%</div> </div>
1	E	702	<div> <div>23%</div> <div>29% 21% 50%</div> </div>
1	F	702	<div> <div>23%</div> <div>30% 20% 50%</div> </div>
1	G	702	<div> <div>23%</div> <div>30% 20% 50%</div> </div>
1	H	702	<div> <div>23%</div> <div>30% 20% 50%</div> </div>

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Mol	Chain	Length	Quality of chain
1	I	702	<div><div><div></div><div></div><div></div></div><div>23%30%21%50%</div></div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 24849 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 Flagellar biosynthesis protein FlhA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	B	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	C	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	D	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	E	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	F	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	G	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	H	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		
1	I	353	Total	C	N	O	S	0	0
			2761	1750	480	523	8		

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	696	GLU	-	expression tag	UNP A0A0F5SXE4
A	697	ASN	-	expression tag	UNP A0A0F5SXE4
A	698	LEU	-	expression tag	UNP A0A0F5SXE4
A	699	TYR	-	expression tag	UNP A0A0F5SXE4
A	700	PHE	-	expression tag	UNP A0A0F5SXE4
A	701	GLN	-	expression tag	UNP A0A0F5SXE4
B	696	GLU	-	expression tag	UNP A0A0F5SXE4
B	697	ASN	-	expression tag	UNP A0A0F5SXE4
B	698	LEU	-	expression tag	UNP A0A0F5SXE4
B	699	TYR	-	expression tag	UNP A0A0F5SXE4
B	700	PHE	-	expression tag	UNP A0A0F5SXE4
B	701	GLN	-	expression tag	UNP A0A0F5SXE4

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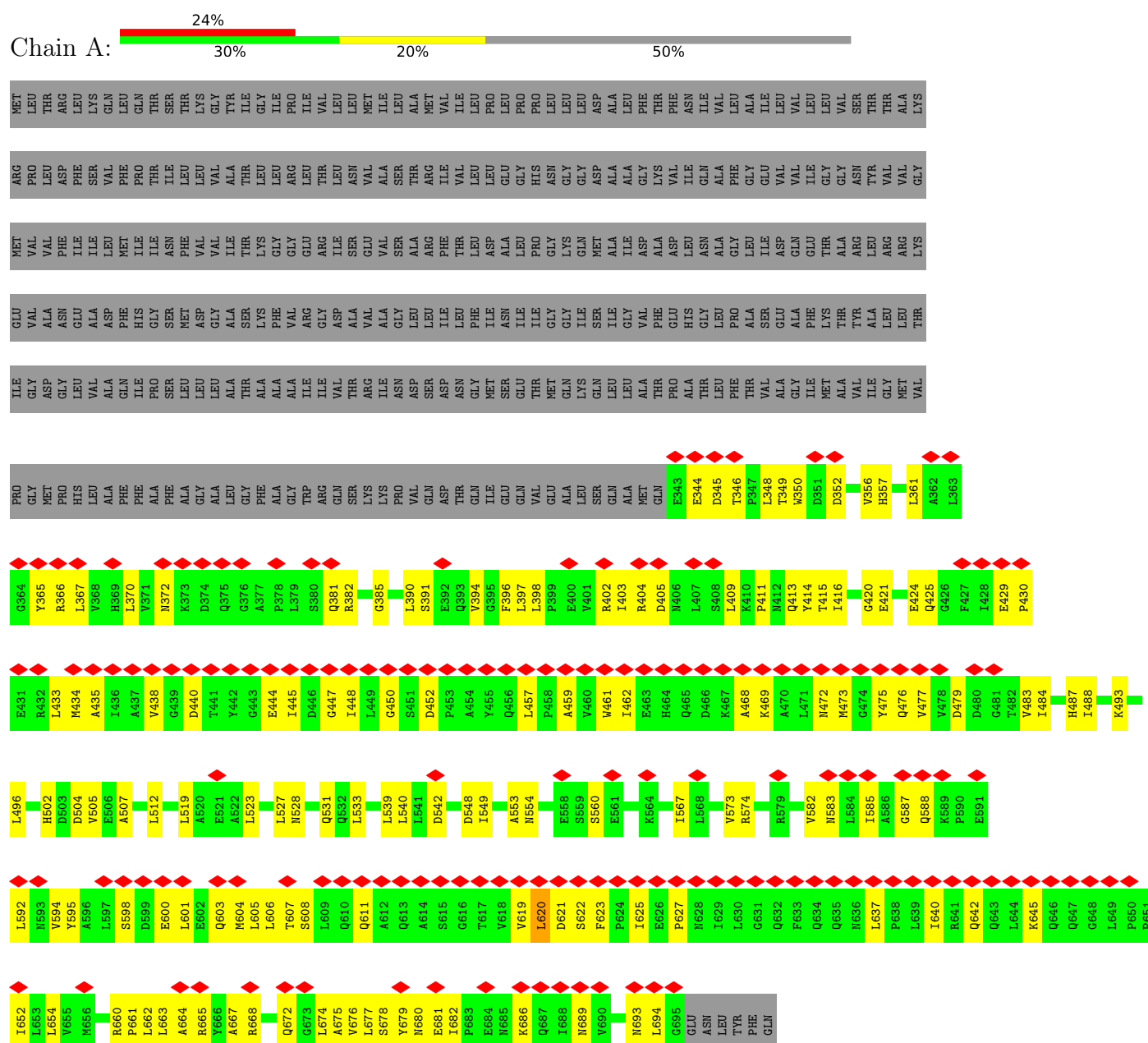
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Chain	Residue	Modelled	Actual	Comment	Reference
C	696	GLU	-	expression tag	UNP A0A0F5SXE4
C	697	ASN	-	expression tag	UNP A0A0F5SXE4
C	698	LEU	-	expression tag	UNP A0A0F5SXE4
C	699	TYR	-	expression tag	UNP A0A0F5SXE4
C	700	PHE	-	expression tag	UNP A0A0F5SXE4
C	701	GLN	-	expression tag	UNP A0A0F5SXE4
D	696	GLU	-	expression tag	UNP A0A0F5SXE4
D	697	ASN	-	expression tag	UNP A0A0F5SXE4
D	698	LEU	-	expression tag	UNP A0A0F5SXE4
D	699	TYR	-	expression tag	UNP A0A0F5SXE4
D	700	PHE	-	expression tag	UNP A0A0F5SXE4
D	701	GLN	-	expression tag	UNP A0A0F5SXE4
E	696	GLU	-	expression tag	UNP A0A0F5SXE4
E	697	ASN	-	expression tag	UNP A0A0F5SXE4
E	698	LEU	-	expression tag	UNP A0A0F5SXE4
E	699	TYR	-	expression tag	UNP A0A0F5SXE4
E	700	PHE	-	expression tag	UNP A0A0F5SXE4
E	701	GLN	-	expression tag	UNP A0A0F5SXE4
F	696	GLU	-	expression tag	UNP A0A0F5SXE4
F	697	ASN	-	expression tag	UNP A0A0F5SXE4
F	698	LEU	-	expression tag	UNP A0A0F5SXE4
F	699	TYR	-	expression tag	UNP A0A0F5SXE4
F	700	PHE	-	expression tag	UNP A0A0F5SXE4
F	701	GLN	-	expression tag	UNP A0A0F5SXE4
G	696	GLU	-	expression tag	UNP A0A0F5SXE4
G	697	ASN	-	expression tag	UNP A0A0F5SXE4
G	698	LEU	-	expression tag	UNP A0A0F5SXE4
G	699	TYR	-	expression tag	UNP A0A0F5SXE4
G	700	PHE	-	expression tag	UNP A0A0F5SXE4
G	701	GLN	-	expression tag	UNP A0A0F5SXE4
H	696	GLU	-	expression tag	UNP A0A0F5SXE4
H	697	ASN	-	expression tag	UNP A0A0F5SXE4
H	698	LEU	-	expression tag	UNP A0A0F5SXE4
H	699	TYR	-	expression tag	UNP A0A0F5SXE4
H	700	PHE	-	expression tag	UNP A0A0F5SXE4
H	701	GLN	-	expression tag	UNP A0A0F5SXE4
I	696	GLU	-	expression tag	UNP A0A0F5SXE4
I	697	ASN	-	expression tag	UNP A0A0F5SXE4
I	698	LEU	-	expression tag	UNP A0A0F5SXE4
I	699	TYR	-	expression tag	UNP A0A0F5SXE4
I	700	PHE	-	expression tag	UNP A0A0F5SXE4
I	701	GLN	-	expression tag	UNP A0A0F5SXE4

### 3 Residue-property plots

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

#### • Molecule 1: Flagellar biosynthesis protein FlhA



Chain B:

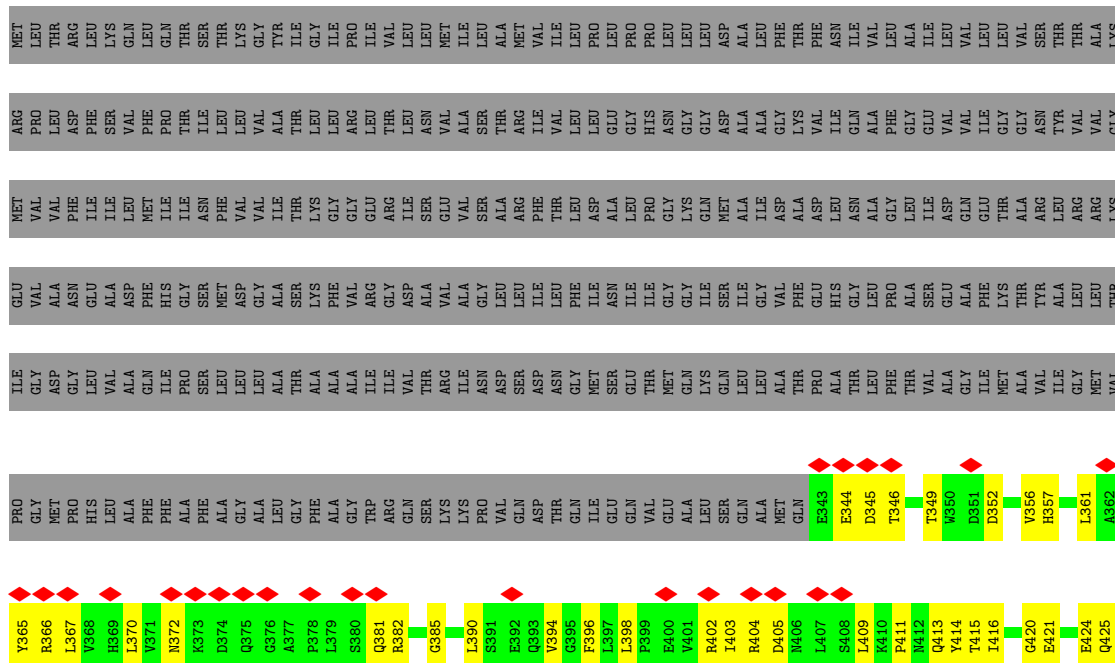
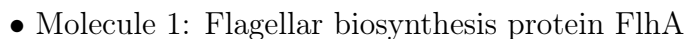


ILE	GLY	ASP	ALA	VAL	GLU	MET	VAL	VAL	MET	ARG	MET
GLY	ASP	ASN	ALA	VAL	VAL	VAL	PHE	ASP	LEU	PRO	LEU
GLY	LEU	LEU	ALA	ASP	GLU	ILE	ILE	PHE	GLY	ASP	ARG
VAL	VAL	ALA	ASP	ALA	ALA	ILE	ILE	SER	LYS	GLN	LYS
ALA	GLN	ILE	PHE	PHE	ASP	LEU	LEU	VAL	LEU	LEU	LEU
ILE	ILE	PRO	HIS	GLY	GLY	ILE	ILE	PRO	GLN	GLN	GLN
THR	THR	SER	SER	SER	ASP	ASN	ILE	THR	THR	SER	THR
LEU	LEU	LEU	MET	MET	VAL	PHE	PHE	LEU	LEU	LYS	LYS
LEU	LEU	GLY	ASP	VAL	VAL	VAL	VAL	VAL	VAL	VAL	GLY
LEU	ALA	ALA	ALA	ALA	LYS	LYS	THR	THR	ALA	ALA	ILE
ALA	ALA	VAL	VAL	VAL	PHE	GLY	GLY	LEU	LEU	PRO	ILE
ILE	ILE	ARG	ARG	ARG	VAL	VAL	VAL	SER	ILE	ARG	PRO
ILE	ILE	ASP	ASP	LEU	ASP	GLY	GLY	LEU	GLY	LEU	ILE
VAL	VAL	SER	LEU	LEU	VAL	ALA	ALA	THR	THR	THR	VAL
THR	THR	ASN	ILE	ILE	THR	PHE	PHE	VAL	ILE	VAL	ILE
THR	THR	ASN	ILE	ILE	THR	ARG	ARG	ARG	ILE	THR	VAL
LEU	LEU	LEU	THR	THR	THR	THR	THR	VAL	VAL	VAL	ILE
LEU	LEU	ASN	GLY	PHE	LEU	LEU	LEU	LEU	GLY	PRO	ILE
PRO	MET	MET	ILE	ILE	ASP	ASP	ASP	LEU	GLY	ASP	LEU
SER	SER	ASN	ASN	ASN	ALA	ALA	ALA	GLY	GLY	LEU	LEU
GLU	GLU	GLU	ILE	ILE	ILE	LEU	GLY	GLY	HIS	PRO	PRO
THR	THR	THR	ILE	ILE	GLY	GLY	GLY	ASN	ASN	LEU	LEU
GLN	GLN	THR	GLY	ILE	LYS	GLN	LYS	GLY	GLY	LEU	LEU
LEU	LEU	THR	VAL	THR	THR	THR	THR	ASP	ASP	LEU	LEU
ALA	ALA	ALA	GLU	GLU	ASP	ASP	ILE	VAL	ILE	ALA	ALA
THR	THR	THR	HIS	ASP	GLU	ASP	ASP	VAL	THR	THR	PHE
ALA	ALA	LEU	GLY	GLY	GLY	ALA	ALA	GLY	GLY	GLY	GLY
LEU	LEU	THR	PRO	PRO	ALA	ALA	ALA	GLY	ILE	ILE	ILE
THR	THR	THR	VAL	VAL	THR	THR	THR	VAL	THR	THR	THR
VAL	VAL	VAL	THR	THR	THR	ALA	ALA	GLY	GLY	VAL	VAL
VAL	VAL	VAL	TYR	TYR	ALA	ALA	ALA	ASN	ASN	SER	SER
GLY	GLY	GLY	ALA	ARG	THR	ARG	ARG	VAL	TYR	THR	THR
MET	MET	MET	LEU	THR	LEU	ARG	VAL	VAL	VAL	VAL	VAL
VAL	VAL	VAL	THR	THR	THR	LYS	GLY	GLY	GLY	GLY	GLY

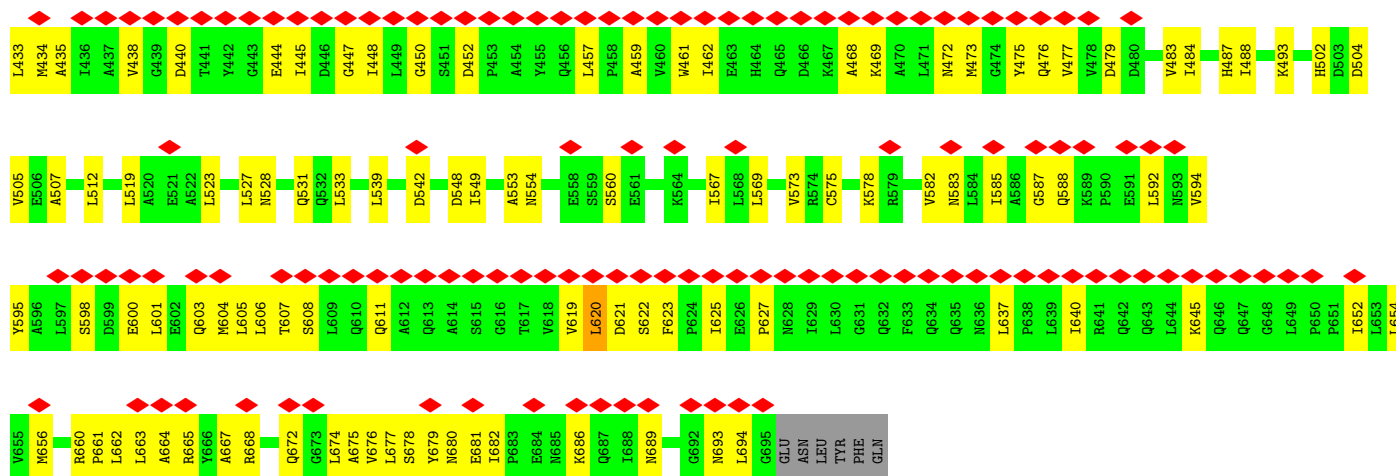
PRO	GLY	MET	PRO	HIS	LEU	PHE	PHE	PHE	ALA	GLY	ALA	LEU	GLY	PHE	ALA	GLY	TRP	ARG	GLN	SER	LYS	PRO	VAL	VAL	GLN	THR	ILE	GLU	GLN	VAL	GLU	ALA	LEU	SER	GLN	ALA	MET	GLN	E343	E344	D345	T346	P347	L348	T349	W350	D351	D352	V356	H357	L361	A362	L363																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																												
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## Chain C:

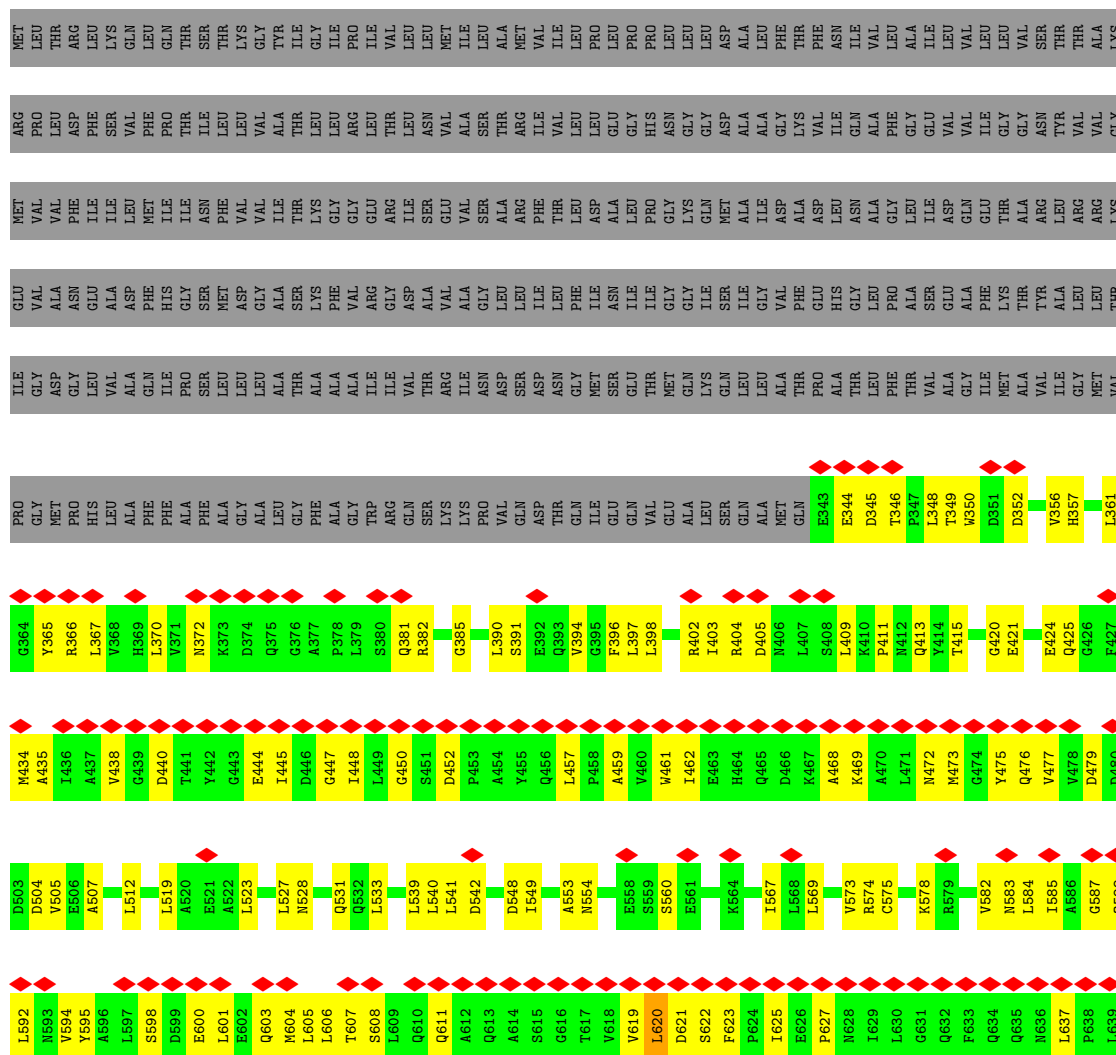
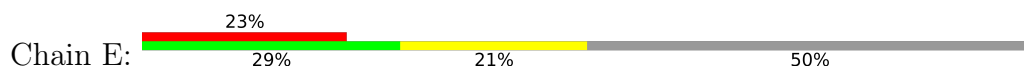
[illegible]

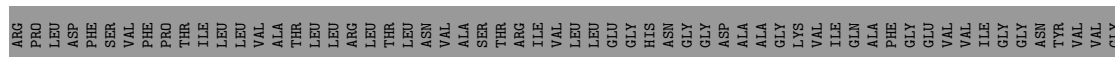


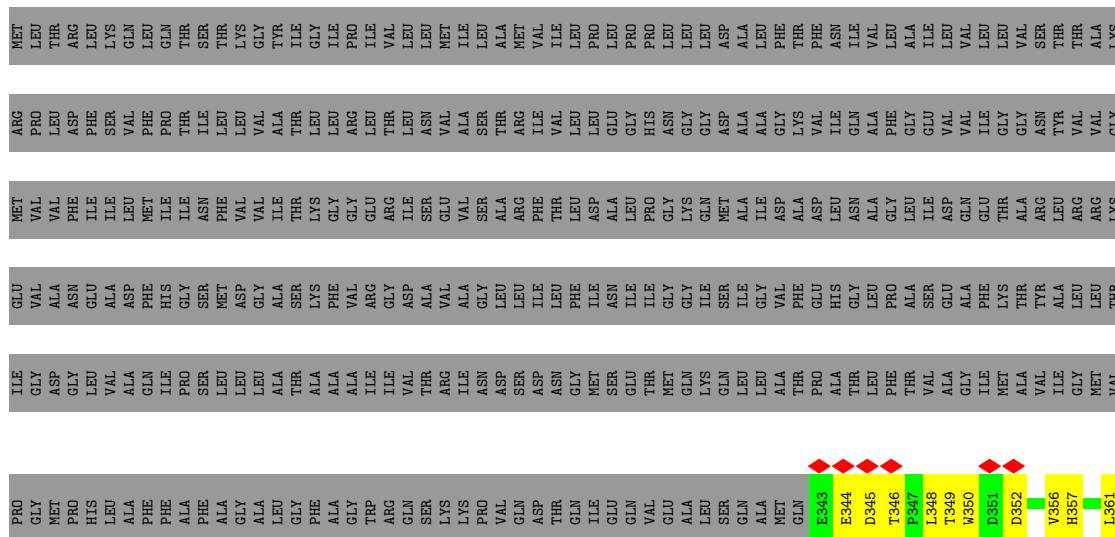
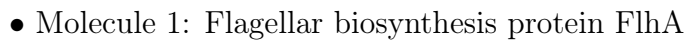


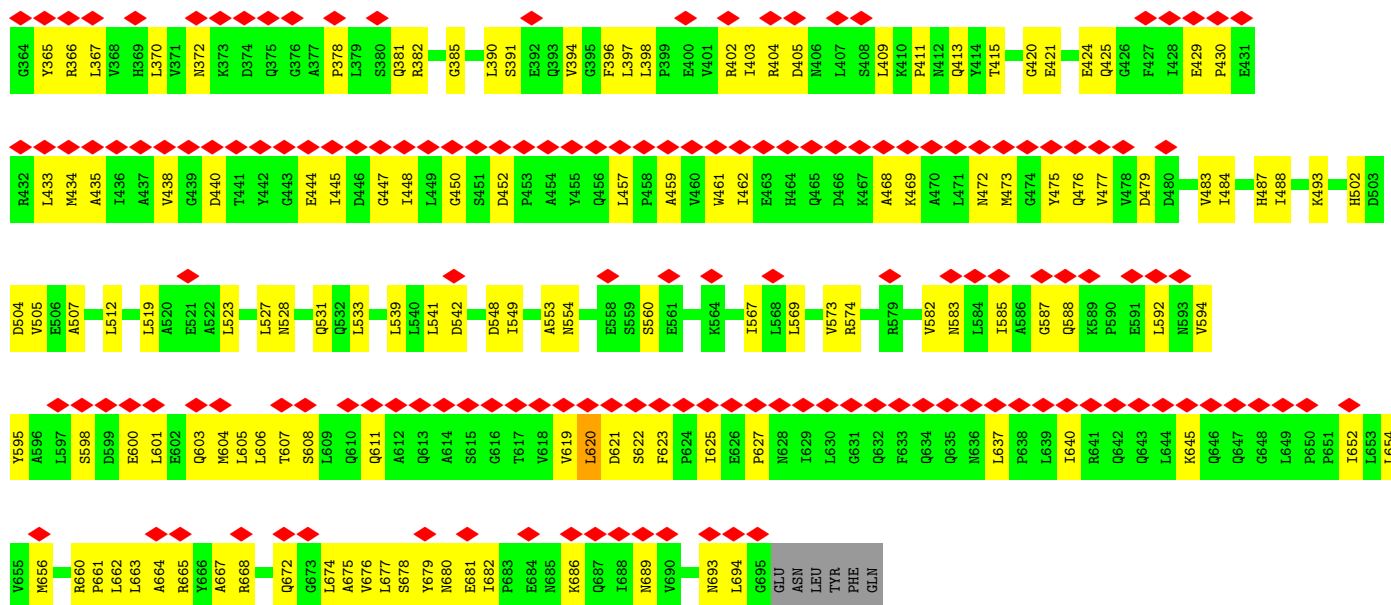


• Molecule 1: Flagellar biosynthesis protein FlhA

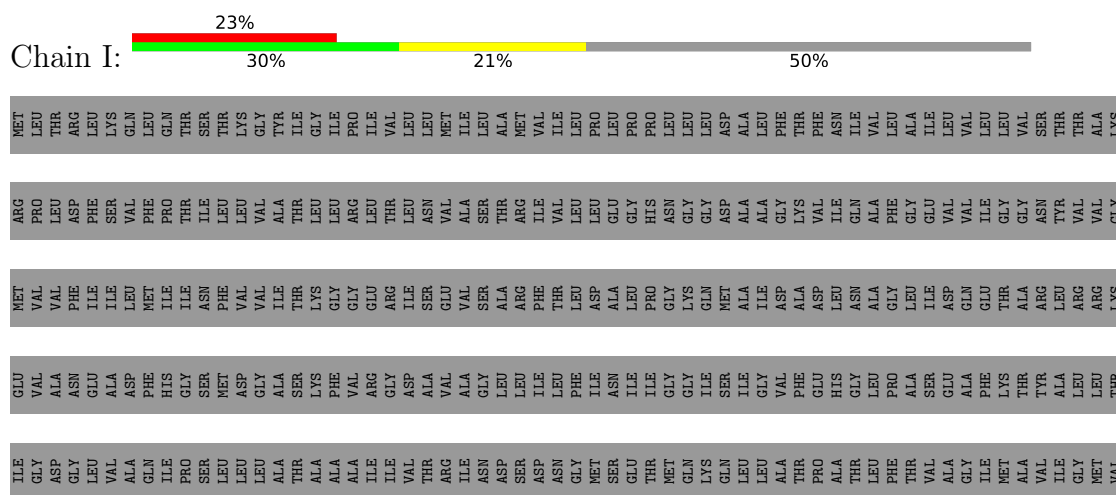








• Molecule 1: Flagellar biosynthesis protein FlhA





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C9	Depositor
Number of particles used	9756	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	48	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.048	Depositor
Minimum map value	-0.022	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.0146	Depositor
Map size ( $\text{\AA}$ )	348.528, 348.528, 348.528	wwPDB
Map dimensions	424, 424, 424	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.822, 0.822, 0.822	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	0.43	0/2806	0.50	1/3819 (0.0%)
1	B	0.43	0/2806	0.50	1/3819 (0.0%)
1	C	0.43	0/2806	0.50	1/3819 (0.0%)
1	D	0.43	0/2806	0.50	1/3819 (0.0%)
1	E	0.43	0/2806	0.50	1/3819 (0.0%)
1	F	0.43	0/2806	0.50	1/3819 (0.0%)
1	G	0.43	0/2806	0.50	1/3819 (0.0%)
1	H	0.43	0/2806	0.50	1/3819 (0.0%)
1	I	0.43	0/2806	0.50	1/3819 (0.0%)
All	All	0.43	0/25254	0.50	9/34371 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	345	ASP	N-CA-C	-6.45	106.14	112.97
1	G	345	ASP	N-CA-C	-6.44	106.14	112.97
1	I	345	ASP	N-CA-C	-6.42	106.17	112.97
1	B	345	ASP	N-CA-C	-6.42	106.17	112.97
1	E	345	ASP	N-CA-C	-6.40	106.19	112.97
1	A	345	ASP	N-CA-C	-6.38	106.20	112.97
1	C	345	ASP	N-CA-C	-6.38	106.21	112.97
1	H	345	ASP	N-CA-C	-6.38	106.20	112.97
1	D	345	ASP	N-CA-C	-6.38	106.21	112.97

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	2761	0	2832	120	0
1	B	2761	0	2832	117	0
1	C	2761	0	2832	113	0
1	D	2761	0	2832	111	0
1	E	2761	0	2832	123	0
1	F	2761	0	2832	115	0
1	G	2761	0	2832	118	0
1	H	2761	0	2832	118	0
1	I	2761	0	2832	118	0
All	All	24849	0	25488	1020	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 20.

All (1020) 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:409:LEU:HB2	1:A:413:GLN:HE21	1.34	0.93
1:B:409:LEU:HB2	1:B:413:GLN:HE21	1.34	0.92
1:I:409:LEU:HB2	1:I:413:GLN:HE21	1.34	0.91
1:C:409:LEU:HB2	1:C:413:GLN:HE21	1.34	0.91
1:E:409:LEU:HB2	1:E:413:GLN:HE21	1.34	0.91
1:H:409:LEU:HB2	1:H:413:GLN:HE21	1.34	0.91
1:D:409:LEU:HB2	1:D:413:GLN:HE21	1.34	0.91
1:F:409:LEU:HB2	1:F:413:GLN:HE21	1.34	0.91
1:G:409:LEU:HB2	1:G:413:GLN:HE21	1.34	0.90
1:A:620:LEU:HD12	1:A:621:ASP:H	1.37	0.90
1:F:620:LEU:HD12	1:F:621:ASP:H	1.37	0.90
1:G:620:LEU:HD12	1:G:621:ASP:H	1.37	0.89
1:I:620:LEU:HD12	1:I:621:ASP:H	1.37	0.89
1:D:620:LEU:HD12	1:D:621:ASP:H	1.37	0.89
1:B:620:LEU:HD12	1:B:621:ASP:H	1.37	0.88
1:H:620:LEU:HD12	1:H:621:ASP:H	1.37	0.88
1:E:620:LEU:HD12	1:E:621:ASP:H	1.37	0.87
1:C:620:LEU:HD12	1:C:621:ASP:H	1.37	0.86
1:B:567:ILE:HG12	1:B:620:LEU:HD13	1.57	0.86
1:I:567:ILE:HG12	1:I:620:LEU:HD13	1.57	0.86
1:E:567:ILE:HG12	1:E:620:LEU:HD13	1.58	0.84
1:A:567:ILE:HG12	1:A:620:LEU:HD13	1.58	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:567:ILE:HG12	1:F:620:LEU:HD13	1.58	0.84
1:D:567:ILE:HG12	1:D:620:LEU:HD13	1.58	0.84
1:D:390:LEU:HD11	1:D:493:LYS:HG3	1.60	0.84
1:G:567:ILE:HG12	1:G:620:LEU:HD13	1.58	0.83
1:C:567:ILE:HG12	1:C:620:LEU:HD13	1.57	0.83
1:E:390:LEU:HD11	1:E:493:LYS:HG3	1.60	0.83
1:F:390:LEU:HD11	1:F:493:LYS:HG3	1.60	0.83
1:H:567:ILE:HG12	1:H:620:LEU:HD13	1.58	0.83
1:C:390:LEU:HD11	1:C:493:LYS:HG3	1.60	0.83
1:B:390:LEU:HD11	1:B:493:LYS:HG3	1.60	0.83
1:G:390:LEU:HD11	1:G:493:LYS:HG3	1.60	0.83
1:I:390:LEU:HD11	1:I:493:LYS:HG3	1.60	0.82
1:A:390:LEU:HD11	1:A:493:LYS:HG3	1.60	0.82
1:H:390:LEU:HD11	1:H:493:LYS:HG3	1.60	0.81
1:F:619:VAL:HG21	1:F:623:PHE:H	1.46	0.81
1:D:619:VAL:HG21	1:D:623:PHE:H	1.46	0.81
1:A:619:VAL:HG21	1:A:623:PHE:H	1.46	0.81
1:C:619:VAL:HG21	1:C:623:PHE:H	1.46	0.81
1:E:619:VAL:HG21	1:E:623:PHE:H	1.46	0.80
1:B:619:VAL:HG21	1:B:623:PHE:H	1.46	0.80
1:H:619:VAL:HG21	1:H:623:PHE:H	1.46	0.80
1:I:619:VAL:HG21	1:I:623:PHE:H	1.46	0.80
1:G:619:VAL:HG21	1:G:623:PHE:H	1.46	0.80
1:C:502:HIS:O	1:C:502:HIS:ND1	2.18	0.77
1:D:502:HIS:O	1:D:502:HIS:ND1	2.18	0.77
1:F:502:HIS:O	1:F:502:HIS:ND1	2.18	0.77
1:B:502:HIS:ND1	1:B:502:HIS:O	2.18	0.77
1:H:502:HIS:O	1:H:502:HIS:ND1	2.18	0.76
1:E:502:HIS:O	1:E:502:HIS:ND1	2.18	0.76
1:A:502:HIS:O	1:A:502:HIS:ND1	2.18	0.76
1:I:502:HIS:O	1:I:502:HIS:ND1	2.18	0.75
1:G:502:HIS:ND1	1:G:502:HIS:O	2.18	0.75
1:B:366:ARG:N	1:B:411:PRO:O	2.20	0.75
1:C:366:ARG:N	1:C:411:PRO:O	2.20	0.75
1:D:366:ARG:N	1:D:411:PRO:O	2.20	0.74
1:A:390:LEU:HD22	1:A:398:LEU:HD11	1.70	0.74
1:G:366:ARG:N	1:G:411:PRO:O	2.20	0.74
1:B:390:LEU:HD22	1:B:398:LEU:HD11	1.70	0.73
1:I:366:ARG:N	1:I:411:PRO:O	2.20	0.73
1:I:390:LEU:HD22	1:I:398:LEU:HD11	1.70	0.73
1:H:390:LEU:HD22	1:H:398:LEU:HD11	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:366:ARG:N	1:F:411:PRO:O	2.20	0.73
1:C:390:LEU:HD22	1:C:398:LEU:HD11	1.70	0.73
1:H:366:ARG:N	1:H:411:PRO:O	2.20	0.73
1:G:390:LEU:HD22	1:G:398:LEU:HD11	1.70	0.73
1:A:366:ARG:N	1:A:411:PRO:O	2.20	0.72
1:E:366:ARG:N	1:E:411:PRO:O	2.20	0.72
1:B:623:PHE:HB3	1:B:665:ARG:HH21	1.55	0.72
1:D:390:LEU:HD22	1:D:398:LEU:HD11	1.70	0.72
1:D:623:PHE:HB3	1:D:665:ARG:HH21	1.55	0.72
1:A:623:PHE:HB3	1:A:665:ARG:HH21	1.55	0.71
1:F:390:LEU:HD22	1:F:398:LEU:HD11	1.70	0.71
1:F:623:PHE:HB3	1:F:665:ARG:HH21	1.55	0.71
1:G:623:PHE:HB3	1:G:665:ARG:HH21	1.55	0.71
1:E:390:LEU:HD22	1:E:398:LEU:HD11	1.70	0.71
1:E:623:PHE:HB3	1:E:665:ARG:HH21	1.55	0.71
1:H:623:PHE:HB3	1:H:665:ARG:HH21	1.55	0.71
1:I:623:PHE:HB3	1:I:665:ARG:HH21	1.55	0.70
1:C:623:PHE:HB3	1:C:665:ARG:HH21	1.55	0.70
1:A:361:LEU:HB2	1:A:403:ILE:HG12	1.73	0.70
1:A:424:GLU:OE1	1:A:487:HIS:NE2	2.26	0.69
1:D:361:LEU:HB2	1:D:403:ILE:HG12	1.74	0.69
1:C:424:GLU:OE1	1:C:487:HIS:NE2	2.26	0.69
1:E:424:GLU:OE1	1:E:487:HIS:NE2	2.26	0.69
1:H:424:GLU:OE1	1:H:487:HIS:NE2	2.26	0.69
1:C:361:LEU:HB2	1:C:403:ILE:HG12	1.74	0.69
1:D:424:GLU:OE1	1:D:487:HIS:NE2	2.26	0.69
1:F:361:LEU:HB2	1:F:403:ILE:HG12	1.74	0.69
1:I:361:LEU:HB2	1:I:403:ILE:HG12	1.73	0.69
1:B:424:GLU:OE1	1:B:487:HIS:NE2	2.26	0.69
1:B:361:LEU:HB2	1:B:403:ILE:HG12	1.74	0.68
1:F:424:GLU:OE1	1:F:487:HIS:NE2	2.26	0.68
1:E:452:ASP:HB3	1:E:457:LEU:H	1.59	0.68
1:G:361:LEU:HB2	1:G:403:ILE:HG12	1.74	0.68
1:H:452:ASP:HB3	1:H:457:LEU:H	1.59	0.68
1:I:424:GLU:OE1	1:I:487:HIS:NE2	2.26	0.68
1:G:424:GLU:OE1	1:G:487:HIS:NE2	2.26	0.68
1:G:592:LEU:O	1:G:689:ASN:N	2.27	0.68
1:E:361:LEU:HB2	1:E:403:ILE:HG12	1.74	0.67
1:A:592:LEU:O	1:A:689:ASN:N	2.28	0.67
1:A:654:LEU:HB3	1:A:679:TYR:HE2	1.60	0.67
1:B:592:LEU:O	1:B:689:ASN:N	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:592:LEU:O	1:C:689:ASN:N	2.27	0.67
1:D:654:LEU:HB3	1:D:679:TYR:HE2	1.60	0.67
1:H:592:LEU:O	1:H:689:ASN:N	2.28	0.67
1:I:592:LEU:O	1:I:689:ASN:N	2.28	0.67
1:E:592:LEU:O	1:E:689:ASN:N	2.28	0.67
1:H:361:LEU:HB2	1:H:403:ILE:HG12	1.74	0.67
1:D:592:LEU:O	1:D:689:ASN:N	2.28	0.67
1:G:654:LEU:HB3	1:G:679:TYR:HE2	1.60	0.67
1:A:452:ASP:HB3	1:A:457:LEU:H	1.59	0.67
1:F:592:LEU:O	1:F:689:ASN:N	2.28	0.67
1:F:654:LEU:HB3	1:F:679:TYR:HE2	1.60	0.67
1:B:452:ASP:HB3	1:B:457:LEU:H	1.59	0.67
1:F:452:ASP:HB3	1:F:457:LEU:H	1.59	0.67
1:I:654:LEU:HB3	1:I:679:TYR:HE2	1.60	0.67
1:D:452:ASP:HB3	1:D:457:LEU:H	1.59	0.67
1:C:452:ASP:HB3	1:C:457:LEU:H	1.59	0.66
1:F:587:GLY:O	1:F:686:LYS:NZ	2.29	0.66
1:G:452:ASP:HB3	1:G:457:LEU:H	1.59	0.66
1:H:587:GLY:O	1:H:686:LYS:NZ	2.29	0.66
1:I:452:ASP:HB3	1:I:457:LEU:H	1.59	0.66
1:B:654:LEU:HB3	1:B:679:TYR:HE2	1.60	0.66
1:G:587:GLY:O	1:G:686:LYS:NZ	2.29	0.66
1:E:587:GLY:O	1:E:686:LYS:NZ	2.29	0.66
1:I:587:GLY:O	1:I:686:LYS:NZ	2.29	0.66
1:H:654:LEU:HB3	1:H:679:TYR:HE2	1.60	0.66
1:D:587:GLY:O	1:D:686:LYS:NZ	2.29	0.66
1:A:587:GLY:O	1:A:686:LYS:NZ	2.29	0.66
1:E:654:LEU:HB3	1:E:679:TYR:HE2	1.60	0.66
1:C:654:LEU:HB3	1:C:679:TYR:HE2	1.60	0.66
1:C:587:GLY:O	1:C:686:LYS:NZ	2.29	0.65
1:B:587:GLY:O	1:B:686:LYS:NZ	2.29	0.65
1:E:433:LEU:N	1:E:462:ILE:O	2.30	0.65
1:F:433:LEU:N	1:F:462:ILE:O	2.30	0.65
1:G:433:LEU:N	1:G:462:ILE:O	2.30	0.65
1:I:433:LEU:N	1:I:462:ILE:O	2.30	0.65
1:H:433:LEU:N	1:H:462:ILE:O	2.30	0.65
1:A:433:LEU:N	1:A:462:ILE:O	2.30	0.64
1:D:433:LEU:N	1:D:462:ILE:O	2.30	0.64
1:B:433:LEU:N	1:B:462:ILE:O	2.30	0.63
1:C:433:LEU:N	1:C:462:ILE:O	2.30	0.63
1:E:608:SER:HB3	1:E:625:ILE:HB	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:421:GLU:N	1:C:421:GLU:OE1	2.31	0.63
1:F:421:GLU:N	1:F:421:GLU:OE1	2.31	0.63
1:D:608:SER:HB3	1:D:625:ILE:HB	1.80	0.63
1:F:608:SER:HB3	1:F:625:ILE:HB	1.80	0.63
1:D:421:GLU:N	1:D:421:GLU:OE1	2.31	0.63
1:E:421:GLU:N	1:E:421:GLU:OE1	2.31	0.63
1:A:421:GLU:N	1:A:421:GLU:OE1	2.31	0.63
1:B:421:GLU:OE1	1:B:421:GLU:N	2.31	0.63
1:H:421:GLU:N	1:H:421:GLU:OE1	2.31	0.63
1:A:429:GLU:HG3	1:A:430:PRO:HD2	1.80	0.63
1:C:608:SER:HB3	1:C:625:ILE:HB	1.80	0.63
1:G:608:SER:HB3	1:G:625:ILE:HB	1.80	0.63
1:B:608:SER:HB3	1:B:625:ILE:HB	1.80	0.63
1:I:421:GLU:N	1:I:421:GLU:OE1	2.31	0.62
1:G:421:GLU:OE1	1:G:421:GLU:N	2.31	0.62
1:H:429:GLU:HG3	1:H:430:PRO:HD2	1.80	0.62
1:H:608:SER:HB3	1:H:625:ILE:HB	1.80	0.62
1:I:608:SER:HB3	1:I:625:ILE:HB	1.80	0.62
1:I:429:GLU:HG3	1:I:430:PRO:HD2	1.81	0.62
1:H:512:LEU:HD12	1:H:549:ILE:HG23	1.82	0.62
1:I:512:LEU:HD12	1:I:549:ILE:HG23	1.82	0.62
1:E:429:GLU:HG3	1:E:430:PRO:HD2	1.80	0.62
1:G:512:LEU:HD12	1:G:549:ILE:HG23	1.82	0.62
1:B:429:GLU:HG3	1:B:430:PRO:HD2	1.81	0.62
1:E:512:LEU:HD12	1:E:549:ILE:HG23	1.82	0.62
1:D:598:SER:HB3	1:D:693:ASN:HB3	1.82	0.62
1:F:429:GLU:HG3	1:F:430:PRO:HD2	1.81	0.61
1:A:512:LEU:HD12	1:A:549:ILE:HG23	1.82	0.61
1:G:382:ARG:HH12	1:G:479:ASP:HB3	1.66	0.61
1:G:429:GLU:HG3	1:G:430:PRO:HD2	1.81	0.61
1:E:598:SER:HB3	1:E:693:ASN:HB3	1.82	0.61
1:F:598:SER:HB3	1:F:693:ASN:HB3	1.82	0.61
1:C:598:SER:HB3	1:C:693:ASN:HB3	1.82	0.61
1:C:429:GLU:HG3	1:C:430:PRO:HD2	1.81	0.61
1:F:512:LEU:HD12	1:F:549:ILE:HG23	1.82	0.61
1:C:604:MET:O	1:C:607:THR:OG1	2.18	0.61
1:D:429:GLU:HG3	1:D:430:PRO:HD2	1.81	0.61
1:E:382:ARG:HH12	1:E:479:ASP:HB3	1.66	0.61
1:A:608:SER:HB3	1:A:625:ILE:HB	1.80	0.61
1:E:604:MET:O	1:E:607:THR:OG1	2.18	0.61
1:G:604:MET:O	1:G:607:THR:OG1	2.18	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:598:SER:HB3	1:B:693:ASN:HB3	1.82	0.61
1:B:604:MET:O	1:B:607:THR:OG1	2.18	0.61
1:H:382:ARG:HH12	1:H:479:ASP:HB3	1.66	0.61
1:B:382:ARG:HH12	1:B:479:ASP:HB3	1.66	0.60
1:B:512:LEU:HD12	1:B:549:ILE:HG23	1.82	0.60
1:C:665:ARG:HG3	1:C:668:ARG:HH22	1.67	0.60
1:F:604:MET:O	1:F:607:THR:OG1	2.18	0.60
1:G:356:VAL:HG11	1:H:548:ASP:HB3	1.84	0.60
1:G:598:SER:HB3	1:G:693:ASN:HB3	1.82	0.60
1:H:604:MET:O	1:H:607:THR:OG1	2.18	0.60
1:A:382:ARG:HH12	1:A:479:ASP:HB3	1.66	0.60
1:C:382:ARG:HH12	1:C:479:ASP:HB3	1.66	0.60
1:D:512:LEU:HD12	1:D:549:ILE:HG23	1.82	0.60
1:D:604:MET:O	1:D:607:THR:OG1	2.18	0.60
1:D:665:ARG:HG3	1:D:668:ARG:HH22	1.67	0.60
1:H:598:SER:HB3	1:H:693:ASN:HB3	1.82	0.60
1:A:356:VAL:HG11	1:B:548:ASP:HB3	1.83	0.60
1:A:548:ASP:HB3	1:I:356:VAL:HG11	1.84	0.60
1:C:598:SER:OG	1:C:694:LEU:O	2.20	0.60
1:C:512:LEU:HD12	1:C:549:ILE:HG23	1.82	0.60
1:I:598:SER:HB3	1:I:693:ASN:HB3	1.82	0.60
1:A:598:SER:HB3	1:A:693:ASN:HB3	1.82	0.60
1:D:382:ARG:HH12	1:D:479:ASP:HB3	1.66	0.59
1:I:604:MET:O	1:I:607:THR:OG1	2.18	0.59
1:B:356:VAL:HG11	1:C:548:ASP:HB3	1.84	0.59
1:F:382:ARG:HH12	1:F:479:ASP:HB3	1.66	0.59
1:D:598:SER:OG	1:D:694:LEU:O	2.20	0.59
1:A:604:MET:O	1:A:607:THR:OG1	2.18	0.59
1:B:665:ARG:HG3	1:B:668:ARG:HH22	1.67	0.59
1:E:356:VAL:HG11	1:G:548:ASP:HB3	1.84	0.59
1:I:382:ARG:HH12	1:I:479:ASP:HB3	1.66	0.59
1:I:598:SER:OG	1:I:694:LEU:O	2.20	0.59
1:I:665:ARG:HG3	1:I:668:ARG:HH22	1.67	0.59
1:G:598:SER:OG	1:G:694:LEU:O	2.20	0.59
1:H:356:VAL:HG11	1:I:548:ASP:HB3	1.84	0.59
1:G:665:ARG:HG3	1:G:668:ARG:HH22	1.67	0.59
1:E:678:SER:HB3	1:E:681:GLU:HG3	1.84	0.59
1:F:665:ARG:HG3	1:F:668:ARG:HH22	1.67	0.59
1:F:678:SER:HB3	1:F:681:GLU:HG3	1.84	0.59
1:E:665:ARG:HG3	1:E:668:ARG:HH22	1.67	0.59
1:B:448:ILE:HG23	1:B:461:TRP:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:665:ARG:HG3	1:H:668:ARG:HH22	1.67	0.59
1:I:448:ILE:HG23	1:I:461:TRP:HB2	1.85	0.59
1:C:448:ILE:HG23	1:C:461:TRP:HB2	1.85	0.58
1:D:448:ILE:HG23	1:D:461:TRP:HB2	1.85	0.58
1:C:678:SER:HB3	1:C:681:GLU:HG3	1.84	0.58
1:A:448:ILE:HG23	1:A:461:TRP:HB2	1.85	0.58
1:E:448:ILE:HG23	1:E:461:TRP:HB2	1.85	0.58
1:G:678:SER:HB3	1:G:681:GLU:HG3	1.84	0.58
1:H:448:ILE:HG23	1:H:461:TRP:HB2	1.85	0.58
1:E:527:LEU:HD11	1:E:560:SER:HB2	1.86	0.58
1:I:678:SER:HB3	1:I:681:GLU:HG3	1.84	0.58
1:A:665:ARG:HG3	1:A:668:ARG:HH22	1.67	0.58
1:E:598:SER:OG	1:E:694:LEU:O	2.19	0.58
1:F:527:LEU:HD11	1:F:560:SER:HB2	1.86	0.58
1:H:678:SER:HB3	1:H:681:GLU:HG3	1.84	0.58
1:H:598:SER:OG	1:H:694:LEU:O	2.20	0.58
1:G:527:LEU:HD11	1:G:560:SER:HB2	1.86	0.58
1:C:433:LEU:O	1:C:462:ILE:N	2.31	0.58
1:D:527:LEU:HD11	1:D:560:SER:HB2	1.86	0.58
1:D:678:SER:HB3	1:D:681:GLU:HG3	1.84	0.57
1:F:448:ILE:HG23	1:F:461:TRP:HB2	1.85	0.57
1:F:598:SER:OG	1:F:694:LEU:O	2.20	0.57
1:C:356:VAL:HG11	1:D:548:ASP:HB3	1.86	0.57
1:G:448:ILE:HG23	1:G:461:TRP:HB2	1.85	0.57
1:B:678:SER:HB3	1:B:681:GLU:HG3	1.84	0.57
1:E:433:LEU:O	1:E:462:ILE:N	2.31	0.57
1:A:527:LEU:HD11	1:A:560:SER:HB2	1.86	0.57
1:A:678:SER:HB3	1:A:681:GLU:HG3	1.84	0.57
1:H:527:LEU:HD11	1:H:560:SER:HB2	1.86	0.57
1:C:527:LEU:HD11	1:C:560:SER:HB2	1.86	0.57
1:E:548:ASP:HB3	1:F:356:VAL:HG11	1.84	0.57
1:I:433:LEU:O	1:I:462:ILE:N	2.31	0.56
1:I:527:LEU:HD11	1:I:560:SER:HB2	1.86	0.56
1:H:621:ASP:HB3	1:H:661:PRO:HB2	1.88	0.56
1:G:621:ASP:HB3	1:G:661:PRO:HB2	1.88	0.56
1:A:621:ASP:HB3	1:A:661:PRO:HB2	1.88	0.56
1:B:527:LEU:HD11	1:B:560:SER:HB2	1.86	0.56
1:I:621:ASP:HB3	1:I:661:PRO:HB2	1.88	0.56
1:D:356:VAL:HG11	1:F:548:ASP:HB3	1.88	0.55
1:H:625:ILE:HG22	1:H:627:PRO:HD3	1.88	0.55
1:I:625:ILE:HG22	1:I:627:PRO:HD3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:583:ASN:OD1	1:A:588:GLN:NE2	2.40	0.55
1:E:588:GLN:N	1:E:588:GLN:OE1	2.40	0.55
1:B:621:ASP:HB3	1:B:661:PRO:HB2	1.88	0.55
1:E:621:ASP:HB3	1:E:661:PRO:HB2	1.88	0.55
1:G:588:GLN:N	1:G:588:GLN:OE1	2.40	0.55
1:C:625:ILE:HG22	1:C:627:PRO:HD3	1.88	0.55
1:I:583:ASN:OD1	1:I:588:GLN:NE2	2.40	0.55
1:A:433:LEU:O	1:A:462:ILE:N	2.31	0.55
1:C:588:GLN:OE1	1:C:588:GLN:N	2.39	0.55
1:D:625:ILE:HG22	1:D:627:PRO:HD3	1.88	0.55
1:I:620:LEU:CD1	1:I:621:ASP:H	2.16	0.55
1:F:588:GLN:OE1	1:F:588:GLN:N	2.40	0.55
1:G:625:ILE:HG22	1:G:627:PRO:HD3	1.88	0.55
1:F:583:ASN:OD1	1:F:588:GLN:NE2	2.40	0.55
1:G:583:ASN:OD1	1:G:588:GLN:NE2	2.40	0.55
1:I:588:GLN:N	1:I:588:GLN:OE1	2.40	0.55
1:A:625:ILE:HG22	1:A:627:PRO:HD3	1.88	0.55
1:B:583:ASN:OD1	1:B:588:GLN:NE2	2.40	0.55
1:H:583:ASN:OD1	1:H:588:GLN:NE2	2.40	0.55
1:D:621:ASP:HB3	1:D:661:PRO:HB2	1.88	0.54
1:G:433:LEU:O	1:G:462:ILE:N	2.31	0.54
1:B:405:ASP:OD1	1:B:405:ASP:N	2.40	0.54
1:A:435:ALA:HA	1:A:477:VAL:HA	1.89	0.54
1:C:621:ASP:HB3	1:C:661:PRO:HB2	1.88	0.54
1:E:583:ASN:OD1	1:E:588:GLN:NE2	2.40	0.54
1:H:588:GLN:OE1	1:H:588:GLN:N	2.40	0.54
1:F:621:ASP:HB3	1:F:661:PRO:HB2	1.88	0.54
1:B:625:ILE:HG22	1:B:627:PRO:HD3	1.88	0.54
1:D:583:ASN:OD1	1:D:588:GLN:NE2	2.40	0.54
1:F:625:ILE:HG22	1:F:627:PRO:HD3	1.88	0.54
1:C:405:ASP:OD1	1:C:405:ASP:N	2.40	0.54
1:D:433:LEU:O	1:D:462:ILE:N	2.31	0.54
1:D:588:GLN:N	1:D:588:GLN:OE1	2.40	0.54
1:A:405:ASP:OD1	1:A:405:ASP:N	2.40	0.54
1:A:588:GLN:OE1	1:A:588:GLN:N	2.40	0.54
1:C:583:ASN:OD1	1:C:588:GLN:NE2	2.40	0.54
1:H:585:ILE:HD13	1:H:675:ALA:HB1	1.90	0.54
1:B:588:GLN:N	1:B:588:GLN:OE1	2.40	0.54
1:C:620:LEU:CD1	1:C:621:ASP:H	2.16	0.54
1:D:620:LEU:CD1	1:D:621:ASP:H	2.16	0.54
1:E:585:ILE:HD13	1:E:675:ALA:HB1	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:620:LEU:CD1	1:G:621:ASP:H	2.16	0.54
1:C:585:ILE:HD13	1:C:675:ALA:HB1	1.90	0.54
1:D:585:ILE:HD13	1:D:675:ALA:HB1	1.90	0.54
1:B:585:ILE:HD13	1:B:675:ALA:HB1	1.90	0.53
1:F:585:ILE:HD13	1:F:675:ALA:HB1	1.90	0.53
1:I:585:ILE:HD13	1:I:675:ALA:HB1	1.90	0.53
1:B:344:GLU:C	1:B:346:THR:H	2.16	0.53
1:D:344:GLU:C	1:D:346:THR:H	2.16	0.53
1:E:435:ALA:HA	1:E:477:VAL:HA	1.90	0.53
1:E:625:ILE:HG22	1:E:627:PRO:HD3	1.88	0.53
1:G:585:ILE:HD13	1:G:675:ALA:HB1	1.90	0.53
1:H:344:GLU:C	1:H:346:THR:H	2.16	0.53
1:H:435:ALA:HA	1:H:477:VAL:HA	1.89	0.53
1:B:402:ARG:HH12	1:B:404:ARG:HD3	1.74	0.53
1:C:435:ALA:HA	1:C:477:VAL:HA	1.89	0.53
1:D:405:ASP:OD1	1:D:405:ASP:N	2.40	0.53
1:F:435:ALA:HA	1:F:477:VAL:HA	1.89	0.53
1:A:585:ILE:HD13	1:A:675:ALA:HB1	1.90	0.53
1:I:405:ASP:OD1	1:I:405:ASP:N	2.40	0.53
1:C:344:GLU:C	1:C:346:THR:H	2.16	0.53
1:C:402:ARG:HH12	1:C:404:ARG:HD3	1.74	0.53
1:G:344:GLU:C	1:G:346:THR:H	2.16	0.53
1:D:402:ARG:HH12	1:D:404:ARG:HD3	1.74	0.53
1:E:445:ILE:HG23	1:E:447:GLY:H	1.74	0.53
1:F:402:ARG:HH12	1:F:404:ARG:HD3	1.74	0.53
1:D:435:ALA:HA	1:D:477:VAL:HA	1.90	0.53
1:G:435:ALA:HA	1:G:477:VAL:HA	1.90	0.53
1:H:402:ARG:HH12	1:H:404:ARG:HD3	1.74	0.53
1:B:435:ALA:HA	1:B:477:VAL:HA	1.89	0.53
1:F:344:GLU:C	1:F:346:THR:H	2.16	0.53
1:F:445:ILE:HG23	1:F:447:GLY:H	1.74	0.53
1:I:435:ALA:HA	1:I:477:VAL:HA	1.89	0.53
1:I:344:GLU:C	1:I:346:THR:H	2.16	0.53
1:I:402:ARG:HH12	1:I:404:ARG:HD3	1.74	0.53
1:A:402:ARG:HH12	1:A:404:ARG:HD3	1.74	0.52
1:E:344:GLU:C	1:E:346:THR:H	2.16	0.52
1:G:402:ARG:HH12	1:G:404:ARG:HD3	1.74	0.52
1:G:445:ILE:HG23	1:G:447:GLY:H	1.74	0.52
1:E:620:LEU:CD1	1:E:621:ASP:H	2.16	0.52
1:A:445:ILE:HG23	1:A:447:GLY:H	1.74	0.52
1:B:445:ILE:HG23	1:B:447:GLY:H	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:405:ASP:OD1	1:F:405:ASP:N	2.40	0.52
1:H:405:ASP:OD1	1:H:405:ASP:N	2.40	0.52
1:H:620:LEU:CD1	1:H:621:ASP:H	2.16	0.52
1:B:662:LEU:HD22	1:B:663:LEU:HD22	1.92	0.52
1:B:433:LEU:O	1:B:462:ILE:N	2.31	0.52
1:C:662:LEU:HD22	1:C:663:LEU:HD22	1.92	0.52
1:E:405:ASP:N	1:E:405:ASP:OD1	2.40	0.52
1:B:598:SER:OG	1:B:694:LEU:O	2.20	0.52
1:D:662:LEU:HD22	1:D:663:LEU:HD22	1.92	0.52
1:G:349:THR:N	1:G:352:ASP:OD1	2.34	0.52
1:G:405:ASP:OD1	1:G:405:ASP:N	2.40	0.52
1:C:445:ILE:HG23	1:C:447:GLY:H	1.74	0.52
1:H:429:GLU:CG	1:H:430:PRO:HD2	2.40	0.52
1:A:598:SER:OG	1:A:694:LEU:O	2.20	0.51
1:A:662:LEU:HD22	1:A:663:LEU:HD22	1.92	0.51
1:B:429:GLU:CG	1:B:430:PRO:HD2	2.41	0.51
1:E:402:ARG:HH12	1:E:404:ARG:HD3	1.74	0.51
1:F:662:LEU:HD22	1:F:663:LEU:HD22	1.92	0.51
1:G:429:GLU:CG	1:G:430:PRO:HD2	2.41	0.51
1:H:349:THR:N	1:H:352:ASP:OD1	2.34	0.51
1:A:344:GLU:C	1:A:346:THR:H	2.16	0.51
1:C:429:GLU:CG	1:C:430:PRO:HD2	2.41	0.51
1:E:662:LEU:HD22	1:E:663:LEU:HD22	1.92	0.51
1:I:349:THR:N	1:I:352:ASP:OD1	2.34	0.51
1:D:445:ILE:HG23	1:D:447:GLY:H	1.74	0.51
1:I:445:ILE:HG23	1:I:447:GLY:H	1.74	0.51
1:E:450:GLY:N	1:E:459:ALA:O	2.43	0.51
1:G:592:LEU:HD23	1:G:682:ILE:HD11	1.93	0.51
1:D:450:GLY:N	1:D:459:ALA:O	2.42	0.51
1:G:662:LEU:HD22	1:G:663:LEU:HD22	1.92	0.51
1:H:445:ILE:HG23	1:H:447:GLY:H	1.74	0.51
1:H:592:LEU:HD23	1:H:682:ILE:HD11	1.93	0.51
1:E:429:GLU:CG	1:E:430:PRO:HD2	2.40	0.51
1:I:662:LEU:HD22	1:I:663:LEU:HD22	1.92	0.51
1:D:429:GLU:CG	1:D:430:PRO:HD2	2.40	0.51
1:E:592:LEU:HD23	1:E:682:ILE:HD11	1.93	0.51
1:A:429:GLU:CG	1:A:430:PRO:HD2	2.40	0.51
1:C:450:GLY:N	1:C:459:ALA:O	2.42	0.51
1:H:433:LEU:O	1:H:462:ILE:N	2.31	0.51
1:G:450:GLY:N	1:G:459:ALA:O	2.43	0.51
1:A:450:GLY:N	1:A:459:ALA:O	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:LEU:CD1	1:A:621:ASP:H	2.16	0.50
1:F:450:GLY:N	1:F:459:ALA:O	2.43	0.50
1:I:429:GLU:CG	1:I:430:PRO:HD2	2.41	0.50
1:I:592:LEU:HD23	1:I:682:ILE:HD11	1.93	0.50
1:H:413:GLN:CD	1:H:425:GLN:HE21	2.19	0.50
1:H:662:LEU:HD22	1:H:663:LEU:HD22	1.92	0.50
1:F:429:GLU:CG	1:F:430:PRO:HD2	2.40	0.50
1:F:592:LEU:HD23	1:F:682:ILE:HD11	1.93	0.50
1:F:680:ASN:N	1:F:680:ASN:OD1	2.45	0.50
1:G:413:GLN:OE1	1:G:425:GLN:NE2	2.45	0.50
1:E:413:GLN:CD	1:E:425:GLN:HE21	2.19	0.50
1:B:504:ASP:O	1:B:507:ALA:N	2.45	0.50
1:G:504:ASP:O	1:G:507:ALA:N	2.45	0.50
1:H:504:ASP:O	1:H:507:ALA:N	2.45	0.50
1:C:413:GLN:CD	1:C:425:GLN:HE21	2.19	0.49
1:C:504:ASP:O	1:C:507:ALA:N	2.45	0.49
1:C:539:LEU:HD11	1:C:573:VAL:HG11	1.94	0.49
1:C:660:ARG:NE	1:C:676:VAL:O	2.45	0.49
1:D:660:ARG:NE	1:D:676:VAL:O	2.45	0.49
1:E:504:ASP:O	1:E:507:ALA:N	2.45	0.49
1:H:450:GLY:HA3	1:H:461:TRP:HE1	1.77	0.49
1:I:539:LEU:HD11	1:I:573:VAL:HG11	1.94	0.49
1:A:504:ASP:O	1:A:507:ALA:N	2.45	0.49
1:A:539:LEU:HD11	1:A:573:VAL:HG11	1.94	0.49
1:C:592:LEU:HD23	1:C:682:ILE:HD11	1.93	0.49
1:D:484:ILE:O	1:D:487:HIS:N	2.45	0.49
1:F:413:GLN:CD	1:F:425:GLN:HE21	2.19	0.49
1:F:539:LEU:HD11	1:F:573:VAL:HG11	1.94	0.49
1:G:413:GLN:CD	1:G:425:GLN:HE21	2.19	0.49
1:G:539:LEU:HD11	1:G:573:VAL:HG11	1.94	0.49
1:I:413:GLN:OE1	1:I:425:GLN:NE2	2.45	0.49
1:I:504:ASP:O	1:I:507:ALA:N	2.45	0.49
1:A:592:LEU:HD23	1:A:682:ILE:HD11	1.93	0.49
1:B:450:GLY:N	1:B:459:ALA:O	2.43	0.49
1:B:539:LEU:HD11	1:B:573:VAL:HG11	1.94	0.49
1:B:680:ASN:OD1	1:B:680:ASN:N	2.45	0.49
1:C:484:ILE:O	1:C:487:HIS:N	2.46	0.49
1:D:413:GLN:CD	1:D:425:GLN:HE21	2.19	0.49
1:D:504:ASP:O	1:D:507:ALA:N	2.45	0.49
1:H:539:LEU:HD11	1:H:573:VAL:HG11	1.94	0.49
1:B:620:LEU:CD1	1:B:621:ASP:H	2.16	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:539:LEU:HD11	1:D:573:VAL:HG11	1.94	0.49
1:D:592:LEU:HD23	1:D:682:ILE:HD11	1.93	0.49
1:F:450:GLY:HA3	1:F:461:TRP:HE1	1.77	0.49
1:F:504:ASP:O	1:F:507:ALA:N	2.45	0.49
1:G:438:VAL:HB	1:G:476:GLN:HE22	1.77	0.49
1:G:660:ARG:NE	1:G:676:VAL:O	2.45	0.49
1:A:450:GLY:HA3	1:A:461:TRP:HE1	1.77	0.49
1:E:349:THR:N	1:E:352:ASP:OD1	2.34	0.49
1:H:413:GLN:OE1	1:H:425:GLN:NE2	2.45	0.49
1:A:413:GLN:OE1	1:A:425:GLN:NE2	2.45	0.49
1:B:484:ILE:O	1:B:487:HIS:N	2.45	0.49
1:B:592:LEU:HD23	1:B:682:ILE:HD11	1.93	0.49
1:B:660:ARG:NE	1:B:676:VAL:O	2.45	0.49
1:F:484:ILE:O	1:F:487:HIS:N	2.45	0.49
1:G:357:HIS:CG	1:G:402:ARG:HE	2.31	0.49
1:H:450:GLY:N	1:H:459:ALA:O	2.43	0.49
1:I:438:VAL:HB	1:I:476:GLN:HE22	1.78	0.49
1:I:450:GLY:HA3	1:I:461:TRP:HE1	1.77	0.49
1:A:413:GLN:CD	1:A:425:GLN:HE21	2.19	0.49
1:A:438:VAL:HB	1:A:476:GLN:HE22	1.77	0.49
1:D:438:VAL:HB	1:D:476:GLN:HE22	1.78	0.49
1:G:444:GLU:HA	1:G:475:TYR:OH	2.13	0.49
1:A:680:ASN:OD1	1:A:680:ASN:N	2.45	0.49
1:B:413:GLN:CD	1:B:425:GLN:HE21	2.19	0.49
1:C:450:GLY:HA3	1:C:461:TRP:HE1	1.77	0.49
1:E:413:GLN:OE1	1:E:425:GLN:NE2	2.44	0.49
1:E:539:LEU:HD11	1:E:573:VAL:HG11	1.94	0.49
1:E:660:ARG:NE	1:E:676:VAL:O	2.45	0.49
1:G:450:GLY:HA3	1:G:461:TRP:HE1	1.77	0.49
1:H:444:GLU:HA	1:H:475:TYR:OH	2.13	0.49
1:C:438:VAL:HB	1:C:476:GLN:HE22	1.78	0.49
1:D:450:GLY:HA3	1:D:461:TRP:HE1	1.77	0.49
1:E:438:VAL:HB	1:E:476:GLN:HE22	1.78	0.49
1:E:444:GLU:HA	1:E:475:TYR:OH	2.13	0.49
1:E:680:ASN:OD1	1:E:680:ASN:N	2.45	0.49
1:F:438:VAL:HB	1:F:476:GLN:HE22	1.78	0.49
1:G:680:ASN:OD1	1:G:680:ASN:N	2.45	0.49
1:B:357:HIS:CG	1:B:402:ARG:HE	2.31	0.49
1:C:528:ASN:O	1:C:531:GLN:N	2.45	0.49
1:C:680:ASN:OD1	1:C:680:ASN:N	2.45	0.49
1:E:357:HIS:CG	1:E:402:ARG:HE	2.31	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:444:GLU:HA	1:F:475:TYR:OH	2.13	0.49
1:I:413:GLN:CD	1:I:425:GLN:HE21	2.19	0.49
1:I:444:GLU:HA	1:I:475:TYR:OH	2.13	0.49
1:B:438:VAL:HB	1:B:476:GLN:HE22	1.78	0.48
1:E:622:SER:O	1:E:662:LEU:HG	2.13	0.48
1:A:484:ILE:O	1:A:487:HIS:N	2.45	0.48
1:B:622:SER:O	1:B:662:LEU:HG	2.13	0.48
1:E:484:ILE:O	1:E:487:HIS:N	2.45	0.48
1:F:660:ARG:C	1:F:660:ARG:HH11	2.22	0.48
1:A:622:SER:O	1:A:662:LEU:HG	2.13	0.48
1:C:622:SER:O	1:C:662:LEU:HG	2.13	0.48
1:D:622:SER:O	1:D:662:LEU:HG	2.13	0.48
1:E:450:GLY:HA3	1:E:461:TRP:HE1	1.77	0.48
1:H:680:ASN:OD1	1:H:680:ASN:N	2.45	0.48
1:A:444:GLU:HA	1:A:475:TYR:OH	2.13	0.48
1:C:357:HIS:CG	1:C:402:ARG:HE	2.31	0.48
1:D:357:HIS:CG	1:D:402:ARG:HE	2.31	0.48
1:F:413:GLN:OE1	1:F:425:GLN:NE2	2.44	0.48
1:H:357:HIS:CG	1:H:402:ARG:HE	2.31	0.48
1:I:528:ASN:O	1:I:531:GLN:N	2.45	0.48
1:I:622:SER:O	1:I:662:LEU:HG	2.13	0.48
1:A:660:ARG:C	1:A:660:ARG:HH11	2.22	0.48
1:F:620:LEU:CD1	1:F:621:ASP:H	2.16	0.48
1:G:622:SER:O	1:G:662:LEU:HG	2.13	0.48
1:A:349:THR:N	1:A:352:ASP:OD1	2.34	0.48
1:F:357:HIS:CG	1:F:402:ARG:HE	2.31	0.48
1:F:433:LEU:O	1:F:462:ILE:N	2.31	0.48
1:G:484:ILE:O	1:G:487:HIS:N	2.45	0.48
1:G:660:ARG:C	1:G:660:ARG:HH11	2.22	0.48
1:I:450:GLY:N	1:I:459:ALA:O	2.43	0.48
1:I:484:ILE:O	1:I:487:HIS:N	2.45	0.48
1:A:660:ARG:NE	1:A:676:VAL:O	2.45	0.48
1:B:348:LEU:HD23	1:B:348:LEU:HA	1.69	0.48
1:C:372:ASN:OD1	1:C:372:ASN:N	2.46	0.48
1:C:444:GLU:HA	1:C:475:TYR:OH	2.13	0.48
1:D:444:GLU:HA	1:D:475:TYR:OH	2.13	0.48
1:D:623:PHE:HB3	1:D:665:ARG:NH2	2.27	0.48
1:E:660:ARG:C	1:E:660:ARG:HH11	2.22	0.48
1:F:622:SER:O	1:F:662:LEU:HG	2.13	0.48
1:H:672:GLN:OE1	1:H:672:GLN:N	2.43	0.48
1:B:444:GLU:HA	1:B:475:TYR:OH	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:595:TYR:CE2	1:B:640:ILE:HG23	2.49	0.48
1:G:372:ASN:OD1	1:G:372:ASN:N	2.46	0.48
1:H:484:ILE:O	1:H:487:HIS:N	2.45	0.48
1:H:622:SER:O	1:H:662:LEU:HG	2.13	0.48
1:B:450:GLY:HA3	1:B:461:TRP:HE1	1.77	0.48
1:F:468:ALA:O	1:F:472:ASN:N	2.43	0.48
1:G:468:ALA:O	1:G:472:ASN:N	2.43	0.48
1:G:595:TYR:CE2	1:G:640:ILE:HG23	2.49	0.48
1:H:372:ASN:OD1	1:H:372:ASN:N	2.46	0.48
1:H:595:TYR:CE2	1:H:640:ILE:HG23	2.49	0.48
1:A:357:HIS:CG	1:A:402:ARG:HE	2.31	0.48
1:D:660:ARG:HH11	1:D:660:ARG:C	2.22	0.48
1:I:357:HIS:CG	1:I:402:ARG:HE	2.31	0.48
1:E:595:TYR:CE2	1:E:640:ILE:HG23	2.49	0.47
1:F:660:ARG:NE	1:F:676:VAL:O	2.45	0.47
1:G:448:ILE:CG2	1:G:461:TRP:HB2	2.44	0.47
1:H:438:VAL:HB	1:H:476:GLN:HE22	1.78	0.47
1:H:448:ILE:CG2	1:H:461:TRP:HB2	2.44	0.47
1:H:660:ARG:C	1:H:660:ARG:HH11	2.22	0.47
1:I:660:ARG:C	1:I:660:ARG:HH11	2.22	0.47
1:C:595:TYR:CE2	1:C:640:ILE:HG23	2.49	0.47
1:G:623:PHE:HB3	1:G:665:ARG:NH2	2.27	0.47
1:I:623:PHE:HB3	1:I:665:ARG:NH2	2.27	0.47
1:B:413:GLN:OE1	1:B:425:GLN:NE2	2.45	0.47
1:D:528:ASN:O	1:D:531:GLN:N	2.45	0.47
1:E:448:ILE:CG2	1:E:461:TRP:HB2	2.44	0.47
1:A:595:TYR:CE2	1:A:640:ILE:HG23	2.49	0.47
1:A:623:PHE:HB3	1:A:665:ARG:NH2	2.27	0.47
1:C:413:GLN:OE1	1:C:425:GLN:NE2	2.45	0.47
1:I:448:ILE:CG2	1:I:461:TRP:HB2	2.44	0.47
1:E:468:ALA:O	1:E:472:ASN:N	2.43	0.47
1:I:595:TYR:CE2	1:I:640:ILE:HG23	2.49	0.47
1:B:660:ARG:C	1:B:660:ARG:HH11	2.22	0.47
1:D:413:GLN:OE1	1:D:425:GLN:NE2	2.44	0.47
1:D:438:VAL:HB	1:D:476:GLN:NE2	2.30	0.47
1:D:595:TYR:CE2	1:D:640:ILE:HG23	2.49	0.47
1:F:448:ILE:CG2	1:F:461:TRP:HB2	2.44	0.47
1:A:468:ALA:O	1:A:472:ASN:N	2.43	0.47
1:B:349:THR:N	1:B:352:ASP:OD1	2.34	0.47
1:E:415:THR:HG22	1:E:425:GLN:HG3	1.97	0.47
1:F:349:THR:N	1:F:352:ASP:OD1	2.34	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:528:ASN:O	1:F:531:GLN:N	2.45	0.47
1:F:595:TYR:CE2	1:F:640:ILE:HG23	2.49	0.47
1:H:415:THR:HG22	1:H:425:GLN:HG3	1.97	0.47
1:H:468:ALA:O	1:H:472:ASN:N	2.43	0.47
1:H:623:PHE:HB3	1:H:665:ARG:NH2	2.27	0.47
1:I:680:ASN:OD1	1:I:680:ASN:N	2.45	0.47
1:A:415:THR:HG22	1:A:425:GLN:HG3	1.97	0.47
1:A:448:ILE:CG2	1:A:461:TRP:HB2	2.44	0.47
1:I:415:THR:HG22	1:I:425:GLN:HG3	1.97	0.47
1:B:438:VAL:HB	1:B:476:GLN:NE2	2.30	0.47
1:D:680:ASN:N	1:D:680:ASN:OD1	2.45	0.47
1:E:452:ASP:N	1:E:457:LEU:O	2.48	0.47
1:F:623:PHE:HB3	1:F:665:ARG:NH2	2.27	0.47
1:G:390:LEU:O	1:G:394:VAL:HG22	2.15	0.47
1:G:415:THR:HG22	1:G:425:GLN:HG3	1.97	0.47
1:A:438:VAL:HB	1:A:476:GLN:NE2	2.30	0.47
1:C:660:ARG:C	1:C:660:ARG:HH11	2.22	0.47
1:D:372:ASN:N	1:D:372:ASN:OD1	2.46	0.47
1:E:390:LEU:O	1:E:394:VAL:HG22	2.15	0.47
1:E:438:VAL:HB	1:E:476:GLN:NE2	2.30	0.47
1:F:452:ASP:N	1:F:457:LEU:O	2.48	0.47
1:H:660:ARG:NE	1:H:676:VAL:O	2.45	0.47
1:I:390:LEU:O	1:I:394:VAL:HG22	2.15	0.47
1:B:528:ASN:O	1:B:531:GLN:N	2.45	0.46
1:D:468:ALA:O	1:D:472:ASN:N	2.43	0.46
1:F:390:LEU:O	1:F:394:VAL:HG22	2.15	0.46
1:G:438:VAL:HB	1:G:476:GLN:NE2	2.30	0.46
1:B:415:THR:HG22	1:B:425:GLN:HG3	1.97	0.46
1:D:448:ILE:CG2	1:D:461:TRP:HB2	2.44	0.46
1:E:350:TRP:CH2	1:G:396:PHE:HB2	2.50	0.46
1:F:415:THR:HG22	1:F:425:GLN:HG3	1.97	0.46
1:I:660:ARG:NE	1:I:676:VAL:O	2.45	0.46
1:C:415:THR:HG22	1:C:425:GLN:HG3	1.97	0.46
1:D:415:THR:HG22	1:D:425:GLN:HG3	1.97	0.46
1:D:660:ARG:NH1	1:D:664:ALA:HB2	2.31	0.46
1:H:660:ARG:NH1	1:H:664:ALA:HB2	2.31	0.46
1:I:660:ARG:NH1	1:I:664:ALA:HB2	2.31	0.46
1:I:672:GLN:OE1	1:I:672:GLN:N	2.43	0.46
1:B:448:ILE:CG2	1:B:461:TRP:HB2	2.44	0.46
1:C:438:VAL:HB	1:C:476:GLN:NE2	2.30	0.46
1:E:528:ASN:OD1	1:E:528:ASN:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:438:VAL:HB	1:F:476:GLN:NE2	2.30	0.46
1:F:528:ASN:N	1:F:528:ASN:OD1	2.48	0.46
1:F:660:ARG:NH1	1:F:664:ALA:HB2	2.31	0.46
1:G:528:ASN:OD1	1:G:528:ASN:N	2.48	0.46
1:H:390:LEU:O	1:H:394:VAL:HG22	2.15	0.46
1:C:656:MET:HE3	1:C:656:MET:HB3	1.87	0.46
1:C:660:ARG:NH1	1:C:664:ALA:HB2	2.31	0.46
1:D:672:GLN:OE1	1:D:672:GLN:N	2.43	0.46
1:H:434:MET:HG3	1:H:459:ALA:HB1	1.98	0.46
1:I:372:ASN:N	1:I:372:ASN:OD1	2.46	0.46
1:C:448:ILE:CG2	1:C:461:TRP:HB2	2.44	0.46
1:D:390:LEU:O	1:D:394:VAL:HG22	2.15	0.46
1:I:438:VAL:HG13	1:I:440:ASP:H	1.81	0.46
1:A:390:LEU:O	1:A:394:VAL:HG22	2.15	0.46
1:A:434:MET:HG3	1:A:459:ALA:HB1	1.98	0.46
1:A:660:ARG:NH1	1:A:664:ALA:HB2	2.31	0.46
1:B:434:MET:HG3	1:B:459:ALA:HB1	1.98	0.46
1:F:372:ASN:OD1	1:F:372:ASN:N	2.46	0.46
1:G:350:TRP:CH2	1:H:396:PHE:HB2	2.51	0.46
1:G:434:MET:HG3	1:G:459:ALA:HB1	1.98	0.46
1:H:438:VAL:HB	1:H:476:GLN:NE2	2.30	0.46
1:I:434:MET:HG3	1:I:459:ALA:HB1	1.98	0.46
1:B:390:LEU:O	1:B:394:VAL:HG22	2.15	0.46
1:G:660:ARG:NH1	1:G:664:ALA:HB2	2.31	0.46
1:H:438:VAL:HG13	1:H:440:ASP:H	1.81	0.46
1:H:528:ASN:O	1:H:531:GLN:N	2.45	0.46
1:A:438:VAL:HG13	1:A:440:ASP:H	1.81	0.46
1:A:452:ASP:N	1:A:457:LEU:O	2.48	0.46
1:B:452:ASP:N	1:B:457:LEU:O	2.48	0.46
1:C:434:MET:HG3	1:C:459:ALA:HB1	1.98	0.46
1:E:434:MET:HG3	1:E:459:ALA:HB1	1.98	0.46
1:G:541:LEU:HD23	1:G:541:LEU:HA	1.76	0.46
1:B:623:PHE:HB3	1:B:665:ARG:NH2	2.27	0.45
1:F:434:MET:HG3	1:F:459:ALA:HB1	1.98	0.45
1:G:528:ASN:O	1:G:531:GLN:N	2.45	0.45
1:I:438:VAL:HB	1:I:476:GLN:NE2	2.30	0.45
1:B:438:VAL:HG13	1:B:440:ASP:H	1.81	0.45
1:C:350:TRP:CH2	1:D:396:PHE:HB2	2.51	0.45
1:D:528:ASN:OD1	1:D:528:ASN:N	2.48	0.45
1:F:672:GLN:OE1	1:F:672:GLN:N	2.43	0.45
1:B:554:ASN:C	1:B:554:ASN:HD22	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:660:ARG:NH1	1:B:664:ALA:HB2	2.31	0.45
1:C:390:LEU:O	1:C:394:VAL:HG22	2.15	0.45
1:D:434:MET:HG3	1:D:459:ALA:HB1	1.98	0.45
1:E:656:MET:HE3	1:E:656:MET:HB3	1.87	0.45
1:E:660:ARG:NH1	1:E:664:ALA:HB2	2.31	0.45
1:C:438:VAL:HG13	1:C:440:ASP:H	1.81	0.45
1:D:438:VAL:HG13	1:D:440:ASP:H	1.81	0.45
1:E:372:ASN:OD1	1:E:372:ASN:N	2.46	0.45
1:H:452:ASP:N	1:H:457:LEU:O	2.48	0.45
1:H:528:ASN:N	1:H:528:ASN:OD1	2.48	0.45
1:H:637:LEU:HA	1:H:640:ILE:HB	1.99	0.45
1:A:554:ASN:C	1:A:554:ASN:HD22	2.25	0.45
1:B:350:TRP:CH2	1:C:396:PHE:HB2	2.51	0.45
1:B:637:LEU:HA	1:B:640:ILE:HB	1.99	0.45
1:D:349:THR:N	1:D:352:ASP:OD1	2.34	0.45
1:E:396:PHE:HB2	1:F:350:TRP:CH2	2.51	0.45
1:A:567:ILE:HD13	1:A:620:LEU:HB2	1.99	0.45
1:A:637:LEU:HA	1:A:640:ILE:HB	1.99	0.45
1:A:672:GLN:OE1	1:A:672:GLN:N	2.43	0.45
1:F:438:VAL:HG13	1:F:440:ASP:H	1.81	0.45
1:G:438:VAL:HG13	1:G:440:ASP:H	1.81	0.45
1:H:479:ASP:O	1:H:483:VAL:HG23	2.17	0.45
1:A:396:PHE:HB2	1:I:350:TRP:CH2	2.52	0.45
1:B:479:ASP:O	1:B:483:VAL:HG23	2.17	0.45
1:B:600:GLU:O	1:B:604:MET:HG3	2.17	0.45
1:D:452:ASP:N	1:D:457:LEU:O	2.48	0.45
1:E:637:LEU:HA	1:E:640:ILE:HB	1.99	0.45
1:G:600:GLU:O	1:G:604:MET:HG3	2.17	0.45
1:G:637:LEU:HA	1:G:640:ILE:HB	1.99	0.45
1:G:654:LEU:HB3	1:G:679:TYR:CE2	2.48	0.45
1:I:438:VAL:N	1:I:476:GLN:OE1	2.50	0.45
1:I:600:GLU:O	1:I:604:MET:HG3	2.17	0.45
1:A:479:ASP:O	1:A:483:VAL:HG23	2.17	0.45
1:D:567:ILE:HD13	1:D:620:LEU:HB2	1.99	0.45
1:F:438:VAL:N	1:F:476:GLN:OE1	2.50	0.45
1:H:605:LEU:O	1:H:608:SER:N	2.50	0.45
1:I:567:ILE:HD13	1:I:620:LEU:HB2	1.99	0.45
1:A:438:VAL:N	1:A:476:GLN:OE1	2.50	0.45
1:A:600:GLU:O	1:A:604:MET:HG3	2.17	0.45
1:B:672:GLN:OE1	1:B:672:GLN:N	2.43	0.45
1:E:438:VAL:N	1:E:476:GLN:OE1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:603:GLN:HA	1:E:606:LEU:HD12	1.99	0.45
1:G:452:ASP:N	1:G:457:LEU:O	2.48	0.45
1:G:523:LEU:HD22	1:G:553:ALA:HB1	1.99	0.45
1:I:637:LEU:HA	1:I:640:ILE:HB	1.99	0.45
1:A:348:LEU:HD23	1:A:348:LEU:HA	1.69	0.45
1:A:381:GLN:O	1:A:385:GLY:N	2.48	0.45
1:A:548:ASP:OD1	1:A:548:ASP:N	2.50	0.45
1:C:528:ASN:N	1:C:528:ASN:OD1	2.48	0.45
1:E:496:LEU:HA	1:E:496:LEU:HD23	1.74	0.45
1:H:350:TRP:CH2	1:I:396:PHE:HB2	2.52	0.45
1:I:605:LEU:O	1:I:608:SER:N	2.50	0.45
1:B:381:GLN:O	1:B:385:GLY:N	2.48	0.44
1:B:523:LEU:HD22	1:B:553:ALA:HB1	2.00	0.44
1:B:567:ILE:HD13	1:B:620:LEU:HB2	1.99	0.44
1:E:479:ASP:O	1:E:483:VAL:HG23	2.17	0.44
1:E:528:ASN:O	1:E:531:GLN:N	2.45	0.44
1:E:605:LEU:O	1:E:608:SER:N	2.50	0.44
1:G:567:ILE:HD13	1:G:620:LEU:HB2	1.99	0.44
1:I:479:ASP:O	1:I:483:VAL:HG23	2.17	0.44
1:A:350:TRP:CH2	1:B:396:PHE:HB2	2.51	0.44
1:B:548:ASP:OD1	1:B:548:ASP:N	2.50	0.44
1:C:349:THR:N	1:C:352:ASP:OD1	2.34	0.44
1:C:381:GLN:O	1:C:385:GLY:N	2.48	0.44
1:C:637:LEU:HA	1:C:640:ILE:HB	1.99	0.44
1:D:479:ASP:O	1:D:483:VAL:HG23	2.17	0.44
1:D:600:GLU:O	1:D:604:MET:HG3	2.17	0.44
1:D:605:LEU:O	1:D:608:SER:N	2.50	0.44
1:E:582:VAL:HG21	1:E:682:ILE:HG13	2.00	0.44
1:F:567:ILE:HD13	1:F:620:LEU:HB2	1.99	0.44
1:H:603:GLN:HA	1:H:606:LEU:HD12	1.99	0.44
1:F:479:ASP:O	1:F:483:VAL:HG23	2.17	0.44
1:F:502:HIS:HA	1:F:533:LEU:HD13	2.00	0.44
1:F:605:LEU:O	1:F:608:SER:N	2.50	0.44
1:G:582:VAL:HG21	1:G:682:ILE:HG13	2.00	0.44
1:H:438:VAL:N	1:H:476:GLN:OE1	2.50	0.44
1:H:656:MET:HE3	1:H:656:MET:HB3	1.87	0.44
1:I:452:ASP:N	1:I:457:LEU:O	2.48	0.44
1:I:523:LEU:HD22	1:I:553:ALA:HB1	2.00	0.44
1:A:523:LEU:HD22	1:A:553:ALA:HB1	2.00	0.44
1:C:598:SER:HA	1:C:693:ASN:OD1	2.18	0.44
1:D:637:LEU:HA	1:D:640:ILE:HB	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:660:ARG:HA	1:D:660:ARG:HD2	1.87	0.44
1:E:569:LEU:HD23	1:E:569:LEU:HA	1.78	0.44
1:F:603:GLN:HA	1:F:606:LEU:HD12	1.99	0.44
1:H:523:LEU:HD22	1:H:553:ALA:HB1	1.99	0.44
1:I:528:ASN:OD1	1:I:528:ASN:N	2.48	0.44
1:C:523:LEU:HD22	1:C:553:ALA:HB1	2.00	0.44
1:C:605:LEU:O	1:C:608:SER:N	2.50	0.44
1:D:598:SER:HA	1:D:693:ASN:OD1	2.18	0.44
1:E:438:VAL:HG13	1:E:440:ASP:H	1.81	0.44
1:F:554:ASN:C	1:F:554:ASN:HD22	2.25	0.44
1:G:554:ASN:C	1:G:554:ASN:HD22	2.25	0.44
1:I:548:ASP:OD1	1:I:548:ASP:N	2.50	0.44
1:B:605:LEU:O	1:B:608:SER:N	2.50	0.44
1:C:554:ASN:C	1:C:554:ASN:HD22	2.25	0.44
1:F:523:LEU:HD22	1:F:553:ALA:HB1	2.00	0.44
1:G:348:LEU:HA	1:G:348:LEU:HD23	1.69	0.44
1:G:438:VAL:N	1:G:476:GLN:OE1	2.50	0.44
1:B:598:SER:HA	1:B:693:ASN:OD1	2.18	0.44
1:C:479:ASP:O	1:C:483:VAL:HG23	2.17	0.44
1:E:523:LEU:HD22	1:E:553:ALA:HB1	2.00	0.44
1:E:541:LEU:HD23	1:E:541:LEU:HA	1.76	0.44
1:E:598:SER:HA	1:E:693:ASN:OD1	2.18	0.44
1:F:541:LEU:HD23	1:F:541:LEU:HA	1.76	0.44
1:F:598:SER:HA	1:F:693:ASN:OD1	2.18	0.44
1:F:600:GLU:O	1:F:604:MET:HG3	2.17	0.44
1:F:637:LEU:HA	1:F:640:ILE:HB	1.99	0.44
1:G:569:LEU:HD23	1:G:569:LEU:HA	1.78	0.44
1:B:528:ASN:OD1	1:B:528:ASN:N	2.48	0.44
1:C:438:VAL:N	1:C:476:GLN:OE1	2.50	0.44
1:D:502:HIS:HA	1:D:533:LEU:HD13	2.00	0.44
1:D:523:LEU:HD22	1:D:553:ALA:HB1	2.00	0.44
1:I:554:ASN:C	1:I:554:ASN:HD22	2.25	0.44
1:A:372:ASN:OD1	1:A:372:ASN:N	2.46	0.44
1:A:528:ASN:N	1:A:528:ASN:OD1	2.48	0.44
1:B:438:VAL:N	1:B:476:GLN:OE1	2.50	0.44
1:C:468:ALA:O	1:C:472:ASN:N	2.43	0.44
1:E:600:GLU:O	1:E:604:MET:HG3	2.17	0.44
1:G:479:ASP:O	1:G:483:VAL:HG23	2.17	0.44
1:I:381:GLN:O	1:I:385:GLY:N	2.48	0.44
1:I:603:GLN:HA	1:I:606:LEU:HD12	1.99	0.44
1:B:603:GLN:HA	1:B:606:LEU:HD12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:567:ILE:HD13	1:C:620:LEU:HB2	1.99	0.43
1:D:381:GLN:O	1:D:385:GLY:N	2.48	0.43
1:F:582:VAL:HG21	1:F:682:ILE:HG13	2.00	0.43
1:H:600:GLU:O	1:H:604:MET:HG3	2.17	0.43
1:D:438:VAL:N	1:D:476:GLN:OE1	2.50	0.43
1:E:567:ILE:HD13	1:E:620:LEU:HB2	1.99	0.43
1:I:361:LEU:HD23	1:I:361:LEU:HA	1.90	0.43
1:I:502:HIS:HA	1:I:533:LEU:HD13	1.99	0.43
1:A:582:VAL:HG21	1:A:682:ILE:HG13	2.00	0.43
1:A:598:SER:HB3	1:A:694:LEU:H	1.83	0.43
1:A:598:SER:HA	1:A:693:ASN:OD1	2.18	0.43
1:B:594:VAL:HG23	1:B:652:ILE:HB	2.01	0.43
1:C:600:GLU:O	1:C:604:MET:HG3	2.17	0.43
1:C:603:GLN:HA	1:C:606:LEU:HD12	1.99	0.43
1:G:603:GLN:HA	1:G:606:LEU:HD12	1.99	0.43
1:H:502:HIS:HA	1:H:533:LEU:HD13	2.00	0.43
1:H:645:LYS:HD2	1:H:645:LYS:HA	1.76	0.43
1:I:582:VAL:HG21	1:I:682:ILE:HG13	2.00	0.43
1:B:598:SER:HB3	1:B:694:LEU:H	1.84	0.43
1:C:582:VAL:HG21	1:C:682:ILE:HG13	2.00	0.43
1:E:348:LEU:HA	1:E:348:LEU:HD23	1.69	0.43
1:F:598:SER:HB3	1:F:694:LEU:H	1.84	0.43
1:G:598:SER:HA	1:G:693:ASN:OD1	2.18	0.43
1:H:567:ILE:HD13	1:H:620:LEU:HB2	1.99	0.43
1:I:469:LYS:O	1:I:473:MET:HB2	2.19	0.43
1:A:594:VAL:HG23	1:A:652:ILE:HB	2.01	0.43
1:A:603:GLN:HA	1:A:606:LEU:HD12	1.99	0.43
1:D:469:LYS:O	1:D:473:MET:HB2	2.19	0.43
1:D:603:GLN:HA	1:D:606:LEU:HD12	1.99	0.43
1:A:469:LYS:O	1:A:473:MET:HB2	2.19	0.43
1:D:554:ASN:C	1:D:554:ASN:HD22	2.25	0.43
1:D:656:MET:HE3	1:D:656:MET:HB3	1.87	0.43
1:E:502:HIS:HA	1:E:533:LEU:HD13	2.00	0.43
1:E:594:VAL:HG23	1:E:652:ILE:HB	2.01	0.43
1:E:598:SER:HB3	1:E:694:LEU:H	1.84	0.43
1:E:623:PHE:HB3	1:E:665:ARG:NH2	2.27	0.43
1:G:605:LEU:O	1:G:608:SER:N	2.50	0.43
1:H:582:VAL:HG21	1:H:682:ILE:HG13	2.00	0.43
1:H:598:SER:HB3	1:H:694:LEU:H	1.84	0.43
1:B:502:HIS:HA	1:B:533:LEU:HD13	1.99	0.43
1:D:502:HIS:HB3	1:F:519:LEU:HD12	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:469:LYS:O	1:F:473:MET:HB2	2.19	0.43
1:I:598:SER:HA	1:I:693:ASN:OD1	2.18	0.43
1:A:528:ASN:O	1:A:531:GLN:N	2.45	0.43
1:A:542:ASP:N	1:A:542:ASP:OD1	2.52	0.43
1:E:540:LEU:HD23	1:E:540:LEU:HA	1.80	0.43
1:F:594:VAL:HG23	1:F:652:ILE:HB	2.01	0.43
1:G:469:LYS:O	1:G:473:MET:HB2	2.19	0.43
1:H:594:VAL:HG23	1:H:652:ILE:HB	2.01	0.43
1:I:542:ASP:OD1	1:I:542:ASP:N	2.52	0.43
1:C:594:VAL:HG23	1:C:652:ILE:HB	2.01	0.43
1:E:672:GLN:OE1	1:E:672:GLN:N	2.43	0.43
1:H:667:ALA:HB1	1:H:674:LEU:HB3	2.01	0.43
1:B:656:MET:HE3	1:B:656:MET:HB3	1.87	0.43
1:C:598:SER:HB3	1:C:694:LEU:H	1.84	0.43
1:D:582:VAL:HG21	1:D:682:ILE:HG13	2.00	0.43
1:E:554:ASN:C	1:E:554:ASN:HD22	2.25	0.43
1:H:554:ASN:C	1:H:554:ASN:HD22	2.25	0.43
1:I:468:ALA:O	1:I:472:ASN:N	2.43	0.43
1:A:502:HIS:HA	1:A:533:LEU:HD13	1.99	0.42
1:A:605:LEU:O	1:A:608:SER:N	2.50	0.42
1:B:469:LYS:O	1:B:473:MET:HB2	2.19	0.42
1:C:378:PRO:O	1:C:382:ARG:N	2.51	0.42
1:C:540:LEU:HD23	1:C:540:LEU:HA	1.80	0.42
1:G:502:HIS:HB3	1:H:519:LEU:HD12	2.01	0.42
1:H:348:LEU:HA	1:H:348:LEU:HD23	1.69	0.42
1:F:378:PRO:O	1:F:382:ARG:N	2.51	0.42
1:G:667:ALA:HB1	1:G:674:LEU:HB3	2.01	0.42
1:I:594:VAL:HG21	1:I:654:LEU:HG	2.02	0.42
1:A:585:ILE:HD12	1:A:677:LEU:HD11	2.02	0.42
1:B:502:HIS:HB3	1:C:519:LEU:HD12	2.02	0.42
1:C:672:GLN:OE1	1:C:672:GLN:N	2.43	0.42
1:D:548:ASP:OD1	1:D:548:ASP:N	2.50	0.42
1:D:594:VAL:HG21	1:D:654:LEU:HG	2.02	0.42
1:E:350:TRP:HH2	1:G:396:PHE:HB2	1.85	0.42
1:E:594:VAL:HG21	1:E:654:LEU:HG	2.02	0.42
1:E:645:LYS:HD2	1:E:645:LYS:HA	1.77	0.42
1:G:502:HIS:HA	1:G:533:LEU:HD13	2.00	0.42
1:H:585:ILE:HD12	1:H:677:LEU:HD11	2.02	0.42
1:I:667:ALA:HB1	1:I:674:LEU:HB3	2.01	0.42
1:B:496:LEU:HD23	1:B:496:LEU:HA	1.74	0.42
1:B:582:VAL:HG21	1:B:682:ILE:HG13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:623:PHE:HB3	1:C:665:ARG:NH2	2.27	0.42
1:E:469:LYS:O	1:E:473:MET:HB2	2.19	0.42
1:E:654:LEU:HB3	1:E:679:TYR:CE2	2.48	0.42
1:F:367:LEU:O	1:F:370:LEU:HD12	2.19	0.42
1:F:594:VAL:HG21	1:F:654:LEU:HG	2.02	0.42
1:F:645:LYS:HD2	1:F:645:LYS:HA	1.76	0.42
1:G:598:SER:HB3	1:G:694:LEU:H	1.84	0.42
1:H:598:SER:HA	1:H:693:ASN:OD1	2.18	0.42
1:B:585:ILE:HD12	1:B:677:LEU:HD11	2.02	0.42
1:C:469:LYS:O	1:C:473:MET:HB2	2.19	0.42
1:C:585:ILE:HD12	1:C:677:LEU:HD11	2.02	0.42
1:E:502:HIS:HB3	1:G:519:LEU:HD12	2.01	0.42
1:G:585:ILE:HD12	1:G:677:LEU:HD11	2.02	0.42
1:G:645:LYS:HA	1:G:645:LYS:HD2	1.76	0.42
1:H:363:LEU:HD23	1:H:363:LEU:HA	1.86	0.42
1:H:594:VAL:HG21	1:H:654:LEU:HG	2.02	0.42
1:E:367:LEU:O	1:E:370:LEU:HD12	2.20	0.42
1:F:381:GLN:O	1:F:385:GLY:N	2.48	0.42
1:H:381:GLN:O	1:H:385:GLY:N	2.48	0.42
1:H:469:LYS:O	1:H:473:MET:HB2	2.19	0.42
1:H:542:ASP:N	1:H:542:ASP:OD1	2.52	0.42
1:A:519:LEU:HD12	1:I:502:HIS:HB3	2.02	0.42
1:B:505:VAL:HG21	1:B:533:LEU:HB2	2.02	0.42
1:B:594:VAL:HG21	1:B:654:LEU:HG	2.02	0.42
1:C:548:ASP:OD1	1:C:548:ASP:N	2.50	0.42
1:C:594:VAL:HG21	1:C:654:LEU:HG	2.02	0.42
1:F:542:ASP:N	1:F:542:ASP:OD1	2.52	0.42
1:G:594:VAL:HG21	1:G:654:LEU:HG	2.02	0.42
1:I:505:VAL:HG21	1:I:533:LEU:HB2	2.02	0.42
1:A:502:HIS:HB3	1:B:519:LEU:HD12	2.02	0.42
1:A:505:VAL:HG21	1:A:533:LEU:HB2	2.02	0.42
1:A:594:VAL:HG21	1:A:654:LEU:HG	2.02	0.42
1:D:594:VAL:HG23	1:D:652:ILE:HB	2.01	0.42
1:H:505:VAL:HG21	1:H:533:LEU:HB2	2.02	0.42
1:I:594:VAL:HG23	1:I:652:ILE:HB	2.01	0.42
1:I:601:LEU:O	1:I:605:LEU:HG	2.20	0.42
1:A:496:LEU:HA	1:A:496:LEU:HD23	1.74	0.42
1:C:505:VAL:HG21	1:C:533:LEU:HB2	2.02	0.42
1:C:608:SER:HA	1:C:611:GLN:OE1	2.20	0.42
1:D:542:ASP:OD1	1:D:542:ASP:N	2.52	0.42
1:D:585:ILE:HD12	1:D:677:LEU:HD11	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:601:LEU:O	1:E:605:LEU:HG	2.20	0.42
1:G:594:VAL:HG23	1:G:652:ILE:HB	2.01	0.42
1:I:378:PRO:O	1:I:382:ARG:N	2.51	0.42
1:I:585:ILE:HD12	1:I:677:LEU:HD11	2.02	0.42
1:A:361:LEU:HD23	1:A:361:LEU:HA	1.90	0.42
1:B:372:ASN:OD1	1:B:372:ASN:N	2.46	0.42
1:B:608:SER:HA	1:B:611:GLN:OE1	2.20	0.42
1:C:502:HIS:HA	1:C:533:LEU:HD13	2.00	0.42
1:D:598:SER:HB3	1:D:694:LEU:H	1.84	0.42
1:D:601:LEU:O	1:D:605:LEU:HG	2.20	0.42
1:E:585:ILE:HD12	1:E:677:LEU:HD11	2.02	0.42
1:E:667:ALA:HB1	1:E:674:LEU:HB3	2.01	0.42
1:G:505:VAL:HG21	1:G:533:LEU:HB2	2.02	0.42
1:I:541:LEU:HD23	1:I:541:LEU:HA	1.76	0.42
1:I:608:SER:HA	1:I:611:GLN:OE1	2.20	0.42
1:A:667:ALA:HB1	1:A:674:LEU:HB3	2.01	0.41
1:B:367:LEU:O	1:B:370:LEU:HD12	2.19	0.41
1:C:367:LEU:O	1:C:370:LEU:HD12	2.19	0.41
1:C:452:ASP:N	1:C:457:LEU:O	2.48	0.41
1:D:367:LEU:O	1:D:370:LEU:HD12	2.20	0.41
1:D:667:ALA:HB1	1:D:674:LEU:HB3	2.01	0.41
1:E:548:ASP:OD1	1:E:548:ASP:N	2.50	0.41
1:F:585:ILE:HD12	1:F:677:LEU:HD11	2.02	0.41
1:F:601:LEU:O	1:F:605:LEU:HG	2.20	0.41
1:G:361:LEU:HD23	1:G:361:LEU:HA	1.90	0.41
1:G:601:LEU:O	1:G:605:LEU:HG	2.20	0.41
1:I:367:LEU:O	1:I:370:LEU:HD12	2.19	0.41
1:I:598:SER:HB3	1:I:694:LEU:H	1.84	0.41
1:A:608:SER:HA	1:A:611:GLN:OE1	2.20	0.41
1:C:601:LEU:O	1:C:605:LEU:HG	2.20	0.41
1:E:592:LEU:HD12	1:E:592:LEU:HA	1.79	0.41
1:H:502:HIS:HB3	1:I:519:LEU:HD12	2.02	0.41
1:H:608:SER:HA	1:H:611:GLN:OE1	2.20	0.41
1:A:367:LEU:O	1:A:370:LEU:HD12	2.20	0.41
1:C:365:TYR:HB2	1:C:411:PRO:CB	2.50	0.41
1:C:667:ALA:HB1	1:C:674:LEU:HB3	2.01	0.41
1:D:608:SER:HA	1:D:611:GLN:OE1	2.20	0.41
1:E:542:ASP:OD1	1:E:542:ASP:N	2.52	0.41
1:A:365:TYR:HB2	1:A:411:PRO:CB	2.50	0.41
1:C:665:ARG:HA	1:C:668:ARG:CZ	2.51	0.41
1:D:365:TYR:HB2	1:D:411:PRO:CB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:665:ARG:HA	1:F:668:ARG:CZ	2.51	0.41
1:G:608:SER:HA	1:G:611:GLN:OE1	2.20	0.41
1:A:601:LEU:O	1:A:605:LEU:HG	2.20	0.41
1:A:645:LYS:HA	1:A:645:LYS:HD2	1.76	0.41
1:A:665:ARG:HA	1:A:668:ARG:CZ	2.51	0.41
1:B:601:LEU:O	1:B:605:LEU:HG	2.20	0.41
1:D:505:VAL:HG21	1:D:533:LEU:HB2	2.02	0.41
1:D:665:ARG:HA	1:D:668:ARG:CZ	2.51	0.41
1:E:505:VAL:HG21	1:E:533:LEU:HB2	2.02	0.41
1:G:672:GLN:OE1	1:G:672:GLN:N	2.43	0.41
1:H:367:LEU:O	1:H:370:LEU:HD12	2.20	0.41
1:H:378:PRO:O	1:H:382:ARG:N	2.51	0.41
1:H:601:LEU:O	1:H:605:LEU:HG	2.20	0.41
1:D:569:LEU:HA	1:D:569:LEU:HD23	1.78	0.41
1:E:574:ARG:HH12	1:E:660:ARG:NE	2.19	0.41
1:F:608:SER:HA	1:F:611:GLN:OE1	2.20	0.41
1:F:667:ALA:HB1	1:F:674:LEU:HB3	2.01	0.41
1:G:542:ASP:N	1:G:542:ASP:OD1	2.52	0.41
1:G:574:ARG:HH12	1:G:660:ARG:NE	2.19	0.41
1:A:637:LEU:HA	1:A:640:ILE:HD12	2.03	0.41
1:B:541:LEU:HA	1:B:541:LEU:HD23	1.76	0.41
1:C:414:TYR:HE2	1:C:416:ILE:HD11	1.86	0.41
1:E:665:ARG:HA	1:E:668:ARG:CZ	2.51	0.41
1:G:367:LEU:O	1:G:370:LEU:HD12	2.19	0.41
1:H:594:VAL:HG22	1:H:595:TYR:N	2.36	0.41
1:I:391:SER:OG	1:I:397:LEU:HA	2.21	0.41
1:B:365:TYR:HB2	1:B:411:PRO:CB	2.50	0.41
1:C:569:LEU:HD23	1:C:569:LEU:HA	1.78	0.41
1:C:637:LEU:HA	1:C:640:ILE:HD12	2.03	0.41
1:E:396:PHE:HB2	1:F:350:TRP:HH2	1.86	0.41
1:E:519:LEU:HD12	1:F:502:HIS:HB3	2.02	0.41
1:G:350:TRP:HH2	1:H:396:PHE:HB2	1.86	0.41
1:H:365:TYR:HB2	1:H:411:PRO:CB	2.50	0.41
1:I:665:ARG:HA	1:I:668:ARG:CZ	2.51	0.41
1:A:391:SER:OG	1:A:397:LEU:HA	2.21	0.41
1:A:414:TYR:HE2	1:A:416:ILE:HD11	1.86	0.41
1:B:414:TYR:HE2	1:B:416:ILE:HD11	1.86	0.41
1:B:468:ALA:O	1:B:472:ASN:N	2.43	0.41
1:B:637:LEU:HA	1:B:640:ILE:HD12	2.03	0.41
1:B:665:ARG:HA	1:B:668:ARG:CZ	2.51	0.41
1:B:667:ALA:HB1	1:B:674:LEU:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:645:LYS:HD2	1:D:645:LYS:HA	1.76	0.41
1:E:365:TYR:HB2	1:E:411:PRO:CB	2.50	0.41
1:E:594:VAL:HG22	1:E:595:TYR:N	2.36	0.41
1:E:608:SER:HA	1:E:611:GLN:OE1	2.20	0.41
1:F:594:VAL:HG22	1:F:595:TYR:N	2.36	0.41
1:G:381:GLN:O	1:G:385:GLY:N	2.48	0.41
1:G:391:SER:OG	1:G:397:LEU:HA	2.21	0.41
1:G:656:MET:HE3	1:G:656:MET:HB3	1.87	0.41
1:G:665:ARG:HA	1:G:668:ARG:CZ	2.51	0.41
1:H:592:LEU:HD12	1:H:592:LEU:HA	1.79	0.41
1:I:600:GLU:HG3	1:I:601:LEU:H	1.86	0.41
1:I:637:LEU:HA	1:I:640:ILE:HD12	2.03	0.41
1:A:540:LEU:HD23	1:A:540:LEU:HA	1.80	0.41
1:B:350:TRP:CZ3	1:C:396:PHE:HA	2.56	0.41
1:D:414:TYR:HE2	1:D:416:ILE:HD11	1.86	0.41
1:E:350:TRP:CZ3	1:G:396:PHE:HA	2.56	0.41
1:E:381:GLN:O	1:E:385:GLY:N	2.48	0.41
1:A:574:ARG:HH12	1:A:660:ARG:NE	2.19	0.40
1:C:542:ASP:OD1	1:C:542:ASP:N	2.52	0.40
1:F:365:TYR:HB2	1:F:411:PRO:CB	2.50	0.40
1:F:414:TYR:HE2	1:F:416:ILE:HD11	1.86	0.40
1:F:505:VAL:HG21	1:F:533:LEU:HB2	2.02	0.40
1:G:594:VAL:HG22	1:G:595:TYR:N	2.36	0.40
1:H:665:ARG:HA	1:H:668:ARG:CZ	2.51	0.40
1:A:594:VAL:HG22	1:A:595:TYR:N	2.36	0.40
1:B:378:PRO:O	1:B:382:ARG:N	2.51	0.40
1:E:575:CYS:O	1:E:578:LYS:HE3	2.22	0.40
1:F:660:ARG:HA	1:F:660:ARG:HD2	1.87	0.40
1:I:348:LEU:HA	1:I:348:LEU:HD23	1.69	0.40
1:A:421:GLU:OE2	1:B:579:ARG:HB3	2.22	0.40
1:A:600:GLU:HG3	1:A:601:LEU:H	1.87	0.40
1:B:391:SER:OG	1:B:397:LEU:HA	2.21	0.40
1:D:575:CYS:O	1:D:578:LYS:HE3	2.22	0.40
1:H:541:LEU:HA	1:H:541:LEU:HD23	1.76	0.40
1:H:569:LEU:HD23	1:H:569:LEU:HA	1.78	0.40
1:H:637:LEU:HA	1:H:640:ILE:HD12	2.03	0.40
1:I:363:LEU:HD23	1:I:363:LEU:HA	1.86	0.40
1:I:574:ARG:HH12	1:I:660:ARG:NE	2.19	0.40
1:I:594:VAL:HG22	1:I:595:TYR:N	2.36	0.40
1:E:391:SER:OG	1:E:397:LEU:HA	2.21	0.40
1:E:584:LEU:HA	1:E:584:LEU:HD23	1.73	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:592:LEU:HD12	1:F:592:LEU:HA	1.79	0.40
1:F:600:GLU:HG3	1:F:601:LEU:H	1.87	0.40
1:G:548:ASP:N	1:G:548:ASP:OD1	2.50	0.40
1:H:391:SER:OG	1:H:397:LEU:HA	2.21	0.40
1:I:365:TYR:HB2	1:I:411:PRO:CB	2.50	0.40
1:A:350:TRP:HH2	1:B:396:PHE:HB2	1.86	0.40
1:A:642:GLN:HA	1:A:645:LYS:HB2	2.03	0.40
1:B:575:CYS:O	1:B:578:LYS:HE3	2.22	0.40
1:C:502:HIS:HB3	1:D:519:LEU:HD12	2.04	0.40
1:F:574:ARG:HH12	1:F:660:ARG:NE	2.19	0.40
1:F:584:LEU:HA	1:F:584:LEU:HD23	1.73	0.40
1:H:574:ARG:HH12	1:H:660:ARG:NE	2.19	0.40
1:I:414:TYR:HE2	1:I:416:ILE:HD11	1.86	0.40
1:I:642:GLN:HA	1:I:645:LYS:HB2	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	351/702 (50%)	302 (86%)	48 (14%)	1 (0%)	37	68
1	B	351/702 (50%)	301 (86%)	49 (14%)	1 (0%)	37	68
1	C	351/702 (50%)	301 (86%)	49 (14%)	1 (0%)	37	68
1	D	351/702 (50%)	302 (86%)	48 (14%)	1 (0%)	37	68
1	E	351/702 (50%)	302 (86%)	48 (14%)	1 (0%)	37	68
1	F	351/702 (50%)	302 (86%)	48 (14%)	1 (0%)	37	68
1	G	351/702 (50%)	302 (86%)	48 (14%)	1 (0%)	37	68
1	H	351/702 (50%)	301 (86%)	49 (14%)	1 (0%)	37	68
1	I	351/702 (50%)	302 (86%)	48 (14%)	1 (0%)	37	68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	3159/6318 (50%)	2715 (86%)	435 (14%)	9 (0%)	38	68

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	420	GLY
1	B	420	GLY
1	C	420	GLY
1	D	420	GLY
1	E	420	GLY
1	F	420	GLY
1	G	420	GLY
1	H	420	GLY
1	I	420	GLY

### 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	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	B	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	C	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	D	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	E	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	F	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	G	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	H	305/584 (52%)	303 (99%)	2 (1%)	81	88
1	I	305/584 (52%)	303 (99%)	2 (1%)	81	88
All	All	2745/5256 (52%)	2727 (99%)	18 (1%)	80	88

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	488	ILE
1	A	620	LEU
1	B	488	ILE
1	B	620	LEU
1	C	488	ILE
1	C	620	LEU
1	D	488	ILE
1	D	620	LEU
1	E	488	ILE
1	E	620	LEU
1	F	488	ILE
1	F	620	LEU
1	G	488	ILE
1	G	620	LEU
1	H	488	ILE
1	H	620	LEU
1	I	488	ILE
1	I	620	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	413	GLN
1	A	425	GLN
1	A	472	ASN
1	A	646	GLN
1	B	413	GLN
1	B	425	GLN
1	B	472	ASN
1	B	646	GLN
1	C	413	GLN
1	C	425	GLN
1	C	472	ASN
1	C	646	GLN
1	D	413	GLN
1	D	425	GLN
1	D	472	ASN
1	D	646	GLN
1	E	413	GLN
1	E	425	GLN
1	E	472	ASN
1	E	646	GLN
1	F	413	GLN

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Mol	Chain	Res	Type
1	F	425	GLN
1	F	472	ASN
1	F	646	GLN
1	G	413	GLN
1	G	425	GLN
1	G	472	ASN
1	G	646	GLN
1	H	413	GLN
1	H	425	GLN
1	H	472	ASN
1	H	646	GLN
1	I	413	GLN
1	I	425	GLN
1	I	472	ASN
1	I	646	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

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.

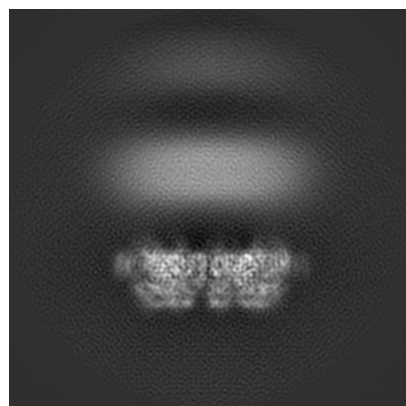
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11827. These allow visual inspection of the internal detail of the map and identification of artifacts.

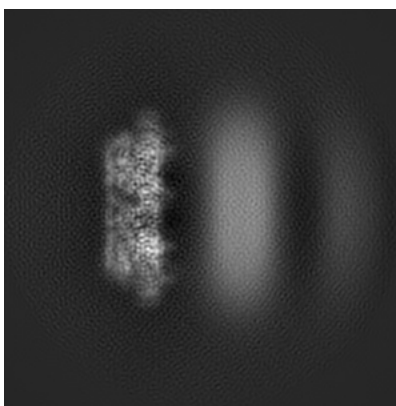
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

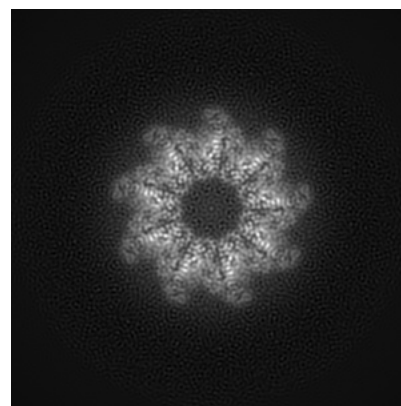
#### 6.1.1 Primary map



X

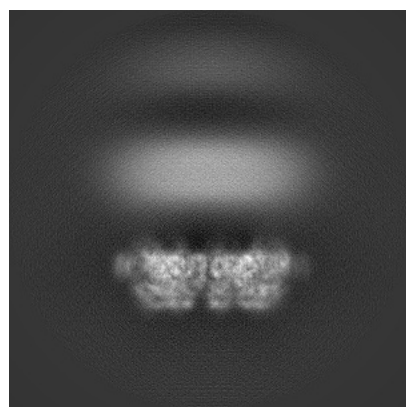


Y

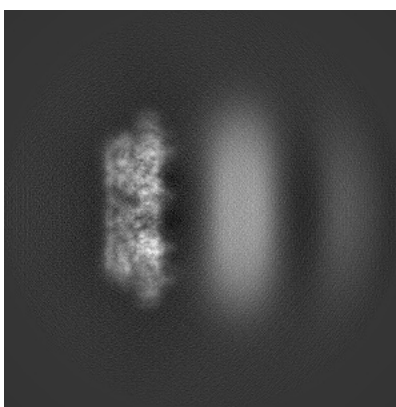


Z

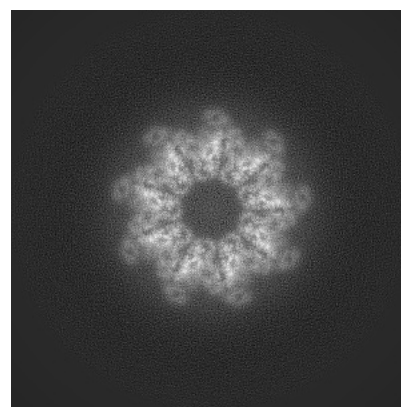
#### 6.1.2 Raw map



X



Y

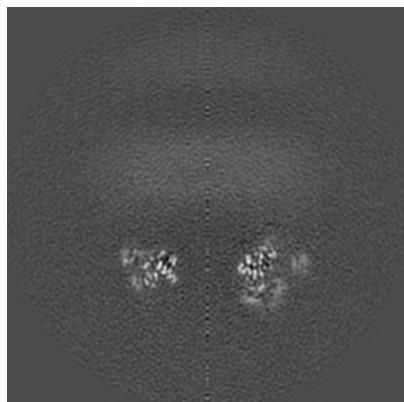


Z

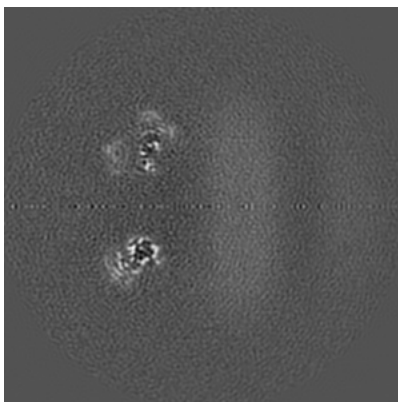
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

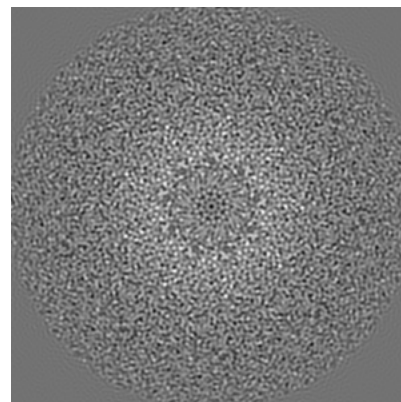
### 6.2.1 Primary map



X Index: 212

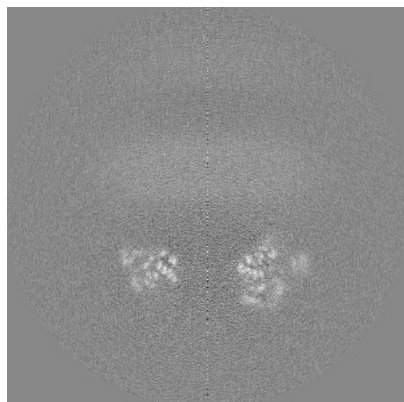


Y Index: 212

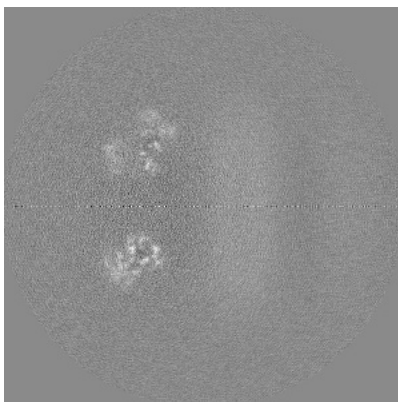


Z Index: 212

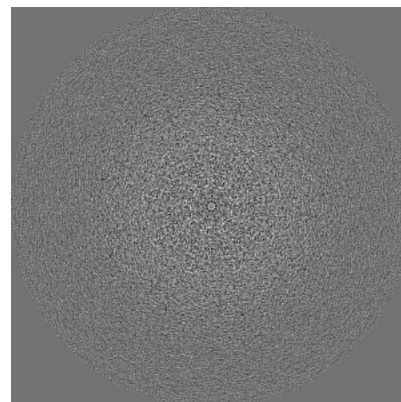
### 6.2.2 Raw map



X Index: 212



Y Index: 212



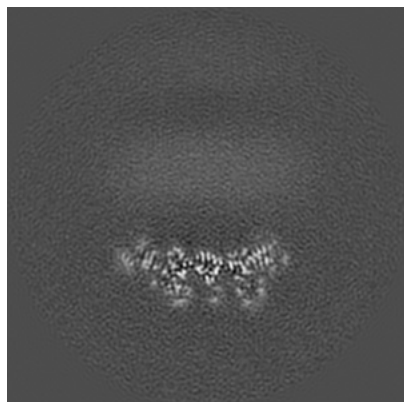
Z Index: 212

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

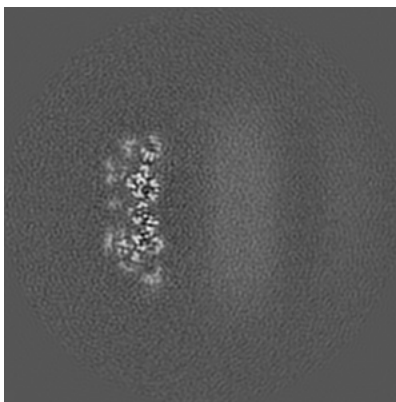


## 6.3 Largest variance slices [i](#)

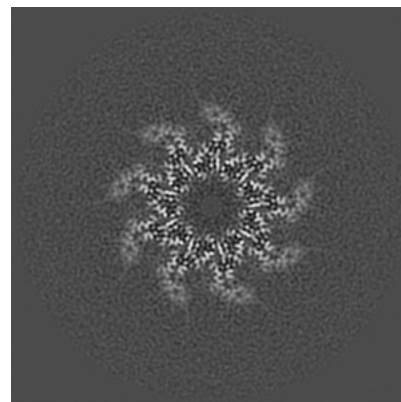
### 6.3.1 Primary map



X Index: 168

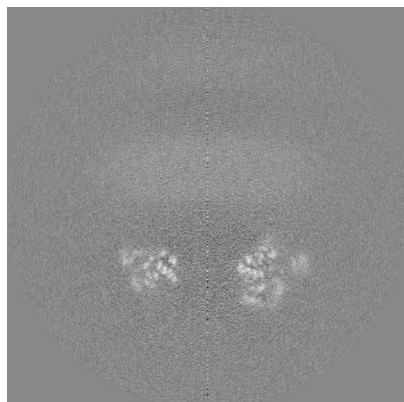


Y Index: 250

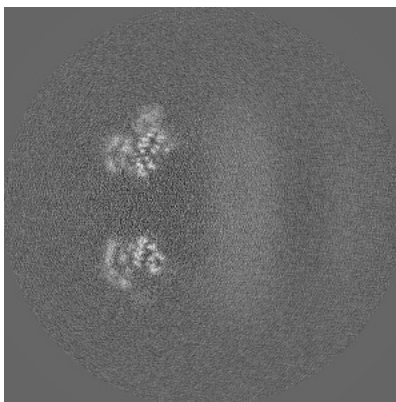


Z Index: 152

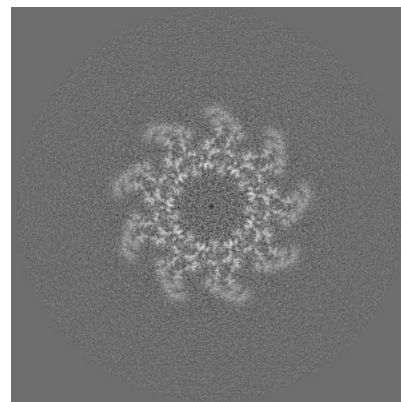
### 6.3.2 Raw map



X Index: 212



Y Index: 219

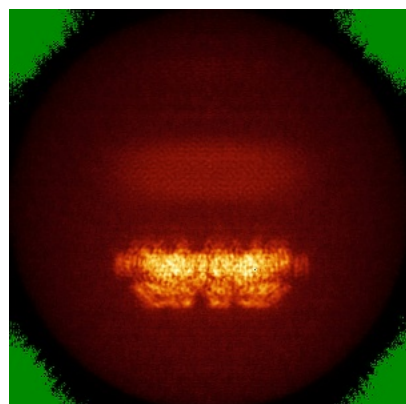


Z Index: 158

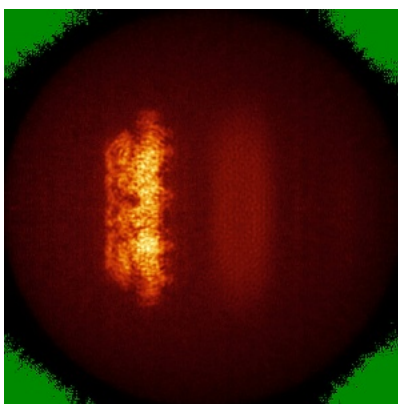
The images above show the largest variance slices of the map in three orthogonal directions.

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

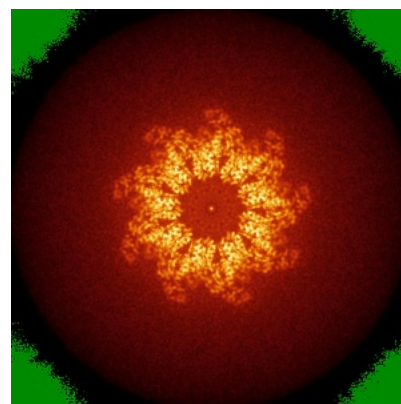
### 6.4.1 Primary map



X

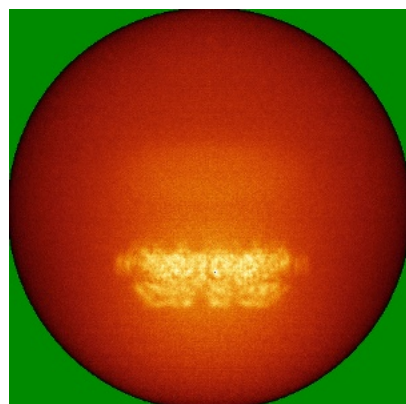


Y

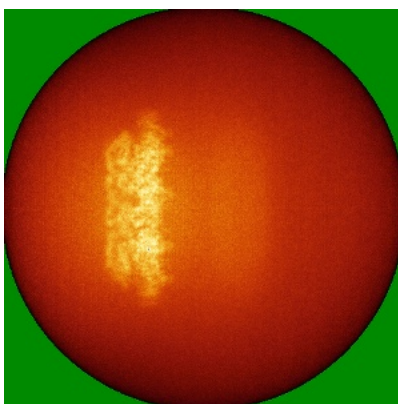


Z

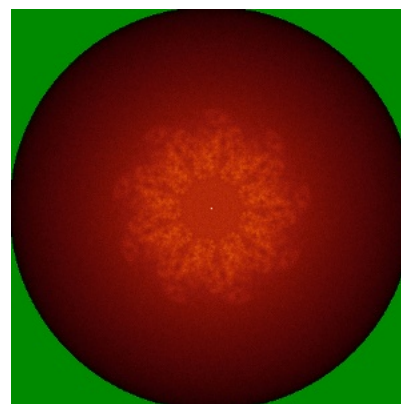
### 6.4.2 Raw map



X



Y



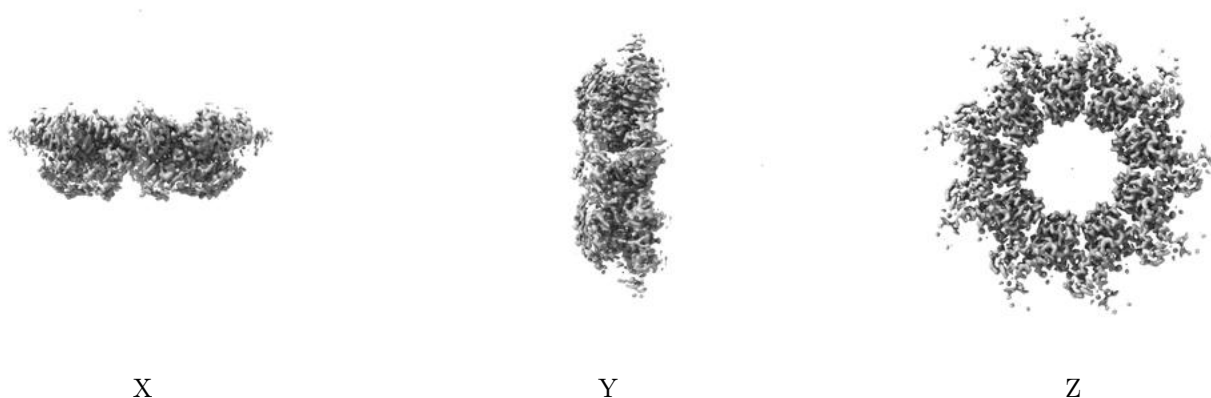
Z

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



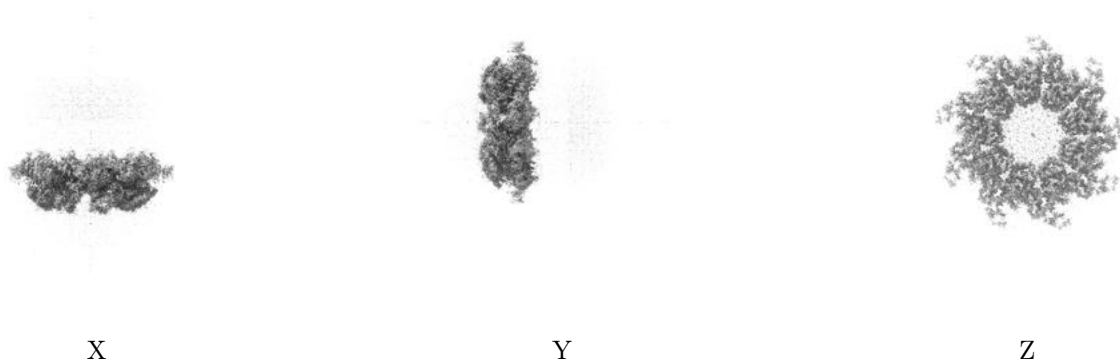
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0146. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.5.2 Raw map



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

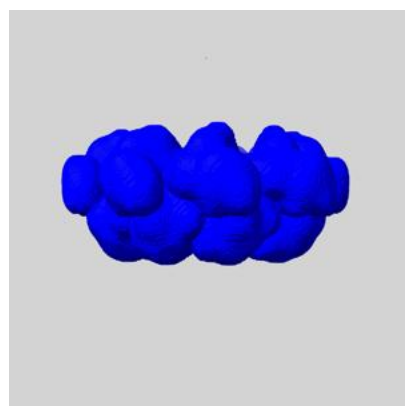
## 6.6 Mask visualisation [i](#)

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

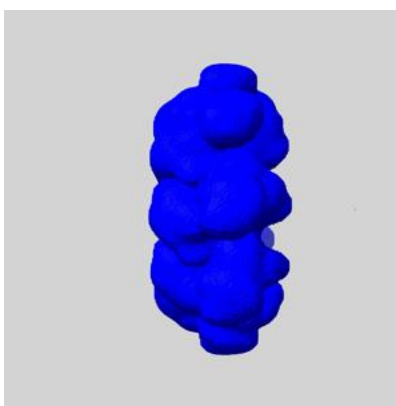
A mask typically either:

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

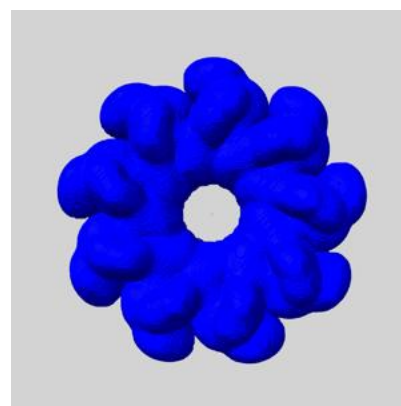
### 6.6.1 emd\_11827\_msk\_1.map [i](#)



X



Y

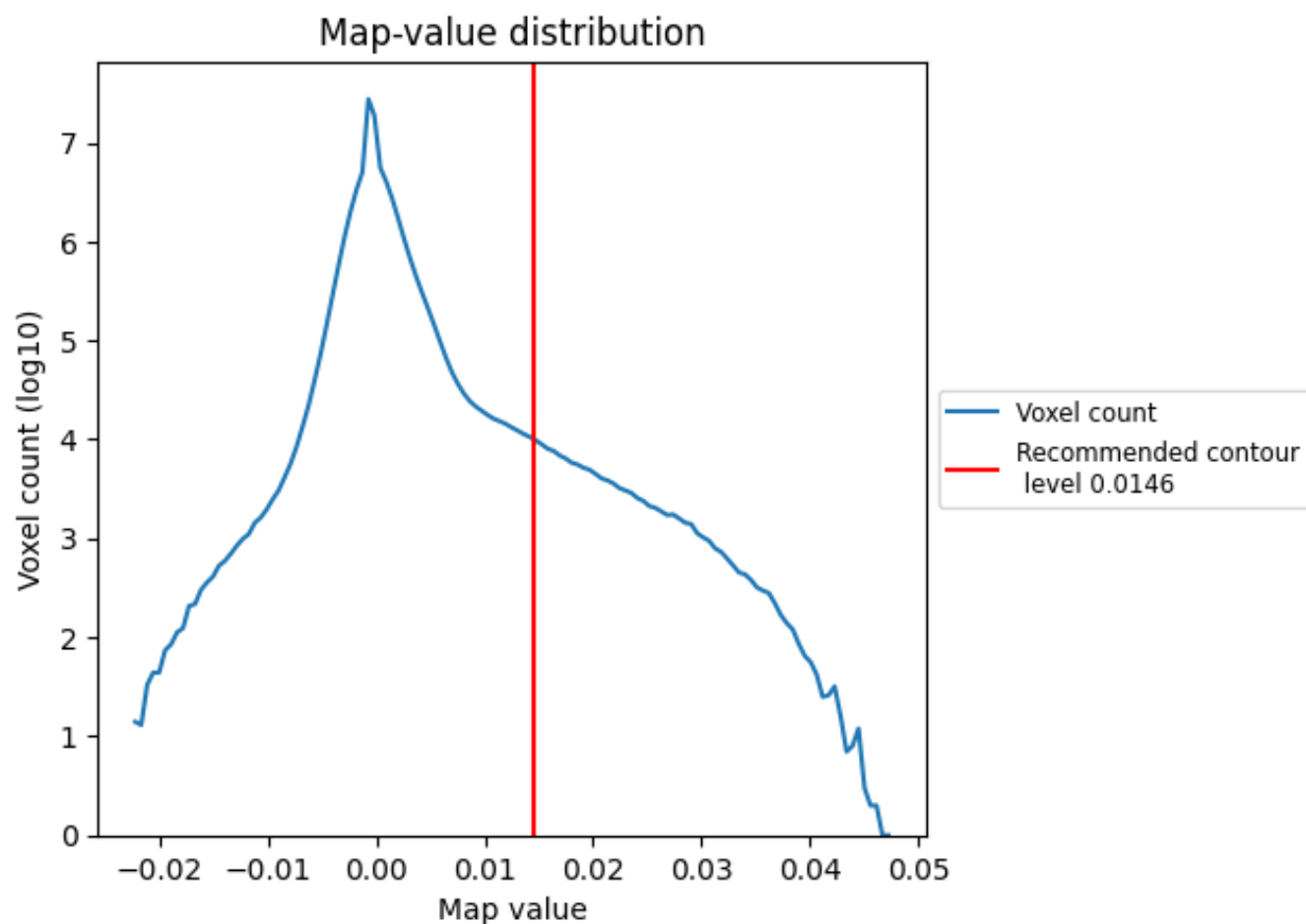


Z

## 7 Map analysis [i](#)

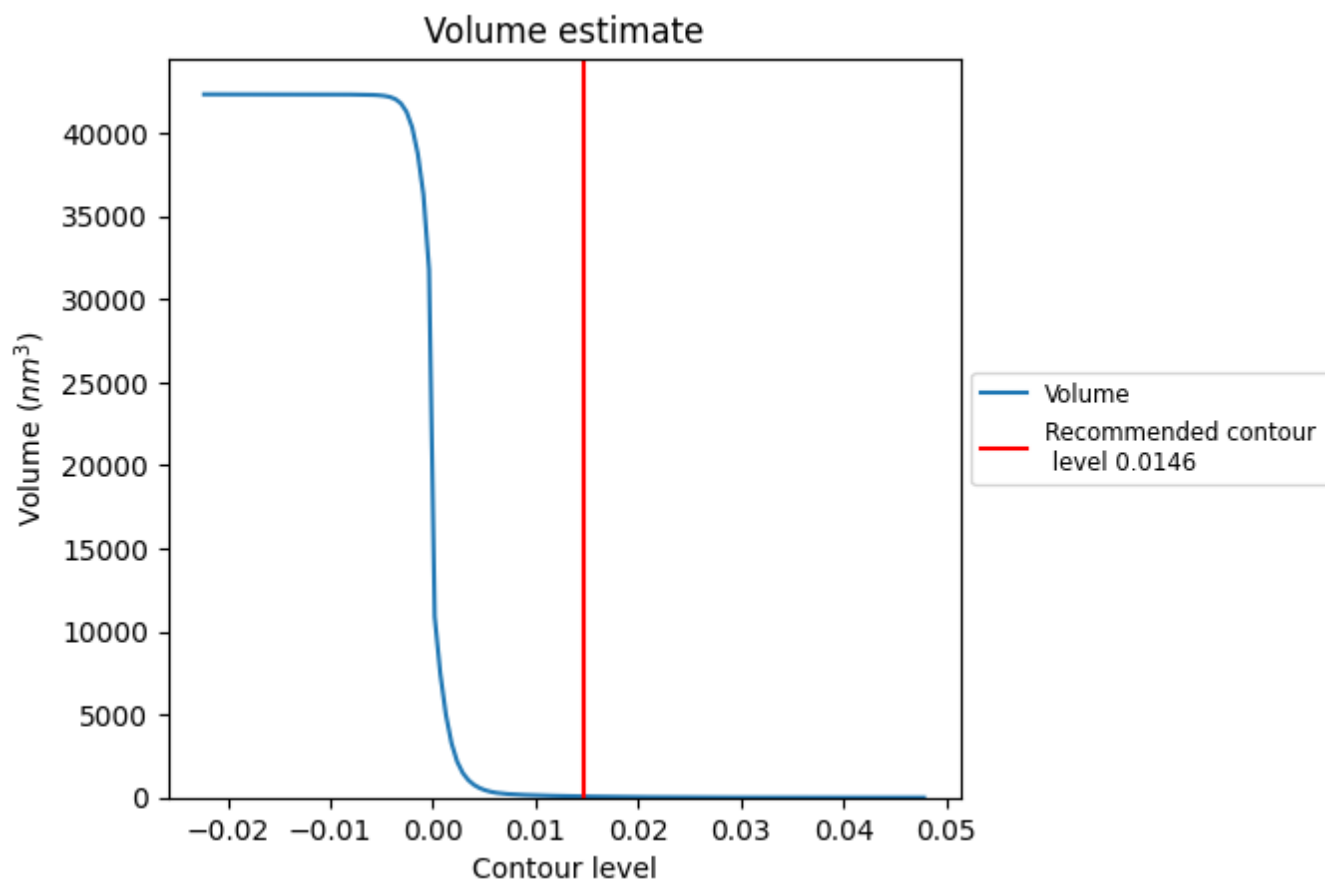
This section contains the results of statistical analysis of the map.

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



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

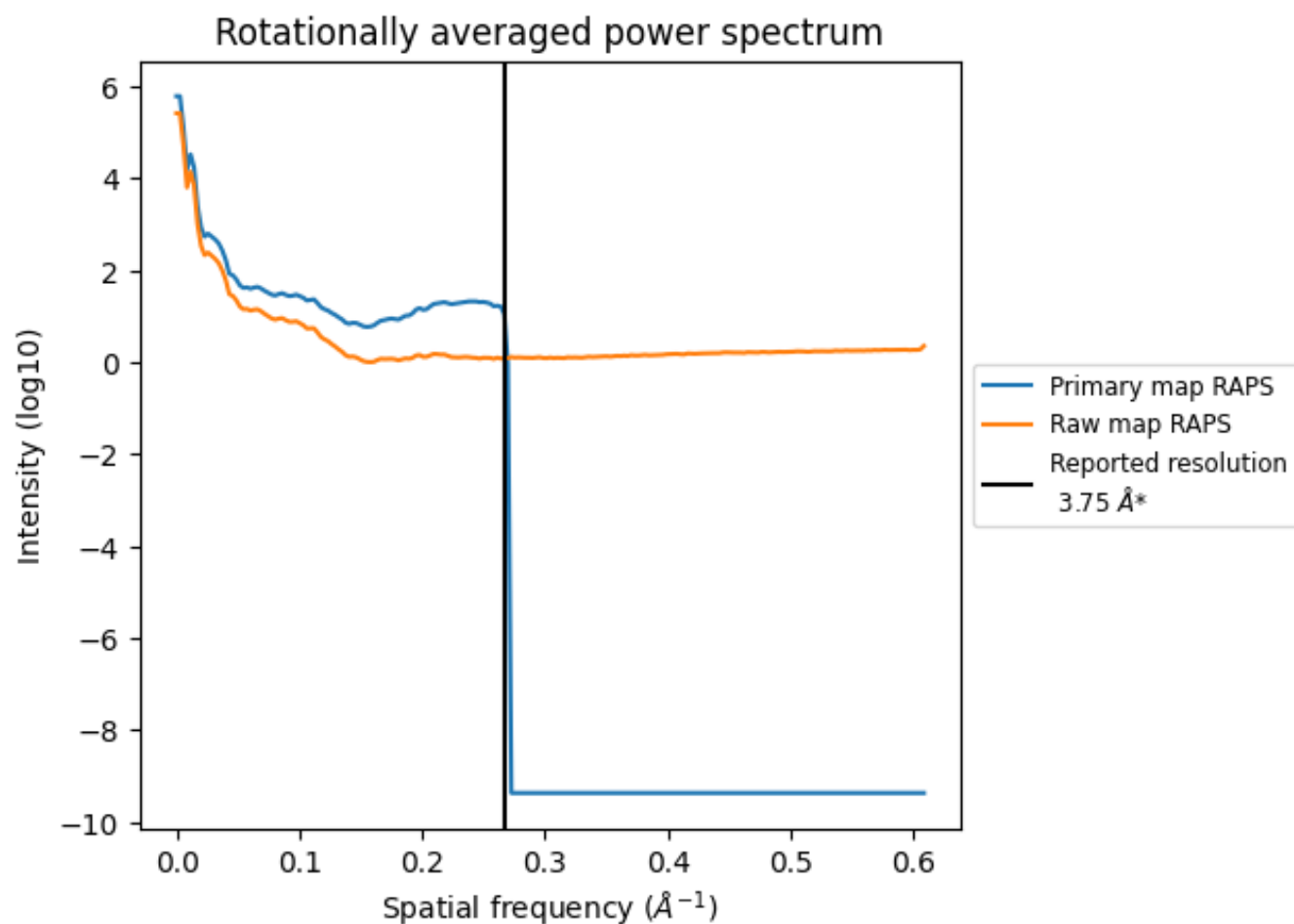
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 69 nm<sup>3</sup>; this corresponds to an approximate mass of 62 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 ⓘ

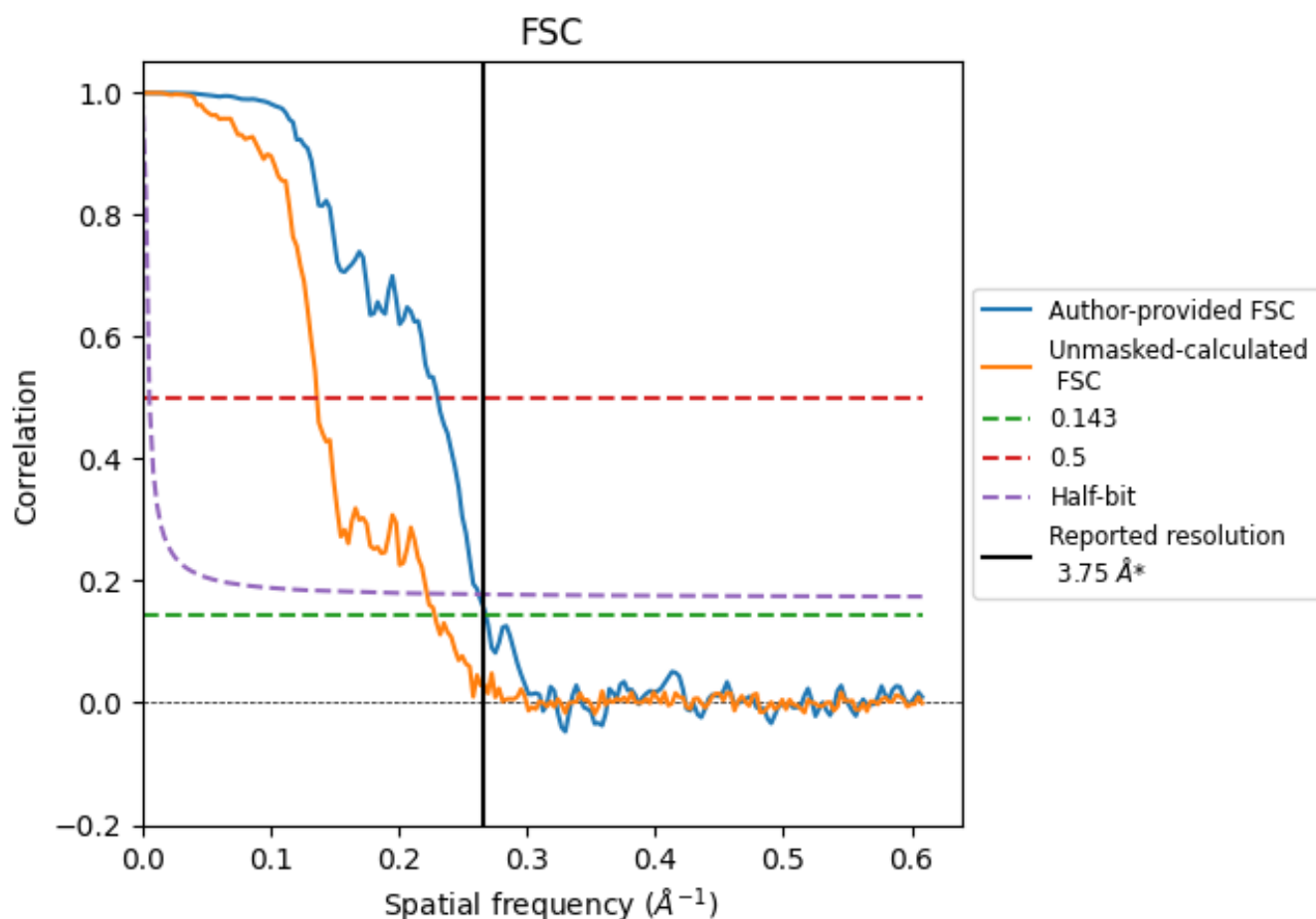


\*Reported resolution corresponds to spatial frequency of 0.267 Å<sup>-1</sup>

## 8 Fourier-Shell correlation [i](#)

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

### 8.1 FSC [i](#)



\*Reported resolution corresponds to spatial frequency of 0.267  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

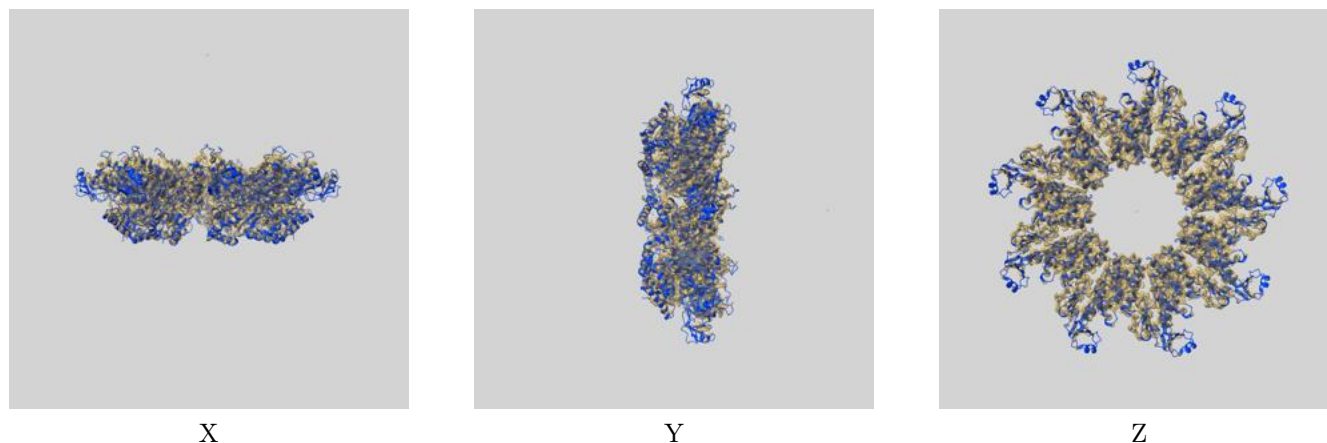
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.75	-	-
Author-provided FSC curve	3.73	4.34	3.81
Unmasked-calculated*	4.38	7.34	4.50

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.38 differs from the reported value 3.75 by more than 10 %

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

This section contains information regarding the fit between EMDB map EMD-11827 and PDB model 7AMY. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

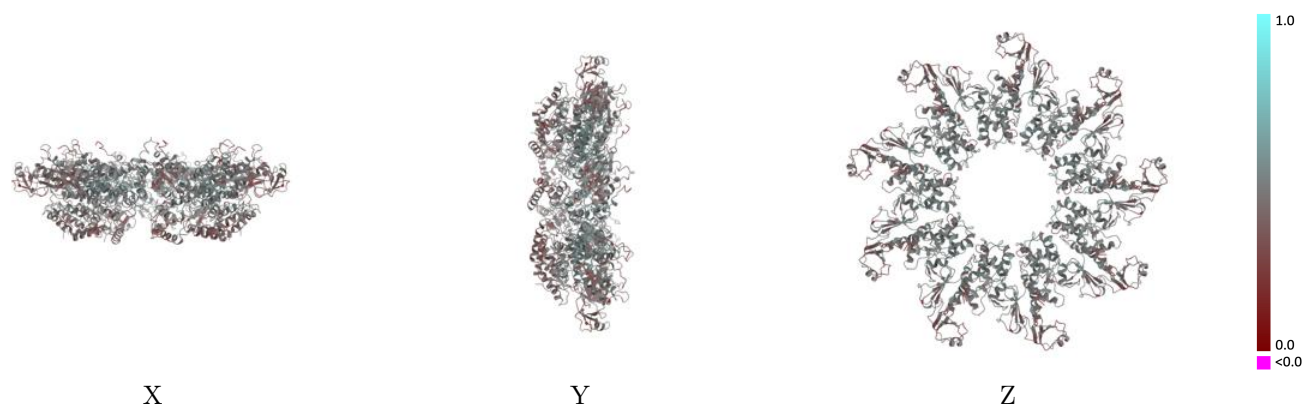
### 9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.0146 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.

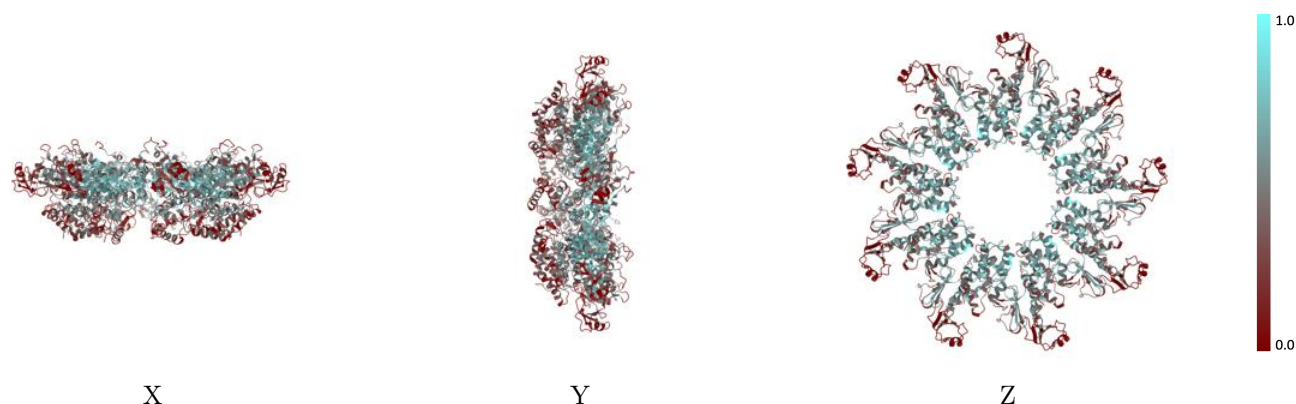


## 9.2 Q-score mapped to coordinate model [i](#)



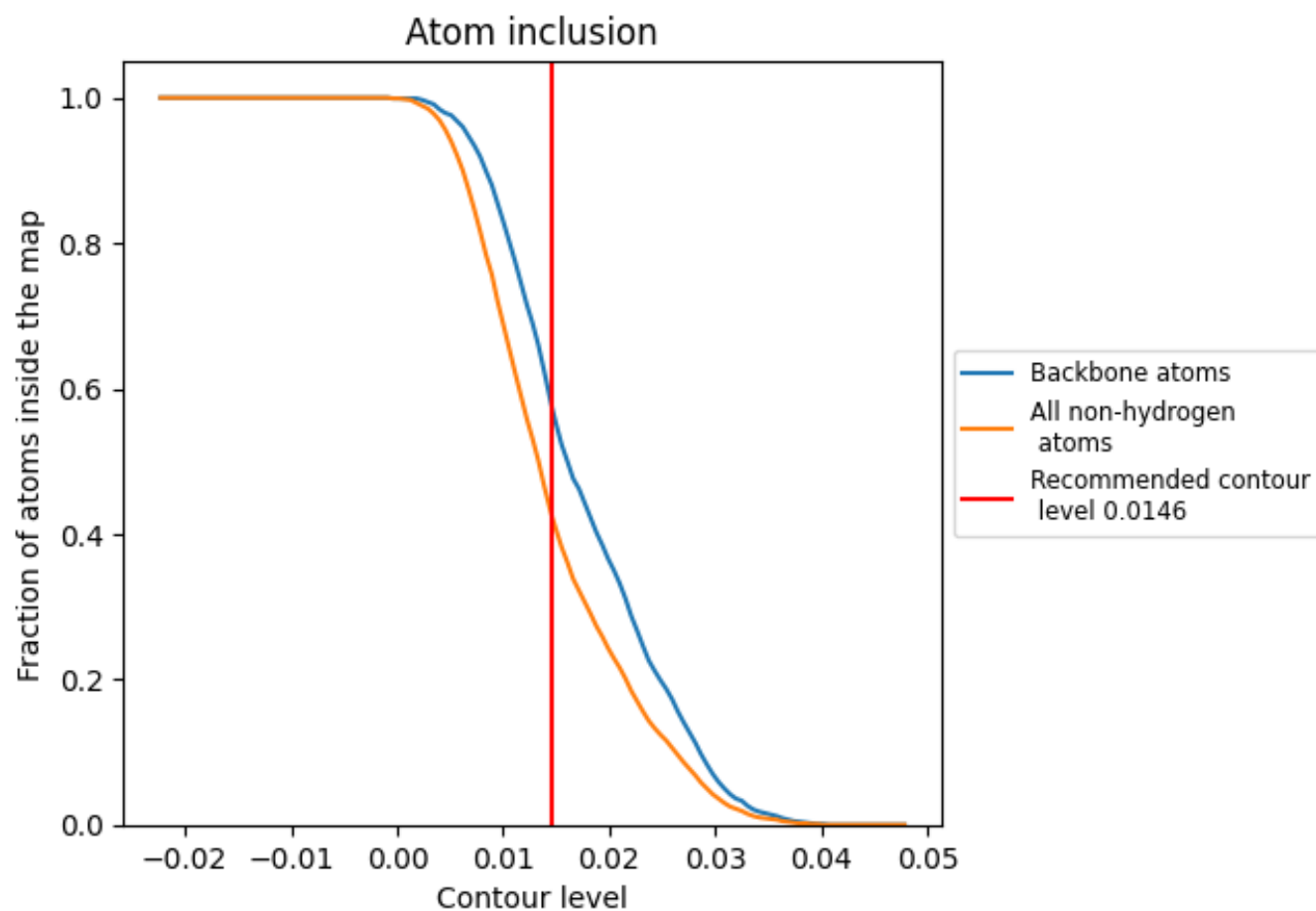
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0146).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 57% of all backbone atoms, 42% 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 (0.0146) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.4240</div>	<div><div></div>0.4560</div>
A	<div><div></div>0.4220</div>	<div><div></div>0.4560</div>
B	<div><div></div>0.4240</div>	<div><div></div>0.4560</div>
C	<div><div></div>0.4220</div>	<div><div></div>0.4550</div>
D	<div><div></div>0.4270</div>	<div><div></div>0.4560</div>
E	<div><div></div>0.4240</div>	<div><div></div>0.4560</div>
F	<div><div></div>0.4220</div>	<div><div></div>0.4560</div>
G	<div><div></div>0.4220</div>	<div><div></div>0.4560</div>
H	<div><div></div>0.4240</div>	<div><div></div>0.4560</div>
I	<div><div></div>0.4230</div>	<div><div></div>0.4570</div>

1.0

0.0

<0.0