



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 08:13 AM UTC

PDB ID : 8AVC / pdb_00008avc
EMDB ID : EMD-15678
Title : Mouse leptin:LEP-R complex cryoEM structure (3:3 model)
Authors : Verstraete, K.; Savvides, S.N.; Verschueren, K.G.; Tsirigotaki, A.
Deposited on : 2022-08-26
Resolution : 4.60 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

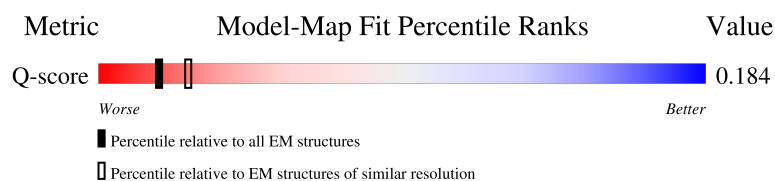
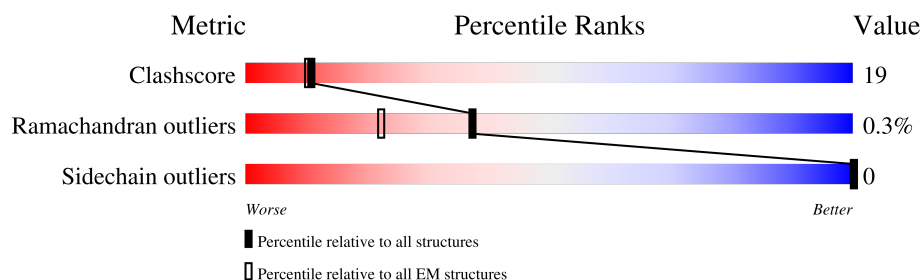
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

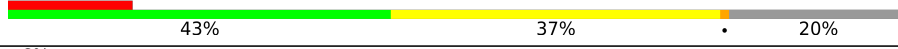
The reported resolution of this entry is 4.60 Å.

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



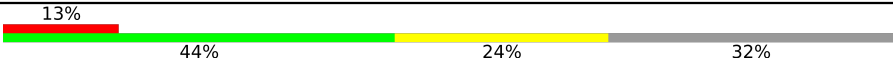
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2407 (4.10 - 5.10)

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

Mol	Chain	Length	Quality of chain
1	A	174	
1	C	174	
1	E	174	
2	B	868	

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Mol	Chain	Length	Quality of chain
2	D	868	
2	F	868	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17505 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 Leptin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	140	Total	C	N	O	S	0	0
			1077	676	182	216	3		
1	C	140	Total	C	N	O	S	0	0
			1077	676	182	216	3		
1	E	140	Total	C	N	O	S	0	0
			1077	676	182	216	3		

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP P41160
A	-5	GLY	-	expression tag	UNP P41160
A	-4	SER	-	expression tag	UNP P41160
A	-3	SER	-	expression tag	UNP P41160
A	-2	HIS	-	expression tag	UNP P41160
A	-1	HIS	-	expression tag	UNP P41160
A	0	HIS	-	expression tag	UNP P41160
A	1	HIS	-	expression tag	UNP P41160
A	2	HIS	-	expression tag	UNP P41160
A	3	HIS	-	expression tag	UNP P41160
A	4	PRO	-	expression tag	UNP P41160
A	5	GLY	-	expression tag	UNP P41160
A	6	GLY	-	expression tag	UNP P41160
A	7	PRO	-	expression tag	UNP P41160
A	8	GLY	-	expression tag	UNP P41160
A	9	SER	-	expression tag	UNP P41160
A	10	GLU	-	expression tag	UNP P41160
A	11	ASN	-	expression tag	UNP P41160
A	12	LEU	-	expression tag	UNP P41160
A	13	TYR	-	expression tag	UNP P41160
A	14	PHE	-	expression tag	UNP P41160
A	15	GLN	-	expression tag	UNP P41160
A	16	GLY	-	expression tag	UNP P41160
A	17	GLY	-	expression tag	UNP P41160

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Chain	Residue	Modelled	Actual	Comment	Reference
A	18	SER	-	expression tag	UNP P41160
A	19	THR	-	expression tag	UNP P41160
A	20	GLY	-	expression tag	UNP P41160
A	21	GLY	ALA	conflict	UNP P41160
C	-6	MET	-	initiating methionine	UNP P41160
C	-5	GLY	-	expression tag	UNP P41160
C	-4	SER	-	expression tag	UNP P41160
C	-3	SER	-	expression tag	UNP P41160
C	-2	HIS	-	expression tag	UNP P41160
C	-1	HIS	-	expression tag	UNP P41160
C	0	HIS	-	expression tag	UNP P41160
C	1	HIS	-	expression tag	UNP P41160
C	2	HIS	-	expression tag	UNP P41160
C	3	HIS	-	expression tag	UNP P41160
C	4	PRO	-	expression tag	UNP P41160
C	5	GLY	-	expression tag	UNP P41160
C	6	GLY	-	expression tag	UNP P41160
C	7	PRO	-	expression tag	UNP P41160
C	8	GLY	-	expression tag	UNP P41160
C	9	SER	-	expression tag	UNP P41160
C	10	GLU	-	expression tag	UNP P41160
C	11	ASN	-	expression tag	UNP P41160
C	12	LEU	-	expression tag	UNP P41160
C	13	TYR	-	expression tag	UNP P41160
C	14	PHE	-	expression tag	UNP P41160
C	15	GLN	-	expression tag	UNP P41160
C	16	GLY	-	expression tag	UNP P41160
C	17	GLY	-	expression tag	UNP P41160
C	18	SER	-	expression tag	UNP P41160
C	19	THR	-	expression tag	UNP P41160
C	20	GLY	-	expression tag	UNP P41160
C	21	GLY	ALA	conflict	UNP P41160
E	-6	MET	-	initiating methionine	UNP P41160
E	-5	GLY	-	expression tag	UNP P41160
E	-4	SER	-	expression tag	UNP P41160
E	-3	SER	-	expression tag	UNP P41160
E	-2	HIS	-	expression tag	UNP P41160
E	-1	HIS	-	expression tag	UNP P41160
E	0	HIS	-	expression tag	UNP P41160
E	1	HIS	-	expression tag	UNP P41160
E	2	HIS	-	expression tag	UNP P41160
E	3	HIS	-	expression tag	UNP P41160

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Chain	Residue	Modelled	Actual	Comment	Reference
E	4	PRO	-	expression tag	UNP P41160
E	5	GLY	-	expression tag	UNP P41160
E	6	GLY	-	expression tag	UNP P41160
E	7	PRO	-	expression tag	UNP P41160
E	8	GLY	-	expression tag	UNP P41160
E	9	SER	-	expression tag	UNP P41160
E	10	GLU	-	expression tag	UNP P41160
E	11	ASN	-	expression tag	UNP P41160
E	12	LEU	-	expression tag	UNP P41160
E	13	TYR	-	expression tag	UNP P41160
E	14	PHE	-	expression tag	UNP P41160
E	15	GLN	-	expression tag	UNP P41160
E	16	GLY	-	expression tag	UNP P41160
E	17	GLY	-	expression tag	UNP P41160
E	18	SER	-	expression tag	UNP P41160
E	19	THR	-	expression tag	UNP P41160
E	20	GLY	-	expression tag	UNP P41160
E	21	GLY	ALA	conflict	UNP P41160

- Molecule 2 is a protein called Leptin receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	593	Total	C	N	O	S	0	0
			4757	3057	798	878	24		
2	D	593	Total	C	N	O	S	0	0
			4757	3057	798	878	24		
2	F	593	Total	C	N	O	S	0	0
			4757	3057	798	878	24		

There are 150 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	840	SER	-	expression tag	UNP P48356
B	841	THR	-	expression tag	UNP P48356
B	842	GLY	-	expression tag	UNP P48356
B	843	GLY	-	expression tag	UNP P48356
B	844	SER	-	expression tag	UNP P48356
B	845	GLY	-	expression tag	UNP P48356
B	846	GLY	-	expression tag	UNP P48356
B	847	SER	-	expression tag	UNP P48356
B	848	GLY	-	expression tag	UNP P48356
B	849	GLY	-	expression tag	UNP P48356

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Chain	Residue	Modelled	Actual	Comment	Reference
B	850	SER	-	expression tag	UNP P48356
B	851	GLY	-	expression tag	UNP P48356
B	852	GLY	-	expression tag	UNP P48356
B	853	SER	-	expression tag	UNP P48356
B	854	GLY	-	expression tag	UNP P48356
B	855	GLY	-	expression tag	UNP P48356
B	856	SER	-	expression tag	UNP P48356
B	857	ARG	-	expression tag	UNP P48356
B	858	MET	-	expression tag	UNP P48356
B	859	LYS	-	expression tag	UNP P48356
B	860	GLN	-	expression tag	UNP P48356
B	861	ILE	-	expression tag	UNP P48356
B	862	GLU	-	expression tag	UNP P48356
B	863	ASP	-	expression tag	UNP P48356
B	864	LYS	-	expression tag	UNP P48356
B	865	ILE	-	expression tag	UNP P48356
B	866	GLU	-	expression tag	UNP P48356
B	867	GLU	-	expression tag	UNP P48356
B	868	ILE	-	expression tag	UNP P48356
B	869	LEU	-	expression tag	UNP P48356
B	870	SER	-	expression tag	UNP P48356
B	871	LYS	-	expression tag	UNP P48356
B	872	ILE	-	expression tag	UNP P48356
B	873	TYR	-	expression tag	UNP P48356
B	874	HIS	-	expression tag	UNP P48356
B	875	ILE	-	expression tag	UNP P48356
B	876	GLU	-	expression tag	UNP P48356
B	877	ASN	-	expression tag	UNP P48356
B	878	GLU	-	expression tag	UNP P48356
B	879	ILE	-	expression tag	UNP P48356
B	880	ALA	-	expression tag	UNP P48356
B	881	ARG	-	expression tag	UNP P48356
B	882	ILE	-	expression tag	UNP P48356
B	883	LYS	-	expression tag	UNP P48356
B	884	LYS	-	expression tag	UNP P48356
B	885	LEU	-	expression tag	UNP P48356
B	886	ILE	-	expression tag	UNP P48356
B	887	GLY	-	expression tag	UNP P48356
B	888	GLU	-	expression tag	UNP P48356
B	889	ARG	-	expression tag	UNP P48356
D	840	SER	-	expression tag	UNP P48356
D	841	THR	-	expression tag	UNP P48356

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Chain	Residue	Modelled	Actual	Comment	Reference
D	842	GLY	-	expression tag	UNP P48356
D	843	GLY	-	expression tag	UNP P48356
D	844	SER	-	expression tag	UNP P48356
D	845	GLY	-	expression tag	UNP P48356
D	846	GLY	-	expression tag	UNP P48356
D	847	SER	-	expression tag	UNP P48356
D	848	GLY	-	expression tag	UNP P48356
D	849	GLY	-	expression tag	UNP P48356
D	850	SER	-	expression tag	UNP P48356
D	851	GLY	-	expression tag	UNP P48356
D	852	GLY	-	expression tag	UNP P48356
D	853	SER	-	expression tag	UNP P48356
D	854	GLY	-	expression tag	UNP P48356
D	855	GLY	-	expression tag	UNP P48356
D	856	SER	-	expression tag	UNP P48356
D	857	ARG	-	expression tag	UNP P48356
D	858	MET	-	expression tag	UNP P48356
D	859	LYS	-	expression tag	UNP P48356
D	860	GLN	-	expression tag	UNP P48356
D	861	ILE	-	expression tag	UNP P48356
D	862	GLU	-	expression tag	UNP P48356
D	863	ASP	-	expression tag	UNP P48356
D	864	LYS	-	expression tag	UNP P48356
D	865	ILE	-	expression tag	UNP P48356
D	866	GLU	-	expression tag	UNP P48356
D	867	GLU	-	expression tag	UNP P48356
D	868	ILE	-	expression tag	UNP P48356
D	869	LEU	-	expression tag	UNP P48356
D	870	SER	-	expression tag	UNP P48356
D	871	LYS	-	expression tag	UNP P48356
D	872	ILE	-	expression tag	UNP P48356
D	873	TYR	-	expression tag	UNP P48356
D	874	HIS	-	expression tag	UNP P48356
D	875	ILE	-	expression tag	UNP P48356
D	876	GLU	-	expression tag	UNP P48356
D	877	ASN	-	expression tag	UNP P48356
D	878	GLU	-	expression tag	UNP P48356
D	879	ILE	-	expression tag	UNP P48356
D	880	ALA	-	expression tag	UNP P48356
D	881	ARG	-	expression tag	UNP P48356
D	882	ILE	-	expression tag	UNP P48356
D	883	LYS	-	expression tag	UNP P48356

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Chain	Residue	Modelled	Actual	Comment	Reference
D	884	LYS	-	expression tag	UNP P48356
D	885	LEU	-	expression tag	UNP P48356
D	886	ILE	-	expression tag	UNP P48356
D	887	GLY	-	expression tag	UNP P48356
D	888	GLU	-	expression tag	UNP P48356
D	889	ARG	-	expression tag	UNP P48356
F	840	SER	-	expression tag	UNP P48356
F	841	THR	-	expression tag	UNP P48356
F	842	GLY	-	expression tag	UNP P48356
F	843	GLY	-	expression tag	UNP P48356
F	844	SER	-	expression tag	UNP P48356
F	845	GLY	-	expression tag	UNP P48356
F	846	GLY	-	expression tag	UNP P48356
F	847	SER	-	expression tag	UNP P48356
F	848	GLY	-	expression tag	UNP P48356
F	849	GLY	-	expression tag	UNP P48356
F	850	SER	-	expression tag	UNP P48356
F	851	GLY	-	expression tag	UNP P48356
F	852	GLY	-	expression tag	UNP P48356
F	853	SER	-	expression tag	UNP P48356
F	854	GLY	-	expression tag	UNP P48356
F	855	GLY	-	expression tag	UNP P48356
F	856	SER	-	expression tag	UNP P48356
F	857	ARG	-	expression tag	UNP P48356
F	858	MET	-	expression tag	UNP P48356
F	859	LYS	-	expression tag	UNP P48356
F	860	GLN	-	expression tag	UNP P48356
F	861	ILE	-	expression tag	UNP P48356
F	862	GLU	-	expression tag	UNP P48356
F	863	ASP	-	expression tag	UNP P48356
F	864	LYS	-	expression tag	UNP P48356
F	865	ILE	-	expression tag	UNP P48356
F	866	GLU	-	expression tag	UNP P48356
F	867	GLU	-	expression tag	UNP P48356
F	868	ILE	-	expression tag	UNP P48356
F	869	LEU	-	expression tag	UNP P48356
F	870	SER	-	expression tag	UNP P48356
F	871	LYS	-	expression tag	UNP P48356
F	872	ILE	-	expression tag	UNP P48356
F	873	TYR	-	expression tag	UNP P48356
F	874	HIS	-	expression tag	UNP P48356
F	875	ILE	-	expression tag	UNP P48356

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Chain	Residue	Modelled	Actual	Comment	Reference
F	876	GLU	-	expression tag	UNP P48356
F	877	ASN	-	expression tag	UNP P48356
F	878	GLU	-	expression tag	UNP P48356
F	879	ILE	-	expression tag	UNP P48356
F	880	ALA	-	expression tag	UNP P48356
F	881	ARG	-	expression tag	UNP P48356
F	882	ILE	-	expression tag	UNP P48356
F	883	LYS	-	expression tag	UNP P48356
F	884	LYS	-	expression tag	UNP P48356
F	885	LEU	-	expression tag	UNP P48356
F	886	ILE	-	expression tag	UNP P48356
F	887	GLY	-	expression tag	UNP P48356
F	888	GLU	-	expression tag	UNP P48356
F	889	ARG	-	expression tag	UNP P48356

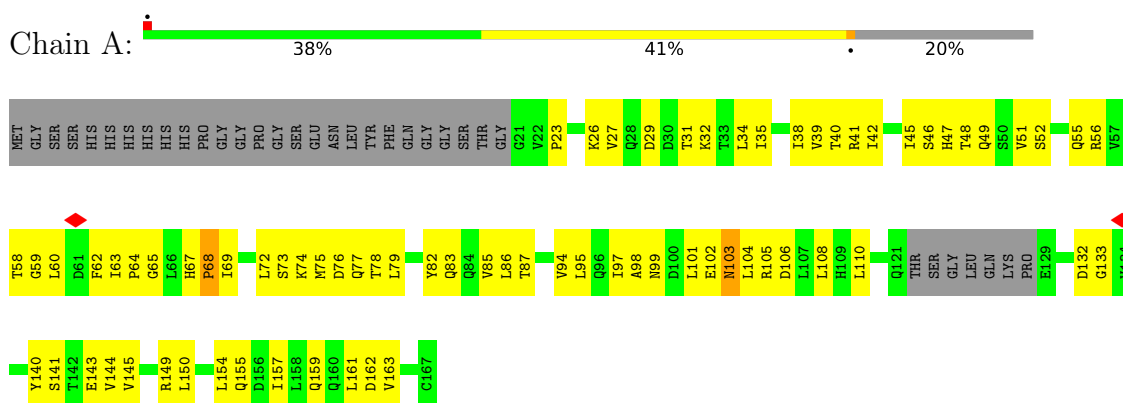
- Molecule 3 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	AltConf
3	B	1	Total Ni 1 1	0
3	D	1	Total Ni 1 1	0
3	F	1	Total Ni 1 1	0

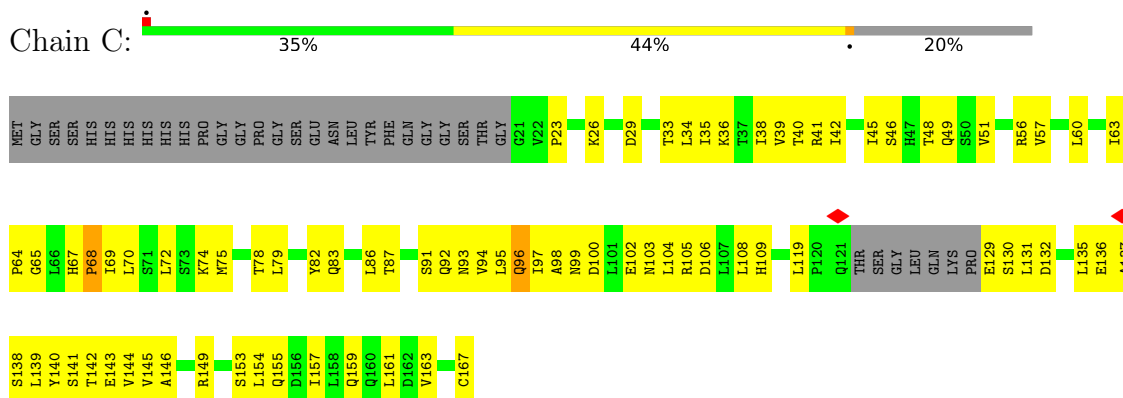
3 Residue-property plots [i](#)

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.

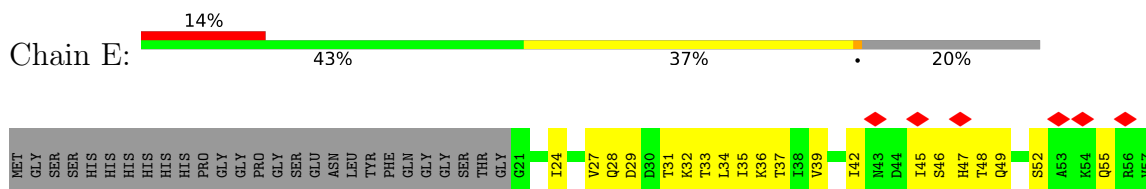
• Molecule 1: Leptin

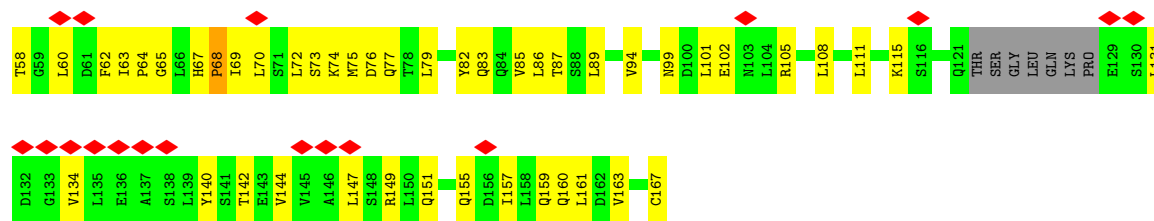


• Molecule 1: Leptin

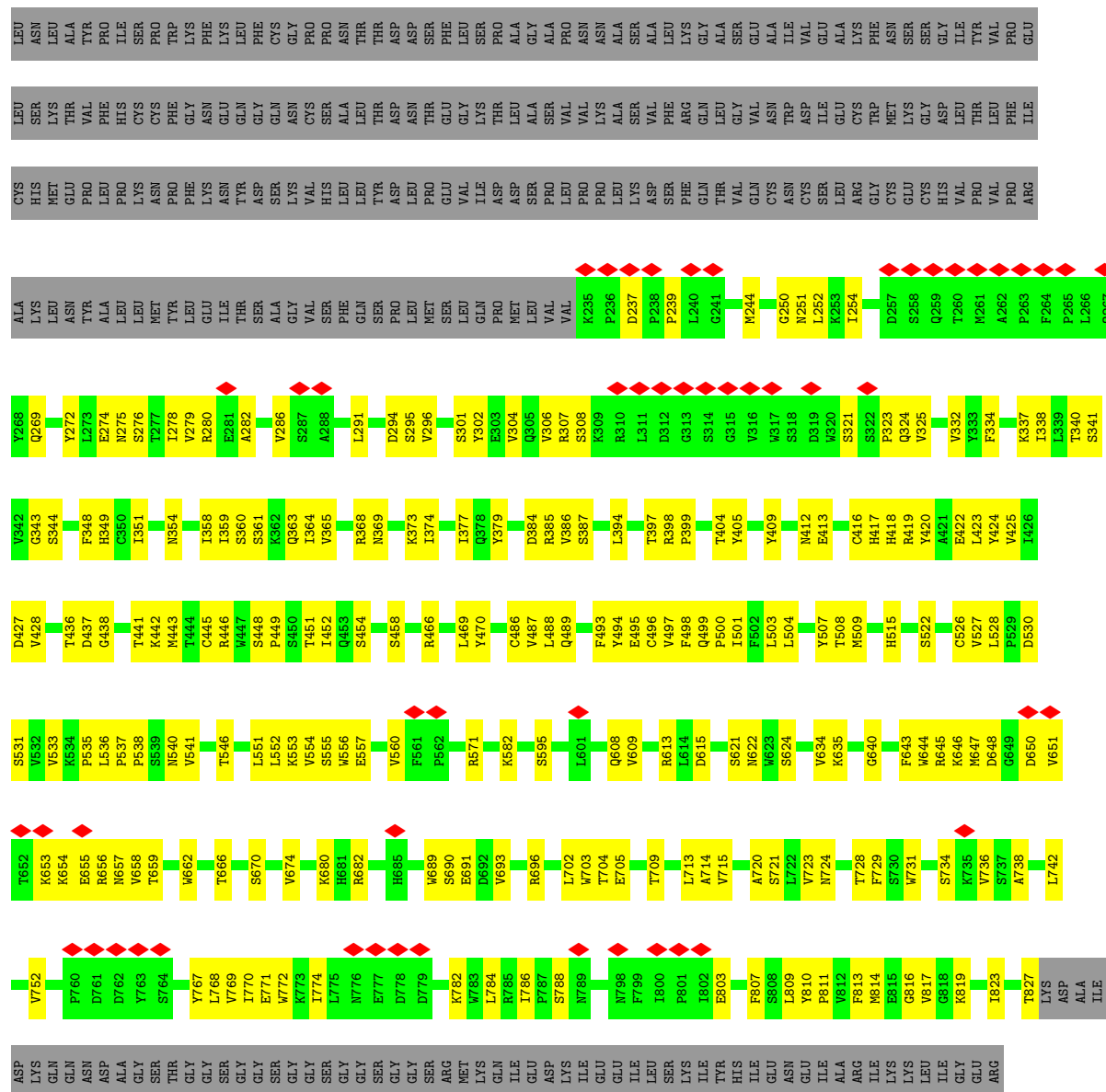


• Molecule 1: Leptin

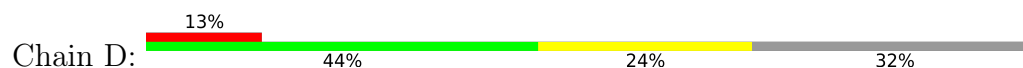


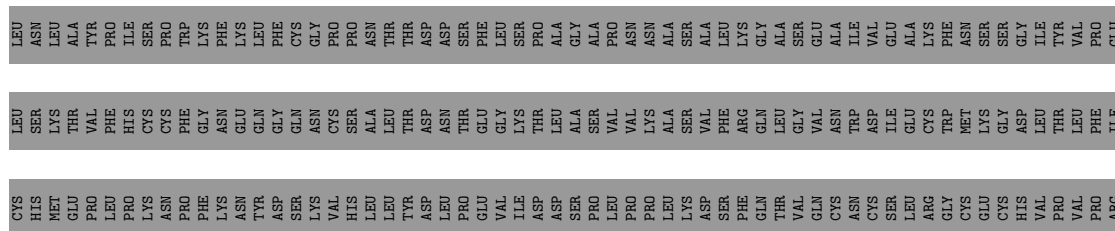


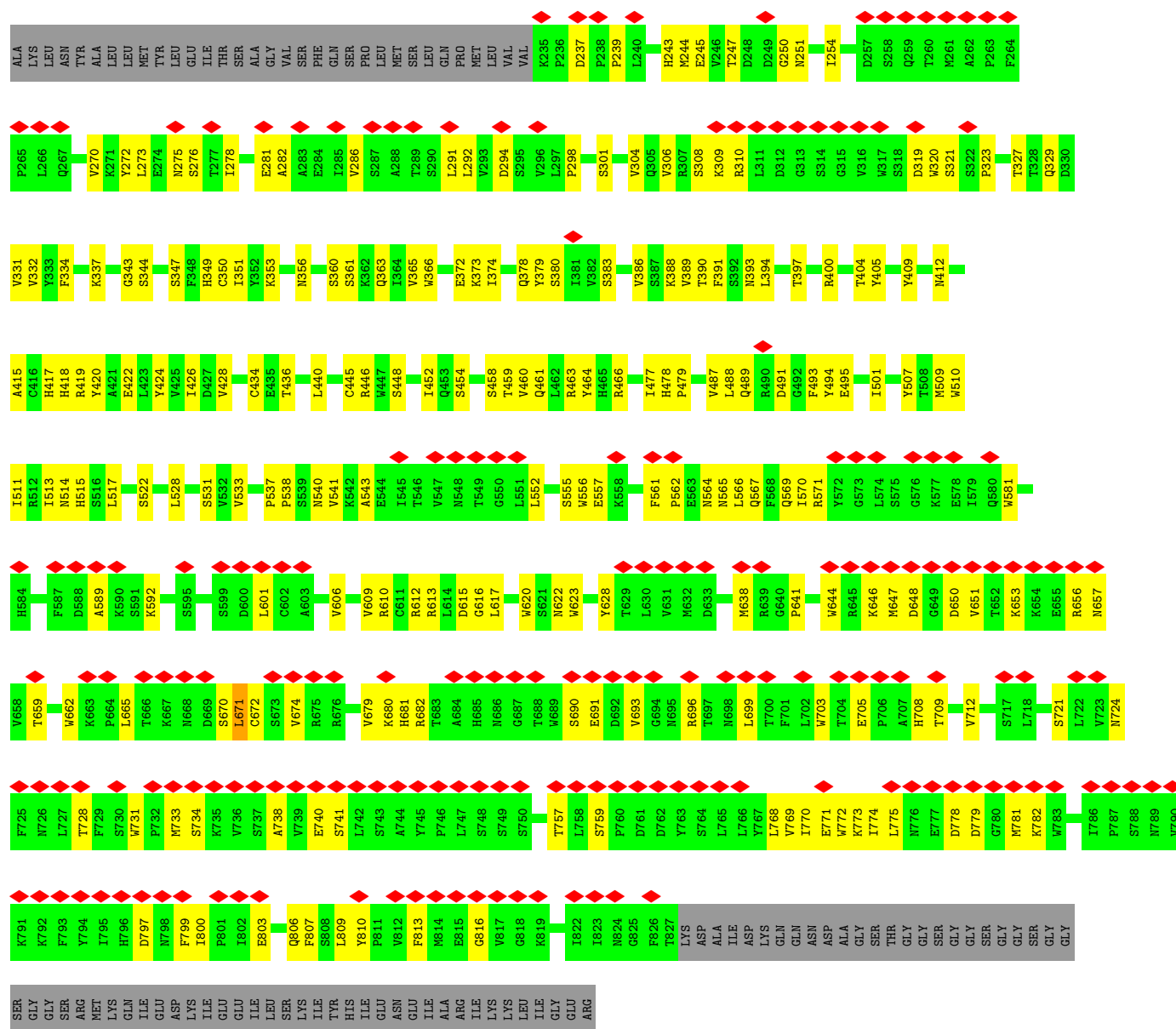
• Molecule 2: Leptin receptor



• Molecule 2: Leptin receptor







4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	37530	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	45	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.927	Depositor
Minimum map value	-0.383	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.025	Depositor
Recommended contour level	0.178	Depositor
Map size (Å)	371.392, 371.392, 371.392	wwPDB
Map dimensions	224, 224, 224	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.658, 1.658, 1.658	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NI

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	0.30	0/1088	0.83	2/1477 (0.1%)
1	C	0.30	0/1088	0.84	2/1477 (0.1%)
1	E	0.21	0/1088	0.66	0/1477
2	B	0.14	0/4890	0.37	0/6668
2	D	0.14	0/4890	0.38	2/6668 (0.0%)
2	F	0.12	0/4890	0.33	0/6668
All	All	0.17	0/17934	0.47	6/24435 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	0	3

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	96	GLN	CB-CG-CD	-10.38	94.95	112.60
2	D	434	CYS	N-CA-C	5.96	116.42	108.38
1	C	96	GLN	N-CA-C	-5.96	104.87	111.36
1	A	103	ASN	CB-CA-C	5.50	120.80	110.63
1	A	103	ASN	CA-CB-CG	5.49	118.09	112.60
2	D	434	CYS	CA-CB-SG	5.17	126.30	114.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	415	ALA	Peptide
2	D	467	ARG	Peptide
2	D	501	ILE	Peptide

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	1077	0	1109	74	0
1	C	1077	0	1109	84	0
1	E	1077	0	1109	50	0
2	B	4757	0	4708	167	0
2	D	4757	0	4709	157	0
2	F	4757	0	4709	154	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
3	F	1	0	0	0	0
All	All	17505	0	17453	648	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 19.

All (648) 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:103:ASN:ND2	2:B:504:LEU:H	1.25	1.30
2:B:409:TYR:OH	1:C:141:SER:HB3	1.34	1.21
1:A:103:ASN:ND2	2:B:504:LEU:N	1.99	1.11
1:A:34:LEU:HD11	1:A:104:LEU:HD13	1.39	1.03
1:C:83:GLN:HE22	1:C:102:GLU:HG2	1.27	0.98
2:B:409:TYR:OH	1:C:141:SER:CB	2.12	0.95
2:D:434:CYS:HB3	2:D:445:CYS:HA	1.44	0.95
1:C:60:LEU:HD22	1:C:145:VAL:HG11	1.52	0.90
1:C:36:LYS:HE2	2:D:563:GLU:HB3	1.54	0.89
1:A:103:ASN:HD21	2:B:504:LEU:H	1.19	0.88
2:B:541:VAL:HA	2:B:555:SER:O	1.75	0.87
1:C:96:GLN:OE1	2:D:501:ILE:O	1.93	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:GLN:NE2	1:C:98:ALA:O	2.11	0.83
2:F:329:GLN:HG3	2:F:419:ARG:HH12	1.44	0.83
1:C:155:GLN:O	1:C:159:GLN:NE2	2.13	0.81
1:A:73:SER:O	1:A:77:GLN:NE2	2.13	0.81
1:C:99:ASN:ND2	2:D:502:PHE:N	2.30	0.79
1:A:23:PRO:HG2	1:A:26:LYS:HG2	1.66	0.78
1:C:96:GLN:NE2	1:C:100:ASP:OD2	2.16	0.78
1:E:63:ILE:O	1:E:149:ARG:NH2	2.17	0.77
1:A:99:ASN:HB3	2:B:501:ILE:O	1.85	0.77
1:E:69:ILE:HD11	1:E:75:MET:HA	1.66	0.76
2:F:412:ASN:HB3	2:F:415:ALA:HB3	1.66	0.76
1:C:65:GLY:HA3	1:C:149:ARG:NH1	1.99	0.76
2:D:434:CYS:CB	2:D:445:CYS:HA	2.15	0.75
2:D:400:ARG:HH12	1:E:47:HIS:HB3	1.53	0.74
2:B:436:THR:HB	2:B:533:VAL:HA	1.70	0.74
2:F:298:PRO:HG2	2:F:353:LYS:HB2	1.70	0.73
1:E:155:GLN:O	1:E:159:GLN:NE2	2.21	0.73
2:B:373:LYS:NZ	2:B:374:ILE:O	2.19	0.72
2:D:638:MET:HE1	2:D:671:LEU:HD21	1.71	0.72
2:F:487:VAL:O	2:F:489:GLN:NE2	2.23	0.72
2:F:571:ARG:HB2	2:F:581:TRP:HB3	1.72	0.72
2:D:487:VAL:O	2:D:489:GLN:NE2	2.23	0.71
2:D:358:ILE:HG21	2:D:385:ARG:HG2	1.71	0.71
1:E:45:ILE:HD13	1:E:147:LEU:HD13	1.73	0.71
1:C:99:ASN:ND2	2:D:502:PHE:H	1.87	0.71
2:D:334:PHE:HB2	2:D:349:HIS:HB2	1.73	0.71
2:F:446:ARG:NH1	2:F:495:GLU:OE2	2.23	0.71
2:B:446:ARG:HG2	2:B:495:GLU:HG2	1.73	0.71
1:E:52:SER:HB3	1:E:55:GLN:HB2	1.73	0.70
2:B:351:ILE:HD11	2:B:386:VAL:HG22	1.74	0.70
2:B:648:ASP:HB2	2:B:657:ASN:HB2	1.74	0.70
1:A:83:GLN:O	1:A:87:THR:HG23	1.91	0.70
2:F:459:THR:HB	2:F:461:GLN:HE22	1.56	0.70
2:B:469:LEU:HB3	2:B:470:TYR:HD1	1.55	0.69
2:B:486:CYS:HA	2:B:496:CYS:HB3	1.73	0.69
2:D:802:ILE:HG13	2:D:803:GLU:HG2	1.76	0.68
2:B:332:VAL:HB	2:B:351:ILE:HB	1.75	0.68
2:D:237:ASP:O	2:D:259:GLN:NE2	2.26	0.68
2:B:768:LEU:HD13	2:B:809:LEU:HD21	1.76	0.68
2:F:703:TRP:NE1	2:F:705:GLU:O	2.26	0.68
2:B:535:PRO:HB2	2:B:621:SER:HB3	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:535:PRO:HB2	2:D:621:SER:HB3	1.75	0.67
2:B:446:ARG:NH1	2:B:495:GLU:OE2	2.28	0.67
2:B:405:TYR:HA	2:B:423:LEU:O	1.95	0.67
2:D:703:TRP:NE1	2:D:705:GLU:O	2.27	0.67
2:B:703:TRP:NE1	2:B:705:GLU:O	2.27	0.67
1:C:60:LEU:HB2	1:C:63:ILE:HD13	1.77	0.67
1:A:99:ASN:CB	2:B:501:ILE:O	2.43	0.67
1:E:72:LEU:HA	1:E:75:MET:HE2	1.77	0.67
1:A:72:LEU:HA	1:A:75:MET:HE2	1.77	0.67
2:F:680:LYS:HA	2:F:690:SER:HA	1.77	0.67
1:A:141:SER:O	1:A:145:VAL:HG12	1.95	0.66
1:C:138:SER:O	1:C:142:THR:HG23	1.95	0.66
1:C:41:ARG:HH22	2:D:440:LEU:HB2	1.60	0.66
1:A:74:LYS:O	1:A:78:THR:HG23	1.95	0.66
1:C:79:LEU:C	1:C:105:ARG:HH22	2.04	0.66
1:A:63:ILE:HD11	1:A:85:VAL:HG21	1.77	0.66
2:B:361:SER:N	2:B:384:ASP:O	2.27	0.66
2:D:415:ALA:HA	1:E:55:GLN:HE21	1.61	0.66
1:C:96:GLN:O	1:C:99:ASN:HB3	1.94	0.66
2:B:470:TYR:HA	2:B:528:LEU:HD21	1.76	0.65
1:C:69:ILE:HA	1:C:74:LYS:HE3	1.79	0.65
1:C:83:GLN:O	1:C:87:THR:HG23	1.96	0.65
2:F:360:SER:HB3	2:F:363:GLN:HG2	1.79	0.64
1:C:56:ARG:HH22	1:C:149:ARG:HH11	1.44	0.64
1:E:72:LEU:HB2	1:E:167:CYS:HA	1.80	0.64
1:A:69:ILE:HG23	1:A:74:LYS:HG3	1.77	0.64
2:D:644:TRP:NE1	2:D:771:GLU:OE2	2.24	0.64
1:A:79:LEU:HD12	1:A:101:LEU:HD22	1.78	0.63
1:A:103:ASN:CG	2:B:504:LEU:H	2.01	0.63
1:A:145:VAL:O	1:A:149:ARG:HG3	1.99	0.63
2:F:351:ILE:HD11	2:F:386:VAL:HG22	1.79	0.63
1:C:23:PRO:HB2	1:C:26:LYS:HE2	1.80	0.63
2:D:774:ILE:HG23	2:D:803:GLU:HB3	1.81	0.63
1:E:157:ILE:O	1:E:161:LEU:HG	1.98	0.63
2:F:452:ILE:HD13	2:F:460:VAL:HB	1.81	0.63
2:F:682:ARG:HB3	2:F:709:THR:HB	1.81	0.63
1:A:103:ASN:HD21	2:B:504:LEU:N	1.85	0.62
2:F:418:HIS:HE1	2:F:419:ARG:HH21	1.46	0.62
2:F:428:VAL:HG22	2:F:522:SER:HA	1.81	0.62
1:A:38:ILE:O	1:A:42:ILE:HD12	1.99	0.62
2:D:397:THR:HG23	2:D:425:VAL:HB	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:60:LEU:HD11	1:E:142:THR:HA	1.81	0.62
2:B:275:ASN:HA	2:B:279:VAL:HG22	1.81	0.61
2:B:571:ARG:NH2	2:B:608:GLN:OE1	2.33	0.61
2:B:644:TRP:NE1	2:B:771:GLU:OE2	2.26	0.61
2:D:304:VAL:O	2:D:323:PRO:HA	1.99	0.61
2:B:438:GLY:HA3	2:B:536:LEU:HD21	1.82	0.61
2:F:648:ASP:HB2	2:F:657:ASN:HB2	1.82	0.61
2:B:466:ARG:HD3	2:B:507:TYR:CZ	2.35	0.61
1:C:157:ILE:O	1:C:161:LEU:HG	2.00	0.61
2:B:398:ARG:NH1	2:B:399:PRO:O	2.34	0.61
1:A:65:GLY:HA3	1:A:149:ARG:HH21	1.66	0.61
1:C:132:ASP:HA	1:C:135:LEU:HD12	1.82	0.61
1:E:79:LEU:HD12	1:E:101:LEU:HD22	1.81	0.61
2:F:337:LYS:HB2	2:F:517:LEU:HD11	1.83	0.61
1:A:59:GLY:HA3	2:F:418:HIS:HD2	1.65	0.60
2:D:608:GLN:NE2	2:D:624:SER:O	2.23	0.60
2:F:458:SER:HB2	2:F:515:HIS:HD2	1.66	0.60
2:F:566:LEU:H	2:F:589:ALA:HB2	1.65	0.60
1:C:72:LEU:HD23	1:C:75:MET:HE3	1.83	0.60
2:D:508:THR:HA	2:D:526:CYS:HA	1.83	0.60
2:B:358:ILE:HG13	2:B:385:ARG:HH22	1.66	0.60
1:A:103:ASN:HD21	2:B:503:LEU:N	1.99	0.60
2:B:643:PHE:HB3	2:B:662:TRP:HB3	1.84	0.60
1:A:23:PRO:O	1:A:26:LYS:HG3	2.02	0.59
2:B:613:ARG:NH2	2:B:615:ASP:OD2	2.35	0.59
1:C:42:ILE:O	1:C:49:GLN:NE2	2.35	0.59
2:F:638:MET:HE3	2:F:671:LEU:HD11	1.85	0.59
2:B:344:SER:HB2	2:B:394:LEU:HD12	1.84	0.59
2:F:562:PRO:HG2	2:F:566:LEU:HD21	1.84	0.59
2:D:344:SER:HB2	2:D:394:LEU:HD12	1.83	0.59
2:F:373:LYS:NZ	2:F:374:ILE:O	2.35	0.59
2:D:446:ARG:HG2	2:D:495:GLU:HG2	1.84	0.59
2:D:770:ILE:HD11	2:D:807:PHE:HB3	1.84	0.59
2:B:680:LYS:HA	2:B:690:SER:HA	1.85	0.59
1:C:141:SER:O	1:C:145:VAL:HG12	2.03	0.59
2:B:428:VAL:HG22	2:B:522:SER:HA	1.82	0.58
2:D:275:ASN:HB2	2:D:301:SER:HB3	1.84	0.58
1:E:157:ILE:HA	1:E:160:GLN:HE21	1.68	0.58
2:F:778:ASP:OD1	2:F:779:ASP:N	2.34	0.58
1:C:33:THR:HG22	1:C:36:LYS:HZ3	1.67	0.58
2:D:465:HIS:ND1	2:D:480:THR:O	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:612:ARG:HE	2:F:616:GLY:HA2	1.68	0.58
1:C:91:SER:HB3	1:C:94:VAL:HG22	1.84	0.58
2:D:647:MET:HE3	2:D:656:ARG:HD2	1.84	0.58
2:F:418:HIS:CE1	2:F:419:ARG:HH21	2.21	0.58
2:F:693:VAL:HG12	2:F:696:ARG:HB2	1.85	0.58
2:B:443:MET:O	2:B:497:VAL:HA	2.04	0.58
2:D:339:LEU:HA	2:D:424:TYR:HB2	1.84	0.58
1:E:42:ILE:HA	1:E:45:ILE:HD12	1.85	0.58
2:B:422:GLU:OE1	2:B:424:TYR:OH	2.19	0.58
2:B:651:VAL:O	2:B:656:ARG:NH2	2.37	0.58
1:C:65:GLY:HA3	1:C:149:ARG:HH12	1.68	0.57
2:B:286:VAL:HG11	2:B:291:LEU:HD13	1.86	0.57
1:E:83:GLN:O	1:E:87:THR:HG23	2.04	0.57
2:B:360:SER:HB3	2:B:363:GLN:HE22	1.70	0.57
2:D:436:THR:HB	2:D:533:VAL:HA	1.86	0.57
2:D:650:ASP:HB2	2:D:653:LYS:HG3	1.85	0.57
2:F:298:PRO:HB3	2:F:332:VAL:HG23	1.86	0.57
2:B:634:VAL:HG12	2:B:635:LYS:HG3	1.85	0.57
2:F:446:ARG:HG2	2:F:495:GLU:HG2	1.85	0.57
2:F:464:TYR:HB3	2:F:509:MET:HG3	1.86	0.57
1:C:99:ASN:HD21	2:D:502:PHE:H	1.52	0.57
2:B:360:SER:HB3	2:B:363:GLN:NE2	2.20	0.57
2:B:650:ASP:HB3	2:B:653:LYS:HG3	1.86	0.56
1:E:69:ILE:HG12	1:E:75:MET:SD	2.45	0.56
2:D:434:CYS:HB2	2:D:444:THR:O	2.05	0.56
2:F:665:LEU:HD12	2:F:674:VAL:HG21	1.87	0.56
1:C:106:ASP:HB3	2:D:469:LEU:HD11	1.87	0.56
2:D:351:ILE:HD12	2:D:385:ARG:HD2	1.86	0.56
2:D:643:PHE:HB3	2:D:662:TRP:HB3	1.86	0.56
2:D:380:SER:HB2	2:D:388:LYS:HB3	1.87	0.56
2:D:564:ASN:OD1	2:D:565:ASN:N	2.39	0.56
1:A:83:GLN:HE22	1:A:102:GLU:HG3	1.70	0.56
2:F:651:VAL:O	2:F:656:ARG:NH2	2.39	0.56
2:B:458:SER:HB2	2:B:515:HIS:HD2	1.71	0.56
2:F:543:ALA:HB1	2:F:552:LEU:HD11	1.88	0.56
2:F:286:VAL:HG11	2:F:291:LEU:HD13	1.87	0.56
2:F:564:ASN:OD1	2:F:565:ASN:N	2.38	0.56
1:A:60:LEU:HD22	1:A:145:VAL:CG1	2.36	0.55
2:B:276:SER:OG	2:B:278:ILE:O	2.24	0.55
1:C:83:GLN:OE1	1:C:105:ARG:NE	2.39	0.55
1:E:86:LEU:HD22	1:E:94:VAL:HG13	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:679:VAL:HG22	2:D:712:VAL:HG22	1.89	0.55
2:D:752:VAL:HG21	2:D:827:THR:HG22	1.89	0.55
2:B:728:THR:HB	2:B:816:GLY:HA2	1.89	0.55
2:F:446:ARG:NH2	2:F:491:ASP:OD2	2.39	0.55
2:F:298:PRO:HG3	2:F:351:ILE:HG22	1.88	0.55
2:B:721:SER:OG	2:B:724:ASN:OD1	2.24	0.55
2:B:368:ARG:HB2	2:B:374:ILE:HD11	1.89	0.55
1:C:48:THR:O	1:C:51:VAL:HG22	2.06	0.55
2:F:244:MET:HG3	2:F:254:ILE:HG12	1.87	0.55
1:E:102:GLU:O	1:E:105:ARG:HG2	2.07	0.55
2:F:650:ASP:HB2	2:F:653:LYS:HG3	1.88	0.55
2:B:412:ASN:ND2	2:B:413:GLU:OE1	2.40	0.55
2:B:609:VAL:N	2:B:624:SER:OG	2.38	0.54
2:D:302:TYR:O	2:D:325:VAL:HA	2.07	0.54
1:A:60:LEU:HD22	1:A:145:VAL:HG13	1.89	0.54
2:D:286:VAL:HG11	2:D:291:LEU:HD13	1.89	0.54
2:D:337:LYS:NZ	2:D:421:ALA:HA	2.22	0.54
2:D:570:ILE:HG23	2:D:609:VAL:HG22	1.89	0.54
1:A:35:ILE:O	1:A:39:VAL:HG23	2.07	0.54
2:B:772:TRP:HH2	2:B:784:LEU:HG	1.72	0.54
2:F:409:TYR:CZ	2:F:420:TYR:HB3	2.43	0.54
2:F:728:THR:HB	2:F:816:GLY:HA2	1.90	0.54
1:A:69:ILE:HG22	1:A:75:MET:SD	2.48	0.54
2:F:452:ILE:HD12	2:F:513:ILE:HD12	1.89	0.54
2:D:373:LYS:NZ	2:D:374:ILE:O	2.40	0.54
1:A:31:THR:O	1:A:35:ILE:HG12	2.08	0.54
2:B:546:THR:HG22	2:B:553:LYS:NZ	2.22	0.54
1:C:153:SER:O	1:C:157:ILE:HG12	2.08	0.54
2:B:655:GLU:HA	2:B:703:TRP:O	2.08	0.54
2:D:666:THR:HG23	2:D:669:ASP:H	1.72	0.54
2:B:409:TYR:CE1	2:B:420:TYR:HB3	2.43	0.54
2:B:693:VAL:HG22	2:B:696:ARG:HB2	1.89	0.54
2:F:334:PHE:HB2	2:F:349:HIS:HB2	1.89	0.54
2:F:770:ILE:HD11	2:F:807:PHE:HB3	1.90	0.54
2:F:391:PHE:CE1	2:F:394:LEU:HA	2.43	0.53
1:A:40:THR:HG21	2:B:560:VAL:HG13	1.90	0.53
1:C:75:MET:SD	1:C:157:ILE:HD12	2.48	0.53
2:B:365:VAL:HG13	2:B:373:LYS:HE2	1.90	0.53
2:D:337:LYS:HZ3	2:D:421:ALA:HA	1.72	0.53
2:D:638:MET:HG3	2:D:722:LEU:HD12	1.91	0.53
2:D:715:VAL:HA	2:D:720:ALA:HA	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:741:SER:OG	2:F:757:THR:OG1	2.20	0.53
2:B:348:PHE:HE2	2:B:423:LEU:HD11	1.73	0.53
2:D:746:PRO:HA	2:D:752:VAL:HG22	1.90	0.53
2:B:244:MET:HG3	2:B:254:ILE:HG12	1.90	0.53
1:C:41:ARG:HH22	2:D:440:LEU:CB	2.21	0.53
2:F:458:SER:HB2	2:F:515:HIS:CD2	2.43	0.53
2:B:239:PRO:HG2	2:B:306:VAL:HG12	1.91	0.53
1:A:157:ILE:O	1:A:161:LEU:HG	2.09	0.52
1:C:102:GLU:O	1:C:105:ARG:HG2	2.10	0.52
2:D:564:ASN:HD22	2:D:613:ARG:HE	1.55	0.52
1:E:89:LEU:HB2	1:E:94:VAL:HG11	1.91	0.52
2:B:689:TRP:NE1	2:B:691:GLU:OE2	2.37	0.52
2:D:242:LEU:HD21	2:D:304:VAL:HG12	1.92	0.52
2:D:374:ILE:HD13	2:D:391:PHE:HZ	1.73	0.52
2:D:440:LEU:HD23	2:D:503:LEU:HD13	1.92	0.52
1:C:92:GLN:HA	1:C:95:LEU:HG	1.92	0.52
2:D:276:SER:OG	2:D:278:ILE:O	2.27	0.52
2:B:537:PRO:HG3	2:B:622:ASN:HB2	1.92	0.52
1:A:154:LEU:O	1:A:157:ILE:HG22	2.10	0.52
2:B:397:THR:HG23	2:B:425:VAL:HB	1.92	0.52
2:B:552:LEU:O	2:B:595:SER:HA	2.10	0.52
1:A:150:LEU:HD23	1:A:154:LEU:HD23	1.91	0.52
2:B:540:ASN:HB3	2:B:557:GLU:HG2	1.92	0.52
1:C:35:ILE:O	1:C:39:VAL:HG23	2.09	0.52
2:D:409:TYR:CZ	2:D:420:TYR:HB3	2.45	0.52
1:E:69:ILE:HA	1:E:74:LYS:HD2	1.91	0.52
2:F:606:VAL:HG13	2:F:628:TYR:HE1	1.75	0.52
2:F:644:TRP:NE1	2:F:771:GLU:OE2	2.39	0.52
1:E:99:ASN:HB3	2:F:501:ILE:O	2.10	0.52
2:F:445:CYS:O	2:F:495:GLU:HA	2.10	0.52
2:F:613:ARG:NH2	2:F:615:ASP:OD2	2.42	0.52
2:F:641:PRO:HB2	2:F:662:TRP:HB2	1.92	0.52
1:C:72:LEU:HB2	1:C:167:CYS:HA	1.91	0.51
2:D:309:LYS:HB2	2:D:317:TRP:CD2	2.45	0.51
2:D:728:THR:HB	2:D:816:GLY:HA2	1.92	0.51
2:B:304:VAL:O	2:B:323:PRO:HA	2.10	0.51
2:D:348:PHE:HZ	2:D:423:LEU:HD13	1.75	0.51
2:D:446:ARG:NH1	2:D:495:GLU:OE2	2.42	0.51
2:D:605:TYR:N	2:D:629:THR:OG1	2.39	0.51
2:F:733:MET:HE3	2:F:733:MET:O	2.11	0.51
2:B:645:ARG:HH21	2:B:817:VAL:HG11	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:635:LYS:O	2:D:716:ASN:ND2	2.43	0.51
1:C:92:GLN:O	1:C:95:LEU:HG	2.10	0.51
2:B:670:SER:HB2	2:B:674:VAL:HG23	1.91	0.51
2:F:571:ARG:HG2	2:F:623:TRP:CZ3	2.46	0.51
2:B:409:TYR:HH	1:C:141:SER:CB	2.19	0.51
2:D:665:LEU:HB2	2:D:670:SER:HB3	1.93	0.51
2:D:693:VAL:HG22	2:D:699:LEU:HD22	1.92	0.51
2:B:554:VAL:HG23	2:B:556:TRP:HZ3	1.76	0.51
2:B:644:TRP:CD2	2:B:769:VAL:HG11	2.46	0.51
2:D:814:MET:HE3	2:D:814:MET:O	2.11	0.51
2:F:537:PRO:HG3	2:F:622:ASN:HB2	1.93	0.50
2:D:299:GLY:N	2:D:328:THR:OG1	2.44	0.50
1:A:26:LYS:HE3	1:A:27:VAL:HG23	1.92	0.50
2:B:442:LYS:HB3	2:B:499:GLN:HB2	1.93	0.50
2:B:448:SER:HB2	2:B:493:PHE:CZ	2.47	0.50
2:F:740:GLU:HG2	2:F:759:SER:HB3	1.93	0.50
1:A:105:ARG:O	1:A:108:LEU:HG	2.12	0.50
2:D:445:CYS:O	2:D:495:GLU:HA	2.11	0.50
1:C:96:GLN:HE22	1:C:100:ASP:CG	2.18	0.50
2:B:770:ILE:HD11	2:B:807:PHE:HB3	1.94	0.50
2:D:399:PRO:HB3	2:D:404:THR:HG22	1.94	0.50
2:B:731:TRP:O	2:B:734:SER:OG	2.20	0.50
1:A:86:LEU:HD22	1:A:94:VAL:HG13	1.93	0.50
2:B:546:THR:OG1	2:B:551:LEU:N	2.45	0.50
2:B:654:LYS:NZ	2:B:704:THR:O	2.35	0.50
1:C:72:LEU:HA	1:C:75:MET:HE3	1.92	0.50
2:D:675:ARG:N	2:D:715:VAL:O	2.45	0.50
1:A:132:ASP:OD1	1:A:133:GLY:N	2.44	0.49
1:C:132:ASP:O	1:C:136:GLU:OE1	2.30	0.49
2:D:584:HIS:CD2	2:D:596:LEU:HD21	2.46	0.49
2:D:611:CYS:H	2:D:621:SER:HG	1.56	0.49
2:F:365:VAL:HG13	2:F:373:LYS:HE2	1.94	0.49
1:A:140:TYR:O	1:A:143:GLU:HG3	2.13	0.49
2:F:538:PRO:HG2	2:F:609:VAL:HG12	1.94	0.49
2:B:269:GLN:OE1	2:B:307:ARG:NH2	2.43	0.49
2:B:814:MET:HE3	2:B:814:MET:HA	1.94	0.49
2:D:537:PRO:HG3	2:D:622:ASN:HB2	1.95	0.49
1:E:70:LEU:HD23	1:E:70:LEU:H	1.77	0.49
1:A:157:ILE:HG13	1:A:161:LEU:HG	1.95	0.49
2:B:458:SER:HB2	2:B:515:HIS:CD2	2.47	0.49
2:F:237:ASP:O	2:F:308:SER:OG	2.29	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:489:GLN:NE2	2:B:495:GLU:HB2	2.27	0.49
1:C:57:VAL:HB	1:C:60:LEU:HD21	1.95	0.49
1:C:109:HIS:CD2	1:C:119:LEU:HD11	2.47	0.49
1:A:76:ASP:OD1	1:A:105:ARG:NH2	2.45	0.49
1:A:52:SER:HB3	1:A:55:GLN:HB2	1.93	0.49
2:B:715:VAL:HA	2:B:720:ALA:HA	1.95	0.49
2:D:270:VAL:HG22	2:D:306:VAL:HG22	1.93	0.49
2:D:571:ARG:HD2	2:D:581:TRP:CD2	2.48	0.49
1:E:24:ILE:HA	1:E:27:VAL:HG12	1.95	0.49
1:E:48:THR:HG22	1:E:144:VAL:HA	1.94	0.49
1:A:154:LEU:HA	1:A:157:ILE:HG22	1.95	0.49
2:B:274:GLU:HB3	2:B:280:ARG:HD3	1.94	0.49
1:C:140:TYR:HA	1:C:143:GLU:OE2	2.13	0.49
2:B:334:PHE:HB2	2:B:349:HIS:HB2	1.95	0.49
1:C:129:GLU:HG2	1:C:130:SER:H	1.77	0.49
2:D:319:ASP:OD1	2:D:320:TRP:N	2.46	0.49
2:F:768:LEU:HD13	2:F:809:LEU:HD21	1.94	0.49
2:B:244:MET:HB3	2:B:252:LEU:HD11	1.94	0.48
2:D:772:TRP:NE1	2:D:797:ASP:OD2	2.35	0.48
1:E:35:ILE:O	1:E:39:VAL:HG23	2.13	0.48
2:D:644:TRP:CG	2:D:645:ARG:H	2.31	0.48
2:F:693:VAL:HG22	2:F:699:LEU:HD22	1.95	0.48
2:B:752:VAL:HG21	2:B:827:THR:HG22	1.95	0.48
2:D:767:TYR:HA	2:D:788:SER:HB3	1.94	0.48
2:B:644:TRP:CG	2:B:645:ARG:H	2.31	0.48
1:C:83:GLN:NE2	1:C:102:GLU:HG2	2.10	0.48
2:D:458:SER:HB2	2:D:515:HIS:HD2	1.77	0.48
2:F:679:VAL:N	2:F:691:GLU:O	2.42	0.48
1:A:60:LEU:HG	1:A:60:LEU:O	2.13	0.48
2:D:348:PHE:HB2	2:D:389:VAL:HG12	1.95	0.48
2:D:571:ARG:NH2	2:D:608:GLN:OE1	2.47	0.48
2:F:434:CYS:HA	2:F:445:CYS:HA	1.94	0.48
2:F:541:VAL:HA	2:F:555:SER:O	2.13	0.48
2:F:613:ARG:NH1	2:F:617:LEU:O	2.46	0.48
2:B:343:GLY:N	2:B:394:LEU:O	2.32	0.48
2:D:511:ILE:N	2:D:522:SER:OG	2.29	0.48
2:B:774:ILE:HG23	2:B:803:GLU:HB3	1.95	0.48
2:D:651:VAL:O	2:D:656:ARG:NH2	2.47	0.48
1:E:65:GLY:O	1:E:67:HIS:ND1	2.34	0.48
2:B:449:PRO:HB2	2:B:452:ILE:HG22	1.94	0.48
2:D:544:GLU:HG3	2:D:553:LYS:HB3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:ARG:HE	1:A:56:ARG:HB3	1.36	0.48
2:B:738:ALA:HB2	2:B:813:PHE:HZ	1.79	0.48
2:D:349:HIS:HA	2:D:387:SER:O	2.13	0.48
2:D:417:HIS:HA	1:E:58:THR:OG1	2.14	0.48
2:D:634:VAL:HG12	2:D:635:LYS:HG3	1.96	0.48
1:E:76:ASP:OD1	1:E:108:LEU:HD13	2.14	0.48
2:F:452:ILE:C	2:F:454:SER:H	2.20	0.48
2:B:337:LYS:HG2	2:B:422:GLU:HB2	1.95	0.47
2:B:416:CYS:O	1:C:57:VAL:HA	2.14	0.47
2:D:541:VAL:HA	2:D:555:SER:O	2.14	0.47
2:B:449:PRO:HG3	2:B:494:TYR:CD2	2.49	0.47
2:D:417:HIS:HD2	2:D:419:ARG:HB2	1.78	0.47
2:B:640:GLY:N	2:B:723:VAL:O	2.39	0.47
2:D:738:ALA:HB2	2:D:813:PHE:HZ	1.78	0.47
2:F:477:ILE:HG23	2:F:510:TRP:CZ2	2.50	0.47
2:F:679:VAL:HG22	2:F:712:VAL:HG22	1.96	0.47
2:F:254:ILE:HG21	2:F:304:VAL:HG11	1.96	0.47
2:F:306:VAL:O	2:F:321:SER:N	2.44	0.47
2:F:463:ARG:HB2	2:F:510:TRP:NE1	2.29	0.47
2:F:540:ASN:HB2	2:F:557:GLU:HG2	1.97	0.47
2:B:769:VAL:O	2:B:810:TYR:N	2.29	0.47
1:C:99:ASN:ND2	2:D:501:ILE:C	2.72	0.47
2:B:647:MET:HE1	2:B:658:VAL:HG22	1.97	0.47
2:F:397:THR:HG21	2:F:405:TYR:HA	1.96	0.47
1:C:140:TYR:HA	1:C:143:GLU:CD	2.40	0.47
2:D:422:GLU:OE1	2:D:424:TYR:OH	2.25	0.47
2:F:738:ALA:HB2	2:F:813:PHE:HZ	1.79	0.47
2:B:538:PRO:HG2	2:B:609:VAL:HG12	1.96	0.47
1:C:34:LEU:O	1:C:38:ILE:HG12	2.14	0.47
1:C:67:HIS:O	1:C:68:PRO:C	2.56	0.47
2:F:646:LYS:HB3	2:F:659:THR:HB	1.97	0.47
1:A:39:VAL:HG22	1:A:154:LEU:CD1	2.45	0.47
1:A:98:ALA:O	1:A:102:GLU:OE1	2.33	0.47
2:D:488:LEU:HB2	2:D:494:TYR:CE1	2.50	0.47
2:D:682:ARG:HB3	2:D:709:THR:HB	1.97	0.47
2:B:337:LYS:HE2	2:B:422:GLU:HG3	1.97	0.47
1:C:92:GLN:OE1	1:C:92:GLN:N	2.39	0.47
2:F:612:ARG:HB2	2:F:620:TRP:CD2	2.49	0.47
2:B:368:ARG:HG3	2:B:369:ASN:HD22	1.79	0.46
2:B:442:LYS:HA	2:B:498:PHE:O	2.15	0.46
1:C:143:GLU:HG2	1:C:144:VAL:N	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:309:LYS:HB2	2:D:317:TRP:CE3	2.50	0.46
2:D:768:LEU:HD13	2:D:809:LEU:HD21	1.96	0.46
2:F:366:TRP:CE2	2:F:389:VAL:HB	2.50	0.46
2:F:478:HIS:CG	2:F:479:PRO:HD2	2.50	0.46
2:F:275:ASN:HB2	2:F:301:SER:HB3	1.97	0.46
2:B:427:ASP:HB3	2:B:451:THR:HG21	1.98	0.46
2:D:612:ARG:HB2	2:D:620:TRP:CE3	2.51	0.46
1:A:95:LEU:HB3	2:B:500:PRO:HG3	1.97	0.46
2:F:613:ARG:CZ	2:F:617:LEU:HB2	2.45	0.46
1:C:70:LEU:H	1:C:74:LYS:HE3	1.81	0.46
2:D:271:LYS:HG3	2:D:320:TRP:CH2	2.50	0.46
1:E:27:VAL:O	1:E:31:THR:HG23	2.15	0.46
2:F:424:TYR:HB3	2:F:426:ILE:HD11	1.97	0.46
2:B:244:MET:HE1	2:B:324:GLN:HB3	1.97	0.46
2:D:351:ILE:HD11	2:D:386:VAL:HG22	1.97	0.46
2:D:439:TYR:O	2:D:441:THR:HG23	2.16	0.46
2:F:422:GLU:OE1	2:F:424:TYR:OH	2.23	0.46
1:A:48:THR:HG22	1:A:144:VAL:HG23	1.97	0.46
2:B:682:ARG:HB3	2:B:709:THR:HB	1.97	0.46
1:E:33:THR:O	1:E:37:THR:HG23	2.15	0.46
2:F:243:HIS:NE2	2:F:245:GLU:HB3	2.30	0.46
2:F:601:LEU:HD12	2:F:671:LEU:HD13	1.98	0.46
2:B:442:LYS:HB2	2:B:497:VAL:CG1	2.46	0.46
2:B:526:CYS:SG	2:B:527:VAL:N	2.89	0.46
2:B:655:GLU:OE1	2:B:702:LEU:HD11	2.16	0.46
1:C:145:VAL:HG13	1:C:146:ALA:N	2.31	0.46
2:F:769:VAL:O	2:F:810:TYR:N	2.26	0.46
2:B:508:THR:HG22	2:B:526:CYS:HB2	1.98	0.46
2:F:276:SER:OG	2:F:278:ILE:O	2.31	0.46
2:B:294:ASP:OD2	2:B:295:SER:OG	2.20	0.46
2:B:571:ARG:HE	2:B:608:GLN:HB2	1.80	0.46
1:C:69:ILE:HG23	1:C:74:LYS:HG2	1.96	0.46
2:F:378:GLN:NE2	2:F:390:THR:O	2.41	0.46
1:C:99:ASN:HD22	2:D:501:ILE:C	2.24	0.45
2:D:526:CYS:SG	2:D:527:VAL:N	2.89	0.45
2:D:773:LYS:HD3	2:D:781:MET:HE1	1.98	0.45
1:E:82:TYR:O	1:E:86:LEU:HG	2.17	0.45
2:F:251:ASN:ND2	2:F:294:ASP:OD1	2.49	0.45
1:E:32:LYS:O	1:E:35:ILE:HG22	2.17	0.45
2:B:377:ILE:H	2:B:377:ILE:HD12	1.81	0.45
2:D:503:LEU:HG	2:D:504:LEU:HD22	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:247:THR:HG21	2:F:292:LEU:HD21	1.98	0.45
2:B:254:ILE:HG21	2:B:304:VAL:HG11	1.98	0.45
2:D:458:SER:HB2	2:D:515:HIS:CD2	2.51	0.45
2:F:721:SER:OG	2:F:724:ASN:OD1	2.27	0.45
2:F:404:THR:HG21	2:F:454:SER:HB2	1.97	0.45
2:F:570:ILE:HG12	2:F:609:VAL:HG22	1.97	0.45
1:A:45:ILE:HG23	1:A:47:HIS:CE1	2.51	0.45
2:B:571:ARG:HH21	2:B:608:GLN:CD	2.24	0.45
2:D:405:TYR:HB3	2:D:424:TYR:CE2	2.52	0.45
1:E:28:GLN:O	1:E:31:THR:OG1	2.29	0.45
2:F:772:TRP:CE2	2:F:782:LYS:HB2	2.52	0.45
2:D:509:MET:HE1	2:D:527:VAL:HG12	1.98	0.45
1:E:46:SER:O	1:E:49:GLN:HG3	2.15	0.45
2:F:306:VAL:N	2:F:321:SER:OG	2.46	0.45
2:B:736:VAL:O	2:B:819:LYS:N	2.32	0.45
2:F:304:VAL:O	2:F:323:PRO:HA	2.17	0.45
2:F:775:LEU:HD13	2:F:806:GLN:HB2	1.98	0.45
2:B:302:TYR:O	2:B:325:VAL:HA	2.17	0.45
2:D:613:ARG:NH1	2:D:617:LEU:O	2.50	0.45
2:F:272:TYR:CE1	2:F:282:ALA:HB3	2.52	0.45
2:F:436:THR:HB	2:F:533:VAL:HG22	1.98	0.45
1:A:82:TYR:O	1:A:86:LEU:HG	2.17	0.45
1:A:106:ASP:O	1:A:110:LEU:HD23	2.17	0.45
1:A:51:VAL:HG21	1:A:144:VAL:HG21	1.98	0.44
2:D:303:GLU:HA	2:D:324:GLN:O	2.17	0.44
2:D:714:ALA:O	2:D:721:SER:N	2.50	0.44
2:F:361:SER:OG	2:F:383:SER:O	2.22	0.44
2:B:546:THR:HG1	2:B:551:LEU:H	1.65	0.44
1:A:46:SER:O	1:A:49:GLN:HG2	2.16	0.44
2:B:404:THR:HG21	2:B:454:SER:HB2	1.99	0.44
2:B:487:VAL:O	2:B:489:GLN:NE2	2.49	0.44
1:E:105:ARG:O	1:E:108:LEU:HG	2.18	0.44
2:F:644:TRP:CE2	2:F:769:VAL:HG11	2.53	0.44
1:A:67:HIS:O	1:A:68:PRO:C	2.60	0.44
2:D:721:SER:OG	2:D:724:ASN:OD1	2.29	0.44
1:A:23:PRO:HG2	1:A:26:LYS:CG	2.43	0.44
1:A:56:ARG:HH12	2:F:415:ALA:HA	1.82	0.44
2:B:608:GLN:NE2	2:B:624:SER:O	2.30	0.44
2:D:543:ALA:HB1	2:D:552:LEU:HD11	2.00	0.44
2:F:347:SER:HB3	2:F:390:THR:HG23	2.00	0.44
2:F:418:HIS:CE1	2:F:419:ARG:HE	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:272:TYR:CZ	2:B:282:ALA:HB3	2.53	0.44
2:B:417:HIS:CG	2:B:418:HIS:H	2.35	0.44
2:D:438:GLY:HA3	2:D:536:LEU:HD21	1.98	0.44
2:F:372:GLU:N	2:F:372:GLU:OE1	2.51	0.44
1:A:63:ILE:HA	1:A:64:PRO:HD3	1.78	0.44
2:B:239:PRO:HG3	2:B:308:SER:HB3	2.00	0.44
2:B:666:THR:O	2:B:670:SER:N	2.38	0.44
2:D:341:SER:C	2:D:425:VAL:HG13	2.43	0.44
2:D:797:ASP:OD1	2:D:798:ASN:N	2.51	0.44
2:F:301:SER:HA	2:F:327:THR:HA	2.00	0.44
2:F:511:ILE:N	2:F:522:SER:OG	2.37	0.44
2:F:612:ARG:HH11	2:F:620:TRP:CD1	2.36	0.44
2:F:270:VAL:HG22	2:F:306:VAL:HG13	1.98	0.44
2:F:773:LYS:HD3	2:F:781:MET:CE	2.48	0.44
2:B:306:VAL:O	2:B:321:SER:N	2.49	0.43
2:F:478:HIS:CD2	2:F:479:PRO:HD2	2.53	0.43
2:F:564:ASN:HD22	2:F:613:ARG:HE	1.65	0.43
2:F:570:ILE:HG23	2:F:609:VAL:HG22	1.99	0.43
2:D:268:TYR:HE2	2:D:289:THR:HG22	1.83	0.43
2:D:567:GLN:HG2	2:D:614:LEU:HA	2.00	0.43
2:F:458:SER:HA	2:F:514:ASN:O	2.18	0.43
2:F:647:MET:HA	2:F:647:MET:HE2	1.99	0.43
2:B:470:TYR:HA	2:B:528:LEU:HD11	2.00	0.43
1:C:33:THR:HA	1:C:36:LYS:NZ	2.33	0.43
2:D:441:THR:C	2:D:501:ILE:HD13	2.43	0.43
2:D:644:TRP:CE2	2:D:769:VAL:HG11	2.53	0.43
2:F:612:ARG:HG2	2:F:613:ARG:N	2.33	0.43
2:F:670:SER:O	2:F:672:CYS:N	2.50	0.43
2:B:509:MET:HE1	2:B:527:VAL:HG12	2.01	0.43
2:B:546:THR:HG22	2:B:553:LYS:HZ3	1.81	0.43
2:B:646:LYS:HB3	2:B:659:THR:HB	1.99	0.43
2:D:272:TYR:CE1	2:D:282:ALA:HB3	2.52	0.43
2:D:654:LYS:NZ	2:D:704:THR:O	2.34	0.43
2:B:237:ASP:O	2:B:308:SER:OG	2.36	0.43
2:B:645:ARG:HD3	2:B:729:PHE:CE1	2.52	0.43
2:D:488:LEU:HB2	2:D:494:TYR:CZ	2.54	0.43
2:B:419:ARG:HA	1:C:138:SER:OG	2.18	0.43
1:C:96:GLN:HB2	2:D:501:ILE:H	1.84	0.43
2:D:612:ARG:HB2	2:D:620:TRP:CD2	2.54	0.43
1:E:67:HIS:O	1:E:68:PRO:C	2.61	0.43
2:F:571:ARG:HG2	2:F:623:TRP:CE3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:405:TYR:HB3	2:B:424:TYR:CE2	2.53	0.43
1:C:139:LEU:O	1:C:143:GLU:OE1	2.36	0.43
2:D:242:LEU:HD13	2:D:256:TRP:HB3	2.00	0.43
2:F:681:HIS:HB3	2:F:708:HIS:NE2	2.34	0.43
1:A:99:ASN:HA	1:A:102:GLU:OE1	2.19	0.43
2:B:251:ASN:OD1	2:B:294:ASP:HA	2.19	0.43
2:D:361:SER:N	2:D:384:ASP:O	2.50	0.43
1:E:65:GLY:H	1:E:149:ARG:HH21	1.67	0.43
1:E:85:VAL:HA	1:E:131:LEU:HD21	2.00	0.43
2:F:738:ALA:HB2	2:F:813:PHE:CZ	2.54	0.43
2:F:774:ILE:HG23	2:F:803:GLU:HB3	2.00	0.43
1:A:29:ASP:HA	1:A:32:LYS:NZ	2.33	0.43
2:B:767:TYR:HA	2:B:788:SER:HB3	2.01	0.43
2:D:330:ASP:HB2	2:D:419:ARG:HH21	1.84	0.43
1:E:29:ASP:HA	1:E:32:LYS:HG2	2.00	0.43
1:A:47:HIS:CG	2:F:400:ARG:HD3	2.54	0.43
1:A:155:GLN:O	1:A:159:GLN:OE1	2.37	0.43
1:E:73:SER:O	1:E:77:GLN:OE1	2.37	0.43
2:B:419:ARG:HE	1:C:137:ALA:HB1	1.84	0.42
2:B:437:ASP:OD2	2:B:441:THR:OG1	2.21	0.42
2:B:541:VAL:HG22	2:B:556:TRP:HB3	2.01	0.42
1:C:95:LEU:HD12	1:C:96:GLN:N	2.34	0.42
1:E:29:ASP:O	1:E:32:LYS:HG2	2.19	0.42
2:F:488:LEU:HD13	2:F:494:TYR:HE1	1.84	0.42
2:B:275:ASN:HB2	2:B:301:SER:HB3	2.00	0.42
2:F:731:TRP:O	2:F:734:SER:OG	2.31	0.42
2:B:354:ASN:ND2	2:B:412:ASN:OD1	2.51	0.42
2:D:800:ILE:HG22	2:D:802:ILE:HG12	2.01	0.42
2:F:239:PRO:HG2	2:F:306:VAL:HG12	1.99	0.42
2:F:250:GLY:HA2	2:F:334:PHE:CE1	2.53	0.42
2:F:378:GLN:HB3	2:F:390:THR:O	2.18	0.42
2:F:448:SER:HB2	2:F:493:PHE:CE1	2.54	0.42
2:D:641:PRO:HD3	2:D:714:ALA:HB2	2.00	0.42
2:F:733:MET:HG2	2:F:816:GLY:HA3	2.01	0.42
2:B:530:ASP:OD1	2:B:531:SER:N	2.53	0.42
2:B:742:LEU:HG	2:B:823:ILE:HD13	2.00	0.42
1:C:42:ILE:HA	1:C:45:ILE:CD1	2.49	0.42
2:F:353:LYS:HD3	2:F:356:ASN:H	1.84	0.42
2:F:466:ARG:HD3	2:F:507:TYR:CE2	2.55	0.42
1:C:29:ASP:O	1:C:33:THR:HG23	2.19	0.42
1:C:131:LEU:HD12	1:C:132:ASP:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:239:PRO:HG2	2:D:306:VAL:HG12	1.99	0.42
2:B:374:ILE:HB	2:B:379:TYR:HE1	1.84	0.42
1:C:104:LEU:O	1:C:108:LEU:HG	2.19	0.42
2:B:644:TRP:CG	2:B:645:ARG:N	2.87	0.42
1:C:99:ASN:O	1:C:102:GLU:HB2	2.20	0.42
2:D:459:THR:HG1	2:D:514:ASN:HB2	1.84	0.42
2:D:553:LYS:NZ	2:D:593:SER:HB2	2.35	0.42
2:D:637:PRO:HD3	2:D:716:ASN:CG	2.45	0.42
1:E:33:THR:HA	1:E:36:LYS:HG2	2.01	0.42
1:C:41:ARG:HG2	1:C:97:ILE:HD11	2.00	0.42
2:D:569:GLN:HG2	2:D:623:TRP:HH2	1.85	0.42
2:F:774:ILE:HD11	2:F:800:ILE:HD12	2.02	0.42
1:A:29:ASP:OD1	1:A:32:LYS:NZ	2.49	0.42
1:A:63:ILE:HA	1:A:63:ILE:HD13	1.88	0.42
2:B:349:HIS:HA	2:B:387:SER:O	2.20	0.42
2:B:466:ARG:HD3	2:B:507:TYR:CE1	2.55	0.42
2:B:488:LEU:HB2	2:B:494:TYR:HE1	1.85	0.42
2:B:274:GLU:O	2:B:279:VAL:HA	2.20	0.41
2:B:713:LEU:HB3	2:B:720:ALA:HB1	2.01	0.41
1:C:56:ARG:NH2	1:C:149:ARG:HH11	2.15	0.41
2:D:334:PHE:HA	2:D:335:PRO:HA	1.90	0.41
2:F:309:LYS:HG2	2:F:310:ARG:N	2.35	0.41
2:F:556:TRP:O	2:F:592:LYS:HB3	2.19	0.41
2:D:241:GLY:O	2:D:256:TRP:HA	2.20	0.41
2:D:706:PRO:O	2:D:730:SER:HA	2.20	0.41
2:F:331:VAL:HG13	2:F:350:CYS:SG	2.60	0.41
2:F:343:GLY:N	2:F:394:LEU:O	2.38	0.41
1:C:140:TYR:O	1:C:144:VAL:HG22	2.20	0.41
2:D:309:LYS:HD3	2:D:317:TRP:CG	2.55	0.41
2:D:675:ARG:H	2:D:716:ASN:HA	1.84	0.41
2:D:731:TRP:O	2:D:734:SER:OG	2.28	0.41
1:A:140:TYR:O	1:A:144:VAL:HG12	2.20	0.41
1:C:93:ASN:O	1:C:97:ILE:HG13	2.20	0.41
1:C:96:GLN:CD	1:C:96:GLN:C	2.88	0.41
2:F:374:ILE:O	2:F:379:TYR:OH	2.29	0.41
1:A:31:THR:O	1:A:34:LEU:HG	2.21	0.41
1:A:58:THR:OG1	2:F:417:HIS:HA	2.20	0.41
2:B:436:THR:HG21	2:B:533:VAL:HG22	2.02	0.41
2:D:420:TYR:CZ	1:E:140:TYR:HB3	2.55	0.41
2:B:338:ILE:HD12	2:B:348:PHE:CE2	2.55	0.41
2:B:768:LEU:HB3	2:B:809:LEU:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:452:ILE:C	2:D:454:SER:H	2.29	0.41
2:D:675:ARG:HG3	2:D:716:ASN:C	2.46	0.41
1:E:62:PHE:CZ	1:E:134:VAL:HG21	2.56	0.41
1:E:63:ILE:HA	1:E:64:PRO:HD3	1.82	0.41
2:F:366:TRP:CD2	2:F:389:VAL:HB	2.55	0.41
2:F:567:GLN:NE2	2:F:612:ARG:HD3	2.35	0.41
1:A:154:LEU:HD12	1:A:155:GLN:N	2.36	0.41
2:B:340:THR:OG1	2:B:341:SER:N	2.52	0.41
2:B:644:TRP:CE2	2:B:769:VAL:HG11	2.56	0.41
2:B:767:TYR:HB2	2:B:786:ILE:O	2.20	0.41
2:B:772:TRP:CE2	2:B:782:LYS:HB2	2.55	0.41
2:B:650:ASP:OD1	2:B:651:VAL:N	2.54	0.41
2:B:714:ALA:HB3	2:B:721:SER:HB3	2.02	0.41
1:C:103:ASN:ND2	2:D:502:PHE:HD1	2.19	0.41
2:D:348:PHE:O	2:D:388:LYS:HA	2.21	0.41
2:D:459:THR:OG1	2:D:514:ASN:HB2	2.20	0.41
2:D:681:HIS:N	2:D:689:TRP:O	2.42	0.41
1:A:162:ASP:OD1	1:A:162:ASP:N	2.53	0.41
2:B:445:CYS:O	2:B:495:GLU:HA	2.20	0.41
1:C:46:SER:HA	1:C:49:GLN:OE1	2.20	0.41
1:C:64:PRO:HB2	1:C:78:THR:HG23	2.03	0.41
1:C:82:TYR:O	1:C:86:LEU:HG	2.21	0.41
2:D:417:HIS:CD2	2:D:419:ARG:HB2	2.55	0.41
2:D:552:LEU:O	2:D:595:SER:HA	2.20	0.41
2:D:608:GLN:HB3	2:D:623:TRP:HB3	2.02	0.41
1:E:105:ARG:HA	1:E:108:LEU:HG	2.03	0.41
2:F:344:SER:O	2:F:393:ASN:N	2.54	0.41
2:F:409:TYR:CE2	2:F:420:TYR:HB3	2.55	0.41
2:F:612:ARG:HD2	2:F:620:TRP:NE1	2.35	0.41
2:F:797:ASP:HB3	2:F:799:PHE:CZ	2.56	0.41
2:B:250:GLY:O	2:B:296:VAL:HG22	2.21	0.41
1:E:32:LYS:HA	1:E:35:ILE:HG22	2.03	0.41
2:F:273:LEU:HB2	2:F:281:GLU:OE2	2.20	0.41
2:F:380:SER:HB2	2:F:388:LYS:HB3	2.02	0.41
2:F:419:ARG:HD3	2:F:419:ARG:HA	1.89	0.41
1:A:47:HIS:HB3	2:F:400:ARG:HD3	2.03	0.40
1:C:36:LYS:O	1:C:40:THR:HG23	2.21	0.40
2:D:382:VAL:HG23	2:D:388:LYS:HB2	2.02	0.40
2:F:319:ASP:OD1	2:F:320:TRP:N	2.55	0.40
2:F:528:LEU:O	2:F:531:SER:OG	2.39	0.40
2:F:569:GLN:OE1	2:F:610:ARG:NE	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:797:ASP:HB3	2:F:799:PHE:CE2	2.56	0.40
1:A:29:ASP:HA	1:A:32:LYS:HG2	2.02	0.40
1:A:105:ARG:HA	1:A:108:LEU:HG	2.02	0.40
2:D:345:ASN:HA	2:D:391:PHE:O	2.21	0.40
2:D:466:ARG:HA	2:D:506:GLY:O	2.21	0.40
2:D:775:LEU:HD13	2:D:806:GLN:HB2	2.04	0.40
2:F:440:LEU:HD13	2:F:561:PHE:CD2	2.56	0.40
1:A:41:ARG:O	1:A:45:ILE:HD12	2.22	0.40
2:B:571:ARG:HA	2:B:582:LYS:O	2.22	0.40
2:B:768:LEU:HD23	2:B:811:PRO:HA	2.03	0.40
2:B:784:LEU:HD23	2:B:784:LEU:HA	1.90	0.40
1:C:39:VAL:HG22	1:C:154:LEU:CD1	2.51	0.40
2:D:449:PRO:HB2	2:D:452:ILE:HG22	2.03	0.40
1:E:34:LEU:O	1:E:37:THR:OG1	2.33	0.40
1:E:42:ILE:HG21	1:E:151:GLN:HB2	2.03	0.40
1:E:111:LEU:O	1:E:115:LYS:HG2	2.21	0.40
1:A:41:ARG:HD2	1:A:97:ILE:HD13	2.02	0.40
1:A:60:LEU:HD12	1:A:62:PHE:CZ	2.56	0.40
2:B:359:ILE:HD11	2:B:364:ILE:HG12	2.03	0.40
1:C:92:GLN:HG2	1:C:93:ASN:N	2.37	0.40
2:F:610:ARG:HG2	2:F:623:TRP:CZ3	2.56	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	136/174 (78%)	122 (90%)	12 (9%)	2 (2%)	8	38
1	C	136/174 (78%)	122 (90%)	12 (9%)	2 (2%)	8	38
1	E	136/174 (78%)	122 (90%)	12 (9%)	2 (2%)	8	38
2	B	591/868 (68%)	570 (96%)	21 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	591/868 (68%)	571 (97%)	20 (3%)	0	100	100
2	F	591/868 (68%)	573 (97%)	17 (3%)	1 (0%)	43	77
All	All	2181/3126 (70%)	2080 (95%)	94 (4%)	7 (0%)	37	71

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	PRO
1	C	68	PRO
1	E	68	PRO
2	F	671	LEU
1	A	163	VAL
1	C	163	VAL
1	E	163	VAL

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	129/155 (83%)	129 (100%)	0	100	100
1	C	129/155 (83%)	129 (100%)	0	100	100
1	E	129/155 (83%)	129 (100%)	0	100	100
2	B	546/781 (70%)	546 (100%)	0	100	100
2	D	546/781 (70%)	546 (100%)	0	100	100
2	F	546/781 (70%)	546 (100%)	0	100	100
All	All	2025/2808 (72%)	2025 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	83	GLN
1	A	103	ASN
2	B	354	ASN
2	B	363	GLN
2	B	461	GLN
2	B	489	GLN
2	B	668	ASN
1	C	99	ASN
1	C	159	GLN
2	D	354	ASN
1	E	55	GLN
1	E	99	ASN
1	E	159	GLN
2	F	378	GLN
2	F	461	GLN
2	F	478	HIS
2	F	548	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

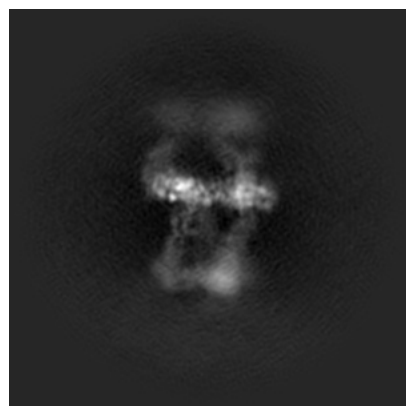
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-15678. These allow visual inspection of the internal detail of the map and identification of artifacts.

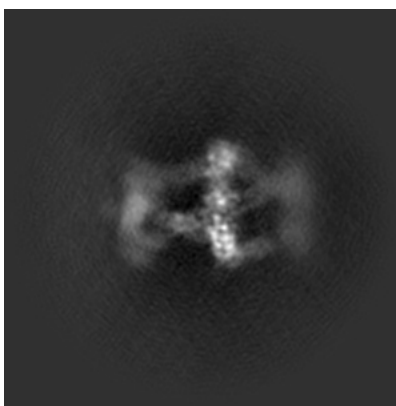
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

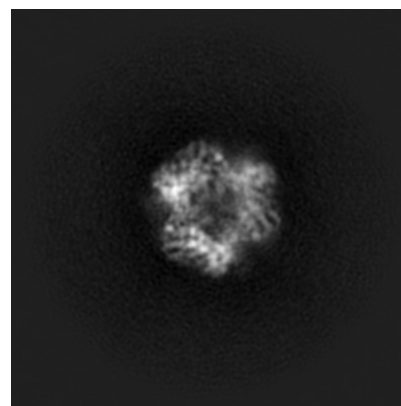
6.1.1 Primary map



X

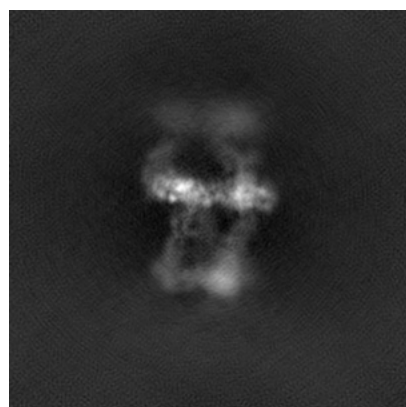


Y

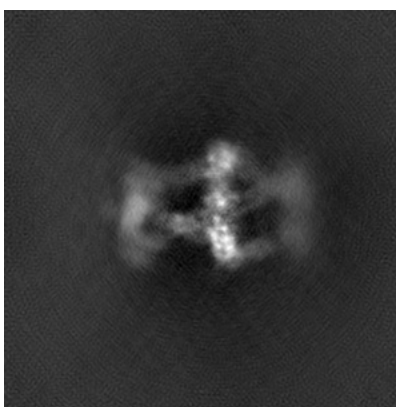


Z

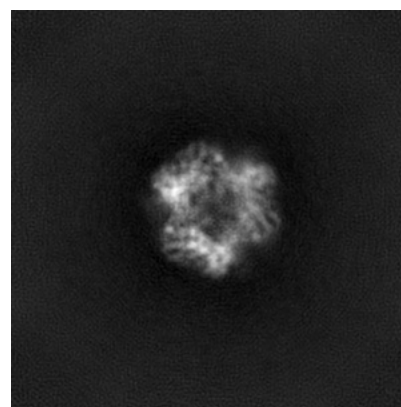
6.1.2 Raw map



X



Y

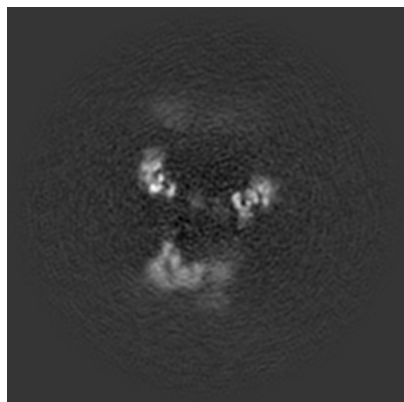


Z

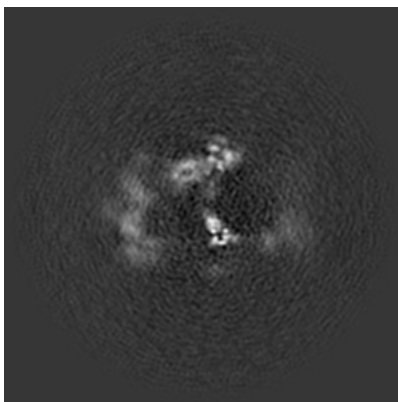
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

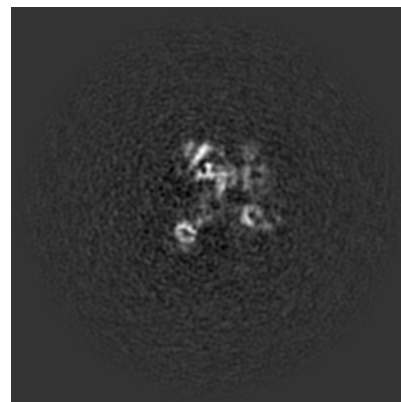
6.2.1 Primary map



X Index: 112

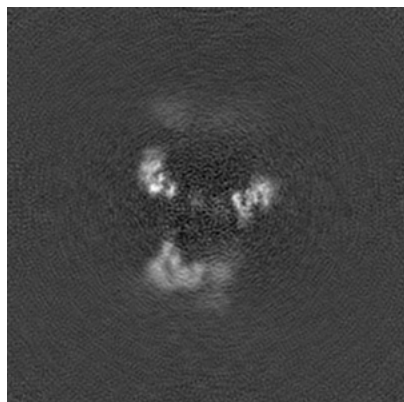


Y Index: 112

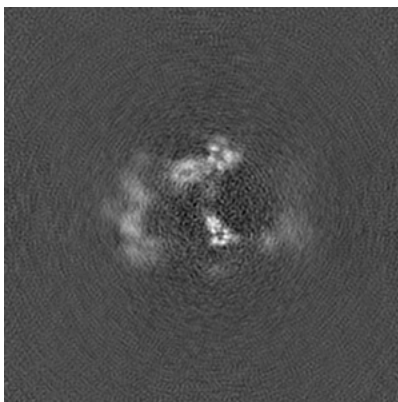


Z Index: 112

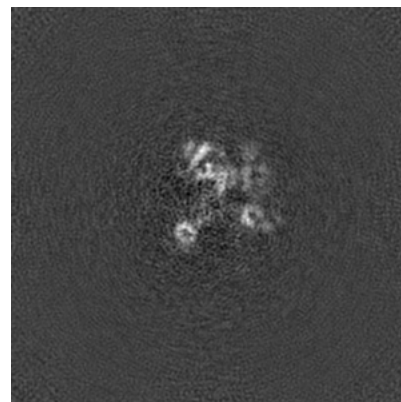
6.2.2 Raw map



X Index: 112



Y Index: 112

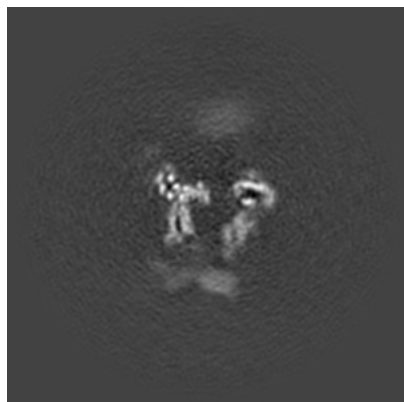


Z Index: 112

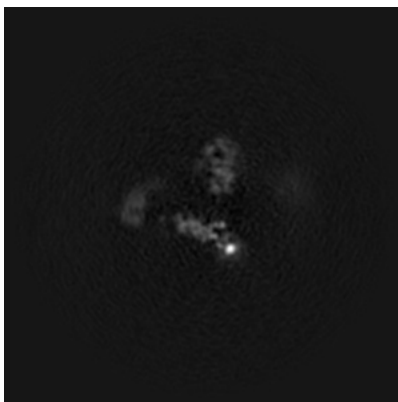
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

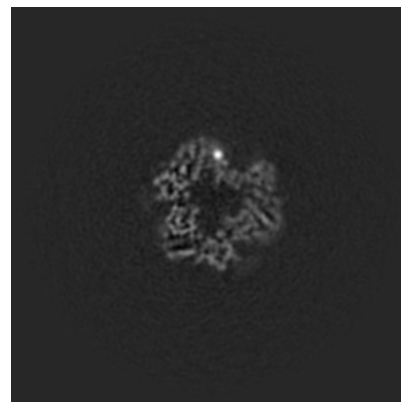
6.3.1 Primary map



X Index: 101

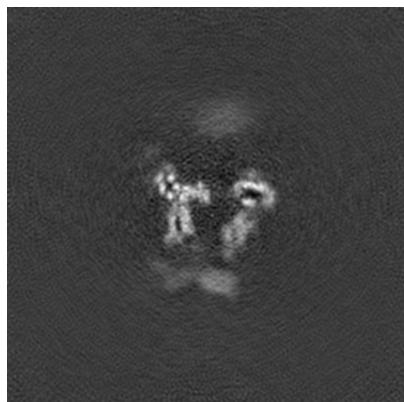


Y Index: 100

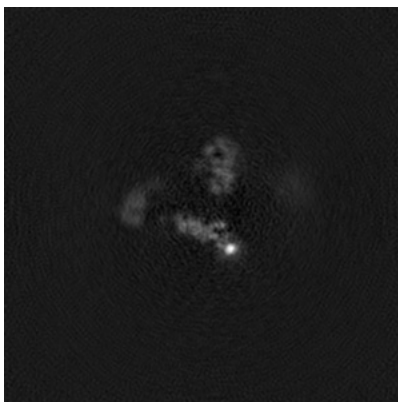


Z Index: 121

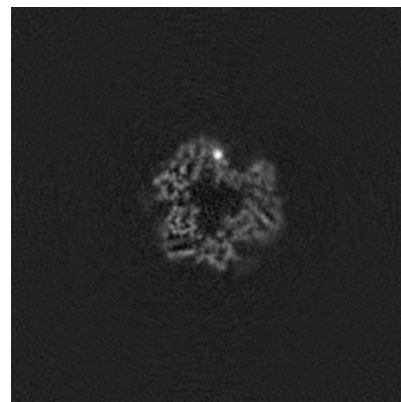
6.3.2 Raw map



X Index: 101



Y Index: 100

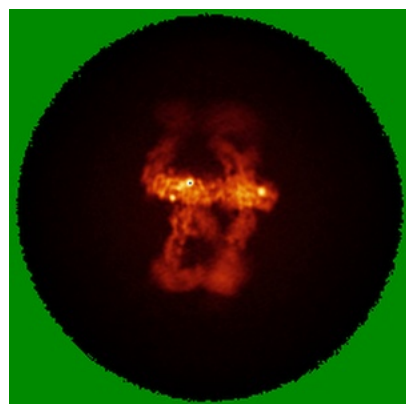


Z Index: 121

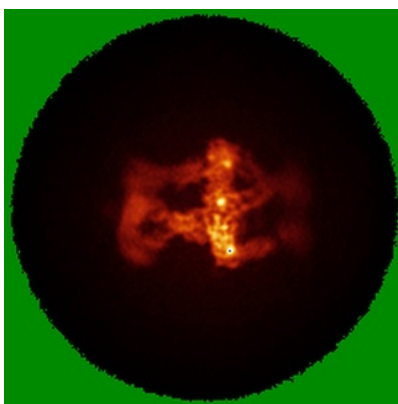
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

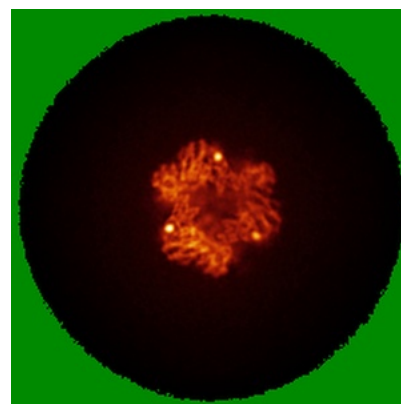
6.4.1 Primary map



X

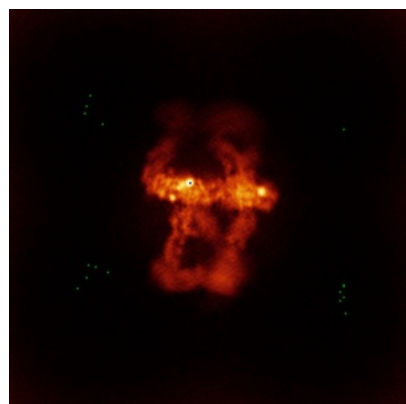


Y

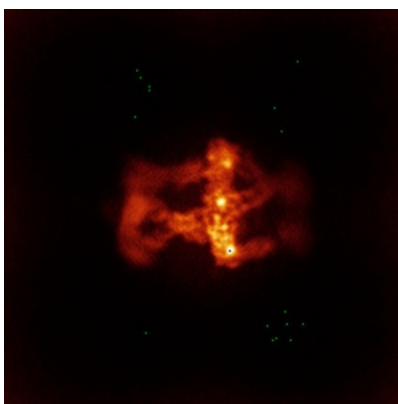


Z

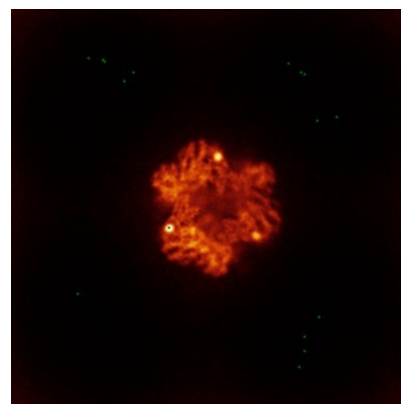
6.4.2 Raw map



X



Y

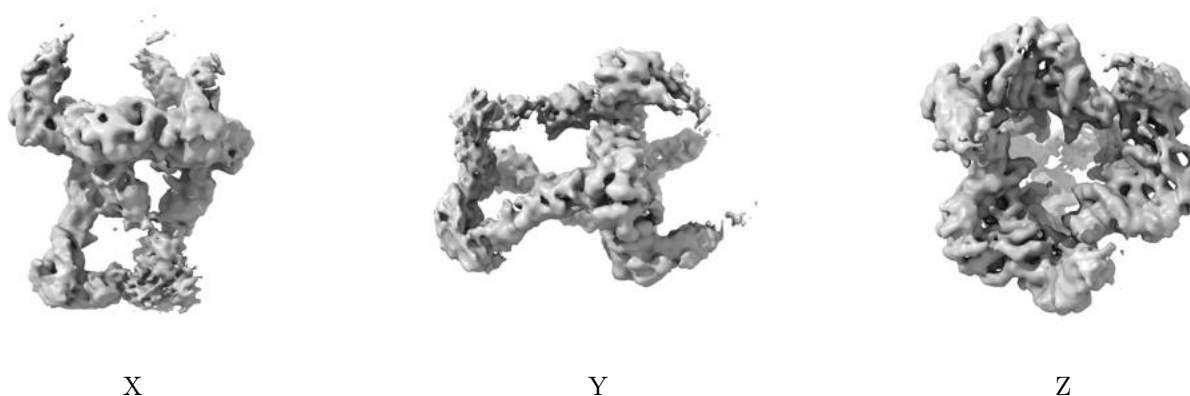


Z

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 0.178. 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.

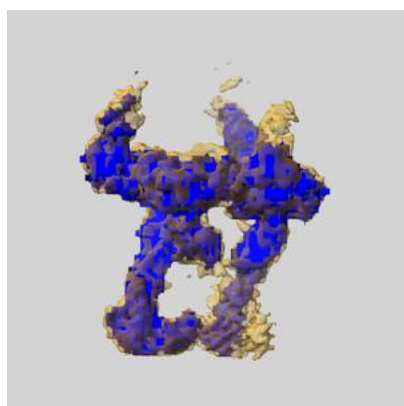
6.6 Mask visualisation [i](#)

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

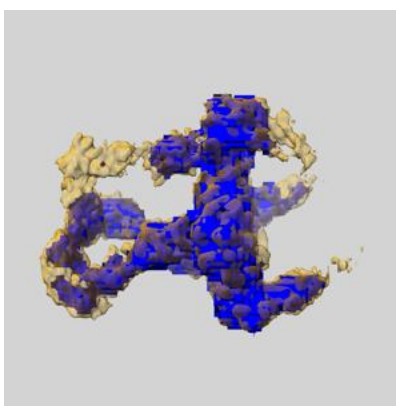
A mask typically either:

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

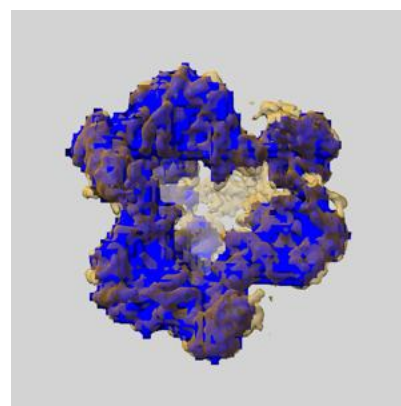
6.6.1 emd_15678_msk_1.map [i](#)



X



Y

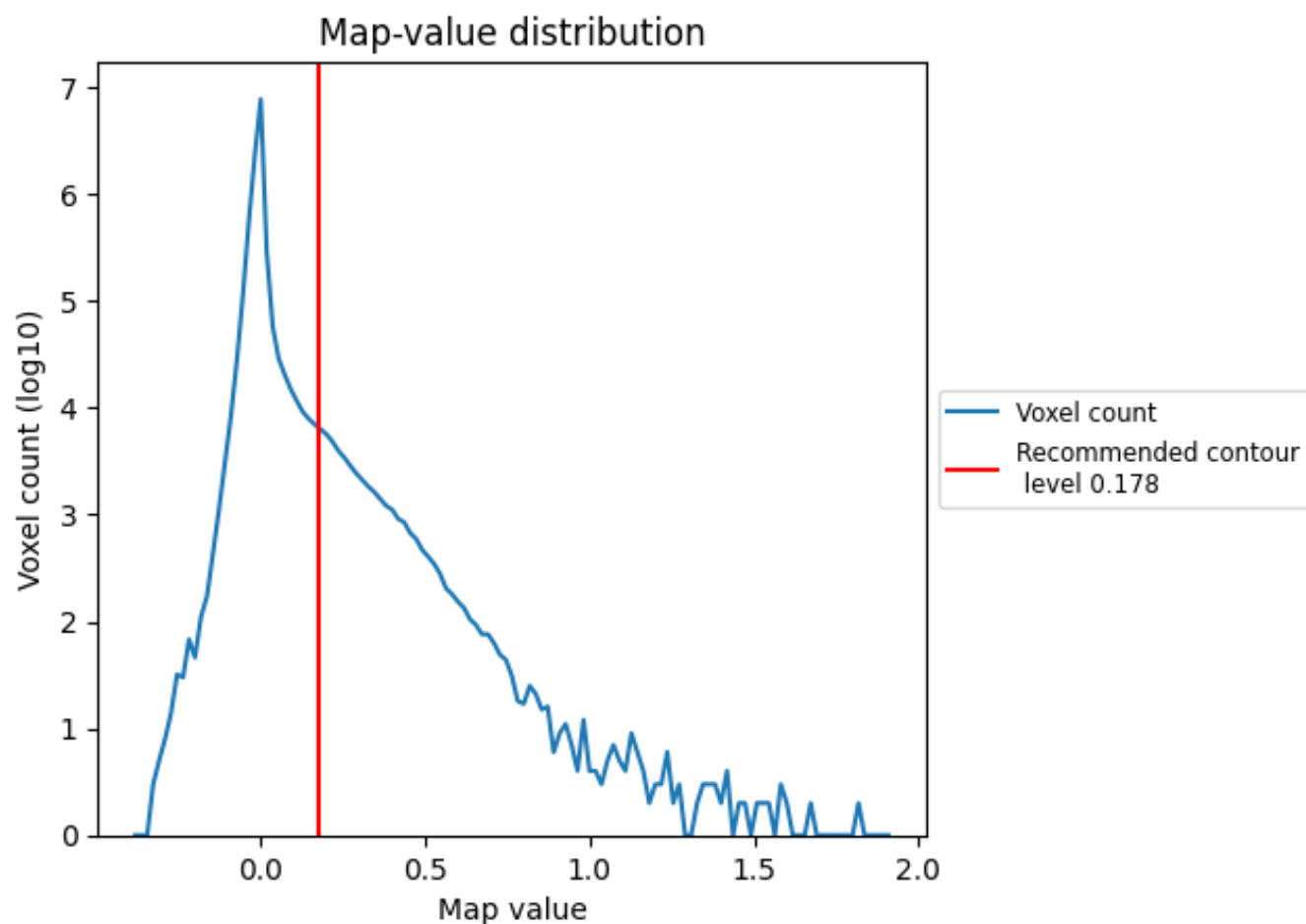


Z

7 Map analysis [i](#)

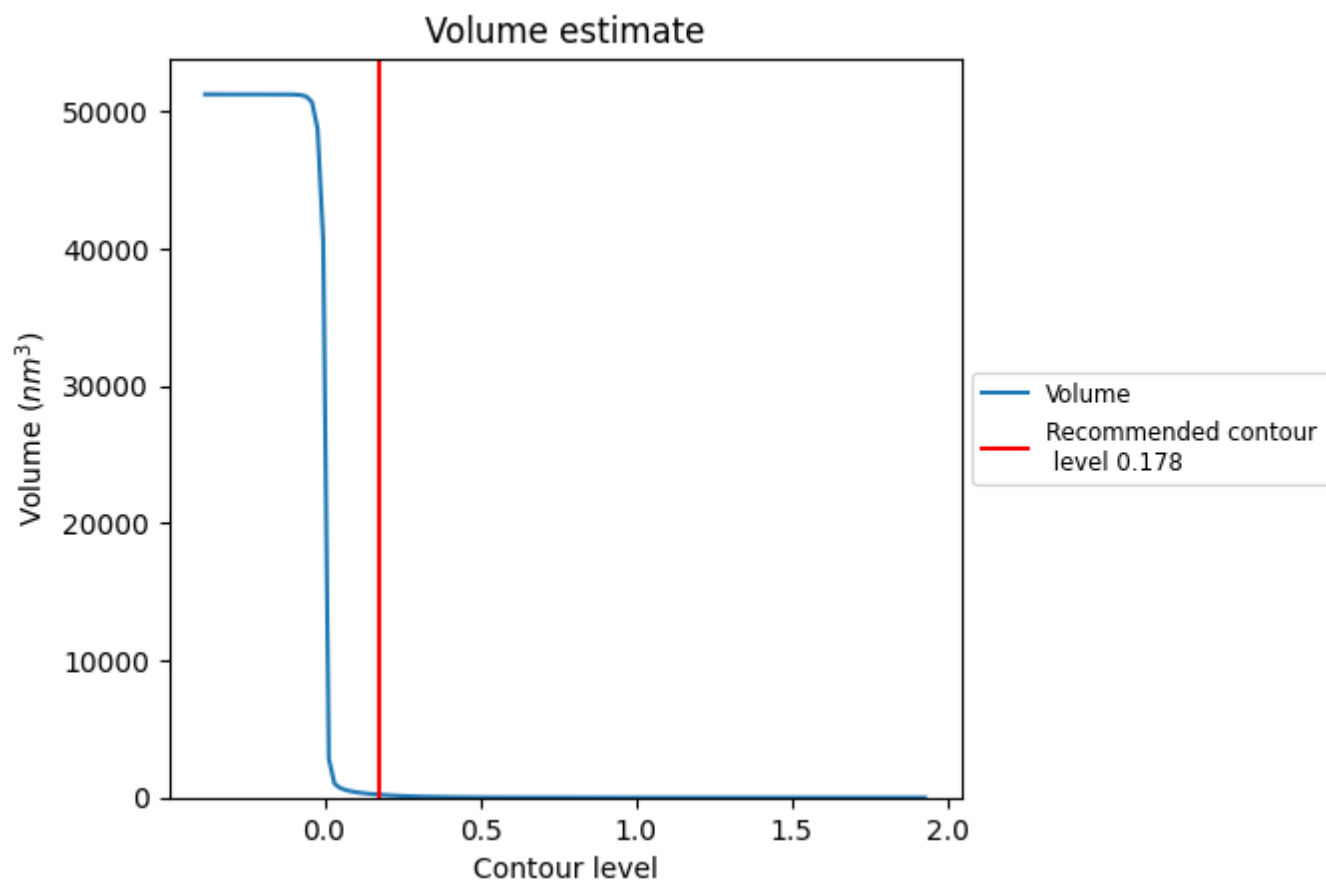
This section contains the results of statistical analysis of the map.

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.

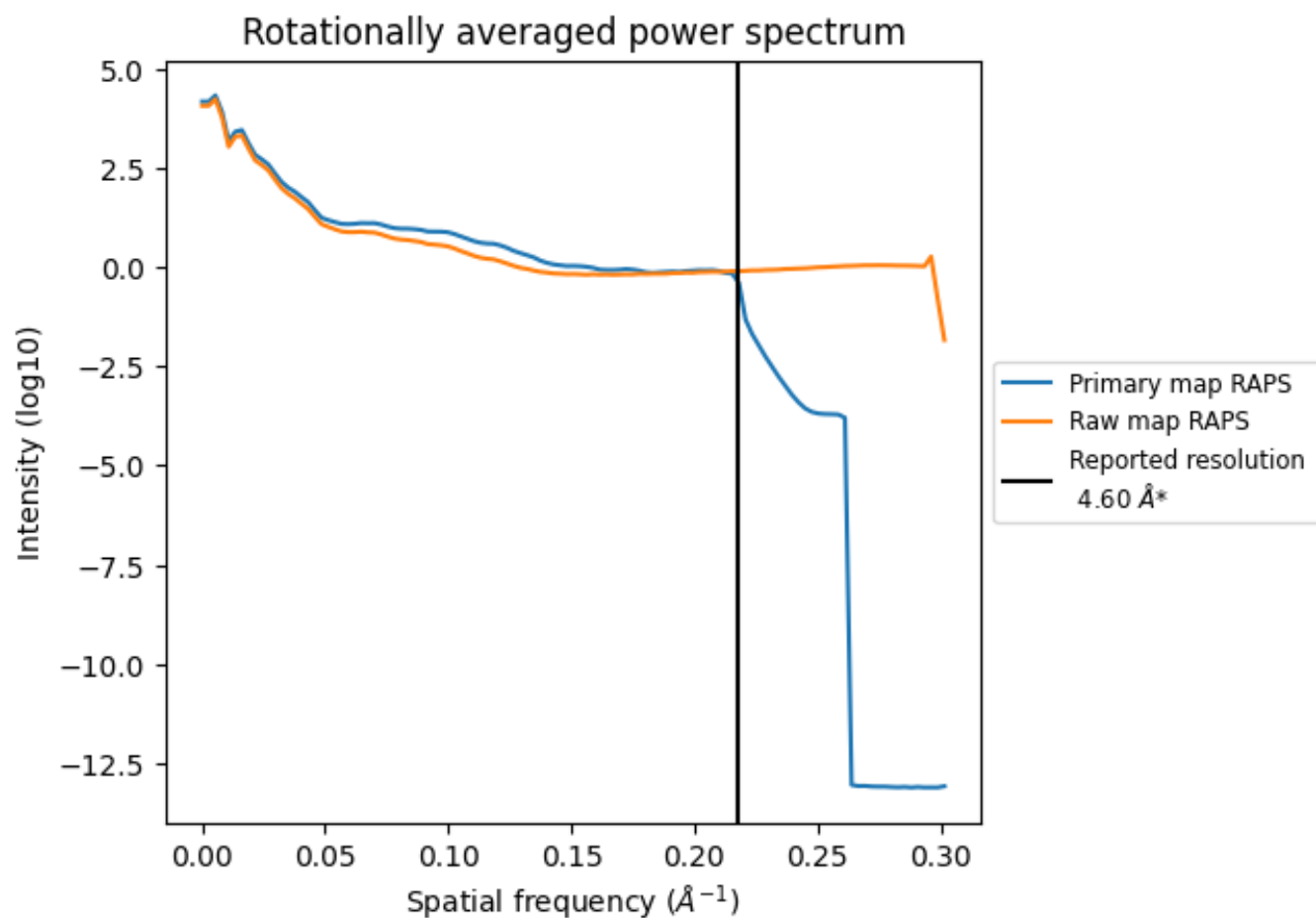
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 209 nm³; this corresponds to an approximate mass of 189 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 ⓘ

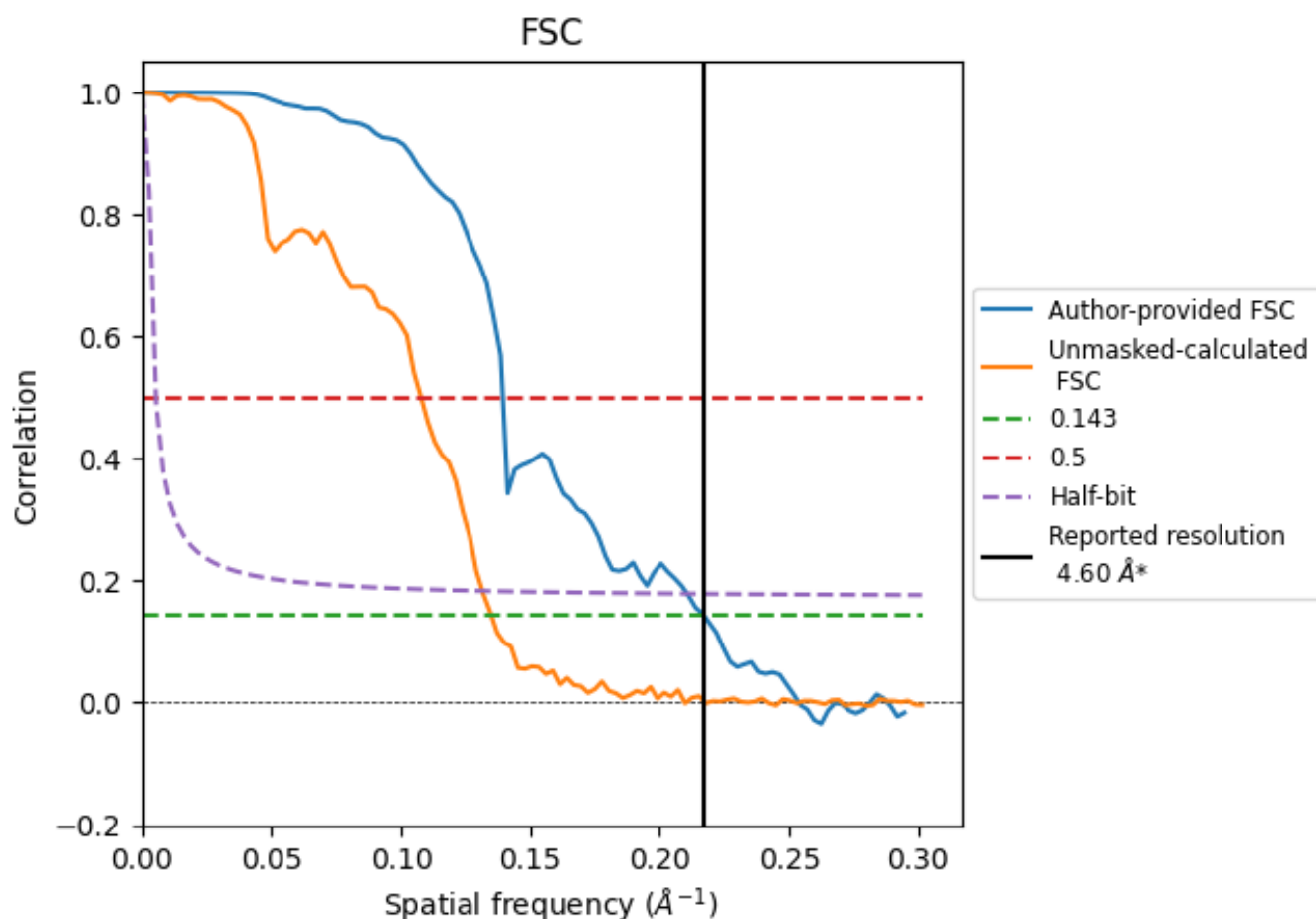


*Reported resolution corresponds to spatial frequency of 0.217 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.217 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

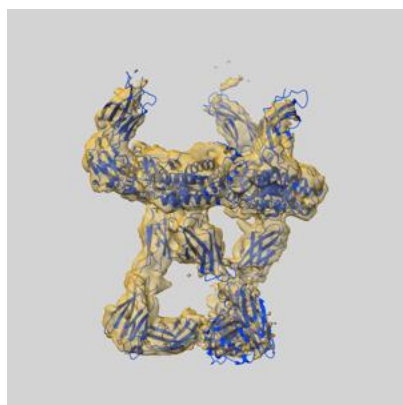
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.60	-	-
Author-provided FSC curve	4.60	7.17	4.75
Unmasked-calculated*	7.40	9.28	7.60

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 7.40 differs from the reported value 4.6 by more than 10 %

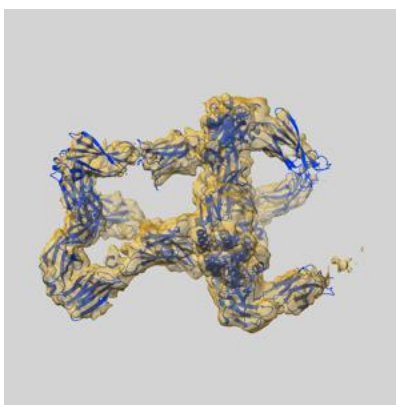
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-15678 and PDB model 8AVC. Per-residue inclusion information can be found in section 3 on page 11.

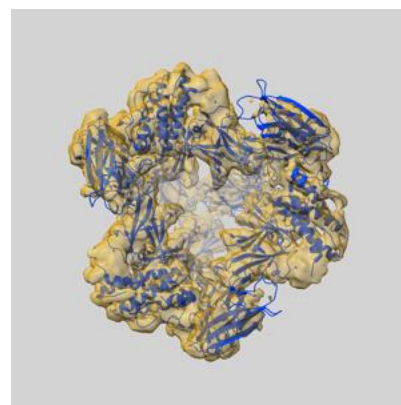
9.1 Map-model overlay [i](#)



X



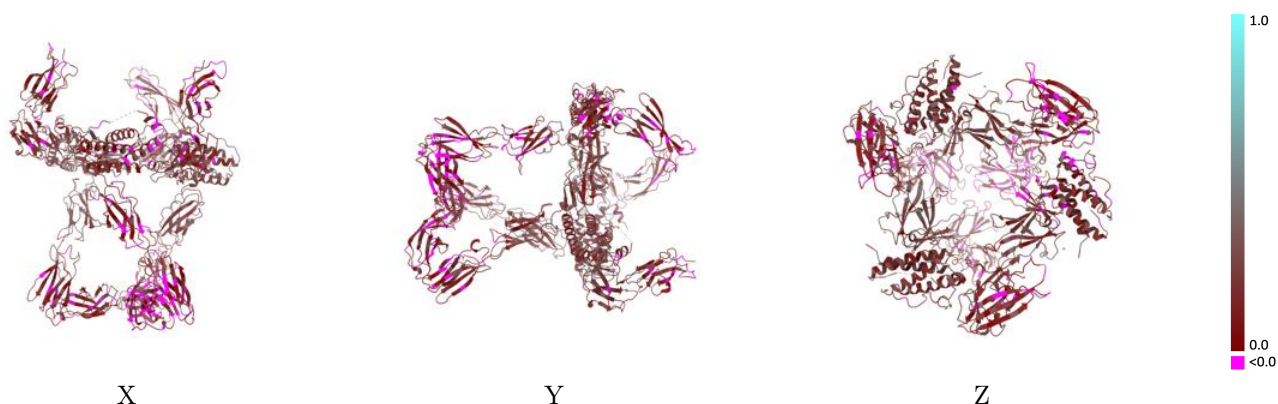
Y



Z

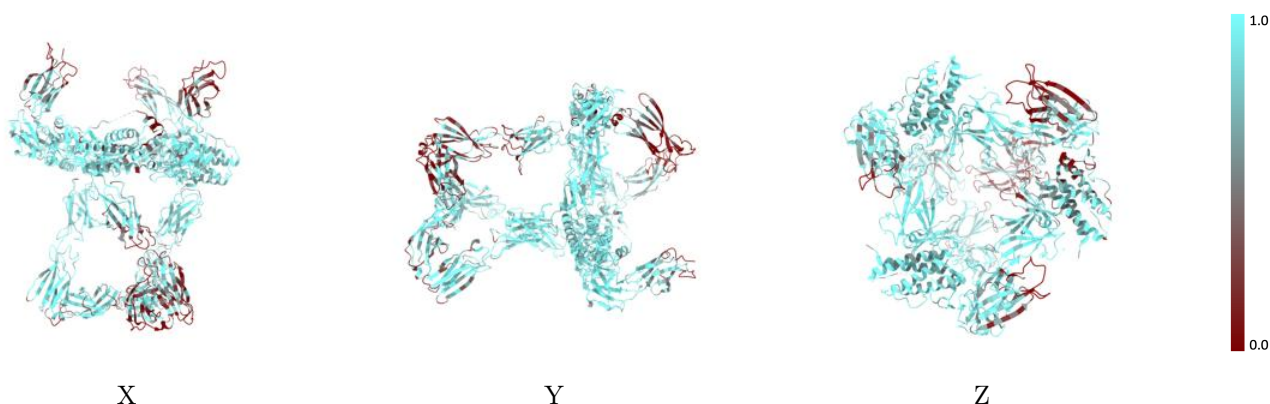
The images above show the 3D surface view of the map at the recommended contour level 0.178 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



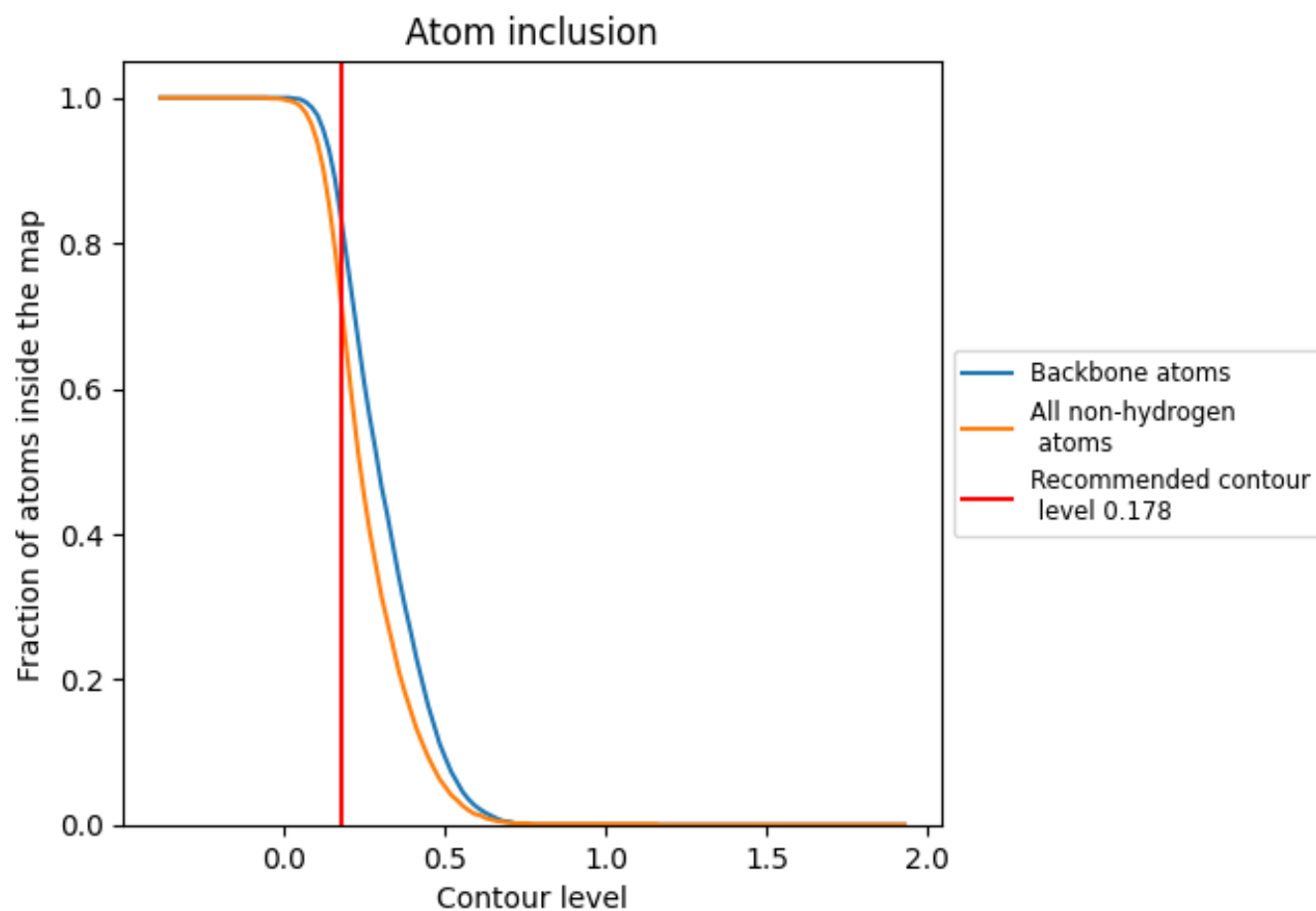
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.178).

9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 72% 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 (0.178) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7190	0.1840
A	0.8280	0.2210
B	0.8120	0.2110
C	0.8190	0.2270
D	0.7060	0.1760
E	0.6780	0.1760
F	0.6010	0.1500

1.0

0.0

<0.0