



Full wwPDB EM Validation Report ⓘ

Jun 27, 2024 – 12:22 AM JST

PDB ID : 8XUS
EMDB ID : EMD-38682
Title : JN.1 Spike Trimer in complex with heparan sulfate
Authors : Yue, C.; Liu, P.
Deposited on : 2024-01-14
Resolution : 3.18 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.dev92
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

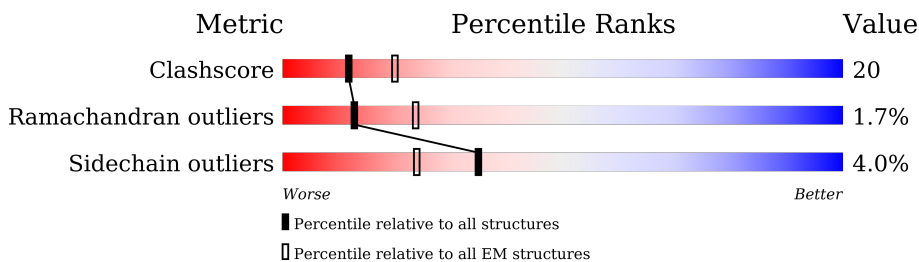
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.18 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	1206	
1	B	1206	
1	C	1206	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 25713 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	C	1063	8314	5319	1383	1574	38	0	0
1	A	1063	8314	5319	1383	1574	38	0	0
1	B	1063	8314	5319	1383	1574	38	0	0

There are 222 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	ALA	-	expression tag	UNP P0DTC2
C	-1	THR	-	expression tag	UNP P0DTC2
C	16	MET	-	insertion	UNP P0DTC2
C	17	PRO	-	insertion	UNP P0DTC2
C	18	LEU	-	insertion	UNP P0DTC2
C	19	PHE	-	insertion	UNP P0DTC2
C	22	ILE	THR	variant	UNP P0DTC2
C	24	THR	ARG	variant	UNP P0DTC2
C	?	-	LEU	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	27	SER	ALA	variant	UNP P0DTC2
C	50	LEU	SER	conflict	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	127	PHE	VAL	conflict	UNP P0DTC2
C	143	ASP	GLY	variant	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	157	SER	PHE	conflict	UNP P0DTC2
C	158	GLY	ARG	variant	UNP P0DTC2
C	?	-	ASN	deletion	UNP P0DTC2
C	212	ILE	LEU	variant	UNP P0DTC2
C	213	GLY	VAL	variant	UNP P0DTC2
C	216	PHE	LEU	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	245	ASN	HIS	conflict	UNP P0DTC2
C	264	ASP	ALA	conflict	UNP P0DTC2
C	332	VAL	ILE	conflict	UNP P0DTC2
C	339	HIS	GLY	variant	UNP P0DTC2
C	356	THR	LYS	conflict	UNP P0DTC2
C	371	PHE	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	376	ALA	THR	variant	UNP P0DTC2
C	403	LYS	ARG	conflict	UNP P0DTC2
C	405	ASN	ASP	variant	UNP P0DTC2
C	408	SER	ARG	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	445	HIS	VAL	conflict	UNP P0DTC2
C	446	SER	GLY	variant	UNP P0DTC2
C	450	ASP	ASN	conflict	UNP P0DTC2
C	452	TRP	LEU	conflict	UNP P0DTC2
C	455	SER	LEU	conflict	UNP P0DTC2
C	460	LYS	ASN	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2
C	481	LYS	ASN	conflict	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	484	LYS	GLU	variant	UNP P0DTC2
C	486	PRO	PHE	variant	UNP P0DTC2
C	498	ARG	GLN	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	505	HIS	TYR	variant	UNP P0DTC2
C	554	LYS	GLU	conflict	UNP P0DTC2
C	570	VAL	ALA	conflict	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	621	SER	PRO	conflict	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	681	ARG	PRO	variant	UNP P0DTC2
C	683	ALA	ARG	conflict	UNP P0DTC2
C	685	ALA	ARG	conflict	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	817	PRO	PHE	conflict	UNP P0DTC2
C	892	PRO	ALA	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	899	PRO	ALA	conflict	UNP P0DTC2
C	939	PHE	SER	conflict	UNP P0DTC2
C	942	PRO	ALA	conflict	UNP P0DTC2
C	954	HIS	GLN	variant	UNP P0DTC2
C	969	LYS	ASN	variant	UNP P0DTC2
C	986	PRO	LYS	variant	UNP P0DTC2
C	987	PRO	VAL	variant	UNP P0DTC2
C	1143	LEU	PRO	conflict	UNP P0DTC2
A	-2	ALA	-	expression tag	UNP P0DTC2
A	-1	THR	-	expression tag	UNP P0DTC2
A	16	MET	-	insertion	UNP P0DTC2
A	17	PRO	-	insertion	UNP P0DTC2
A	18	LEU	-	insertion	UNP P0DTC2
A	19	PHE	-	insertion	UNP P0DTC2
A	22	ILE	THR	variant	UNP P0DTC2
A	24	THR	ARG	variant	UNP P0DTC2
A	?	-	LEU	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	27	SER	ALA	variant	UNP P0DTC2
A	50	LEU	SER	conflict	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	127	PHE	VAL	conflict	UNP P0DTC2
A	143	ASP	GLY	variant	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	157	SER	PHE	conflict	UNP P0DTC2
A	158	GLY	ARG	variant	UNP P0DTC2
A	?	-	ASN	deletion	UNP P0DTC2
A	212	ILE	LEU	variant	UNP P0DTC2
A	213	GLY	VAL	variant	UNP P0DTC2
A	216	PHE	LEU	conflict	UNP P0DTC2
A	245	ASN	HIS	conflict	UNP P0DTC2
A	264	ASP	ALA	conflict	UNP P0DTC2
A	332	VAL	ILE	conflict	UNP P0DTC2
A	339	HIS	GLY	variant	UNP P0DTC2
A	356	THR	LYS	conflict	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	403	LYS	ARG	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	445	HIS	VAL	conflict	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	450	ASP	ASN	conflict	UNP P0DTC2
A	452	TRP	LEU	conflict	UNP P0DTC2
A	455	SER	LEU	conflict	UNP P0DTC2
A	460	LYS	ASN	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	481	LYS	ASN	conflict	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	484	LYS	GLU	variant	UNP P0DTC2
A	486	PRO	PHE	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	554	LYS	GLU	conflict	UNP P0DTC2
A	570	VAL	ALA	conflict	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	621	SER	PRO	conflict	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	ARG	PRO	variant	UNP P0DTC2
A	683	ALA	ARG	conflict	UNP P0DTC2
A	685	ALA	ARG	conflict	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	conflict	UNP P0DTC2
A	892	PRO	ALA	conflict	UNP P0DTC2
A	899	PRO	ALA	conflict	UNP P0DTC2
A	939	PHE	SER	conflict	UNP P0DTC2
A	942	PRO	ALA	conflict	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2
A	986	PRO	LYS	variant	UNP P0DTC2
A	987	PRO	VAL	variant	UNP P0DTC2
A	1143	LEU	PRO	conflict	UNP P0DTC2
B	-2	ALA	-	expression tag	UNP P0DTC2
B	-1	THR	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	16	MET	-	insertion	UNP P0DTC2
B	17	PRO	-	insertion	UNP P0DTC2
B	18	LEU	-	insertion	UNP P0DTC2
B	19	PHE	-	insertion	UNP P0DTC2
B	22	ILE	THR	variant	UNP P0DTC2
B	24	THR	ARG	variant	UNP P0DTC2
B	?	-	LEU	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	27	SER	ALA	variant	UNP P0DTC2
B	50	LEU	SER	conflict	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	127	PHE	VAL	conflict	UNP P0DTC2
B	143	ASP	GLY	variant	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	157	SER	PHE	conflict	UNP P0DTC2
B	158	GLY	ARG	variant	UNP P0DTC2
B	?	-	ASN	deletion	UNP P0DTC2
B	212	ILE	LEU	variant	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	216	PHE	LEU	conflict	UNP P0DTC2
B	245	ASN	HIS	conflict	UNP P0DTC2
B	264	ASP	ALA	conflict	UNP P0DTC2
B	332	VAL	ILE	conflict	UNP P0DTC2
B	339	HIS	GLY	variant	UNP P0DTC2
B	356	THR	LYS	conflict	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	403	LYS	ARG	conflict	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	445	HIS	VAL	conflict	UNP P0DTC2
B	446	SER	GLY	variant	UNP P0DTC2
B	450	ASP	ASN	conflict	UNP P0DTC2
B	452	TRP	LEU	conflict	UNP P0DTC2
B	455	SER	LEU	conflict	UNP P0DTC2
B	460	LYS	ASN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	481	LYS	ASN	conflict	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	484	LYS	GLU	variant	UNP P0DTC2
B	486	PRO	PHE	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	554	LYS	GLU	conflict	UNP P0DTC2
B	570	VAL	ALA	conflict	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	621	SER	PRO	conflict	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	ARG	PRO	variant	UNP P0DTC2
B	683	ALA	ARG	conflict	UNP P0DTC2
B	685	ALA	ARG	conflict	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	conflict	UNP P0DTC2
B	892	PRO	ALA	conflict	UNP P0DTC2
B	899	PRO	ALA	conflict	UNP P0DTC2
B	939	PHE	SER	conflict	UNP P0DTC2
B	942	PRO	ALA	conflict	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	986	PRO	LYS	variant	UNP P0DTC2
B	987	PRO	VAL	variant	UNP P0DTC2
B	1143	LEU	PRO	conflict	UNP P0DTC2

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



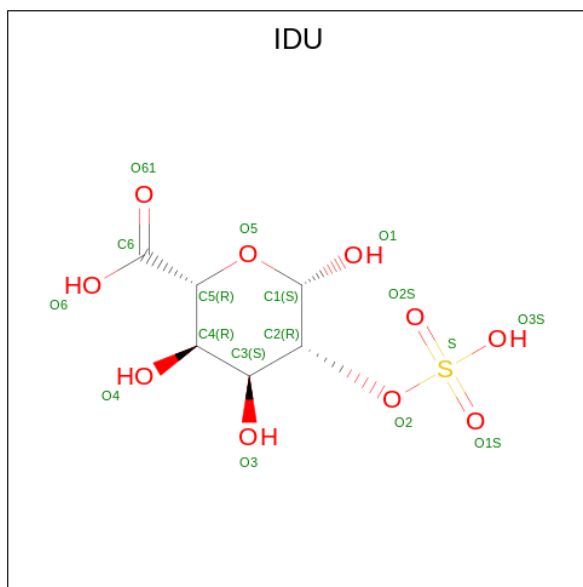
Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	
2	C	1	Total	C	N	O	0
			14	8	1	5	

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Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
2	A	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0
2	B	1	Total 14	C 8	N 1	O 5	0

- Molecule 3 is 2-O-sulfo-beta-L-altropyranuronic acid (three-letter code: IDU) (formula: C₆H₁₀O₁₀S) (labeled as "Ligand of Interest" by depositor).

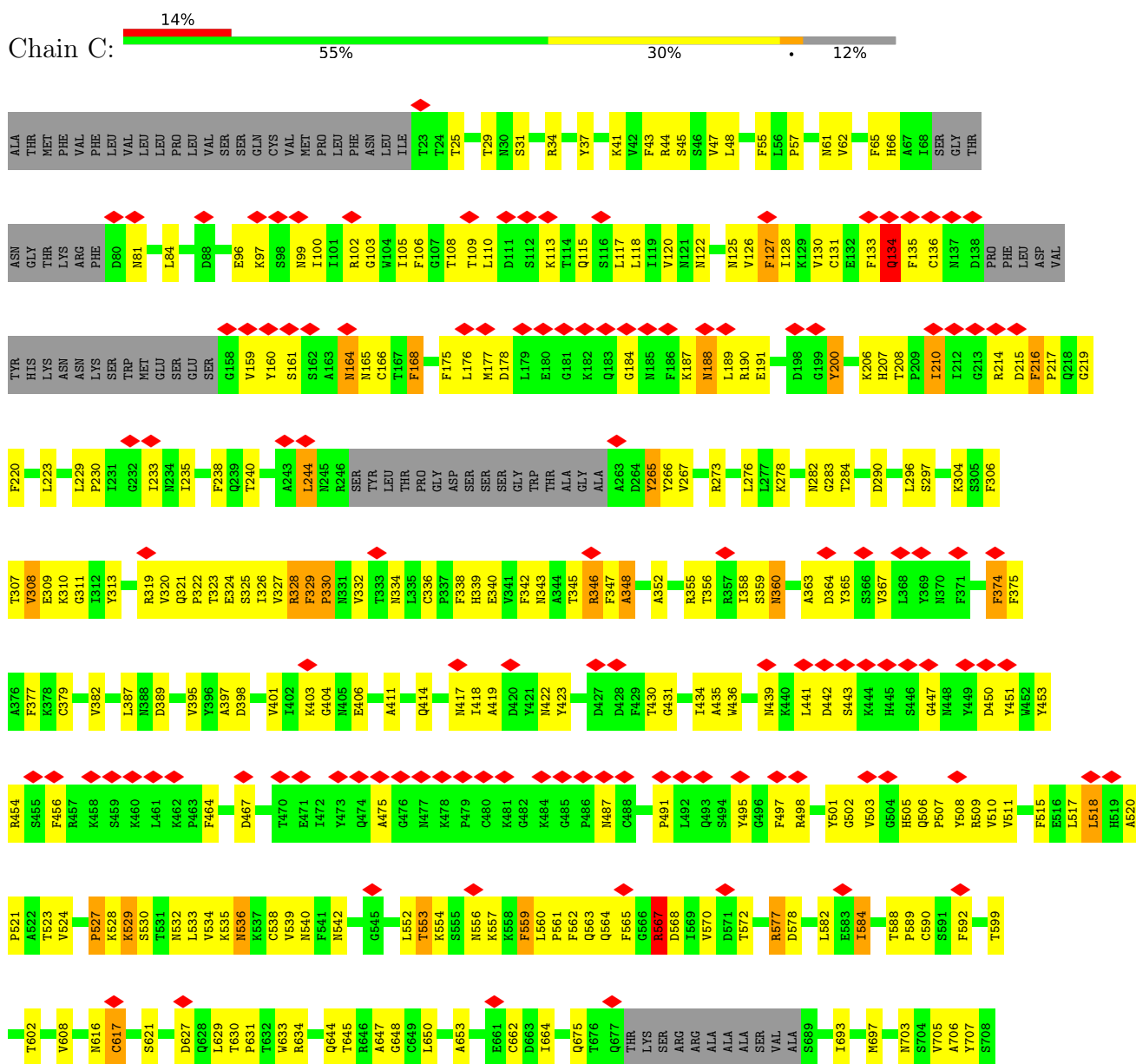


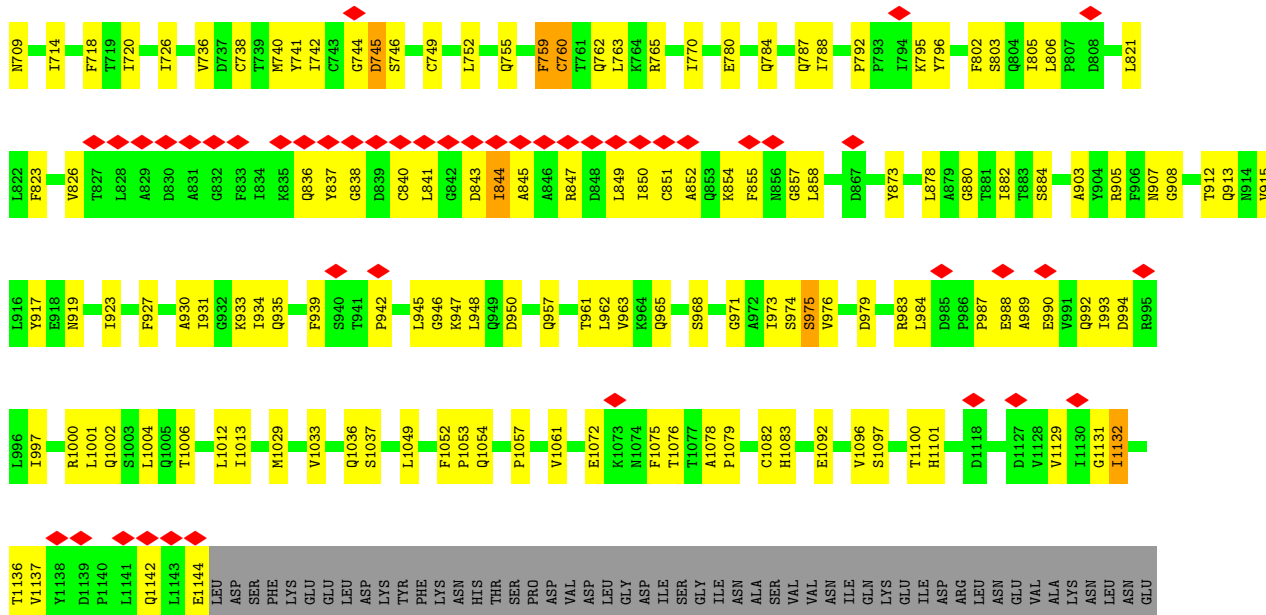
Mol	Chain	Residues	Atoms				AltConf
			Total	C	O	S	
3	B	1	15	6	8	1	0

3 Residue-property plots

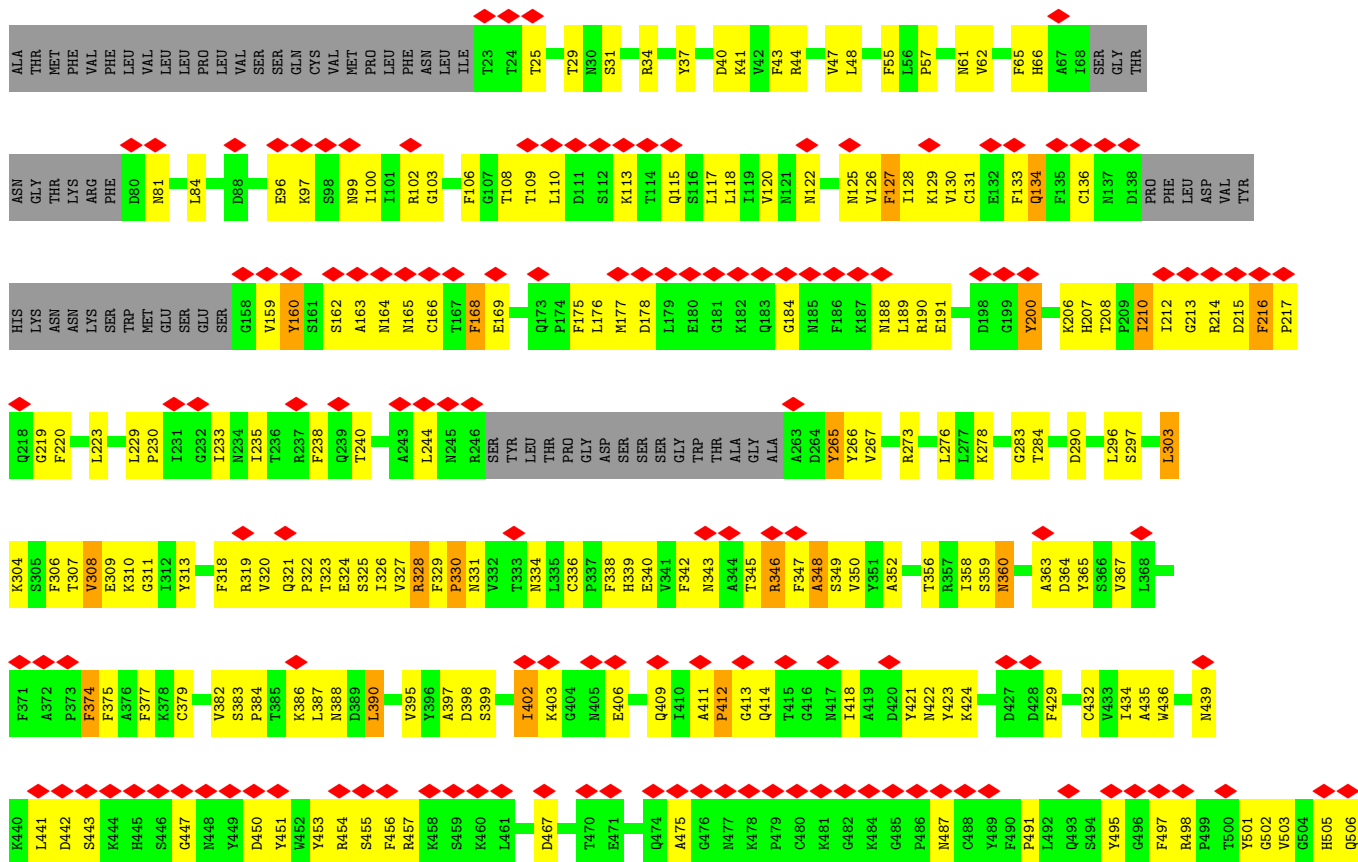
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: Spike glycoprotein





• Molecule 1: Spike glycoprotein



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	301841	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.695	Depositor
Minimum map value	-1.681	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.085	Depositor
Recommended contour level	0.35	Depositor
Map size (\AA)	310.80002, 310.80002, 310.80002	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.036, 1.036, 1.036	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.44	0/8514	0.60	0/11588
1	B	0.45	0/8514	0.60	0/11588
1	C	0.44	0/8514	0.60	0/11588
All	All	0.44	0/25542	0.60	0/34764

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
1	C	0	4
All	All	0	9

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	328	ARG	Sidechain
1	A	346	ARG	Sidechain
1	B	319	ARG	Sidechain
1	B	328	ARG	Sidechain
1	B	346	ARG	Sidechain
1	C	328	ARG	Sidechain
1	C	346	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	C	567	ARG	Sidechain
1	C	577	ARG	Sidechain

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8314	0	8094	383	0
1	B	8314	0	8094	363	0
1	C	8314	0	8094	396	0
2	A	252	0	234	6	0
2	B	252	0	234	5	0
2	C	252	0	234	10	0
3	B	15	0	4	1	0
All	All	25713	0	24988	993	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 (993) 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:C:430:THR:CG2	1:C:517:LEU:HD12	1.63	1.26
1:C:503:VAL:CG2	1:A:505:HIS:CE1	2.20	1.25
1:C:759:PHE:HZ	1:A:1002:GLN:CG	1.49	1.24
1:A:503:VAL:CG2	1:B:505:HIS:CE1	2.21	1.23
1:A:759:PHE:HZ	1:B:1002:GLN:CG	1.56	1.17
1:C:1002:GLN:HG3	1:B:759:PHE:HZ	1.07	1.16
1:C:759:PHE:CZ	1:A:1002:GLN:HG3	1.80	1.16
1:C:503:VAL:HG21	1:A:505:HIS:CE1	1.78	1.15
1:C:1002:GLN:HG3	1:B:759:PHE:CZ	1.81	1.15
1:C:430:THR:HG21	1:C:517:LEU:HD12	1.25	1.15
1:A:759:PHE:CZ	1:B:1002:GLN:HG3	1.82	1.14
1:C:505:HIS:CE1	1:B:503:VAL:CG2	2.29	1.14
1:C:503:VAL:HG21	1:A:505:HIS:HE1	1.00	1.13
1:C:505:HIS:HE1	1:B:503:VAL:HG21	1.06	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:759:PHE:HZ	1:B:1002:GLN:HG3	1.05	1.13
1:A:503:VAL:HG21	1:B:505:HIS:CE1	1.78	1.13
1:A:841:LEU:HB2	1:B:553:THR:HG21	1.22	1.12
1:C:1002:GLN:CG	1:B:759:PHE:HZ	1.63	1.12
1:C:326:ILE:HD11	1:C:534:VAL:HG23	1.19	1.11
1:C:759:PHE:HZ	1:A:1002:GLN:HG3	1.02	1.11
1:C:755:GLN:OE1	1:A:971:GLY:HA2	1.51	1.10
1:C:553:THR:HG21	1:B:841:LEU:HB2	1.13	1.09
1:C:759:PHE:CZ	1:A:1002:GLN:CG	2.33	1.09
1:A:503:VAL:HG21	1:B:505:HIS:HE1	0.97	1.09
1:C:971:GLY:HA2	1:B:755:GLN:OE1	1.51	1.08
1:A:759:PHE:CZ	1:B:1002:GLN:CG	2.38	1.07
1:A:755:GLN:OE1	1:B:971:GLY:HA2	1.54	1.07
1:C:983:ARG:CD	1:A:517:LEU:HD13	1.85	1.06
1:C:505:HIS:HE1	1:B:503:VAL:CG2	1.65	1.06
1:A:503:VAL:CG2	1:B:505:HIS:HE1	1.64	1.06
1:A:841:LEU:HB2	1:B:553:THR:CG2	1.88	1.03
1:C:505:HIS:CE1	1:B:503:VAL:HG21	1.92	1.03
1:A:903:ALA:HB1	1:A:913:GLN:HB3	1.42	1.02
1:C:1083:HIS:ND1	1:C:1137:VAL:HG22	1.74	1.02
1:B:1083:HIS:ND1	1:B:1137:VAL:HG22	1.74	1.02
1:C:553:THR:CG2	1:B:841:LEU:HB2	1.90	1.02
1:C:1002:GLN:CG	1:B:759:PHE:CZ	2.40	1.02
1:A:1083:HIS:ND1	1:A:1137:VAL:HG22	1.74	1.01
1:A:983:ARG:CD	1:B:517:LEU:HD13	1.90	1.00
1:C:1083:HIS:CG	1:C:1137:VAL:HG22	1.97	1.00
1:B:1083:HIS:CG	1:B:1137:VAL:HG22	1.97	0.99
1:C:503:VAL:CG2	1:A:505:HIS:HE1	1.64	0.99
1:A:1083:HIS:CG	1:A:1137:VAL:HG22	1.97	0.98
1:C:326:ILE:HD13	1:C:533:LEU:HA	1.47	0.95
1:C:983:ARG:HD3	1:A:517:LEU:HD13	1.45	0.95
1:C:841:LEU:HB2	1:A:553:THR:HG21	1.48	0.94
1:A:983:ARG:HD3	1:B:517:LEU:HD13	1.50	0.93
1:C:382:VAL:HG11	1:C:387:LEU:HD13	1.50	0.92
1:A:503:VAL:HG23	1:B:505:HIS:CE1	2.06	0.90
1:C:505:HIS:CE1	1:B:503:VAL:HG23	2.04	0.89
1:C:62:VAL:HG11	1:C:216:PHE:HE2	1.36	0.88
1:A:62:VAL:HG11	1:A:216:PHE:HE2	1.36	0.88
1:B:300:LYS:HG2	1:B:308:VAL:HG23	1.55	0.87
1:B:559:PHE:HB2	1:B:584:ILE:HD11	1.56	0.87
1:C:503:VAL:HG23	1:A:505:HIS:CE1	2.07	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:PRO:HG2	1:B:540:ASN:HD21	1.40	0.86
1:C:326:ILE:HD11	1:C:534:VAL:CG2	2.06	0.86
1:C:329:PHE:HB3	1:C:330:PRO:CD	2.06	0.86
1:A:329:PHE:HB3	1:A:330:PRO:HD2	1.58	0.85
1:C:553:THR:HG21	1:B:841:LEU:CB	2.02	0.85
1:A:841:LEU:CB	1:B:553:THR:HG21	2.04	0.85
1:A:322:PRO:HG2	1:A:540:ASN:HD21	1.40	0.85
1:C:322:PRO:HG2	1:C:540:ASN:HD21	1.40	0.85
1:A:326:ILE:HD12	1:A:532:ASN:O	1.76	0.84
1:B:326:ILE:CG2	1:B:532:ASN:O	2.26	0.84
1:B:326:ILE:HD12	1:B:532:ASN:O	1.76	0.84
1:B:557:LYS:HD3	1:B:559:PHE:HE1	1.43	0.83
1:C:1082:CYS:HB2	1:C:1132:ILE:HG13	1.58	0.83
1:B:329:PHE:HB3	1:B:330:PRO:HD2	1.61	0.83
1:A:326:ILE:CG2	1:A:532:ASN:O	2.26	0.83
1:C:759:PHE:CZ	1:A:1002:GLN:HG2	2.14	0.82
1:B:326:ILE:HG23	1:B:532:ASN:O	1.80	0.82
1:C:983:ARG:HD2	1:A:517:LEU:HD13	1.60	0.81
1:A:983:ARG:HD2	1:B:517:LEU:HD13	1.63	0.81
1:A:326:ILE:HG23	1:A:532:ASN:O	1.80	0.81
1:B:325:SER:HA	1:B:539:VAL:HG13	1.63	0.81
1:B:1082:CYS:HB2	1:B:1132:ILE:HG13	1.61	0.81
1:A:325:SER:HA	1:A:539:VAL:HG13	1.63	0.81
1:C:326:ILE:HD13	1:C:533:LEU:CA	2.11	0.80
1:C:325:SER:HA	1:C:539:VAL:HG13	1.63	0.80
1:A:1082:CYS:HB2	1:A:1132:ILE:HG13	1.62	0.80
1:C:838:GLY:HA3	2:A:1309:NAG:H81	1.62	0.79
1:B:1083:HIS:ND1	1:B:1137:VAL:CG2	2.46	0.79
1:C:328:ARG:HG3	1:C:530:SER:HB2	1.66	0.78
1:A:328:ARG:HG3	1:A:530:SER:HB2	1.65	0.78
1:B:328:ARG:HG3	1:B:530:SER:HB2	1.66	0.78
1:C:1083:HIS:ND1	1:C:1137:VAL:CG2	2.46	0.77
1:A:1083:HIS:ND1	1:A:1137:VAL:CG2	2.46	0.77
1:C:178:ASP:HA	1:C:188:ASN:HD21	1.49	0.77
1:A:503:VAL:HG23	1:B:505:HIS:NE2	1.99	0.77
1:C:329:PHE:HB3	1:C:330:PRO:HD3	1.66	0.76
1:A:322:PRO:HG2	1:A:540:ASN:ND2	2.01	0.76
1:C:310:LYS:HG2	1:C:664:ILE:HD11	1.68	0.76
1:C:802:PHE:H	2:C:1313:NAG:H82	1.49	0.76
1:C:322:PRO:HG2	1:C:540:ASN:ND2	2.01	0.76
1:A:310:LYS:HG2	1:A:664:ILE:HD11	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:417:ASN:O	1:C:422:ASN:ND2	2.20	0.75
1:B:310:LYS:HG2	1:B:664:ILE:HD11	1.68	0.75
1:B:322:PRO:HG2	1:B:540:ASN:ND2	2.01	0.75
1:C:788:ILE:HD11	1:A:699:LEU:HB2	1.67	0.75
1:A:62:VAL:HG11	1:A:216:PHE:CE2	2.21	0.75
1:C:503:VAL:HG23	1:A:505:HIS:NE2	2.02	0.75
1:C:707:TYR:HD1	1:B:792:PRO:HG3	1.52	0.75
1:C:62:VAL:HG11	1:C:216:PHE:CE2	2.21	0.75
2:C:1311:NAG:H62	1:B:796:TYR:CD2	2.22	0.75
1:A:37:TYR:HA	1:A:223:LEU:HB2	1.69	0.74
1:A:759:PHE:CZ	1:B:1002:GLN:HG2	2.22	0.73
1:A:43:PHE:HA	1:B:563:GLN:OE1	1.88	0.73
1:A:557:LYS:HD3	1:A:559:PHE:HE1	1.52	0.73
1:C:983:ARG:HD2	1:A:517:LEU:CD1	2.18	0.73
1:C:430:THR:HG22	1:C:517:LEU:HD12	1.64	0.73
1:C:365:TYR:CE2	1:C:387:LEU:HG	2.24	0.73
1:B:577:ARG:HG2	1:B:584:ILE:HG12	1.71	0.73
1:C:430:THR:CG2	1:C:517:LEU:CD1	2.57	0.72
1:C:330:PRO:HG2	1:C:332:VAL:HG22	1.71	0.72
1:C:503:VAL:CB	1:A:505:HIS:CE1	2.72	0.72
1:C:983:ARG:CD	1:A:517:LEU:CD1	2.67	0.72
1:B:65:PHE:HB2	1:B:265:TYR:HB3	1.72	0.71
1:A:65:PHE:HB2	1:A:265:TYR:HB3	1.72	0.71
1:C:65:PHE:HB2	1:C:265:TYR:HB3	1.72	0.71
1:C:971:GLY:CA	1:B:755:GLN:OE1	2.36	0.71
1:B:44:ARG:HB3	1:B:47:VAL:HG11	1.72	0.71
1:C:1083:HIS:CE1	1:C:1137:VAL:HG22	2.26	0.71
1:A:852:ALA:HB1	1:B:570:VAL:HG22	1.72	0.71
1:A:106:PHE:HD1	1:A:235:ILE:HD13	1.56	0.71
1:B:1083:HIS:CE1	1:B:1137:VAL:HG22	2.26	0.71
1:A:503:VAL:CG2	1:B:505:HIS:NE2	2.53	0.71
1:C:44:ARG:HB3	1:C:47:VAL:HG11	1.72	0.70
1:A:1083:HIS:CE1	1:A:1137:VAL:HG22	2.26	0.70
1:B:96:GLU:OE1	1:B:100:ILE:N	2.24	0.70
1:C:43:PHE:HA	1:A:563:GLN:OE1	1.91	0.70
1:C:96:GLU:OE1	1:C:100:ILE:N	2.24	0.70
1:A:44:ARG:HB3	1:A:47:VAL:HG11	1.72	0.70
1:C:503:VAL:CG2	1:A:505:HIS:NE2	2.53	0.70
1:C:987:PRO:HD2	1:C:988:GLU:OE2	1.92	0.70
1:A:96:GLU:OE1	1:A:100:ILE:N	2.24	0.70
1:B:106:PHE:HD1	1:B:235:ILE:HD13	1.56	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:903:ALA:HB1	1:C:913:GLN:HB3	1.72	0.70
1:C:106:PHE:HD1	1:C:235:ILE:HD13	1.56	0.69
1:C:326:ILE:CD1	1:C:534:VAL:HG23	2.11	0.69
1:C:505:HIS:NE2	1:B:503:VAL:HG23	2.08	0.69
1:A:987:PRO:HD2	1:A:988:GLU:OE2	1.92	0.69
1:B:327:VAL:HG11	1:B:528:LYS:CE	2.23	0.69
1:B:770:ILE:HD11	1:B:1012:LEU:HD23	1.75	0.69
1:B:903:ALA:HB1	1:B:913:GLN:HB3	1.75	0.69
1:C:229:LEU:HD23	1:C:229:LEU:H	1.58	0.68
1:C:770:ILE:HD11	1:C:1012:LEU:HD23	1.75	0.68
1:C:454:ARG:NH1	1:C:467:ASP:O	2.27	0.68
1:A:43:PHE:CA	1:B:563:GLN:OE1	2.42	0.68
1:B:128:ILE:HD11	1:B:175:PHE:HZ	1.59	0.68
1:C:456:PHE:HB2	1:C:491:PRO:HA	1.75	0.68
1:C:1129:VAL:HG13	1:B:917:TYR:HB3	1.76	0.68
1:C:326:ILE:CD1	1:C:534:VAL:N	2.57	0.68
1:A:454:ARG:NH1	1:A:467:ASP:O	2.27	0.68
1:A:456:PHE:HB2	1:A:491:PRO:HA	1.75	0.68
1:A:770:ILE:HD11	1:A:1012:LEU:HD23	1.75	0.68
1:A:983:ARG:HD2	1:B:517:LEU:CD1	2.24	0.68
1:C:326:ILE:CD1	1:C:533:LEU:C	2.62	0.67
1:A:327:VAL:HG11	1:A:528:LYS:CE	2.23	0.67
1:B:456:PHE:HB2	1:B:491:PRO:HA	1.75	0.67
1:A:229:LEU:HD23	1:A:229:LEU:H	1.58	0.67
1:C:327:VAL:HG11	1:C:528:LYS:CE	2.23	0.67
1:A:350:VAL:HG21	1:A:418:ILE:HG23	1.74	0.67
1:B:129:LYS:HD2	1:B:160:TYR:CE2	2.30	0.67
1:B:454:ARG:NH1	1:B:467:ASP:O	2.27	0.67
1:C:852:ALA:HB1	1:A:570:VAL:HG22	1.77	0.66
1:A:852:ALA:HB1	1:B:570:VAL:CG2	2.26	0.66
1:B:115:GLN:HE21	1:B:233:ILE:HB	1.61	0.65
1:B:327:VAL:HG11	1:B:528:LYS:HE3	1.78	0.65
1:B:375:PHE:CD1	1:B:508:TYR:OH	2.41	0.65
1:A:379:CYS:HA	1:A:432:CYS:HA	1.79	0.65
1:A:327:VAL:HG11	1:A:528:LYS:HE3	1.78	0.65
1:C:442:ASP:OD1	1:C:509:ARG:NH2	2.30	0.65
1:A:375:PHE:CD1	1:A:508:TYR:OH	2.41	0.65
1:C:115:GLN:HE21	1:C:233:ILE:HB	1.61	0.65
1:C:327:VAL:HG11	1:C:528:LYS:HE3	1.78	0.65
1:C:128:ILE:HD11	1:C:175:PHE:HZ	1.62	0.65
1:B:442:ASP:OD1	1:B:509:ARG:NH2	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:503:VAL:HB	1:A:505:HIS:CE1	2.32	0.64
1:A:128:ILE:HD11	1:A:175:PHE:HZ	1.62	0.64
1:A:796:TYR:CD2	2:B:1311:NAG:H62	2.33	0.64
1:C:206:LYS:HB3	1:C:223:LEU:HD13	1.80	0.64
1:A:442:ASP:OD1	1:A:509:ARG:NH2	2.30	0.64
1:C:406:GLU:OE1	1:C:406:GLU:N	2.30	0.64
1:C:557:LYS:HG3	1:B:844:ILE:O	1.98	0.64
1:C:917:TYR:HB3	1:A:1129:VAL:HG13	1.80	0.64
1:C:326:ILE:CG2	1:C:532:ASN:O	2.45	0.64
1:A:983:ARG:CD	1:B:517:LEU:CD1	2.74	0.64
1:A:200:TYR:HA	1:A:230:PRO:HA	1.80	0.64
1:C:759:PHE:CE2	1:A:1002:GLN:HG2	2.33	0.64
1:C:1079:PRO:HB3	1:B:917:TYR:CE1	2.33	0.64
1:A:118:LEU:HD11	1:A:159:VAL:HB	1.80	0.64
1:C:311:GLY:HA2	1:C:664:ILE:HD12	1.81	0.63
1:A:303:LEU:HD21	1:A:308:VAL:HG12	1.79	0.63
1:C:200:TYR:HA	1:C:230:PRO:HA	1.80	0.63
1:C:852:ALA:HB1	1:A:570:VAL:CG2	2.28	0.63
1:A:503:VAL:CB	1:B:505:HIS:CE1	2.80	0.63
1:C:787:GLN:OE1	1:A:703:ASN:ND2	2.29	0.63
1:B:1076:THR:HB	1:B:1097:SER:HB3	1.81	0.63
1:C:326:ILE:HD11	1:C:533:LEU:C	2.18	0.63
1:B:311:GLY:HA2	1:B:664:ILE:HD12	1.81	0.63
1:A:792:PRO:HG3	1:B:707:TYR:HD1	1.64	0.63
1:C:556:ASN:H	1:B:844:ILE:HG23	1.64	0.63
1:A:1076:THR:HB	1:A:1097:SER:HB3	1.81	0.63
1:B:39:PRO:HG3	1:B:55:PHE:HZ	1.64	0.63
1:C:788:ILE:CD1	1:A:699:LEU:HB2	2.28	0.63
1:B:134:GLN:HG2	1:B:162:SER:HB2	1.80	0.63
1:B:300:LYS:HG2	1:B:308:VAL:CG2	2.29	0.62
1:C:796:TYR:CD2	2:A:1311:NAG:H62	2.34	0.62
1:A:398:ASP:HB2	1:A:512:VAL:HG12	1.81	0.62
1:C:821:LEU:HD11	1:C:935:GLN:HG3	1.82	0.62
1:C:136:CYS:HB2	1:C:159:VAL:HA	1.81	0.62
1:B:1100:THR:HB	2:B:1315:NAG:HN2	1.64	0.62
1:A:311:GLY:HA2	1:A:664:ILE:HD12	1.81	0.62
1:C:189:LEU:HD23	1:C:210:ILE:HA	1.81	0.62
1:A:441:LEU:HD22	1:A:509:ARG:HH12	1.65	0.62
1:A:1100:THR:HB	2:A:1315:NAG:HN2	1.65	0.62
1:C:917:TYR:CE1	1:A:1079:PRO:HB3	2.35	0.62
1:B:326:ILE:HG21	1:B:532:ASN:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1002:GLN:HG2	1:B:759:PHE:CE2	2.34	0.61
1:B:1053:PRO:O	1:B:1054:GLN:NE2	2.28	0.61
1:C:276:LEU:HD11	1:C:304:LYS:HA	1.81	0.61
1:C:662:CYS:HB2	1:C:697:MET:HE3	1.82	0.61
1:A:276:LEU:HD11	1:A:304:LYS:HA	1.82	0.61
1:B:441:LEU:HD22	1:B:509:ARG:HH12	1.65	0.61
1:C:1013:ILE:HD13	1:B:1012:LEU:HB3	1.83	0.61
1:B:323:THR:OG1	1:B:324:GLU:N	2.34	0.61
1:C:441:LEU:HD22	1:C:509:ARG:HH12	1.65	0.60
1:C:1002:GLN:HG2	1:B:759:PHE:CZ	2.32	0.60
1:A:759:PHE:CE2	1:B:1002:GLN:HG2	2.35	0.60
1:B:662:CYS:HB2	1:B:697:MET:HE3	1.82	0.60
1:C:326:ILE:HG23	1:C:532:ASN:O	2.01	0.60
1:C:375:PHE:CD1	1:C:508:TYR:OH	2.42	0.60
1:C:557:LYS:HD3	1:C:559:PHE:HE1	1.65	0.60
1:C:645:THR:HG23	1:C:647:ALA:H	1.65	0.60
1:C:323:THR:OG1	1:C:324:GLU:N	2.34	0.60
1:C:564:GLN:O	1:C:565:PHE:C	2.40	0.60
1:A:34:ARG:NH1	1:A:191:GLU:OE2	2.34	0.60
1:A:323:THR:OG1	1:A:324:GLU:N	2.34	0.60
1:A:356:THR:OG1	1:A:397:ALA:HB3	2.02	0.60
1:C:431:GLY:HA2	1:C:515:PHE:CD2	2.36	0.60
1:A:513:LEU:HG	1:A:515:PHE:CE1	2.37	0.60
1:C:1002:GLN:CG	1:B:759:PHE:CE2	2.84	0.60
1:A:318:PHE:CE1	1:A:593:GLY:HA3	2.37	0.60
1:B:350:VAL:HG21	1:B:418:ILE:HG23	1.84	0.60
1:C:282:ASN:OD1	1:A:558:LYS:HG3	2.02	0.59
1:B:412:PRO:HG3	1:B:429:PHE:HB3	1.84	0.59
1:B:564:GLN:O	1:B:565:PHE:C	2.40	0.59
1:A:567:ARG:HD3	1:A:571:ASP:HA	1.84	0.59
1:A:213:GLY:O	1:A:214:ARG:HG3	2.02	0.59
1:A:564:GLN:O	1:A:565:PHE:C	2.40	0.59
1:C:627:ASP:HA	1:C:634:ARG:HH22	1.68	0.59
1:A:513:LEU:HG	1:A:515:PHE:HE1	1.67	0.59
1:A:206:LYS:HB3	1:A:223:LEU:HD13	1.83	0.59
1:C:359:SER:HA	1:C:524:VAL:HG23	1.85	0.59
1:C:841:LEU:CB	1:A:553:THR:HG21	2.29	0.59
1:C:1053:PRO:O	1:C:1054:GLN:NE2	2.28	0.59
1:B:131:CYS:HB3	1:B:133:PHE:CE2	2.38	0.59
1:B:34:ARG:NH1	1:B:191:GLU:OE2	2.34	0.59
1:B:326:ILE:H	1:B:539:VAL:HG11	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:THR:OG1	1:B:397:ALA:HB3	2.03	0.58
1:B:213:GLY:O	1:B:214:ARG:HG3	2.02	0.58
1:B:334:ASN:ND2	1:B:360:ASN:O	2.35	0.58
1:B:567:ARG:HD3	1:B:571:ASP:HA	1.84	0.58
1:A:326:ILE:H	1:A:539:VAL:HG11	1.68	0.58
1:A:382:VAL:HG23	1:A:517:LEU:HD11	1.84	0.58
1:A:662:CYS:HB2	1:A:697:MET:HE3	1.85	0.58
1:C:334:ASN:ND2	1:C:360:ASN:O	2.35	0.58
1:A:1053:PRO:O	1:A:1054:GLN:NE2	2.28	0.58
1:C:557:LYS:HB2	1:C:584:ILE:HG21	1.84	0.58
1:C:792:PRO:HG3	1:A:707:TYR:HD1	1.68	0.58
1:C:844:ILE:O	1:A:557:LYS:HG3	2.03	0.58
1:C:976:VAL:HG12	1:C:979:ASP:H	1.69	0.58
1:A:326:ILE:HG21	1:A:532:ASN:O	2.00	0.58
1:A:976:VAL:HG12	1:A:979:ASP:H	1.69	0.58
1:B:81:ASN:ND2	1:B:240:THR:O	2.25	0.58
1:B:421:TYR:CD1	1:B:457:ARG:HB3	2.39	0.58
1:C:326:ILE:H	1:C:539:VAL:HG11	1.68	0.58
1:B:62:VAL:HG11	1:B:216:PHE:HE2	1.68	0.58
1:C:742:ILE:HG23	1:C:1000:ARG:HB2	1.86	0.58
1:A:1142:GLN:N	1:A:1142:GLN:OE1	2.36	0.58
1:B:411:ALA:HB3	1:B:414:GLN:HG3	1.85	0.58
1:C:336:CYS:HB2	1:C:338:PHE:CE2	2.39	0.57
1:B:976:VAL:HG12	1:B:979:ASP:H	1.69	0.57
1:A:43:PHE:CB	1:B:563:GLN:OE1	2.53	0.57
1:A:847:ARG:HB2	1:A:851:CYS:CB	2.34	0.57
1:A:102:ARG:HH21	1:A:122:ASN:CA	2.18	0.57
1:B:382:VAL:HG23	1:B:517:LEU:HD11	1.86	0.57
1:C:447:GLY:HA2	1:C:498:ARG:HG2	1.87	0.57
1:C:805:ILE:HB	1:C:878:LEU:HD21	1.85	0.57
1:C:1142:GLN:OE1	1:C:1142:GLN:N	2.36	0.57
1:A:1012:LEU:HB3	1:B:1013:ILE:HD13	1.86	0.57
1:B:965:GLN:O	1:B:968:SER:OG	2.22	0.57
1:A:336:CYS:HB2	1:A:338:PHE:CE2	2.39	0.56
1:A:557:LYS:HD3	1:A:559:PHE:CE1	2.36	0.56
1:A:1144:GLU:N	1:A:1144:GLU:OE2	2.38	0.56
1:B:1142:GLN:N	1:B:1142:GLN:OE1	2.36	0.56
1:C:847:ARG:HB2	1:C:851:CYS:HB2	1.85	0.56
1:A:81:ASN:ND2	1:A:240:THR:O	2.25	0.56
1:A:136:CYS:HB2	1:A:159:VAL:HA	1.86	0.56
1:A:213:GLY:O	1:A:214:ARG:CG	2.54	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:GLY:HA2	1:A:498:ARG:HG2	1.87	0.56
1:B:1144:GLU:OE2	1:B:1144:GLU:N	2.38	0.56
1:C:215:ASP:O	1:C:216:PHE:C	2.44	0.56
1:A:322:PRO:CG	1:A:540:ASN:HD21	2.16	0.56
1:B:39:PRO:HG2	1:B:51:THR:HG21	1.87	0.56
1:C:505:HIS:NE2	1:B:503:VAL:CG2	2.68	0.56
1:A:334:ASN:ND2	1:A:360:ASN:O	2.35	0.56
1:B:102:ARG:HH21	1:B:122:ASN:CA	2.18	0.56
1:B:447:GLY:HA2	1:B:498:ARG:HG2	1.87	0.56
1:C:43:PHE:CA	1:A:563:GLN:OE1	2.53	0.56
1:B:336:CYS:HB2	1:B:338:PHE:CE2	2.39	0.56
1:C:503:VAL:HB	1:A:505:HIS:NE2	2.20	0.56
1:C:736:VAL:HG22	1:C:858:LEU:HD22	1.87	0.56
1:C:1144:GLU:N	1:C:1144:GLU:OE2	2.38	0.56
1:B:429:PHE:HE1	1:B:514:SER:HB2	1.71	0.56
1:C:102:ARG:HH21	1:C:122:ASN:CA	2.18	0.56
1:C:178:ASP:N	1:C:178:ASP:OD1	2.39	0.56
1:A:736:VAL:HG22	1:A:858:LEU:HD22	1.87	0.56
1:B:359:SER:HA	1:B:524:VAL:HG23	1.87	0.56
1:C:404:GLY:HA2	1:C:508:TYR:HD2	1.72	0.55
1:C:1002:GLN:NE2	1:B:759:PHE:CZ	2.74	0.55
1:A:359:SER:HA	1:A:524:VAL:HG23	1.87	0.55
1:C:102:ARG:HH21	1:C:122:ASN:HA	1.71	0.55
1:C:563:GLN:OE1	1:B:43:PHE:HA	2.06	0.55
1:B:213:GLY:O	1:B:214:ARG:CG	2.54	0.55
1:B:215:ASP:O	1:B:216:PHE:C	2.44	0.55
1:C:836:GLN:HG2	1:C:840:CYS:SG	2.47	0.55
1:C:852:ALA:CB	1:A:570:VAL:HG22	2.36	0.55
1:A:215:ASP:O	1:A:216:PHE:C	2.44	0.55
1:A:384:PRO:HA	1:A:387:LEU:HB2	1.87	0.55
1:B:847:ARG:HB2	1:B:851:CYS:HB2	1.88	0.55
1:C:34:ARG:NH1	1:C:191:GLU:OE2	2.34	0.55
1:C:102:ARG:HH12	1:C:176:LEU:HD12	1.72	0.55
1:C:358:ILE:HB	1:C:395:VAL:HB	1.89	0.55
1:C:130:VAL:HG23	1:C:168:PHE:HB3	1.87	0.55
1:A:102:ARG:HH12	1:A:176:LEU:HD12	1.72	0.55
1:A:453:TYR:CD1	1:A:495:TYR:HD1	2.24	0.55
1:A:852:ALA:CB	1:B:570:VAL:HG22	2.36	0.55
1:B:453:TYR:CD1	1:B:495:TYR:HD1	2.24	0.55
1:A:581:THR:C	1:A:583:GLU:H	2.10	0.55
1:C:41:LYS:O	1:A:563:GLN:HA	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:ASP:N	1:A:178:ASP:OD1	2.39	0.55
1:B:102:ARG:HH12	1:B:176:LEU:HD12	1.72	0.55
1:A:189:LEU:HG	1:A:208:THR:O	2.07	0.54
1:A:411:ALA:HB3	1:A:414:GLN:HG3	1.88	0.54
1:A:102:ARG:HH21	1:A:122:ASN:HA	1.71	0.54
1:A:880:GLY:O	1:A:884:SER:OG	2.23	0.54
1:B:570:VAL:HG23	1:B:570:VAL:O	2.08	0.54
1:C:37:TYR:HA	1:C:223:LEU:HB2	1.88	0.54
1:C:326:ILE:HD11	1:C:534:VAL:N	2.21	0.54
1:C:707:TYR:CD1	1:B:792:PRO:HG3	2.39	0.54
1:C:826:VAL:HB	1:C:1057:PRO:HG2	1.90	0.54
1:A:326:ILE:H	1:A:539:VAL:CG1	2.21	0.54
1:A:403:LYS:HD2	1:A:505:HIS:HA	1.89	0.54
1:B:826:VAL:HB	1:B:1057:PRO:HG2	1.90	0.54
1:C:1012:LEU:HB3	1:A:1013:ILE:HD13	1.88	0.54
1:A:570:VAL:O	1:A:570:VAL:HG23	2.08	0.54
1:B:975:SER:O	1:B:975:SER:OG	2.25	0.54
1:C:505:HIS:CE1	1:B:503:VAL:CB	2.91	0.54
1:A:43:PHE:HB2	1:B:563:GLN:OE1	2.08	0.54
1:A:503:VAL:HB	1:B:505:HIS:CE1	2.43	0.54
1:B:329:PHE:HB3	1:B:330:PRO:CD	2.36	0.54
1:C:418:ILE:H	1:C:418:ILE:HD12	1.72	0.54
1:B:736:VAL:HG22	1:B:858:LEU:HD22	1.88	0.54
1:C:720:ILE:HD12	1:C:923:ILE:HD11	1.90	0.54
1:A:742:ILE:HG23	1:A:1000:ARG:HB2	1.88	0.54
1:B:102:ARG:HH21	1:B:122:ASN:HA	1.71	0.54
1:B:475:ALA:HB3	1:B:487:ASN:HB3	1.90	0.54
1:C:62:VAL:CG1	1:C:216:PHE:HE2	2.15	0.54
1:C:453:TYR:CD1	1:C:495:TYR:HD1	2.24	0.54
1:A:560:LEU:O	1:A:562:PHE:N	2.41	0.54
1:B:326:ILE:H	1:B:539:VAL:CG1	2.21	0.54
1:C:326:ILE:H	1:C:539:VAL:CG1	2.21	0.54
1:C:1100:THR:HB	2:C:1315:NAG:HN2	1.73	0.54
1:A:752:LEU:HD12	1:A:993:ILE:HG21	1.90	0.54
1:C:560:LEU:O	1:C:562:PHE:N	2.41	0.53
1:C:975:SER:O	1:C:975:SER:OG	2.25	0.53
1:C:1002:GLN:NE2	1:B:759:PHE:HZ	2.05	0.53
1:A:57:PRO:HG3	1:A:273:ARG:HD2	1.90	0.53
1:A:759:PHE:CE2	1:B:1002:GLN:CG	2.90	0.53
1:B:559:PHE:HB2	1:B:584:ILE:CD1	2.35	0.53
1:C:475:ALA:HB3	1:C:487:ASN:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1002:GLN:HE21	1:B:759:PHE:HZ	1.55	0.53
1:A:117:LEU:HD12	1:A:118:LEU:N	2.24	0.53
1:A:429:PHE:HE1	1:A:514:SER:HB2	1.74	0.53
1:A:826:VAL:HB	1:A:1057:PRO:HG2	1.90	0.53
1:B:200:TYR:HA	1:B:230:PRO:HA	1.91	0.53
1:B:560:LEU:O	1:B:562:PHE:N	2.41	0.53
1:B:720:ILE:HD12	1:B:923:ILE:HD11	1.90	0.53
1:C:57:PRO:HG3	1:C:273:ARG:HD2	1.90	0.53
1:C:375:PHE:CG	1:C:508:TYR:CZ	2.97	0.53
1:C:880:GLY:O	1:C:884:SER:OG	2.23	0.53
1:A:62:VAL:CG1	1:A:216:PHE:HE2	2.15	0.53
1:B:752:LEU:HD12	1:B:993:ILE:HG21	1.90	0.53
1:C:403:LYS:HG2	1:C:505:HIS:HA	1.90	0.53
1:A:129:LYS:HD3	1:A:133:PHE:CZ	2.43	0.53
1:A:965:GLN:O	1:A:968:SER:OG	2.22	0.53
1:A:975:SER:O	1:A:975:SER:OG	2.25	0.53
1:B:129:LYS:HA	1:B:168:PHE:O	2.09	0.53
1:B:439:ASN:O	1:B:443:SER:OG	2.25	0.53
1:B:945:LEU:HD12	1:B:948:LEU:HD12	1.90	0.53
1:C:957:GLN:OE1	1:B:765:ARG:HD3	2.08	0.53
1:B:178:ASP:OD1	1:B:178:ASP:N	2.39	0.53
1:C:765:ARG:HD3	1:A:957:GLN:OE1	2.08	0.53
1:B:847:ARG:HB2	1:B:851:CYS:CB	2.39	0.53
1:C:945:LEU:HD12	1:C:948:LEU:HD12	1.90	0.53
1:B:57:PRO:HG3	1:B:273:ARG:HD2	1.90	0.53
1:B:375:PHE:CG	1:B:508:TYR:CZ	2.97	0.53
1:C:858:LEU:HD12	1:C:963:VAL:HG12	1.91	0.53
1:C:965:GLN:O	1:C:968:SER:OG	2.22	0.53
1:A:421:TYR:CD1	1:A:457:ARG:HB3	2.44	0.53
1:A:720:ILE:HD12	1:A:923:ILE:HD11	1.90	0.53
1:B:880:GLY:O	1:B:884:SER:OG	2.23	0.53
1:C:752:LEU:HD12	1:C:993:ILE:HG21	1.90	0.52
1:B:403:LYS:HD2	1:B:505:HIS:HA	1.90	0.52
1:A:858:LEU:HD12	1:A:963:VAL:HG12	1.91	0.52
1:B:420:ASP:HB3	1:B:460:LYS:HD2	1.91	0.52
1:B:1083:HIS:HB2	1:B:1137:VAL:HG13	1.91	0.52
1:C:993:ILE:O	1:C:997:ILE:HG12	2.10	0.52
1:B:752:LEU:HD12	1:B:993:ILE:CG2	2.40	0.52
1:B:858:LEU:HD12	1:B:963:VAL:HG12	1.91	0.52
1:B:993:ILE:O	1:B:997:ILE:HG12	2.10	0.52
1:C:133:PHE:CD2	1:C:160:TYR:HA	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:GLU:HB2	1:C:223:LEU:HD11	1.91	0.52
1:C:418:ILE:HA	1:C:422:ASN:HB2	1.91	0.52
1:A:945:LEU:HD12	1:A:948:LEU:HD12	1.91	0.52
1:B:37:TYR:HB3	1:B:223:LEU:HB2	1.92	0.52
1:C:803:SER:HB3	2:C:1313:NAG:HN2	1.74	0.52
1:A:375:PHE:CG	1:A:508:TYR:CZ	2.97	0.52
1:A:475:ALA:HB3	1:A:487:ASN:HB3	1.90	0.52
1:B:108:THR:C	1:B:110:LEU:H	2.13	0.52
1:B:912:THR:O	1:B:915:VAL:HG12	2.10	0.52
1:B:559:PHE:CB	1:B:584:ILE:HD11	2.36	0.52
1:B:397:ALA:HA	1:B:512:VAL:O	2.09	0.52
1:C:117:LEU:HD12	1:C:118:LEU:N	2.24	0.52
1:C:755:GLN:OE1	1:A:971:GLY:CA	2.41	0.52
1:A:755:GLN:OE1	1:B:971:GLY:CA	2.42	0.52
1:A:752:LEU:HD12	1:A:993:ILE:CG2	2.40	0.52
1:C:1083:HIS:HB2	1:C:1137:VAL:HG13	1.91	0.51
1:A:108:THR:C	1:A:110:LEU:H	2.13	0.51
1:B:229:LEU:H	1:B:229:LEU:HD23	1.75	0.51
1:B:364:ASP:OD1	1:B:364:ASP:N	2.43	0.51
1:C:1076:THR:HB	1:C:1097:SER:HB3	1.92	0.51
1:A:115:GLN:HE21	1:A:233:ILE:HB	1.76	0.51
1:A:993:ILE:O	1:A:997:ILE:HG12	2.10	0.51
1:C:847:ARG:HB2	1:C:851:CYS:CB	2.41	0.51
1:A:329:PHE:HB3	1:A:330:PRO:CD	2.34	0.51
1:A:1083:HIS:HB2	1:A:1137:VAL:HG13	1.91	0.51
1:B:347:PHE:HD2	1:B:399:SER:HB2	1.75	0.51
1:C:752:LEU:HD12	1:C:993:ILE:CG2	2.40	0.51
1:C:912:THR:O	1:C:915:VAL:HG12	2.11	0.51
1:B:931:ILE:O	1:B:934:ILE:HG22	2.09	0.51
1:A:307:THR:HA	1:A:602:THR:HG21	1.93	0.51
1:B:322:PRO:CG	1:B:540:ASN:ND2	2.74	0.51
1:C:364:ASP:N	1:C:364:ASP:OD1	2.43	0.51
1:C:854:LYS:HE2	1:A:592:PHE:CD2	2.46	0.51
1:C:1006:THR:HG21	1:B:762:GLN:OE1	2.10	0.51
1:B:631:PRO:HB3	1:B:633:TRP:CZ2	2.45	0.51
1:B:742:ILE:O	1:B:1000:ARG:NH1	2.43	0.51
1:B:805:ILE:HB	1:B:878:LEU:HD21	1.93	0.51
1:A:31:SER:HB3	1:A:62:VAL:CG2	2.41	0.50
1:B:559:PHE:HD2	1:B:577:ARG:HD2	1.77	0.50
1:C:631:PRO:HB3	1:C:633:TRP:CZ2	2.46	0.50
1:C:759:PHE:CE2	1:A:1002:GLN:CG	2.90	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:836:GLN:O	1:C:837:TYR:C	2.50	0.50
1:B:382:VAL:CG2	1:B:517:LEU:HD11	2.41	0.50
1:B:557:LYS:HD3	1:B:559:PHE:CE1	2.34	0.50
1:C:276:LEU:CD1	1:C:304:LYS:HA	2.42	0.50
1:C:854:LYS:HE2	1:A:592:PHE:HD2	1.75	0.50
1:A:387:LEU:HD11	1:A:515:PHE:CE2	2.47	0.50
1:C:563:GLN:OE1	1:B:43:PHE:CA	2.60	0.50
1:C:675:GLN:HG3	1:C:693:ILE:HD11	1.94	0.50
1:A:177:MET:H	1:A:207:HIS:CE1	2.30	0.50
1:A:765:ARG:HD3	1:B:957:GLN:OE1	2.11	0.50
1:B:189:LEU:HG	1:B:208:THR:O	2.12	0.50
1:B:27:SER:O	1:B:64:TRP:N	2.45	0.50
1:B:581:THR:C	1:B:583:GLU:H	2.14	0.50
1:A:290:ASP:O	1:A:297:SER:HB3	2.12	0.50
1:C:187:LYS:O	1:C:188:ASN:C	2.50	0.50
1:C:403:LYS:HB3	1:C:406:GLU:OE1	2.12	0.50
1:B:113:LYS:HD2	1:B:113:LYS:O	2.12	0.50
1:B:317:ASN:HB2	1:B:592:PHE:CE1	2.46	0.50
1:C:31:SER:HB3	1:C:62:VAL:CG2	2.41	0.50
1:C:762:GLN:OE1	1:A:1006:THR:HG21	2.12	0.50
1:A:382:VAL:CG2	1:A:517:LEU:HD11	2.41	0.50
1:A:905:ARG:HD2	1:A:1049:LEU:O	2.12	0.50
1:B:215:ASP:HB2	1:B:266:TYR:OH	2.12	0.50
1:C:503:VAL:CB	1:A:505:HIS:NE2	2.75	0.49
1:C:905:ARG:HD2	1:C:1049:LEU:O	2.12	0.49
1:A:364:ASP:OD1	1:A:364:ASP:N	2.43	0.49
1:A:762:GLN:OE1	1:B:1006:THR:HG21	2.12	0.49
1:C:177:MET:H	1:C:207:HIS:CE1	2.30	0.49
1:C:215:ASP:HB2	1:C:266:TYR:OH	2.12	0.49
1:C:290:ASP:O	1:C:297:SER:HB3	2.11	0.49
1:A:360:ASN:H	1:A:523:THR:HB	1.78	0.49
1:B:675:GLN:HG3	1:B:693:ILE:HD11	1.94	0.49
1:C:113:LYS:HD2	1:C:113:LYS:O	2.12	0.49
1:A:276:LEU:CD1	1:A:304:LYS:HA	2.42	0.49
1:A:322:PRO:CG	1:A:540:ASN:ND2	2.74	0.49
1:C:946:GLY:O	1:C:950:ASP:HB2	2.13	0.49
1:A:44:ARG:HB3	1:A:47:VAL:CG1	2.41	0.49
1:B:384:PRO:HA	1:B:387:LEU:HB2	1.94	0.49
1:B:821:LEU:HD11	1:B:935:GLN:HG3	1.94	0.49
1:C:570:VAL:HG23	1:C:572:THR:HG23	1.94	0.49
1:A:113:LYS:HD2	1:A:113:LYS:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:631:PRO:HB3	1:A:633:TRP:CZ2	2.47	0.49
1:A:844:ILE:O	1:B:557:LYS:HG3	2.12	0.49
1:B:131:CYS:SG	1:B:163:ALA:O	2.71	0.49
1:B:946:GLY:O	1:B:950:ASP:HB2	2.13	0.49
1:C:108:THR:C	1:C:110:LEU:H	2.13	0.49
1:C:131:CYS:HB3	1:C:133:PHE:CE2	2.47	0.49
1:A:41:LYS:O	1:B:563:GLN:HA	2.12	0.49
1:A:749:CYS:SG	1:A:997:ILE:HD11	2.53	0.49
1:C:189:LEU:HG	1:C:208:THR:O	2.12	0.49
1:A:191:GLU:HB2	1:A:223:LEU:HD11	1.93	0.49
1:A:347:PHE:CE1	1:A:509:ARG:HD3	2.48	0.49
1:A:412:PRO:HG3	1:A:429:PHE:HB3	1.93	0.49
1:C:360:ASN:H	1:C:523:THR:HB	1.77	0.49
1:A:131:CYS:HB3	1:A:133:PHE:CE2	2.48	0.49
1:A:946:GLY:O	1:A:950:ASP:HB2	2.13	0.49
1:B:177:MET:H	1:B:207:HIS:CE1	2.30	0.49
1:C:44:ARG:HB3	1:C:47:VAL:CG1	2.41	0.48
1:C:307:THR:HA	1:C:602:THR:HG21	1.95	0.48
1:C:326:ILE:HD13	1:C:533:LEU:C	2.31	0.48
1:C:971:GLY:HA2	1:B:755:GLN:CD	2.31	0.48
1:A:497:PHE:HA	1:A:501:TYR:HE2	1.78	0.48
1:A:675:GLN:HG3	1:A:693:ILE:HD11	1.94	0.48
1:B:347:PHE:CE1	1:B:509:ARG:HD3	2.48	0.48
1:B:360:ASN:H	1:B:523:THR:HB	1.77	0.48
1:B:588:THR:HG23	1:B:589:PRO:HD2	1.95	0.48
1:C:578:ASP:O	1:C:582:LEU:N	2.45	0.48
1:C:749:CYS:SG	1:C:997:ILE:HD11	2.53	0.48
1:A:740:MET:HG3	1:B:592:PHE:CZ	2.48	0.48
1:A:973:ILE:HG22	1:A:992:GLN:HG3	1.95	0.48
1:B:556:ASN:O	1:B:557:LYS:C	2.51	0.48
1:B:927:PHE:HZ	1:B:1052:PHE:HE2	1.60	0.48
1:C:568:ASP:HA	1:B:846:ALA:HB1	1.95	0.48
1:C:588:THR:HG22	1:C:589:PRO:O	2.13	0.48
1:C:588:THR:OG1	1:B:841:LEU:HB3	2.13	0.48
1:A:296:LEU:HD11	1:A:308:VAL:HG21	1.95	0.48
1:B:126:VAL:HG11	1:B:175:PHE:CE1	2.49	0.48
1:B:214:ARG:O	1:B:217:PRO:HG3	2.13	0.48
1:B:555:SER:HB3	1:B:584:ILE:HG22	1.95	0.48
1:C:760:CYS:HA	1:C:763:LEU:HG	1.96	0.48
1:A:215:ASP:HB2	1:A:266:TYR:OH	2.12	0.48
1:A:374:PHE:CZ	1:A:434:ILE:HG23	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:927:PHE:HZ	1:A:1052:PHE:HE2	1.61	0.48
1:A:1106:GLN:OE1	1:A:1106:GLN:N	2.47	0.48
1:B:497:PHE:HA	1:B:501:TYR:HE2	1.78	0.48
1:C:214:ARG:O	1:C:217:PRO:HG3	2.13	0.48
1:B:760:CYS:HA	1:B:763:LEU:HG	1.95	0.48
1:B:905:ARG:HD2	1:B:1049:LEU:O	2.12	0.48
1:A:103:GLY:HA3	1:A:120:VAL:HA	1.95	0.48
1:A:106:PHE:CD1	1:A:235:ILE:HG21	2.49	0.48
1:A:802:PHE:O	1:A:804:GLN:N	2.47	0.48
1:C:431:GLY:HA2	1:C:515:PHE:HD2	1.77	0.48
1:B:44:ARG:HB3	1:B:47:VAL:CG1	2.41	0.48
1:B:106:PHE:CD1	1:B:235:ILE:HG21	2.49	0.48
1:B:191:GLU:HB2	1:B:223:LEU:HD11	1.95	0.48
1:C:126:VAL:HG11	1:C:175:PHE:CE1	2.49	0.48
1:A:331:ASN:HB3	1:A:580:GLN:HG2	1.94	0.48
1:B:322:PRO:CG	1:B:540:ASN:HD21	2.16	0.48
1:C:131:CYS:HA	1:C:166:CYS:HA	1.96	0.48
1:A:126:VAL:HG11	1:A:175:PHE:CE1	2.49	0.48
1:B:749:CYS:SG	1:B:997:ILE:HD11	2.53	0.48
1:B:973:ILE:HG22	1:B:992:GLN:HG3	1.95	0.48
1:C:81:ASN:ND2	1:C:240:THR:O	2.25	0.48
1:C:296:LEU:HD11	1:C:308:VAL:HG21	1.95	0.48
1:C:497:PHE:HA	1:C:501:TYR:HE2	1.78	0.48
1:C:843:ASP:O	1:C:845:ALA:N	2.47	0.48
1:C:439:ASN:O	1:C:443:SER:OG	2.25	0.47
1:B:878:LEU:HD13	1:B:1053:PRO:HD2	1.96	0.47
1:C:31:SER:HB3	1:C:62:VAL:HG21	1.96	0.47
1:C:347:PHE:CE1	1:C:509:ARG:HD3	2.48	0.47
1:A:439:ASN:O	1:A:443:SER:OG	2.25	0.47
1:C:106:PHE:HB3	1:C:235:ILE:HG23	1.97	0.47
1:C:557:LYS:HB3	1:C:559:PHE:HD1	1.79	0.47
1:C:973:ILE:HG22	1:C:992:GLN:HG3	1.95	0.47
1:A:760:CYS:HA	1:A:763:LEU:HG	1.96	0.47
1:B:213:GLY:C	1:B:214:ARG:HG3	2.35	0.47
1:B:726:ILE:HG23	1:B:1061:VAL:HG22	1.96	0.47
1:C:103:GLY:HA3	1:C:120:VAL:HA	1.96	0.47
1:C:927:PHE:HZ	1:C:1052:PHE:HE2	1.60	0.47
1:C:1002:GLN:CD	1:B:759:PHE:HZ	2.15	0.47
1:A:520:ALA:O	1:A:521:PRO:C	2.53	0.47
1:A:726:ILE:HG23	1:A:1061:VAL:HG22	1.96	0.47
1:B:342:PHE:CZ	1:B:511:VAL:HG21	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:LEU:HB2	1:C:238:PHE:HD2	1.79	0.47
1:C:329:PHE:HB3	1:C:330:PRO:HD2	1.93	0.47
1:A:1101:HIS:CD2	2:A:1315:NAG:H3	2.49	0.47
1:B:84:LEU:HB2	1:B:238:PHE:HD2	1.79	0.47
1:B:383:SER:HB3	1:B:386:LYS:HG2	1.95	0.47
1:C:322:PRO:CG	1:C:540:ASN:HD21	2.16	0.47
1:C:726:ILE:HG23	1:C:1061:VAL:HG22	1.96	0.47
1:C:850:ILE:O	1:C:854:LYS:HG2	2.15	0.47
1:C:931:ILE:O	1:C:934:ILE:HG22	2.15	0.47
1:A:350:VAL:HG22	1:A:422:ASN:HB3	1.96	0.47
1:C:106:PHE:CD1	1:C:235:ILE:HG21	2.49	0.47
1:C:356:THR:OG1	1:C:397:ALA:HB3	2.15	0.47
1:A:213:GLY:C	1:A:214:ARG:HG3	2.34	0.47
1:A:931:ILE:O	1:A:934:ILE:HG22	2.15	0.47
1:B:117:LEU:HD12	1:B:118:LEU:N	2.30	0.47
1:B:206:LYS:HB3	1:B:223:LEU:HD13	1.97	0.47
1:B:216:PHE:O	1:B:217:PRO:C	2.52	0.47
1:B:1132:ILE:HD12	1:B:1132:ILE:HA	1.78	0.47
1:C:823:PHE:CD1	1:C:1057:PRO:HG3	2.50	0.47
1:A:106:PHE:HB3	1:A:235:ILE:HG23	1.97	0.47
1:A:214:ARG:O	1:A:217:PRO:HG3	2.13	0.47
1:A:906:PHE:CD2	1:A:916:LEU:HB2	2.50	0.47
1:C:177:MET:HB2	1:C:207:HIS:ND1	2.30	0.47
1:C:322:PRO:CG	1:C:540:ASN:ND2	2.74	0.47
1:C:577:ARG:HG2	1:C:584:ILE:HD13	1.97	0.47
1:C:631:PRO:HA	1:C:633:TRP:CZ3	2.50	0.47
1:C:741:TYR:CD2	1:C:1004:LEU:HD22	2.50	0.47
1:A:31:SER:HB3	1:A:62:VAL:HG21	1.96	0.47
1:A:627:ASP:HA	1:A:634:ARG:HH22	1.80	0.47
1:A:823:PHE:CD1	1:A:1057:PRO:HG3	2.50	0.47
1:B:823:PHE:CD1	1:B:1057:PRO:HG3	2.50	0.47
1:A:878:LEU:HD13	1:A:1053:PRO:HD2	1.96	0.46
1:C:175:PHE:CD1	1:C:175:PHE:N	2.82	0.46
1:C:355:ARG:NE	1:C:398:ASP:OD2	2.48	0.46
1:C:742:ILE:HG22	1:C:997:ILE:CD1	2.45	0.46
1:C:983:ARG:HD3	1:A:517:LEU:CD1	2.30	0.46
1:C:1132:ILE:HA	1:C:1132:ILE:HD12	1.77	0.46
1:A:216:PHE:O	1:A:217:PRO:C	2.52	0.46
1:B:204:TYR:CE2	1:B:225:PRO:HG3	2.50	0.46
1:B:350:VAL:HG22	1:B:422:ASN:HB3	1.97	0.46
1:C:25:THR:HG23	1:C:66:HIS:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:878:LEU:HD13	1:C:1053:PRO:HD2	1.96	0.46
1:A:84:LEU:HB2	1:A:238:PHE:HD2	1.79	0.46
1:A:318:PHE:HD2	1:A:629:LEU:HD11	1.81	0.46
1:B:103:GLY:HA3	1:B:120:VAL:HA	1.96	0.46
1:B:559:PHE:CD1	1:B:584:ILE:HD13	2.50	0.46
1:B:718:PHE:HE1	1:B:919:ASN:ND2	2.14	0.46
1:C:755:GLN:CD	1:A:971:GLY:HA2	2.30	0.46
1:A:718:PHE:HE1	1:A:919:ASN:ND2	2.14	0.46
1:B:177:MET:HB2	1:B:207:HIS:ND1	2.31	0.46
1:C:176:LEU:HD22	1:C:190:ARG:HD2	1.98	0.46
1:C:645:THR:HG23	1:C:647:ALA:N	2.31	0.46
1:C:650:LEU:HD23	1:C:653:ALA:HB3	1.98	0.46
1:B:129:LYS:HE2	1:B:169:GLU:HG3	1.98	0.46
1:B:307:THR:HA	1:B:602:THR:HG21	1.96	0.46
1:C:563:GLN:HA	1:B:41:LYS:O	2.16	0.46
1:B:25:THR:HG23	1:B:66:HIS:HB2	1.98	0.46
1:B:454:ARG:HH22	1:B:467:ASP:HB3	1.81	0.46
1:A:133:PHE:CD2	1:A:160:TYR:HA	2.51	0.46
1:B:118:LEU:HD23	1:B:118:LEU:HA	1.81	0.46
1:C:1083:HIS:CE1	1:C:1137:VAL:CG2	2.96	0.46
1:A:175:PHE:CD1	1:A:175:PHE:N	2.82	0.46
1:A:347:PHE:HD2	1:A:399:SER:HB2	1.81	0.46
1:B:627:ASP:HA	1:B:634:ARG:HH22	1.80	0.46
1:B:1092:GLU:N	1:B:1092:GLU:OE2	2.49	0.46
1:C:592:PHE:CD1	1:B:740:MET:HE2	2.51	0.46
1:C:763:LEU:HD13	1:C:1004:LEU:HG	1.98	0.46
1:A:342:PHE:CE2	1:A:436:TRP:CZ3	3.04	0.46
1:C:454:ARG:HH22	1:C:467:ASP:HB3	1.81	0.46
1:C:718:PHE:HE1	1:C:919:ASN:ND2	2.14	0.46
1:A:25:THR:HG23	1:A:66:HIS:HB2	1.98	0.46
1:A:177:MET:HB2	1:A:207:HIS:ND1	2.31	0.46
1:A:439:ASN:OD1	1:A:507:PRO:HD2	2.17	0.46
1:A:631:PRO:HA	1:A:633:TRP:CZ3	2.50	0.46
1:B:575:ALA:HA	1:B:585:LEU:O	2.16	0.46
1:C:423:TYR:OH	1:C:464:PHE:HE1	1.98	0.45
1:A:454:ARG:HH22	1:A:467:ASP:HB3	1.81	0.45
1:B:528:LYS:O	1:B:529:LYS:HB3	2.16	0.45
1:C:105:ILE:HD13	1:C:159:VAL:HG11	1.98	0.45
1:C:439:ASN:OD1	1:C:507:PRO:HD2	2.17	0.45
1:C:528:LYS:O	1:C:529:LYS:HB3	2.16	0.45
1:C:617:CYS:O	1:C:621:SER:N	2.40	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:759:PHE:CZ	1:B:1002:GLN:NE2	2.84	0.45
1:A:1092:GLU:OE2	1:A:1092:GLU:N	2.49	0.45
1:B:106:PHE:HB3	1:B:235:ILE:HG23	1.97	0.45
1:B:162:SER:O	1:B:163:ALA:C	2.54	0.45
1:B:631:PRO:HA	1:B:633:TRP:CZ3	2.51	0.45
1:B:763:LEU:HD13	1:B:1004:LEU:HG	1.98	0.45
1:B:1101:HIS:CD2	2:B:1315:NAG:H3	2.51	0.45
1:C:644:GLN:NE2	1:C:648:GLY:O	2.49	0.45
1:A:528:LYS:O	1:A:529:LYS:HB3	2.16	0.45
1:A:763:LEU:HD13	1:A:1004:LEU:HG	1.98	0.45
1:B:106:PHE:CE2	1:B:238:PHE:HD1	2.34	0.45
1:B:878:LEU:O	1:B:882:ILE:HG12	2.17	0.45
1:C:164:ASN:HB2	2:C:1302:NAG:HN2	1.81	0.45
1:C:343:ASN:OD1	1:C:343:ASN:N	2.50	0.45
1:A:1101:HIS:CG	2:A:1315:NAG:H5	2.52	0.45
1:B:327:VAL:HG11	1:B:528:LYS:CG	2.47	0.45
1:C:1092:GLU:OE2	1:C:1092:GLU:N	2.49	0.45
1:A:176:LEU:HD22	1:A:190:ARG:HD2	1.98	0.45
1:B:520:ALA:O	1:B:521:PRO:C	2.53	0.45
1:B:705:VAL:HG12	1:B:707:TYR:H	1.82	0.45
1:A:131:CYS:HA	1:A:166:CYS:HA	1.99	0.45
1:A:210:ILE:H	1:A:210:ILE:HG13	1.56	0.45
1:A:501:TYR:HB3	1:A:505:HIS:HB2	1.98	0.45
1:A:742:ILE:HD13	1:A:1001:LEU:HD13	1.99	0.45
1:B:363:ALA:O	1:B:527:PRO:HD3	2.17	0.45
1:B:878:LEU:HD12	1:B:878:LEU:HA	1.75	0.45
1:C:505:HIS:CE1	1:B:503:VAL:HB	2.52	0.45
1:C:520:ALA:O	1:C:521:PRO:C	2.53	0.45
1:C:742:ILE:HG22	1:C:997:ILE:HD12	1.99	0.45
1:A:303:LEU:H	1:A:303:LEU:HG	1.67	0.45
1:B:97:LYS:HE2	1:B:184:GLY:HA3	1.99	0.45
1:B:342:PHE:CE2	1:B:436:TRP:CZ3	3.04	0.45
1:A:188:ASN:OD1	1:A:188:ASN:N	2.50	0.45
1:A:343:ASN:N	1:A:343:ASN:OD1	2.50	0.45
1:A:390:LEU:HD13	1:A:390:LEU:HA	1.83	0.45
1:A:503:VAL:HB	1:B:505:HIS:NE2	2.31	0.45
1:B:343:ASN:OD1	1:B:343:ASN:N	2.50	0.45
1:C:41:LYS:HB3	1:A:562:PHE:O	2.17	0.45
1:C:97:LYS:HE2	1:C:184:GLY:HA3	1.98	0.45
1:C:125:ASN:HB2	2:C:1301:NAG:H62	1.99	0.45
1:C:321:GLN:NE2	1:C:630:THR:OG1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:501:TYR:HB3	1:C:505:HIS:HB2	1.98	0.45
1:C:741:TYR:CE2	1:C:1004:LEU:HB2	2.51	0.45
1:C:878:LEU:O	1:C:882:ILE:HG12	2.17	0.45
1:A:129:LYS:HD3	1:A:133:PHE:HZ	1.82	0.45
1:A:327:VAL:HG11	1:A:528:LYS:CG	2.47	0.45
1:A:705:VAL:HG12	1:A:707:TYR:H	1.82	0.45
1:A:744:GLY:O	1:A:746:SER:N	2.50	0.45
1:A:759:PHE:HZ	1:B:1002:GLN:NE2	2.15	0.45
1:C:106:PHE:CE2	1:C:238:PHE:HD1	2.34	0.45
1:C:278:LYS:HB2	1:C:306:PHE:CE1	2.52	0.45
1:C:363:ALA:O	1:C:527:PRO:HD3	2.17	0.45
1:A:106:PHE:CE2	1:A:238:PHE:HD1	2.34	0.45
1:A:363:ALA:O	1:A:527:PRO:HD3	2.17	0.45
1:B:175:PHE:CD1	1:B:175:PHE:N	2.82	0.45
1:C:342:PHE:CE2	1:C:436:TRP:CZ3	3.04	0.44
1:C:1013:ILE:CD1	1:B:1012:LEU:HD13	2.47	0.44
1:B:176:LEU:HD22	1:B:190:ARG:HD2	1.98	0.44
1:C:907:ASN:HB2	1:C:913:GLN:HG2	1.99	0.44
1:B:276:LEU:HD22	1:B:301:CYS:HA	1.99	0.44
1:C:44:ARG:O	1:C:283:GLY:HA2	2.18	0.44
1:C:216:PHE:O	1:C:217:PRO:C	2.52	0.44
1:C:326:ILE:HG12	1:C:532:ASN:O	2.18	0.44
1:A:321:GLN:NE2	1:A:630:THR:OG1	2.49	0.44
1:B:501:TYR:HB3	1:B:505:HIS:HB2	1.98	0.44
1:C:327:VAL:HG11	1:C:528:LYS:CG	2.47	0.44
1:C:780:GLU:O	1:C:784:GLN:NE2	2.49	0.44
1:C:858:LEU:HD13	1:C:962:LEU:HD13	2.00	0.44
1:A:348:ALA:HB1	1:A:352:ALA:O	2.18	0.44
1:B:188:ASN:N	1:B:188:ASN:OD1	2.50	0.44
1:B:450:ASP:OD1	1:B:450:ASP:N	2.50	0.44
1:C:244:LEU:HD23	1:C:244:LEU:HA	1.87	0.44
1:C:567:ARG:HB3	1:C:568:ASP:H	1.67	0.44
1:A:278:LYS:HB2	1:A:306:PHE:CE1	2.52	0.44
1:A:309:GLU:HG2	1:A:313:TYR:OH	2.18	0.44
1:B:34:ARG:NH2	1:B:217:PRO:O	2.51	0.44
1:A:97:LYS:HE2	1:A:184:GLY:HA3	1.98	0.44
1:A:402:ILE:HD11	1:A:510:VAL:HG21	2.00	0.44
1:C:375:PHE:CG	1:C:508:TYR:OH	2.68	0.44
1:C:1002:GLN:HG3	1:B:759:PHE:CE2	2.44	0.44
1:A:125:ASN:HB2	2:A:1301:NAG:H62	1.99	0.44
1:A:742:ILE:HG22	1:A:997:ILE:CD1	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:930:ALA:HA	1:A:933:LYS:HG2	2.00	0.44
1:B:125:ASN:HB2	2:B:1301:NAG:H62	1.99	0.44
1:B:439:ASN:OD1	1:B:507:PRO:HD2	2.17	0.44
1:B:1101:HIS:CG	2:B:1315:NAG:H5	2.53	0.44
1:C:930:ALA:HA	1:C:933:LYS:HG2	2.00	0.44
1:A:878:LEU:O	1:A:882:ILE:HG12	2.17	0.44
1:B:348:ALA:HB1	1:B:352:ALA:O	2.18	0.44
1:C:348:ALA:HB1	1:C:352:ALA:O	2.18	0.44
1:B:195:LYS:HE2	1:B:195:LYS:HB2	1.79	0.44
1:B:358:ILE:HB	1:B:395:VAL:HB	2.00	0.44
1:B:930:ALA:HA	1:B:933:LYS:HG2	2.00	0.44
1:C:322:PRO:CB	1:C:540:ASN:ND2	2.81	0.43
1:C:339:HIS:O	1:C:340:GLU:C	2.53	0.43
1:C:450:ASP:N	1:C:450:ASP:OD1	2.51	0.43
1:C:562:PHE:O	1:B:41:LYS:HD3	2.17	0.43
1:C:705:VAL:C	1:C:707:TYR:H	2.21	0.43
1:A:44:ARG:O	1:A:283:GLY:HA2	2.18	0.43
1:A:581:THR:O	1:A:583:GLU:N	2.50	0.43
1:A:792:PRO:HG3	1:B:707:TYR:CD1	2.50	0.43
1:B:419:ALA:HA	1:B:423:TYR:O	2.18	0.43
1:C:34:ARG:NH2	1:C:217:PRO:O	2.51	0.43
1:C:1001:LEU:HD12	1:C:1001:LEU:HA	1.77	0.43
1:C:1078:ALA:C	1:B:900:MET:HE1	2.39	0.43
1:A:806:LEU:HD23	1:A:878:LEU:HD23	1.99	0.43
1:B:339:HIS:O	1:B:340:GLU:C	2.53	0.43
1:A:413:GLY:HA2	1:A:424:LYS:HD2	1.98	0.43
1:A:857:GLY:O	1:A:858:LEU:HD23	2.18	0.43
3:B:1317:IDU:O1S	3:B:1317:IDU:O3	2.22	0.43
1:A:454:ARG:HH12	1:A:467:ASP:HB3	1.84	0.43
1:C:401:VAL:HG21	1:C:451:TYR:CE2	2.54	0.43
1:A:554:LYS:HZ2	1:A:554:LYS:HA	1.84	0.43
1:B:322:PRO:CB	1:B:540:ASN:ND2	2.81	0.43
1:B:358:ILE:HG22	1:B:524:VAL:HG21	2.00	0.43
1:C:309:GLU:HG2	1:C:313:TYR:OH	2.18	0.43
1:C:742:ILE:HD13	1:C:1001:LEU:HD13	2.00	0.43
1:C:744:GLY:O	1:C:745:ASP:C	2.57	0.43
1:A:108:THR:O	1:A:110:LEU:N	2.50	0.43
1:A:212:ILE:HG13	1:A:212:ILE:O	2.18	0.43
1:A:339:HIS:O	1:A:340:GLU:C	2.53	0.43
1:A:858:LEU:HD13	1:A:962:LEU:HD13	2.00	0.43
1:A:947:LYS:HE3	1:A:947:LYS:HB2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:358:ILE:HG22	1:C:524:VAL:HG21	2.00	0.43
1:C:557:LYS:CD	1:C:559:PHE:HE1	2.32	0.43
1:A:908:GLY:HA3	1:A:1036:GLN:HE22	1.84	0.43
1:A:1072:GLU:OE1	1:A:1072:GLU:N	2.51	0.43
1:B:278:LYS:HB2	1:B:306:PHE:CE1	2.52	0.43
1:B:789:TYR:CE2	1:B:888:PHE:HD2	2.36	0.43
1:C:34:ARG:NH2	1:C:219:GLY:O	2.52	0.43
1:C:382:VAL:HG22	1:C:517:LEU:HD11	2.01	0.43
1:A:164:ASN:O	1:A:166:CYS:N	2.52	0.43
1:B:418:ILE:O	1:B:422:ASN:HB2	2.19	0.43
1:B:454:ARG:HH12	1:B:467:ASP:HB3	1.84	0.43
1:C:118:LEU:HD11	1:C:159:VAL:HB	2.00	0.43
2:C:1311:NAG:H62	1:B:796:TYR:CG	2.54	0.43
1:A:744:GLY:O	1:A:745:ASP:C	2.57	0.43
1:A:841:LEU:HB3	1:B:588:THR:OG1	2.19	0.43
1:A:1083:HIS:CE1	1:A:1137:VAL:CG2	2.96	0.43
1:B:44:ARG:O	1:B:283:GLY:HA2	2.18	0.43
1:B:455:SER:OG	1:B:456:PHE:CD1	2.71	0.43
1:B:908:GLY:HA3	1:B:1036:GLN:HE22	1.84	0.43
1:C:1029:MET:O	1:C:1033:VAL:HB	2.19	0.43
1:A:322:PRO:CB	1:A:540:ASN:ND2	2.81	0.43
1:A:742:ILE:HG22	1:A:997:ILE:HD12	1.99	0.43
1:C:134:GLN:HE21	1:C:135:PHE:H	1.67	0.42
1:A:34:ARG:NH2	1:A:217:PRO:O	2.51	0.42
1:A:34:ARG:NH2	1:A:219:GLY:O	2.52	0.42
1:A:320:VAL:HG22	1:A:629:LEU:HD12	2.01	0.42
1:A:340:GLU:O	1:A:342:PHE:N	2.52	0.42
1:B:303:LEU:H	1:B:303:LEU:HG	1.73	0.42
1:B:375:PHE:CG	1:B:508:TYR:OH	2.68	0.42
1:B:388:ASN:HB3	1:B:527:PRO:HD2	2.00	0.42
1:C:327:VAL:HG13	1:C:542:ASN:HB3	2.01	0.42
1:C:434:ILE:HB	1:C:511:VAL:HG22	2.01	0.42
1:C:744:GLY:O	1:C:746:SER:N	2.52	0.42
1:A:130:VAL:HG23	1:A:168:PHE:HB3	2.00	0.42
1:A:503:VAL:CB	1:B:505:HIS:NE2	2.82	0.42
1:A:847:ARG:HB2	1:A:851:CYS:HB3	2.00	0.42
1:B:299:THR:O	1:B:302:THR:HG22	2.18	0.42
1:B:309:GLU:HG2	1:B:313:TYR:OH	2.19	0.42
1:B:340:GLU:O	1:B:342:PHE:N	2.52	0.42
1:C:62:VAL:HG13	1:C:267:VAL:O	2.20	0.42
1:C:759:PHE:CZ	1:A:1002:GLN:NE2	2.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:908:GLY:HA3	1:C:1036:GLN:HE22	1.84	0.42
1:A:99:ASN:HD21	1:A:102:ARG:HH11	1.68	0.42
1:A:423:TYR:OH	1:A:512:VAL:HG11	2.19	0.42
1:A:780:GLU:O	1:A:784:GLN:NE2	2.49	0.42
1:A:984:LEU:HB3	1:A:989:ALA:HB2	2.02	0.42
1:B:34:ARG:NH2	1:B:219:GLY:O	2.52	0.42
1:B:387:LEU:HD11	1:B:515:PHE:CE2	2.55	0.42
1:B:1029:MET:O	1:B:1033:VAL:HB	2.19	0.42
1:C:454:ARG:HH12	1:C:467:ASP:HB3	1.84	0.42
1:A:847:ARG:HB2	1:A:851:CYS:HB2	2.00	0.42
1:A:878:LEU:HA	1:A:878:LEU:HD12	1.75	0.42
1:B:858:LEU:HD13	1:B:962:LEU:HD13	2.00	0.42
1:C:99:ASN:HD21	1:C:102:ARG:HH11	1.68	0.42
1:C:404:GLY:HA2	1:C:508:TYR:CD2	2.53	0.42
1:A:48:LEU:HD23	1:A:48:LEU:HA	1.73	0.42
1:A:326:ILE:CD1	1:A:532:ASN:O	2.59	0.42
1:A:1114:ILE:HD13	1:A:1114:ILE:HA	1.93	0.42
1:B:320:VAL:HG22	1:B:629:LEU:HD12	2.02	0.42
1:A:574:ASP:O	1:A:587:ILE:N	2.50	0.42
1:A:983:ARG:HD3	1:B:517:LEU:CD1	2.37	0.42
1:B:364:ASP:O	1:B:367:VAL:HG12	2.20	0.42
1:C:379:CYS:HB3	1:C:382:VAL:O	2.19	0.42
1:C:403:LYS:N	1:C:495:TYR:OH	2.51	0.42
1:A:358:ILE:HB	1:A:395:VAL:HB	2.02	0.42
1:C:1101:HIS:CD2	2:C:1315:NAG:H3	2.54	0.42
1:C:984:LEU:HB3	1:C:989:ALA:HB2	2.02	0.42
1:A:738:CYS:O	1:A:742:ILE:HG13	2.19	0.42
1:A:878:LEU:HD11	1:A:1052:PHE:HB3	2.02	0.42
1:B:327:VAL:HG13	1:B:542:ASN:HB3	2.01	0.42
1:B:375:PHE:CG	1:B:508:TYR:CE1	3.08	0.42
1:B:502:GLY:O	1:B:506:GLN:HG2	2.20	0.42
1:C:97:LYS:O	1:C:97:LYS:NZ	2.52	0.42
1:C:419:ALA:HA	1:C:423:TYR:O	2.19	0.42
1:A:129:LYS:HB3	1:A:169:GLU:HG3	2.01	0.42
1:A:387:LEU:HD11	1:A:515:PHE:CZ	2.55	0.42
1:B:559:PHE:CG	1:B:584:ILE:CD1	3.03	0.42
1:B:984:LEU:HB3	1:B:989:ALA:HB2	2.02	0.42
1:C:187:LYS:H	1:C:210:ILE:CG1	2.33	0.41
1:C:340:GLU:O	1:C:342:PHE:N	2.52	0.41
1:C:347:PHE:CE1	1:C:509:ARG:NH1	2.88	0.41
1:C:527:PRO:O	1:C:528:LYS:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:567:ARG:HE	1:B:42:VAL:HG11	1.85	0.41
1:C:616:ASN:HA	1:C:644:GLN:HE22	1.85	0.41
1:C:742:ILE:CD1	1:C:1001:LEU:HD13	2.49	0.41
1:A:62:VAL:HG13	1:A:267:VAL:O	2.20	0.41
1:A:97:LYS:O	1:A:97:LYS:NZ	2.52	0.41
1:A:130:VAL:CG1	1:A:233:ILE:HD13	2.50	0.41
1:A:318:PHE:CD2	1:A:629:LEU:HD11	2.55	0.41
1:A:375:PHE:CG	1:A:508:TYR:OH	2.68	0.41
1:B:878:LEU:HD11	1:B:1052:PHE:HB3	2.02	0.41
1:C:320:VAL:HG22	1:C:629:LEU:HD12	2.02	0.41
1:C:740:MET:HG2	1:C:857:GLY:HA3	2.01	0.41
1:C:806:LEU:HD21	1:C:882:ILE:HD11	2.02	0.41
1:C:878:LEU:HD11	1:C:1052:PHE:HB3	2.02	0.41
1:C:974:SER:OG	1:C:975:SER:N	2.53	0.41
1:A:1029:MET:O	1:A:1033:VAL:HB	2.19	0.41
1:B:99:ASN:HD21	1:B:102:ARG:HH11	1.68	0.41
1:B:974:SER:OG	1:B:975:SER:N	2.53	0.41
1:C:502:GLY:O	1:C:506:GLN:HG2	2.20	0.41
1:C:703:ASN:HB2	1:B:787:GLN:HB3	2.02	0.41
1:C:1101:HIS:HB2	2:C:1315:NAG:H5	2.02	0.41
1:A:327:VAL:HG13	1:A:542:ASN:HB3	2.01	0.41
1:A:347:PHE:CE1	1:A:509:ARG:NH1	2.88	0.41
1:A:450:ASP:OD1	1:A:450:ASP:N	2.50	0.41
1:A:455:SER:OG	1:A:456:PHE:CD1	2.71	0.41
1:B:108:THR:O	1:B:110:LEU:N	2.50	0.41
1:B:527:PRO:O	1:B:528:LYS:HB3	2.20	0.41
1:B:581:THR:O	1:B:583:GLU:N	2.53	0.41
1:A:527:PRO:O	1:A:528:LYS:HB3	2.20	0.41
1:B:322:PRO:HB2	1:B:540:ASN:ND2	2.36	0.41
1:B:347:PHE:CE1	1:B:509:ARG:NH1	2.88	0.41
1:C:629:LEU:HD23	1:C:631:PRO:HG2	2.03	0.41
1:A:581:THR:C	1:A:583:GLU:N	2.74	0.41
1:B:402:ILE:HD11	1:B:510:VAL:HG21	2.01	0.41
1:B:629:LEU:HD23	1:B:631:PRO:HG2	2.03	0.41
1:C:878:LEU:HA	1:C:878:LEU:HD12	1.75	0.41
1:A:388:ASN:HB3	1:A:527:PRO:HD2	2.02	0.41
1:C:322:PRO:HB2	1:C:540:ASN:ND2	2.36	0.41
1:A:502:GLY:O	1:A:506:GLN:HG2	2.20	0.41
1:A:518:LEU:HD12	1:A:518:LEU:HA	1.83	0.41
1:A:697:MET:H	1:A:697:MET:HG2	1.72	0.41
1:A:974:SER:OG	1:A:975:SER:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1001:LEU:HD12	1:A:1001:LEU:HA	1.77	0.41
1:C:375:PHE:CG	1:C:508:TYR:CE1	3.08	0.41
1:C:588:THR:HG23	1:C:589:PRO:HD2	2.03	0.41
1:C:947:LYS:HE3	1:C:947:LYS:HB2	1.83	0.41
1:A:134:GLN:HG2	1:A:162:SER:HB2	2.02	0.41
1:A:345:THR:OG1	1:A:346:ARG:N	2.54	0.41
1:A:375:PHE:CG	1:A:508:TYR:CE1	3.08	0.41
1:A:1012:LEU:HB3	1:B:1013:ILE:CD1	2.51	0.41
1:B:390:LEU:HA	1:B:390:LEU:HD13	1.87	0.41
1:B:411:ALA:HB3	1:B:414:GLN:CG	2.49	0.41
1:B:780:GLU:O	1:B:784:GLN:NE2	2.49	0.41
1:C:120:VAL:HG13	1:C:127:PHE:HB3	2.02	0.41
1:C:364:ASP:O	1:C:367:VAL:HG12	2.20	0.41
1:C:435:ALA:HB2	1:C:510:VAL:HG22	2.03	0.41
1:C:759:PHE:HZ	1:A:1002:GLN:NE2	2.19	0.41
1:C:795:LYS:HB3	2:C:1313:NAG:H83	2.03	0.41
1:C:1012:LEU:HB3	1:A:1013:ILE:CD1	2.50	0.41
1:C:1072:GLU:N	1:C:1072:GLU:OE1	2.54	0.41
1:A:40:ASP:CG	1:A:41:LYS:H	2.24	0.41
1:A:120:VAL:HG13	1:A:127:PHE:HB3	2.02	0.41
1:A:364:ASP:O	1:A:367:VAL:HG12	2.20	0.41
1:A:383:SER:HB3	1:A:386:LYS:HG2	2.02	0.41
1:A:435:ALA:HB2	1:A:510:VAL:HG22	2.03	0.41
1:A:961:THR:O	1:A:965:GLN:HG2	2.21	0.41
1:B:131:CYS:HB3	1:B:133:PHE:CZ	2.56	0.41
1:B:321:GLN:NE2	1:B:630:THR:OG1	2.49	0.41
1:B:435:ALA:HB2	1:B:510:VAL:HG22	2.03	0.41
1:B:802:PHE:HB3	1:B:805:ILE:HG12	2.02	0.41
1:B:961:THR:O	1:B:965:GLN:HG2	2.21	0.41
1:B:1083:HIS:CE1	1:B:1137:VAL:CG2	2.96	0.41
1:C:535:LYS:O	1:C:536:ASN:CB	2.69	0.41
1:C:714:ILE:HB	1:C:1075:PHE:CE2	2.56	0.41
1:C:759:PHE:HZ	1:A:1002:GLN:CD	2.16	0.41
1:C:961:THR:O	1:C:965:GLN:HG2	2.21	0.41
1:A:365:TYR:CE2	1:A:387:LEU:HG	2.56	0.41
1:A:406:GLU:HA	1:A:409:GLN:HB3	2.03	0.41
1:B:130:VAL:HG21	1:B:231:ILE:HD12	2.03	0.41
1:C:345:THR:OG1	1:C:346:ARG:N	2.54	0.40
1:A:535:LYS:O	1:A:536:ASN:CB	2.69	0.40
1:B:40:ASP:CG	1:B:41:LYS:H	2.24	0.40
1:B:743:CYS:HB3	1:B:749:CYS:HB3	1.82	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1079:PRO:HD2	1:C:1131:GLY:O	2.22	0.40
1:A:358:ILE:HG22	1:A:524:VAL:HG21	2.02	0.40
1:A:857:GLY:C	1:A:858:LEU:HD23	2.42	0.40
1:A:913:GLN:H	1:A:913:GLN:HG3	1.74	0.40
1:B:345:THR:OG1	1:B:346:ARG:N	2.54	0.40
1:B:420:ASP:O	1:B:461:LEU:HB2	2.20	0.40
1:C:168:PHE:CZ	1:C:229:LEU:HB2	2.57	0.40
1:C:411:ALA:HB3	1:C:414:GLN:CG	2.51	0.40
1:C:599:THR:HB	1:C:608:VAL:HG12	2.04	0.40
1:A:759:PHE:HZ	1:B:1002:GLN:CD	2.17	0.40
1:A:959:LEU:O	1:A:963:VAL:HG13	2.22	0.40
1:A:1086:LYS:HE3	1:A:1086:LYS:HB2	1.88	0.40
1:B:231:ILE:HG22	1:B:233:ILE:HG23	2.03	0.40
1:B:599:THR:HB	1:B:608:VAL:HG12	2.04	0.40
1:C:48:LEU:HA	1:C:48:LEU:HD23	1.73	0.40
1:C:374:PHE:CZ	1:C:434:ILE:HG23	2.56	0.40
1:C:375:PHE:CD2	1:C:508:TYR:CZ	3.09	0.40
1:A:206:LYS:HB3	1:A:223:LEU:CD1	2.51	0.40
1:A:349:SER:OG	1:A:451:TYR:HA	2.22	0.40
1:A:554:LYS:HA	1:A:554:LYS:NZ	2.36	0.40
1:A:806:LEU:HD21	1:A:882:ILE:HD11	2.02	0.40
1:B:110:LEU:HD13	1:B:110:LEU:HA	1.94	0.40
1:B:326:ILE:CD1	1:B:532:ASN:O	2.59	0.40
1:B:375:PHE:CD2	1:B:508:TYR:CZ	3.09	0.40
1:B:437:ASN:OD1	1:B:439:ASN:ND2	2.54	0.40
1:B:906:PHE:HE1	1:B:1049:LEU:HD11	1.86	0.40
1:C:25:THR:O	1:C:65:PHE:HA	2.21	0.40
1:C:518:LEU:HD12	1:C:518:LEU:HA	1.83	0.40
1:C:806:LEU:HD23	1:C:878:LEU:HD23	2.03	0.40
1:C:855:PHE:HD1	1:A:589:PRO:HG3	1.86	0.40
1:C:990:GLU:O	1:C:993:ILE:HG22	2.22	0.40
1:B:25:THR:O	1:B:65:PHE:HA	2.21	0.40
1:B:244:LEU:HD23	1:B:244:LEU:HA	1.87	0.40
1:B:716:THR:HA	1:B:1110:TYR:HB3	2.04	0.40
1:B:725:GLU:OE1	1:B:1064:HIS:NE2	2.54	0.40
1:B:1083:HIS:CD2	1:B:1137:VAL:H	2.40	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	1053/1206 (87%)	944 (90%)	91 (9%)	18 (2%)	9	40
1	B	1053/1206 (87%)	944 (90%)	92 (9%)	17 (2%)	9	41
1	C	1053/1206 (87%)	942 (90%)	91 (9%)	20 (2%)	8	37
All	All	3159/3618 (87%)	2830 (90%)	274 (9%)	55 (2%)	13	40

All (55) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	329	PHE
1	C	348	ALA
1	C	527	PRO
1	C	536	ASN
1	C	745	ASP
1	A	348	ALA
1	A	527	PRO
1	A	536	ASN
1	A	745	ASP
1	B	348	ALA
1	B	527	PRO
1	B	536	ASN
1	C	109	THR
1	C	559	PHE
1	C	590	CYS
1	A	109	THR
1	A	559	PHE
1	A	803	SER
1	B	109	THR
1	B	213	GLY
1	B	590	CYS
1	C	161	SER
1	C	165	ASN
1	C	188	ASN

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Mol	Chain	Res	Type
1	C	244	LEU
1	C	518	LEU
1	A	165	ASN
1	A	244	LEU
1	A	518	LEU
1	B	244	LEU
1	B	518	LEU
1	B	593	GLY
1	C	529	LYS
1	C	561	PRO
1	A	529	LYS
1	A	561	PRO
1	A	582	LEU
1	B	529	LYS
1	B	557	LYS
1	B	561	PRO
1	C	330	PRO
1	C	706	ALA
1	A	402	ILE
1	B	402	ILE
1	B	582	LEU
1	C	134	GLN
1	C	844	ILE
1	A	163	ALA
1	A	330	PRO
1	B	164	ASN
1	A	412	PRO
1	B	330	PRO
1	C	942	PRO
1	A	942	PRO
1	B	942	PRO

5.3.2 Protein sidechains [i](#)

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	925/1054 (88%)	887 (96%)	38 (4%)	30	64
1	B	925/1054 (88%)	892 (96%)	33 (4%)	35	68
1	C	925/1054 (88%)	885 (96%)	40 (4%)	29	62
All	All	2775/3162 (88%)	2664 (96%)	111 (4%)	35	64

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	29	THR
1	C	45	SER
1	C	55	PHE
1	C	61	ASN
1	C	127	PHE
1	C	134	GLN
1	C	164	ASN
1	C	168	PHE
1	C	200	TYR
1	C	210	ILE
1	C	216	PHE
1	C	220	PHE
1	C	265	TYR
1	C	284	THR
1	C	308	VAL
1	C	319	ARG
1	C	360	ASN
1	C	374	PHE
1	C	377	PHE
1	C	389	ASP
1	C	538	CYS
1	C	552	LEU
1	C	553	THR
1	C	554	LYS
1	C	567	ARG
1	C	584	ILE
1	C	617	CYS
1	C	709	ASN
1	C	738	CYS
1	C	759	PHE
1	C	760	CYS

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Mol	Chain	Res	Type
1	C	849	LEU
1	C	873	TYR
1	C	939	PHE
1	C	975	SER
1	C	994	ASP
1	C	1037	SER
1	C	1096	VAL
1	C	1132	ILE
1	C	1136	THR
1	A	29	THR
1	A	55	PHE
1	A	61	ASN
1	A	127	PHE
1	A	134	GLN
1	A	160	TYR
1	A	168	PHE
1	A	200	TYR
1	A	210	ILE
1	A	216	PHE
1	A	220	PHE
1	A	265	TYR
1	A	284	THR
1	A	303	LEU
1	A	308	VAL
1	A	319	ARG
1	A	360	ASN
1	A	374	PHE
1	A	377	PHE
1	A	390	LEU
1	A	552	LEU
1	A	553	THR
1	A	554	LYS
1	A	582	LEU
1	A	709	ASN
1	A	738	CYS
1	A	759	PHE
1	A	760	CYS
1	A	873	TYR
1	A	916	LEU
1	A	918	GLU
1	A	939	PHE
1	A	975	SER

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Mol	Chain	Res	Type
1	A	994	ASP
1	A	1037	SER
1	A	1096	VAL
1	A	1132	ILE
1	A	1136	THR
1	B	45	SER
1	B	127	PHE
1	B	134	GLN
1	B	168	PHE
1	B	200	TYR
1	B	216	PHE
1	B	220	PHE
1	B	265	TYR
1	B	284	THR
1	B	302	THR
1	B	303	LEU
1	B	360	ASN
1	B	374	PHE
1	B	377	PHE
1	B	390	LEU
1	B	538	CYS
1	B	552	LEU
1	B	553	THR
1	B	554	LYS
1	B	572	THR
1	B	709	ASN
1	B	759	PHE
1	B	760	CYS
1	B	786	LYS
1	B	873	TYR
1	B	939	PHE
1	B	975	SER
1	B	983	ARG
1	B	994	ASP
1	B	1037	SER
1	B	1096	VAL
1	B	1132	ILE
1	B	1136	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	207	HIS
1	C	394	ASN
1	C	505	HIS
1	C	540	ASN
1	C	556	ASN
1	C	804	GLN
1	C	965	GLN
1	C	1101	HIS
1	A	164	ASN
1	A	207	HIS
1	A	540	ASN
1	A	580	GLN
1	A	965	GLN
1	A	1101	HIS
1	B	207	HIS
1	B	394	ASN
1	B	540	ASN
1	B	556	ASN
1	B	804	GLN
1	B	935	GLN
1	B	965	GLN
1	B	1101	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

55 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	1314	1	14,14,15	0.40	0	17,19,21	0.36	0
2	NAG	C	1307	1	14,14,15	0.39	0	17,19,21	0.74	1 (5%)
2	NAG	B	1318	1	14,14,15	0.38	0	17,19,21	0.49	0
2	NAG	A	1316	1	14,14,15	0.39	0	17,19,21	0.64	0
2	NAG	C	1308	1	14,14,15	0.39	0	17,19,21	0.70	0
2	NAG	C	1302	1	14,14,15	0.39	0	17,19,21	0.75	1 (5%)
2	NAG	B	1309	1	14,14,15	0.39	0	17,19,21	0.38	0
2	NAG	A	1315	1	14,14,15	0.40	0	17,19,21	0.38	0
2	NAG	B	1308	1	14,14,15	0.39	0	17,19,21	0.69	0
2	NAG	C	1306	1	14,14,15	0.39	0	17,19,21	0.35	0
2	NAG	B	1314	1	14,14,15	0.39	0	17,19,21	0.36	0
2	NAG	A	1318	1	14,14,15	0.38	0	17,19,21	0.48	0
2	NAG	C	1311	1	14,14,15	0.38	0	17,19,21	0.54	0
2	NAG	A	1305	1	14,14,15	0.41	0	17,19,21	0.69	1 (5%)
2	NAG	B	1307	1	14,14,15	0.40	0	17,19,21	0.74	1 (5%)
2	NAG	C	1309	1	14,14,15	0.40	0	17,19,21	0.36	0
2	NAG	C	1317	1	14,14,15	0.40	0	17,19,21	0.81	0
2	NAG	B	1304	1	14,14,15	0.38	0	17,19,21	0.54	0
2	NAG	C	1301	1	14,14,15	0.40	0	17,19,21	0.77	0
2	NAG	A	1314	1	14,14,15	0.41	0	17,19,21	0.44	0
3	IDU	B	1317	-	15,15,17	0.89	1 (6%)	15,22,26	1.03	1 (6%)
2	NAG	B	1303	1	14,14,15	0.39	0	17,19,21	0.74	0
2	NAG	B	1305	1	14,14,15	0.42	0	17,19,21	0.70	1 (5%)
2	NAG	A	1311	1	14,14,15	0.37	0	17,19,21	0.54	0
2	NAG	A	1313	1	14,14,15	0.40	0	17,19,21	0.81	0
2	NAG	B	1315	1	14,14,15	0.39	0	17,19,21	0.39	0
2	NAG	C	1304	1	14,14,15	0.37	0	17,19,21	0.55	0
2	NAG	B	1312	1	14,14,15	0.40	0	17,19,21	0.81	0
2	NAG	A	1307	1	14,14,15	0.38	0	17,19,21	0.75	1 (5%)
2	NAG	C	1310	1	14,14,15	0.38	0	17,19,21	0.61	0
2	NAG	A	1303	1	14,14,15	0.39	0	17,19,21	0.74	0
2	NAG	B	1316	1	14,14,15	0.39	0	17,19,21	0.64	0
2	NAG	B	1311	1	14,14,15	0.38	0	17,19,21	0.54	0
2	NAG	A	1308	1	14,14,15	0.38	0	17,19,21	0.69	0
2	NAG	A	1317	1	14,14,15	0.41	0	17,19,21	0.81	0
2	NAG	A	1306	1	14,14,15	0.41	0	17,19,21	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	1309	1	14,14,15	0.40	0	17,19,21	0.52	0
2	NAG	C	1316	1	14,14,15	0.39	0	17,19,21	0.64	0
2	NAG	C	1318	1	14,14,15	0.37	0	17,19,21	0.48	0
2	NAG	A	1310	1	14,14,15	0.39	0	17,19,21	0.61	0
2	NAG	A	1312	1	14,14,15	0.39	0	17,19,21	0.81	0
2	NAG	C	1312	1	14,14,15	0.39	0	17,19,21	0.81	0
2	NAG	A	1304	1	14,14,15	0.35	0	17,19,21	0.55	0
2	NAG	B	1302	1	14,14,15	0.39	0	17,19,21	0.66	1 (5%)
2	NAG	B	1306	1	14,14,15	0.41	0	17,19,21	0.70	0
2	NAG	B	1301	1	14,14,15	0.38	0	17,19,21	0.77	0
2	NAG	A	1301	1	14,14,15	0.40	0	17,19,21	0.76	0
2	NAG	C	1303	1	14,14,15	0.39	0	17,19,21	0.74	0
2	NAG	C	1305	1	14,14,15	0.42	0	17,19,21	0.70	1 (5%)
2	NAG	B	1310	1	14,14,15	0.39	0	17,19,21	0.62	0
2	NAG	C	1313	1	14,14,15	0.38	0	17,19,21	0.45	0
2	NAG	A	1302	1	14,14,15	0.37	0	17,19,21	0.97	1 (5%)
2	NAG	C	1315	1	14,14,15	0.40	0	17,19,21	0.41	0
2	NAG	B	1319	1	14,14,15	0.38	0	17,19,21	0.45	0
2	NAG	B	1313	1	14,14,15	0.41	0	17,19,21	0.81	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1314	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1307	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1318	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1316	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1308	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1302	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1309	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1315	1	-	3/6/23/26	0/1/1/1
2	NAG	B	1308	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1306	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1314	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1318	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1311	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1305	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1307	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1309	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1317	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1304	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1301	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1314	1	-	0/6/23/26	0/1/1/1
3	IDU	B	1317	-	-	3/9/22/29	1/1/1/1
2	NAG	B	1303	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1305	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1311	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1313	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1315	1	-	3/6/23/26	0/1/1/1
2	NAG	C	1304	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1312	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1307	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1310	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1303	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1316	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1311	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1308	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1317	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1306	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1309	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1316	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1318	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1310	1	-	1/6/23/26	0/1/1/1
2	NAG	A	1312	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1312	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1304	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1302	1	-	1/6/23/26	0/1/1/1
2	NAG	B	1306	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1301	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1301	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1303	1	-	0/6/23/26	0/1/1/1
2	NAG	C	1305	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1310	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1313	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1302	1	-	1/6/23/26	0/1/1/1
2	NAG	C	1315	1	-	3/6/23/26	0/1/1/1
2	NAG	B	1319	1	-	0/6/23/26	0/1/1/1
2	NAG	B	1313	1	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1317	IDU	O6-C6	-2.89	1.21	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	1302	NAG	C2-N2-C7	3.53	127.92	122.90
2	C	1302	NAG	C2-N2-C7	2.59	126.60	122.90
3	B	1317	IDU	O6-C6-C5	2.36	119.77	113.03
2	C	1307	NAG	C1-O5-C5	2.28	115.28	112.19
2	A	1307	NAG	C1-O5-C5	2.27	115.27	112.19
2	B	1307	NAG	C1-O5-C5	2.26	115.25	112.19
2	B	1305	NAG	C1-O5-C5	2.18	115.15	112.19
2	C	1305	NAG	C1-O5-C5	2.16	115.11	112.19
2	A	1305	NAG	C1-O5-C5	2.14	115.09	112.19
2	B	1302	NAG	C2-N2-C7	2.13	125.93	122.90

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1317	IDU	C3-C2-O2-S
3	B	1317	IDU	O5-C5-C6-O61
2	C	1315	NAG	C8-C7-N2-C2
2	C	1315	NAG	O7-C7-N2-C2
2	A	1315	NAG	C8-C7-N2-C2
2	A	1315	NAG	O7-C7-N2-C2
2	B	1315	NAG	C8-C7-N2-C2
2	B	1315	NAG	O7-C7-N2-C2
2	B	1309	NAG	O5-C5-C6-O6
2	C	1309	NAG	O5-C5-C6-O6
2	C	1315	NAG	O5-C5-C6-O6
2	A	1309	NAG	O5-C5-C6-O6
2	A	1315	NAG	O5-C5-C6-O6
2	B	1315	NAG	O5-C5-C6-O6
2	C	1310	NAG	C1-C2-N2-C7
2	A	1310	NAG	C1-C2-N2-C7
2	B	1310	NAG	C1-C2-N2-C7
2	C	1302	NAG	C3-C2-N2-C7
2	A	1302	NAG	C3-C2-N2-C7
2	B	1302	NAG	C3-C2-N2-C7
3	B	1317	IDU	O5-C5-C6-O6

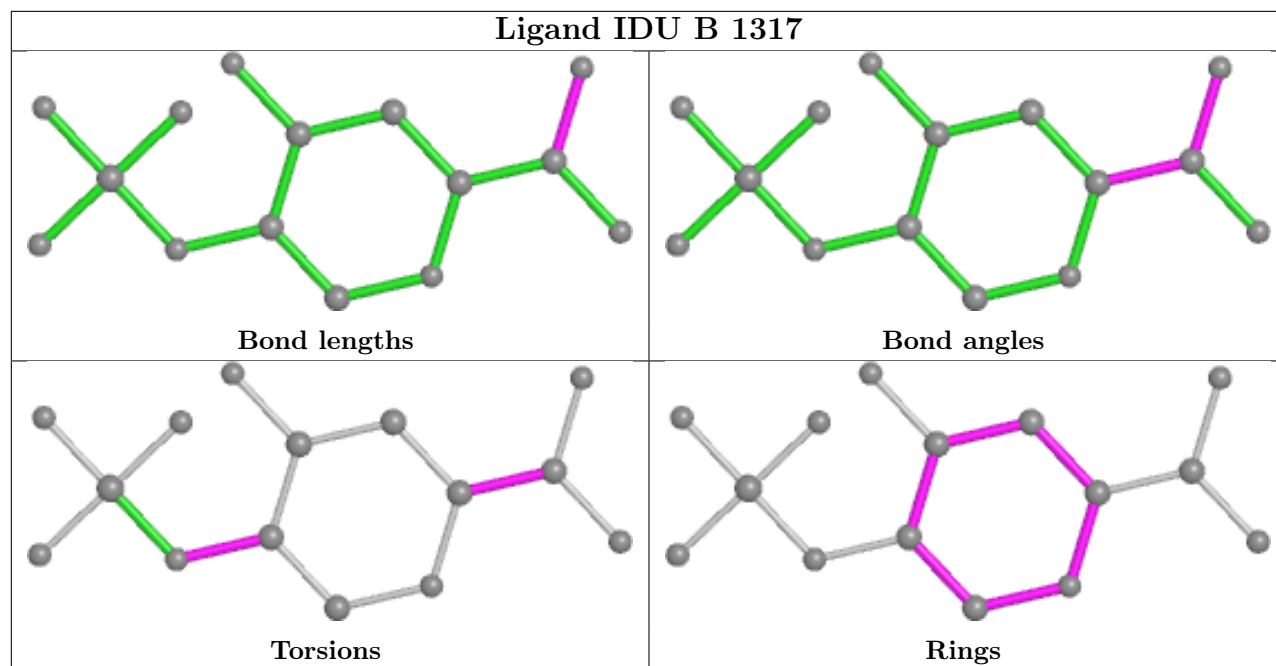
All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1317	IDU	C1-C2-C3-C4-C5-O5

13 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1302	NAG	1	0
2	A	1315	NAG	3	0
2	C	1311	NAG	2	0
2	C	1301	NAG	1	0
3	B	1317	IDU	1	0
2	A	1311	NAG	1	0
2	B	1315	NAG	3	0
2	B	1311	NAG	1	0
2	A	1309	NAG	1	0
2	B	1301	NAG	1	0
2	A	1301	NAG	1	0
2	C	1313	NAG	3	0
2	C	1315	NAG	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

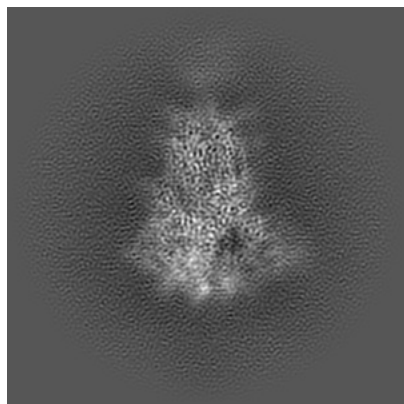
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-38682. These allow visual inspection of the internal detail of the map and identification of artifacts.

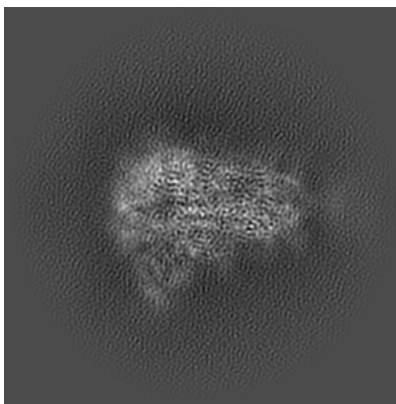
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

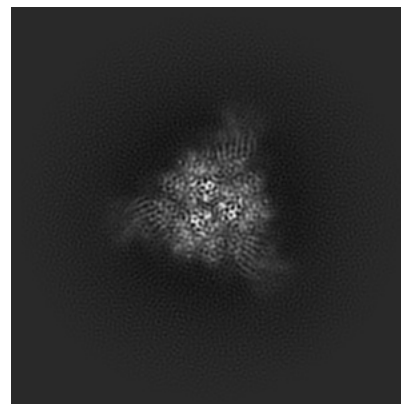
6.1.1 Primary map



X

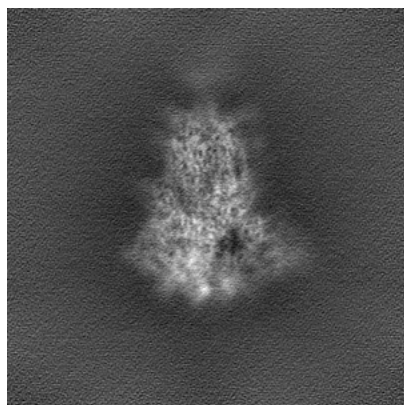


Y

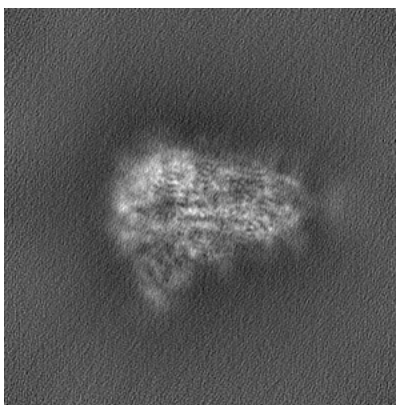


Z

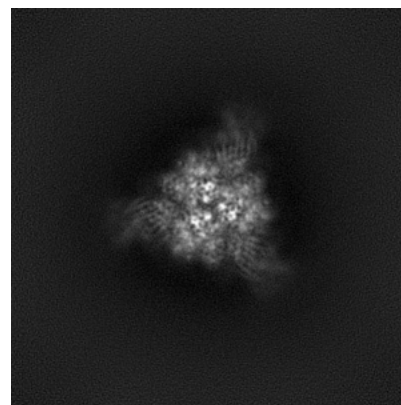
6.1.2 Raw map



X



Y

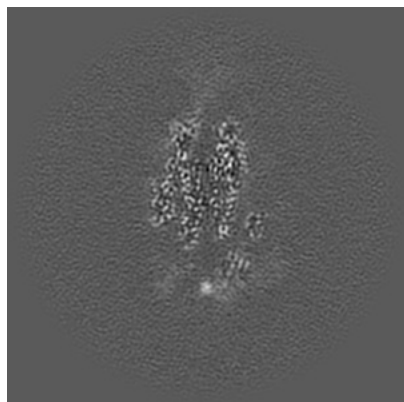


Z

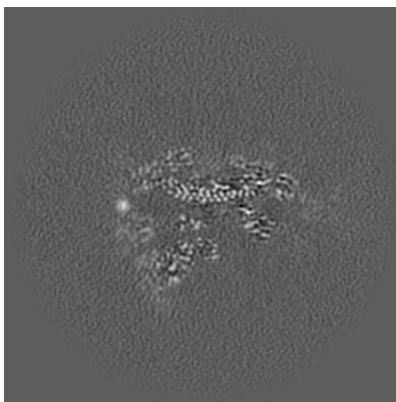
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

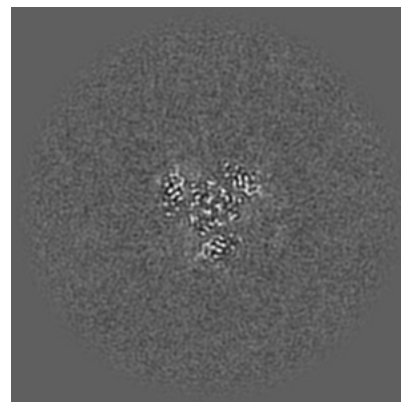
6.2.1 Primary map



X Index: 150

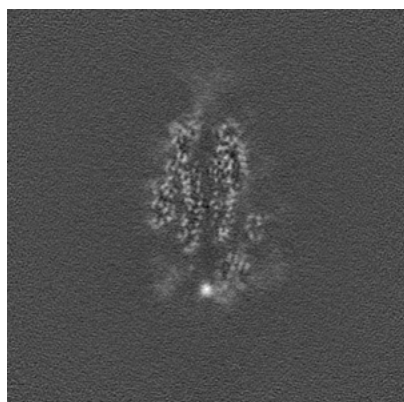


Y Index: 150

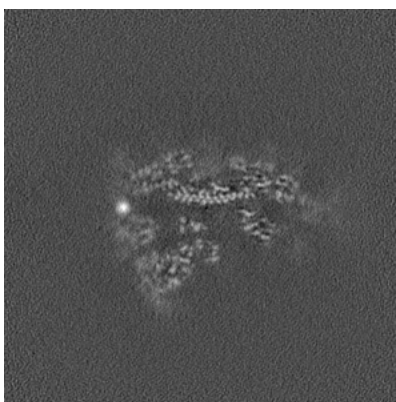


Z Index: 150

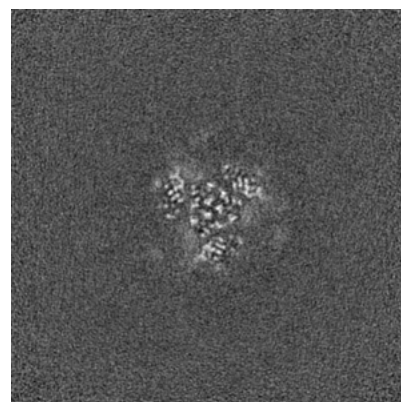
6.2.2 Raw map



X Index: 150



Y Index: 150

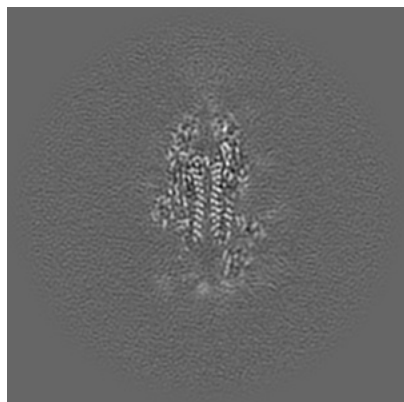


Z Index: 150

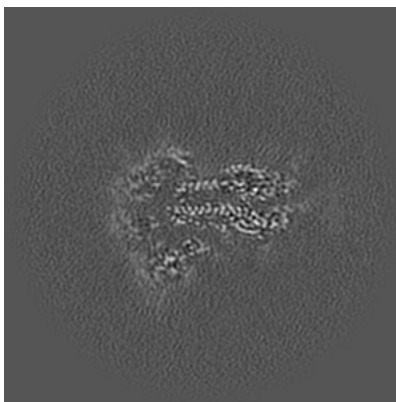
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

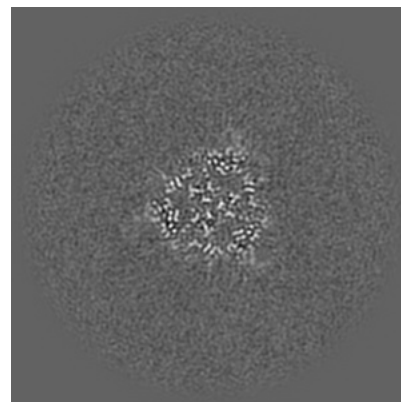
6.3.1 Primary map



X Index: 146

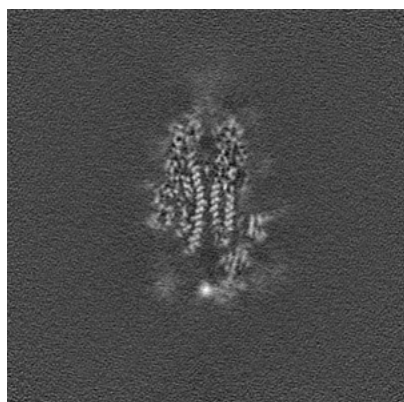


Y Index: 143

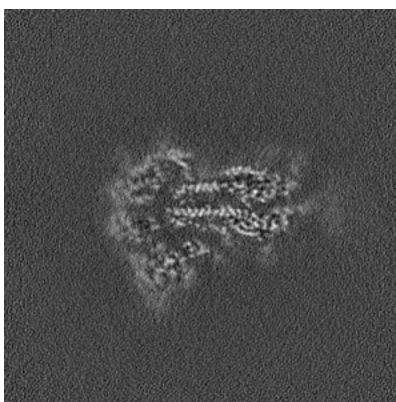


Z Index: 139

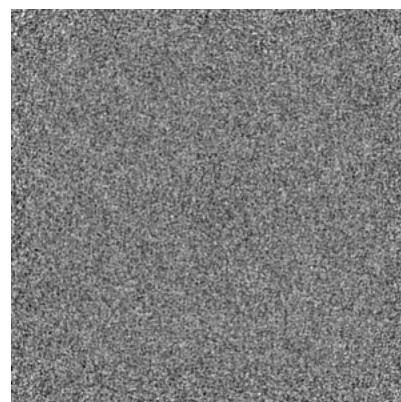
6.3.2 Raw map



X Index: 148



Y Index: 143

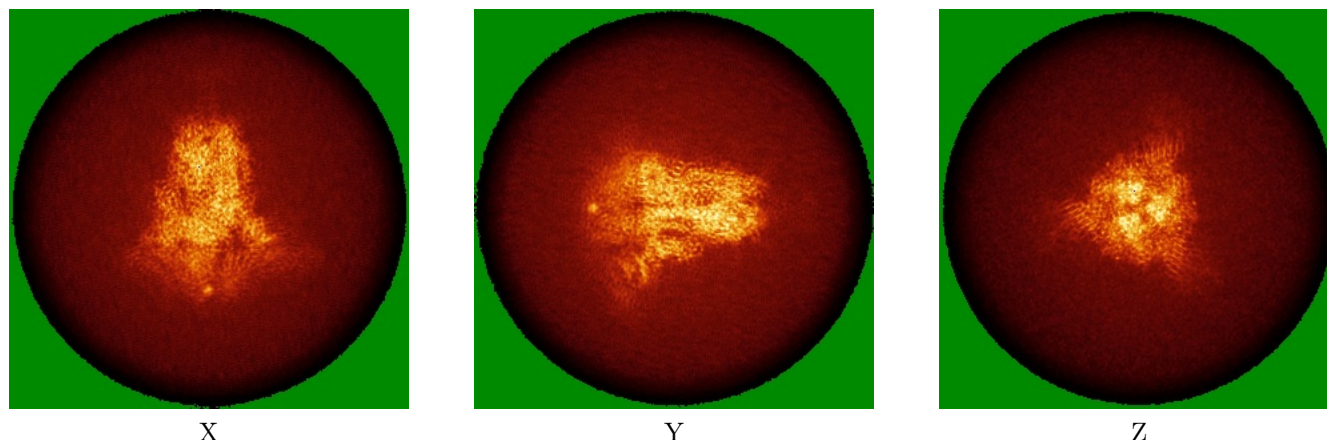


Z Index: 0

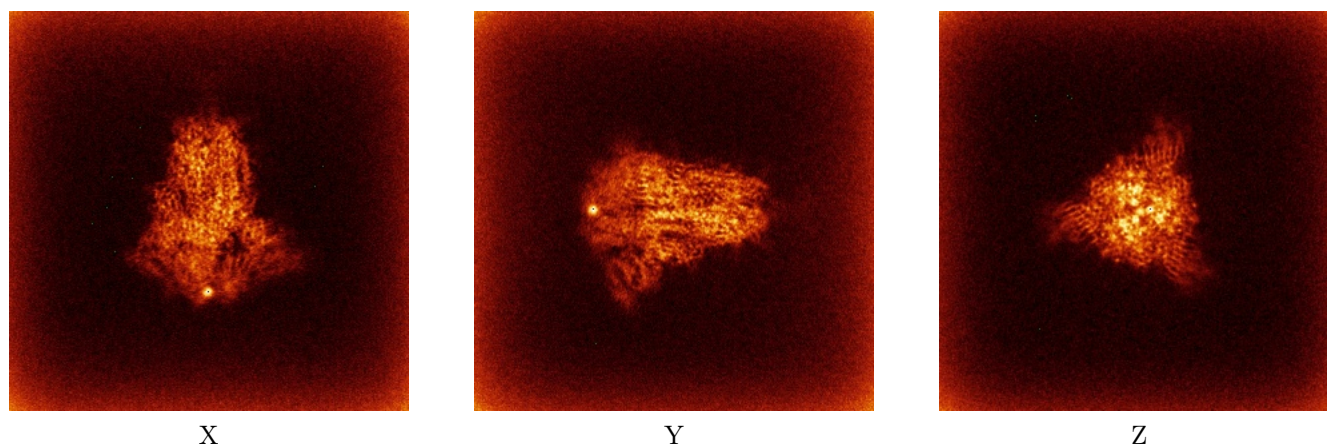
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



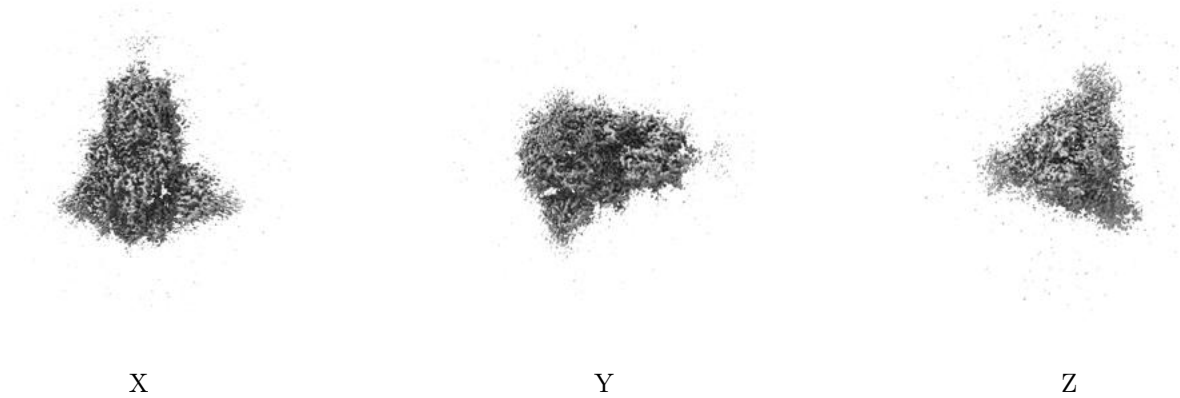
6.4.2 Raw map



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.

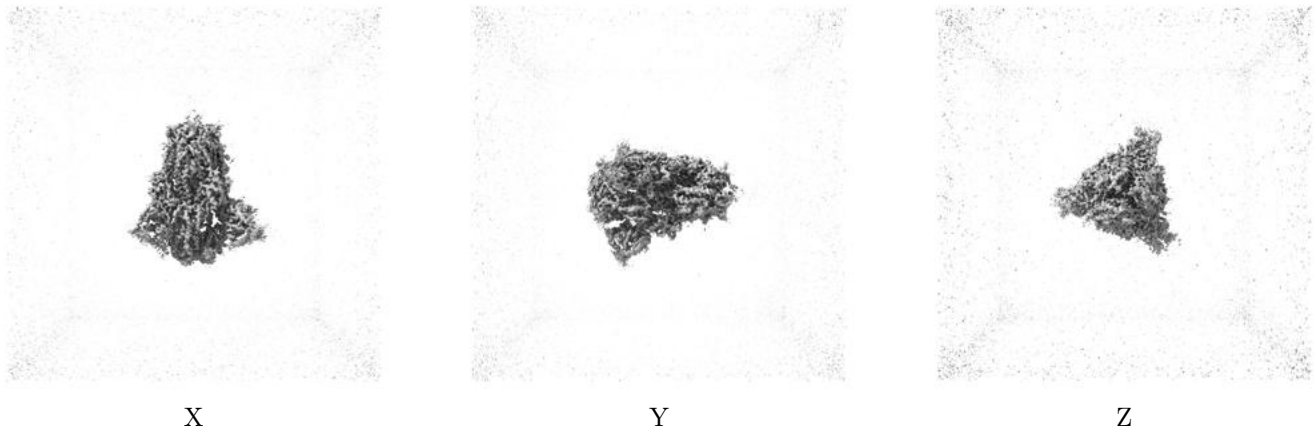
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.35. 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.

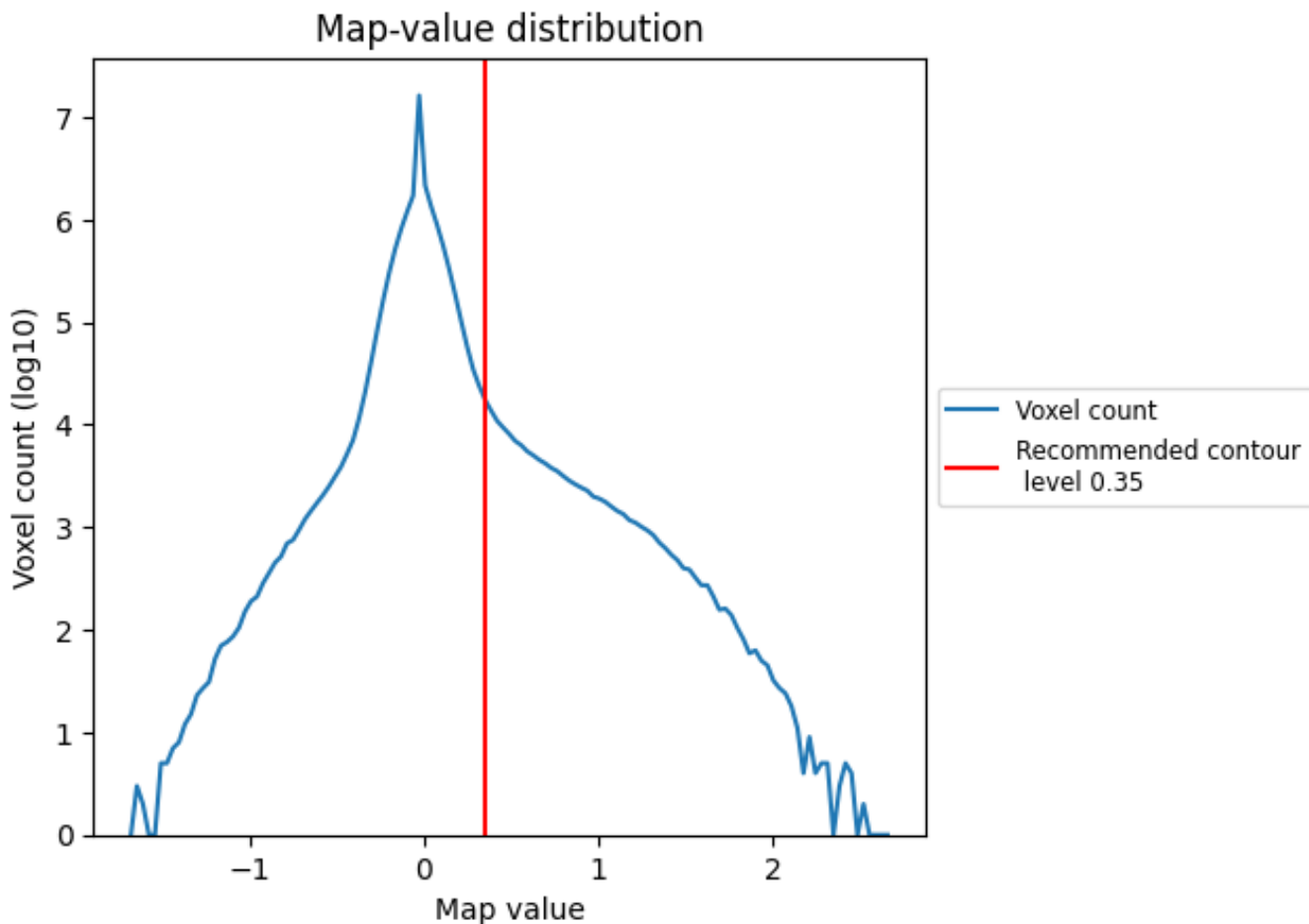
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

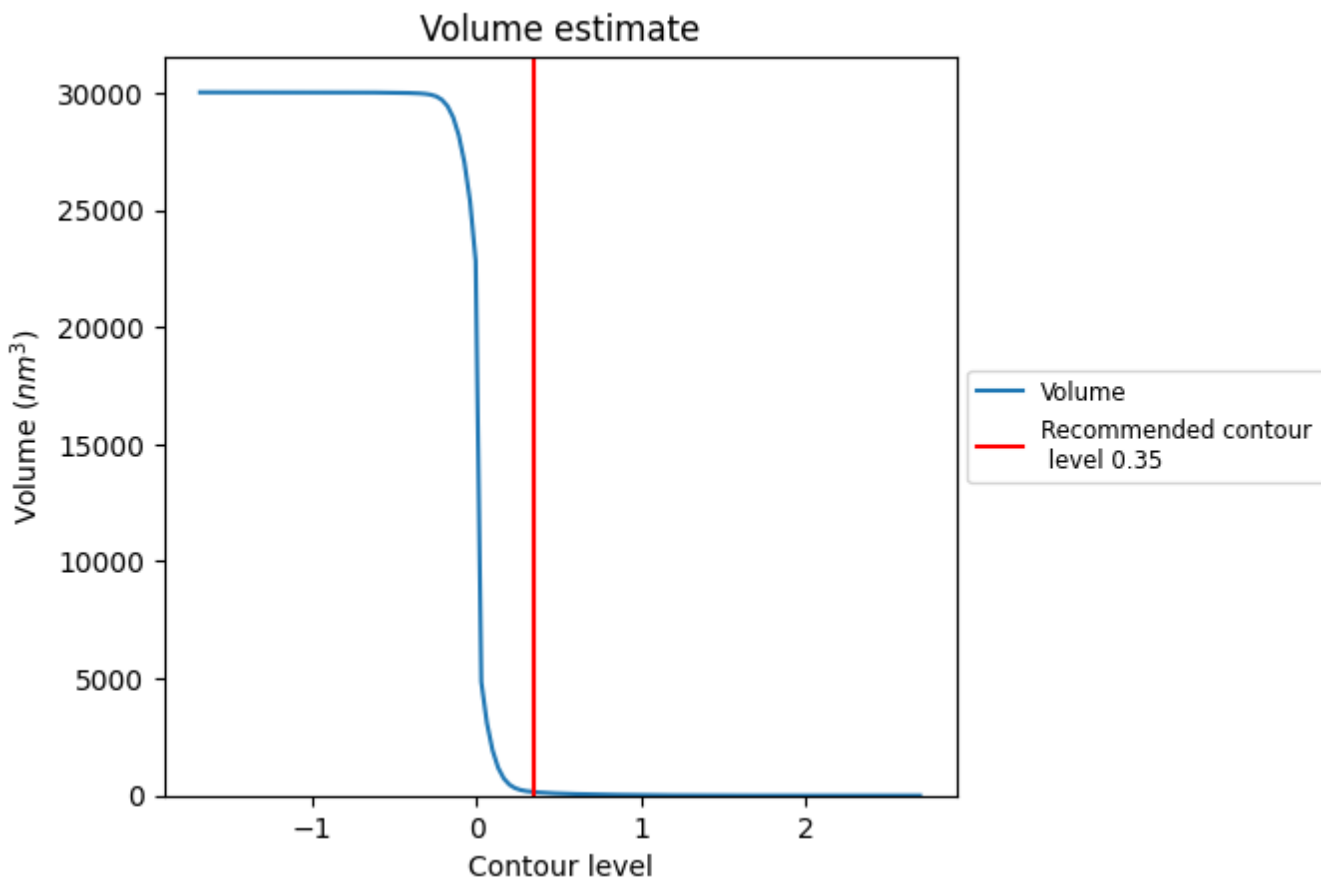
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

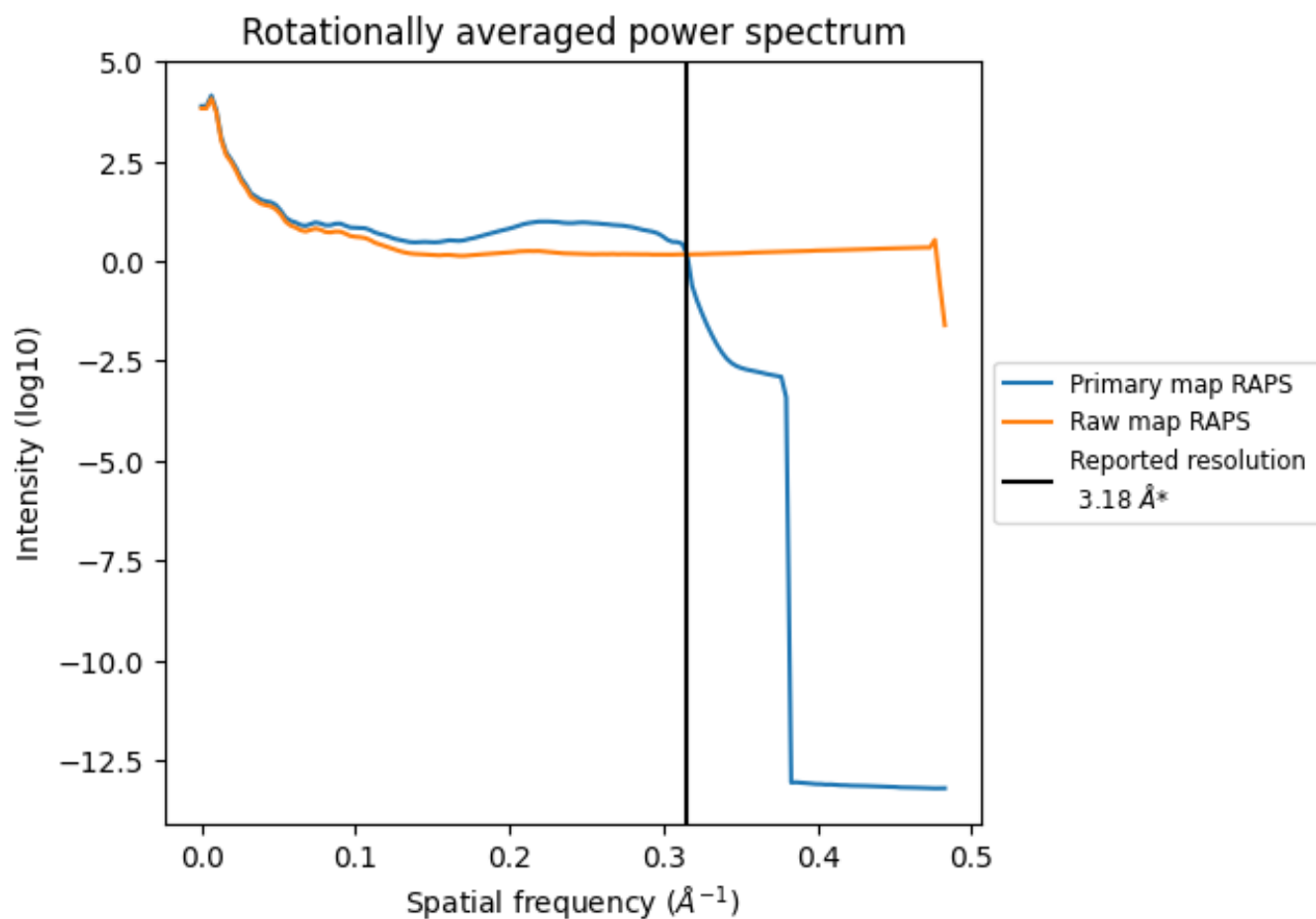
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 150 nm³; this corresponds to an approximate mass of 136 kDa.

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

7.3 Rotationally averaged power spectrum [i](#)

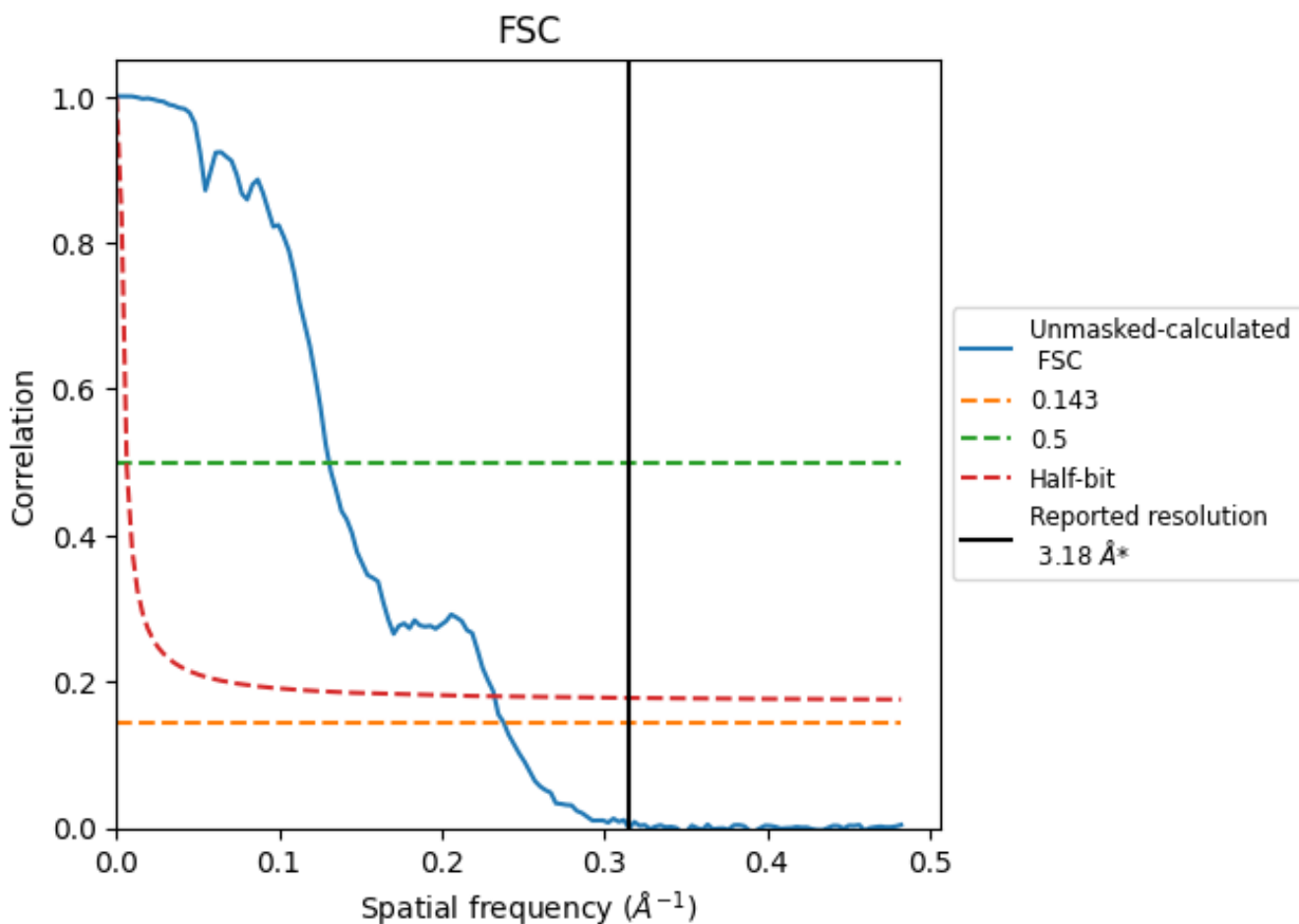


*Reported resolution corresponds to spatial frequency of 0.314 Å⁻¹

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 0.314 Å⁻¹

8.2 Resolution estimates [i](#)

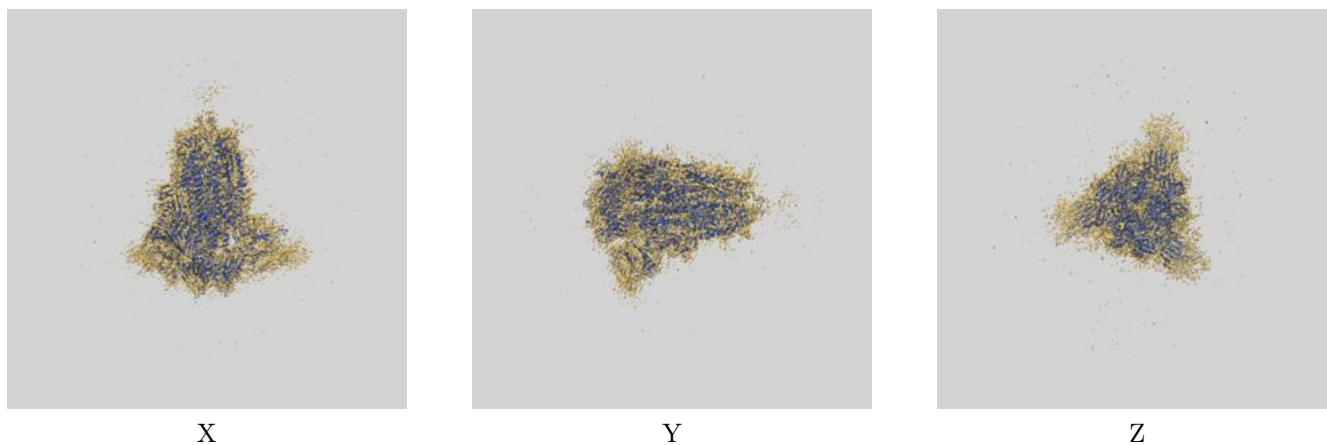
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.18	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.20	7.65	4.30

*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 4.20 differs from the reported value 3.18 by more than 10 %

9 Map-model fit [i](#)

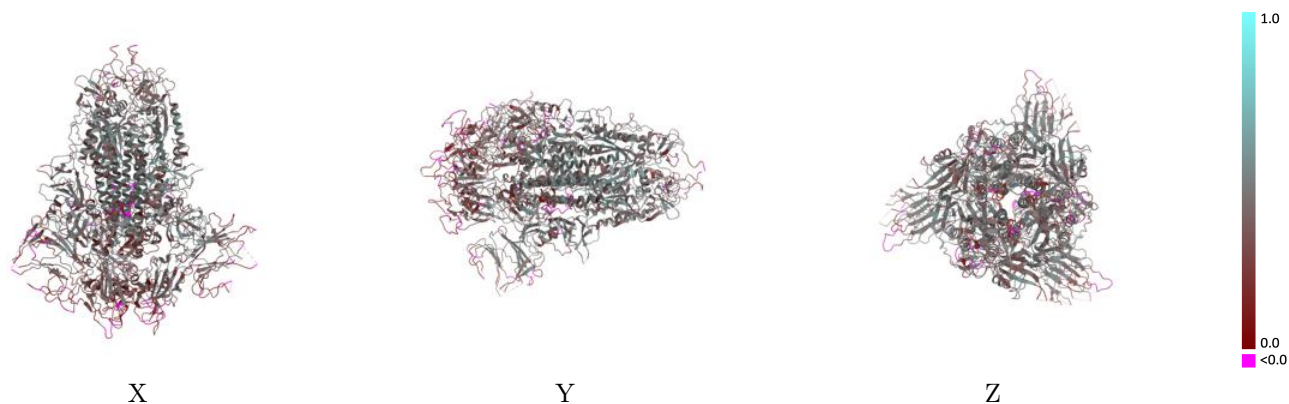
This section contains information regarding the fit between EMDB map EMD-38682 and PDB model 8XUS. Per-residue inclusion information can be found in section 3 on page 13.

9.1 Map-model overlay [i](#)



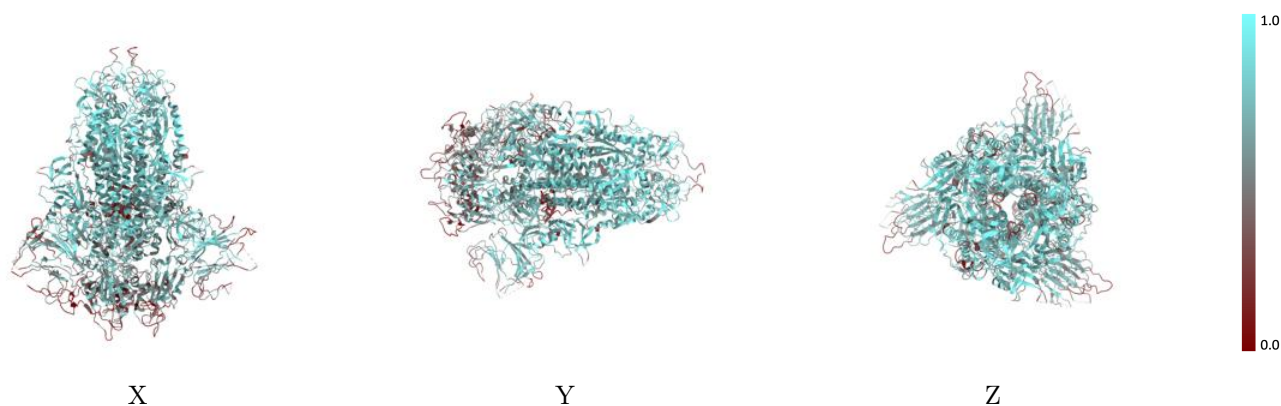
The images above show the 3D surface view of the map at the recommended contour level 0.35 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



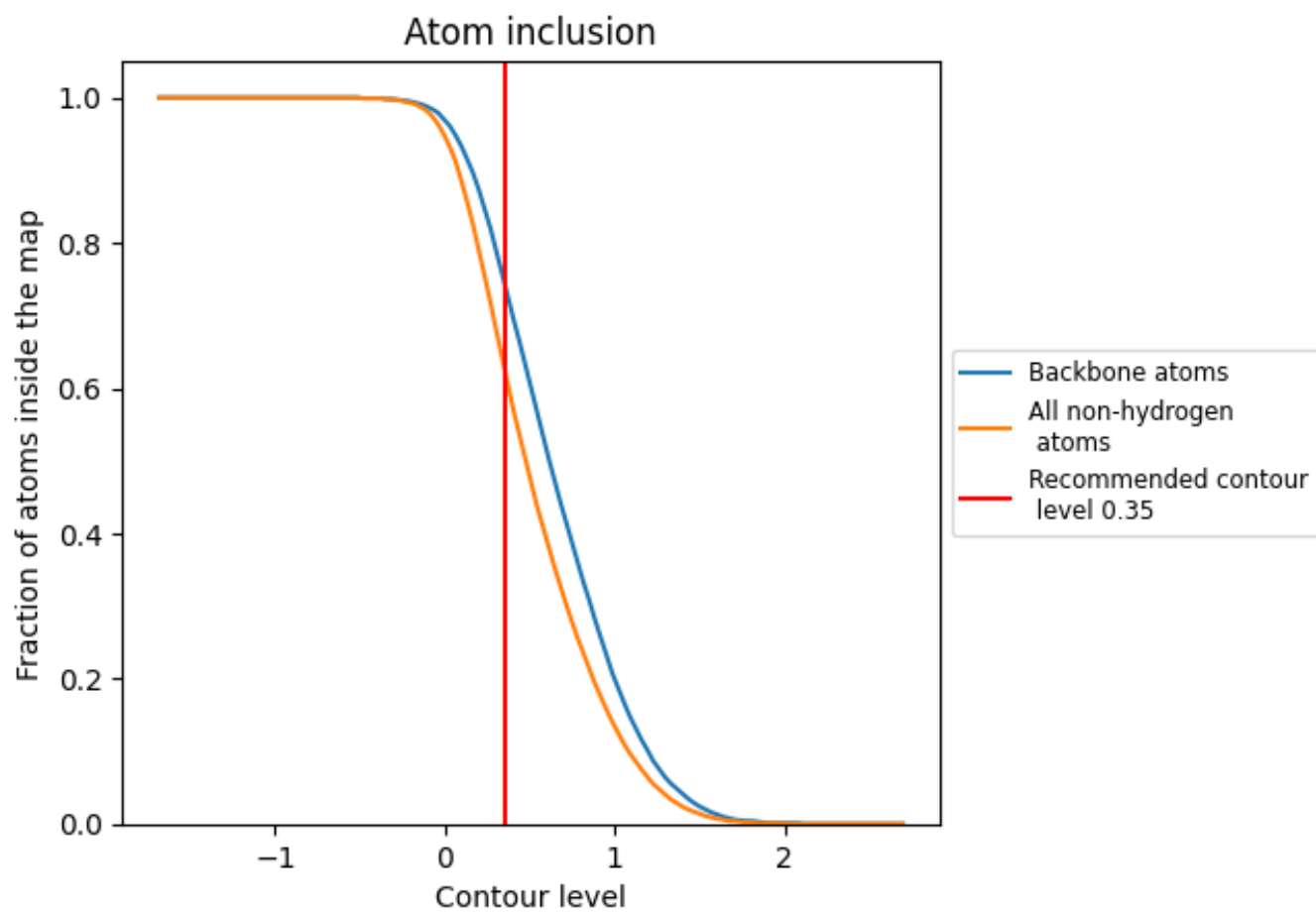
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.35).









9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 63% 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 (0.35) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6270	 0.3720
A	 0.6200	 0.3680
B	 0.6170	 0.3660
C	 0.6440	 0.3820

