



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 12:25 PM UTC

PDB ID : 8T21 / pdb_00008t21
EMDB ID : EMD-40977
Title : Cryo-EM structure of mink variant Y453F trimeric spike protein
Authors : Ahn, H.M.; Calderon, B.; Fan, X.; Gao, Y.; Horgan, N.; Zhou, B.; Liang, B.
Deposited on : 2023-06-05
Resolution : 3.60 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

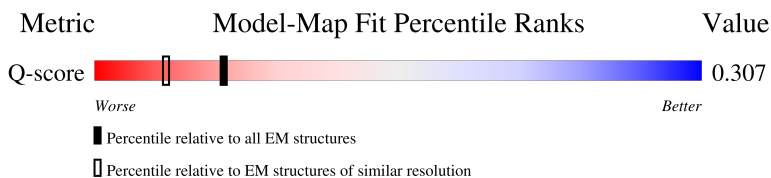
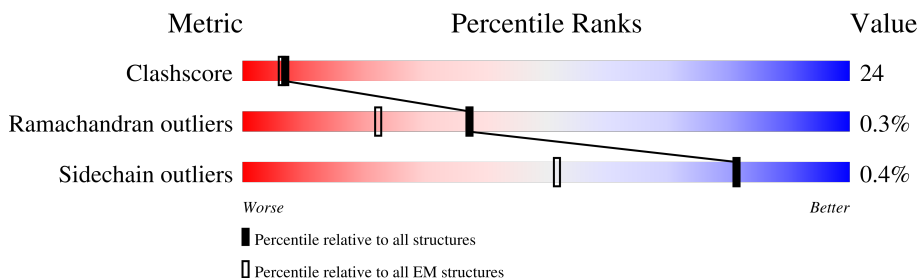
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 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	12797 (3.10 - 4.10)

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	1269	22% 50% 26% •• 21%
1	B	1269	22% 53% 25% • 21%
1	C	1269	11% 55% 24% • 21%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 23430 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
1	A	1001	Total	C	N	O	S	0	0
			7810	4990	1297	1488	35		
1	B	1001	Total	C	N	O	S	0	0
			7810	4990	1297	1488	35		
1	C	1001	Total	C	N	O	S	0	0
			7810	4990	1297	1488	35		

There are 225 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	HIS	deletion	UNP A0A6H1PJZ3
A	?	-	VAL	deletion	UNP A0A6H1PJZ3
A	453	PHE	TYR	variant	UNP A0A6H1PJZ3
A	682	GLY	ARG	engineered mutation	UNP A0A6H1PJZ3
A	683	SER	ARG	engineered mutation	UNP A0A6H1PJZ3
A	685	SER	ARG	engineered mutation	UNP A0A6H1PJZ3
A	817	PRO	PHE	engineered mutation	UNP A0A6H1PJZ3
A	892	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
A	899	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
A	942	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
A	986	PRO	LYS	engineered mutation	UNP A0A6H1PJZ3
A	987	PRO	VAL	engineered mutation	UNP A0A6H1PJZ3
A	1209	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1210	SER	-	expression tag	UNP A0A6H1PJZ3
A	1211	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1212	SER	-	expression tag	UNP A0A6H1PJZ3
A	1213	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1214	SER	-	expression tag	UNP A0A6H1PJZ3
A	1215	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1216	SER	-	expression tag	UNP A0A6H1PJZ3
A	1217	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1218	TYR	-	expression tag	UNP A0A6H1PJZ3
A	1219	ILE	-	expression tag	UNP A0A6H1PJZ3
A	1220	PRO	-	expression tag	UNP A0A6H1PJZ3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1221	GLU	-	expression tag	UNP A0A6H1PJZ3
A	1222	ALA	-	expression tag	UNP A0A6H1PJZ3
A	1223	PRO	-	expression tag	UNP A0A6H1PJZ3
A	1224	ARG	-	expression tag	UNP A0A6H1PJZ3
A	1225	ASP	-	expression tag	UNP A0A6H1PJZ3
A	1226	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1227	GLN	-	expression tag	UNP A0A6H1PJZ3
A	1228	ALA	-	expression tag	UNP A0A6H1PJZ3
A	1229	TYR	-	expression tag	UNP A0A6H1PJZ3
A	1230	VAL	-	expression tag	UNP A0A6H1PJZ3
A	1231	ARG	-	expression tag	UNP A0A6H1PJZ3
A	1232	LYS	-	expression tag	UNP A0A6H1PJZ3
A	1233	ASP	-	expression tag	UNP A0A6H1PJZ3
A	1234	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1235	GLU	-	expression tag	UNP A0A6H1PJZ3
A	1236	TRP	-	expression tag	UNP A0A6H1PJZ3
A	1237	VAL	-	expression tag	UNP A0A6H1PJZ3
A	1238	LEU	-	expression tag	UNP A0A6H1PJZ3
A	1239	LEU	-	expression tag	UNP A0A6H1PJZ3
A	1240	SER	-	expression tag	UNP A0A6H1PJZ3
A	1241	THR	-	expression tag	UNP A0A6H1PJZ3
A	1242	PHE	-	expression tag	UNP A0A6H1PJZ3
A	1243	LEU	-	expression tag	UNP A0A6H1PJZ3
A	1244	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1245	SER	-	expression tag	UNP A0A6H1PJZ3
A	1246	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1247	SER	-	expression tag	UNP A0A6H1PJZ3
A	1248	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1249	SER	-	expression tag	UNP A0A6H1PJZ3
A	1250	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1251	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1252	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1253	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1254	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1255	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1256	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1257	GLY	-	expression tag	UNP A0A6H1PJZ3
A	1258	LEU	-	expression tag	UNP A0A6H1PJZ3
A	1259	ASN	-	expression tag	UNP A0A6H1PJZ3
A	1260	ASP	-	expression tag	UNP A0A6H1PJZ3
A	1261	ILE	-	expression tag	UNP A0A6H1PJZ3
A	1262	PHE	-	expression tag	UNP A0A6H1PJZ3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1263	GLU	-	expression tag	UNP A0A6H1PJZ3
A	1264	ALA	-	expression tag	UNP A0A6H1PJZ3
A	1265	GLN	-	expression tag	UNP A0A6H1PJZ3
A	1266	LYS	-	expression tag	UNP A0A6H1PJZ3
A	1267	ILE	-	expression tag	UNP A0A6H1PJZ3
A	1268	GLU	-	expression tag	UNP A0A6H1PJZ3
A	1269	TRP	-	expression tag	UNP A0A6H1PJZ3
A	1270	HIS	-	expression tag	UNP A0A6H1PJZ3
A	1271	GLU	-	expression tag	UNP A0A6H1PJZ3
B	?	-	HIS	deletion	UNP A0A6H1PJZ3
B	?	-	VAL	deletion	UNP A0A6H1PJZ3
B	453	PHE	TYR	variant	UNP A0A6H1PJZ3
B	682	GLY	ARG	engineered mutation	UNP A0A6H1PJZ3
B	683	SER	ARG	engineered mutation	UNP A0A6H1PJZ3
B	685	SER	ARG	engineered mutation	UNP A0A6H1PJZ3
B	817	PRO	PHE	engineered mutation	UNP A0A6H1PJZ3
B	892	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
B	899	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
B	942	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
B	986	PRO	LYS	engineered mutation	UNP A0A6H1PJZ3
B	987	PRO	VAL	engineered mutation	UNP A0A6H1PJZ3
B	1209	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1210	SER	-	expression tag	UNP A0A6H1PJZ3
B	1211	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1212	SER	-	expression tag	UNP A0A6H1PJZ3
B	1213	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1214	SER	-	expression tag	UNP A0A6H1PJZ3
B	1215	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1216	SER	-	expression tag	UNP A0A6H1PJZ3
B	1217	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1218	TYR	-	expression tag	UNP A0A6H1PJZ3
B	1219	ILE	-	expression tag	UNP A0A6H1PJZ3
B	1220	PRO	-	expression tag	UNP A0A6H1PJZ3
B	1221	GLU	-	expression tag	UNP A0A6H1PJZ3
B	1222	ALA	-	expression tag	UNP A0A6H1PJZ3
B	1223	PRO	-	expression tag	UNP A0A6H1PJZ3
B	1224	ARG	-	expression tag	UNP A0A6H1PJZ3
B	1225	ASP	-	expression tag	UNP A0A6H1PJZ3
B	1226	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1227	GLN	-	expression tag	UNP A0A6H1PJZ3
B	1228	ALA	-	expression tag	UNP A0A6H1PJZ3
B	1229	TYR	-	expression tag	UNP A0A6H1PJZ3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1230	VAL	-	expression tag	UNP A0A6H1PJZ3
B	1231	ARG	-	expression tag	UNP A0A6H1PJZ3
B	1232	LYS	-	expression tag	UNP A0A6H1PJZ3
B	1233	ASP	-	expression tag	UNP A0A6H1PJZ3
B	1234	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1235	GLU	-	expression tag	UNP A0A6H1PJZ3
B	1236	TRP	-	expression tag	UNP A0A6H1PJZ3
B	1237	VAL	-	expression tag	UNP A0A6H1PJZ3
B	1238	LEU	-	expression tag	UNP A0A6H1PJZ3
B	1239	LEU	-	expression tag	UNP A0A6H1PJZ3
B	1240	SER	-	expression tag	UNP A0A6H1PJZ3
B	1241	THR	-	expression tag	UNP A0A6H1PJZ3
B	1242	PHE	-	expression tag	UNP A0A6H1PJZ3
B	1243	LEU	-	expression tag	UNP A0A6H1PJZ3
B	1244	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1245	SER	-	expression tag	UNP A0A6H1PJZ3
B	1246	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1247	SER	-	expression tag	UNP A0A6H1PJZ3
B	1248	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1249	SER	-	expression tag	UNP A0A6H1PJZ3
B	1250	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1251	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1252	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1253	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1254	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1255	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1256	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1257	GLY	-	expression tag	UNP A0A6H1PJZ3
B	1258	LEU	-	expression tag	UNP A0A6H1PJZ3
B	1259	ASN	-	expression tag	UNP A0A6H1PJZ3
B	1260	ASP	-	expression tag	UNP A0A6H1PJZ3
B	1261	ILE	-	expression tag	UNP A0A6H1PJZ3
B	1262	PHE	-	expression tag	UNP A0A6H1PJZ3
B	1263	GLU	-	expression tag	UNP A0A6H1PJZ3
B	1264	ALA	-	expression tag	UNP A0A6H1PJZ3
B	1265	GLN	-	expression tag	UNP A0A6H1PJZ3
B	1266	LYS	-	expression tag	UNP A0A6H1PJZ3
B	1267	ILE	-	expression tag	UNP A0A6H1PJZ3
B	1268	GLU	-	expression tag	UNP A0A6H1PJZ3
B	1269	TRP	-	expression tag	UNP A0A6H1PJZ3
B	1270	HIS	-	expression tag	UNP A0A6H1PJZ3
B	1271	GLU	-	expression tag	UNP A0A6H1PJZ3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	HIS	deletion	UNP A0A6H1PJZ3
C	?	-	VAL	deletion	UNP A0A6H1PJZ3
C	453	PHE	TYR	variant	UNP A0A6H1PJZ3
C	682	GLY	ARG	engineered mutation	UNP A0A6H1PJZ3
C	683	SER	ARG	engineered mutation	UNP A0A6H1PJZ3
C	685	SER	ARG	engineered mutation	UNP A0A6H1PJZ3
C	817	PRO	PHE	engineered mutation	UNP A0A6H1PJZ3
C	892	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
C	899	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
C	942	PRO	ALA	engineered mutation	UNP A0A6H1PJZ3
C	986	PRO	LYS	engineered mutation	UNP A0A6H1PJZ3
C	987	PRO	VAL	engineered mutation	UNP A0A6H1PJZ3
C	1209	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1210	SER	-	expression tag	UNP A0A6H1PJZ3
C	1211	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1212	SER	-	expression tag	UNP A0A6H1PJZ3
C	1213	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1214	SER	-	expression tag	UNP A0A6H1PJZ3
C	1215	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1216	SER	-	expression tag	UNP A0A6H1PJZ3
C	1217	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1218	TYR	-	expression tag	UNP A0A6H1PJZ3
C	1219	ILE	-	expression tag	UNP A0A6H1PJZ3
C	1220	PRO	-	expression tag	UNP A0A6H1PJZ3
C	1221	GLU	-	expression tag	UNP A0A6H1PJZ3
C	1222	ALA	-	expression tag	UNP A0A6H1PJZ3
C	1223	PRO	-	expression tag	UNP A0A6H1PJZ3
C	1224	ARG	-	expression tag	UNP A0A6H1PJZ3
C	1225	ASP	-	expression tag	UNP A0A6H1PJZ3
C	1226	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1227	GLN	-	expression tag	UNP A0A6H1PJZ3
C	1228	ALA	-	expression tag	UNP A0A6H1PJZ3
C	1229	TYR	-	expression tag	UNP A0A6H1PJZ3
C	1230	VAL	-	expression tag	UNP A0A6H1PJZ3
C	1231	ARG	-	expression tag	UNP A0A6H1PJZ3
C	1232	LYS	-	expression tag	UNP A0A6H1PJZ3
C	1233	ASP	-	expression tag	UNP A0A6H1PJZ3
C	1234	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1235	GLU	-	expression tag	UNP A0A6H1PJZ3
C	1236	TRP	-	expression tag	UNP A0A6H1PJZ3
C	1237	VAL	-	expression tag	UNP A0A6H1PJZ3
C	1238	LEU	-	expression tag	UNP A0A6H1PJZ3

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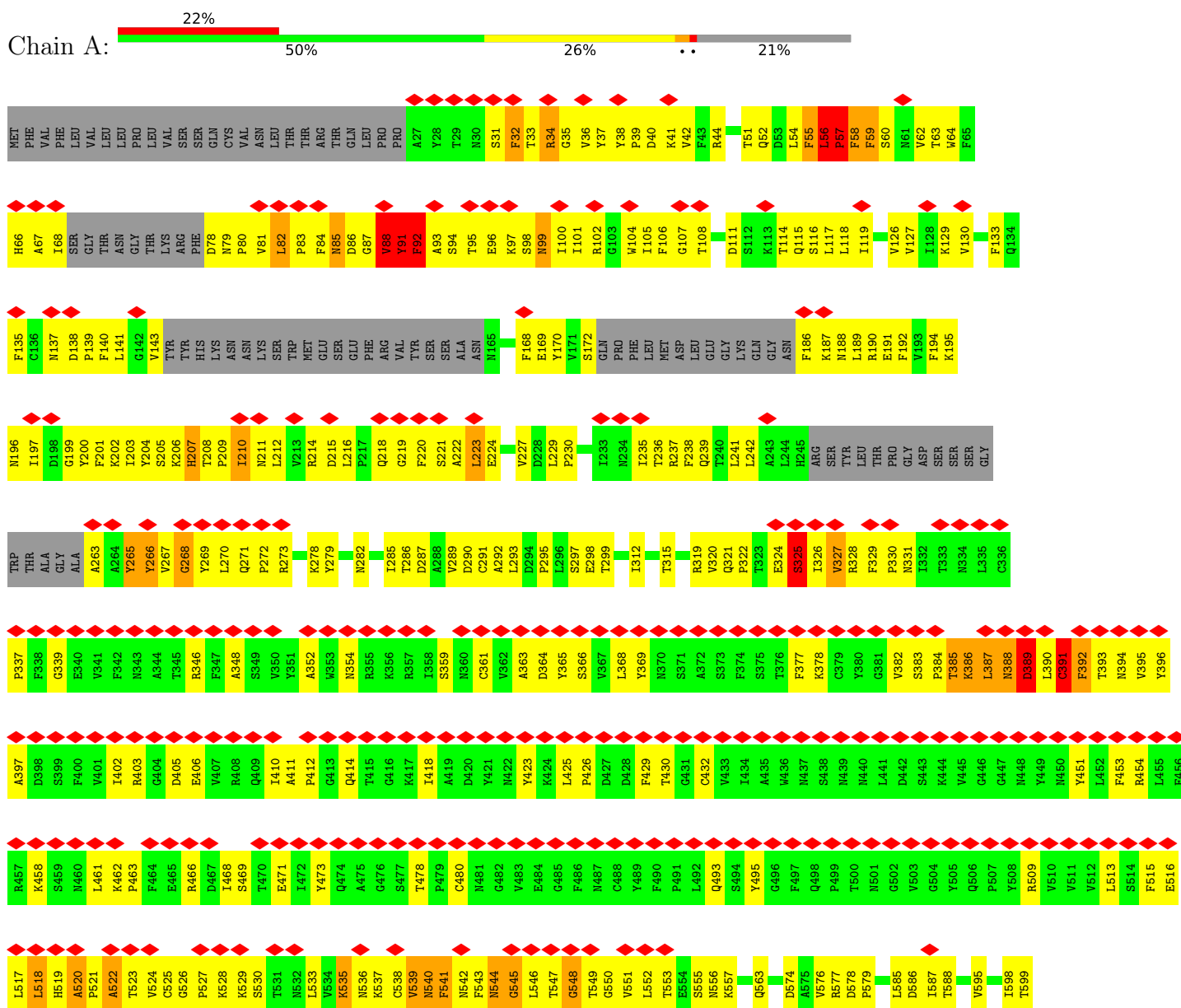
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Chain	Residue	Modelled	Actual	Comment	Reference
C	1239	LEU	-	expression tag	UNP A0A6H1PJZ3
C	1240	SER	-	expression tag	UNP A0A6H1PJZ3
C	1241	THR	-	expression tag	UNP A0A6H1PJZ3
C	1242	PHE	-	expression tag	UNP A0A6H1PJZ3
C	1243	LEU	-	expression tag	UNP A0A6H1PJZ3
C	1244	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1245	SER	-	expression tag	UNP A0A6H1PJZ3
C	1246	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1247	SER	-	expression tag	UNP A0A6H1PJZ3
C	1248	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1249	SER	-	expression tag	UNP A0A6H1PJZ3
C	1250	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1251	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1252	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1253	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1254	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1255	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1256	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1257	GLY	-	expression tag	UNP A0A6H1PJZ3
C	1258	LEU	-	expression tag	UNP A0A6H1PJZ3
C	1259	ASN	-	expression tag	UNP A0A6H1PJZ3
C	1260	ASP	-	expression tag	UNP A0A6H1PJZ3
C	1261	ILE	-	expression tag	UNP A0A6H1PJZ3
C	1262	PHE	-	expression tag	UNP A0A6H1PJZ3
C	1263	GLU	-	expression tag	UNP A0A6H1PJZ3
C	1264	ALA	-	expression tag	UNP A0A6H1PJZ3
C	1265	GLN	-	expression tag	UNP A0A6H1PJZ3
C	1266	LYS	-	expression tag	UNP A0A6H1PJZ3
C	1267	ILE	-	expression tag	UNP A0A6H1PJZ3
C	1268	GLU	-	expression tag	UNP A0A6H1PJZ3
C	1269	TRP	-	expression tag	UNP A0A6H1PJZ3
C	1270	HIS	-	expression tag	UNP A0A6H1PJZ3
C	1271	GLU	-	expression tag	UNP A0A6H1PJZ3

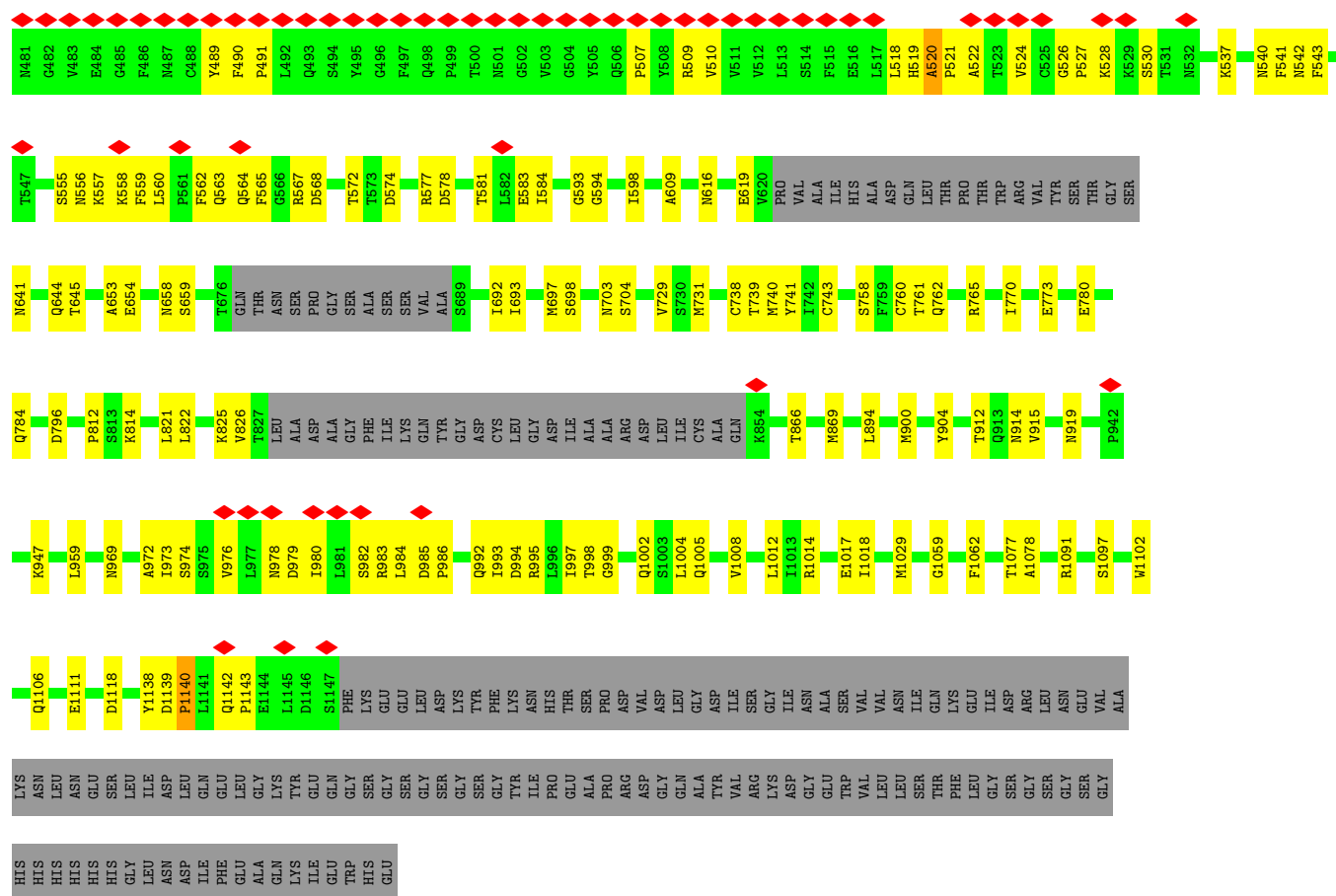
3 Residue-property plots

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.

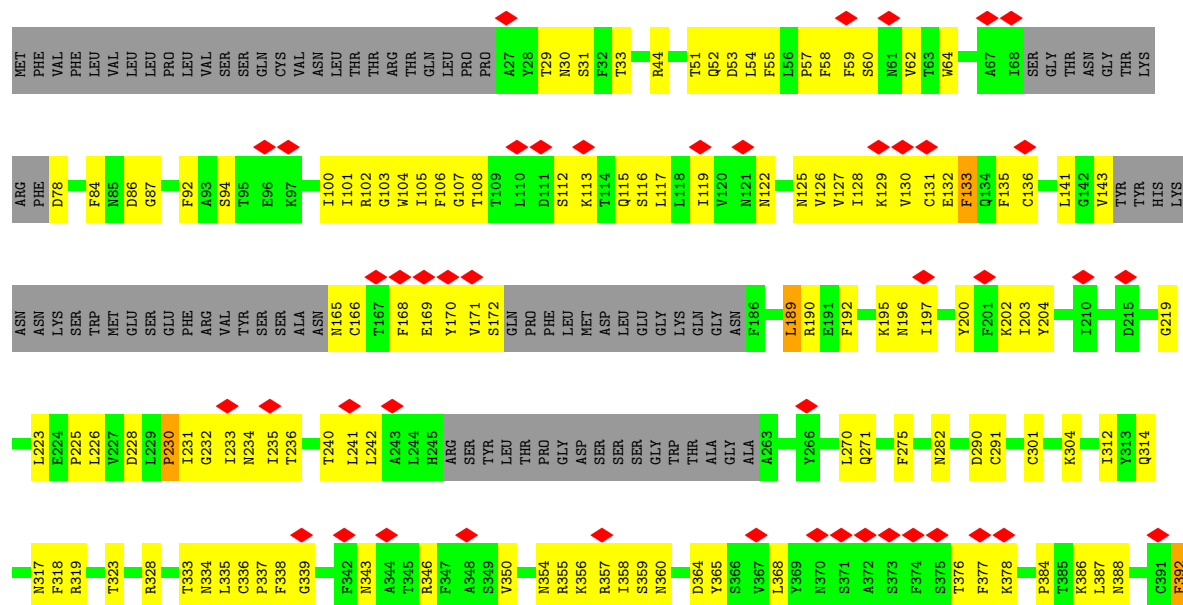
• Molecule 1: Spike glycoprotein







• 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	127065	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 ($e^-/\text{\AA}^2$)	49.98	Depositor
Minimum defocus (nm)	750	Depositor
Maximum defocus (nm)	1750	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.656	Depositor
Minimum map value	-1.052	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.055	Depositor
Recommended contour level	0.415	Depositor
Map size (Å)	568.32, 568.32, 568.32	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.11, 1.11, 1.11	Depositor

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.40	3/7988 (0.0%)	0.78	40/10873 (0.4%)
1	B	0.22	0/7988	0.54	6/10873 (0.1%)
1	C	0.19	0/7988	0.48	3/10873 (0.0%)
All	All	0.28	3/23964 (0.0%)	0.62	49/32619 (0.2%)

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	22
1	B	0	7
1	C	0	4
All	All	0	33

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	88	VAL	CA-C	5.87	1.60	1.52
1	A	91	TYR	N-CA	5.57	1.53	1.46
1	A	59	PHE	N-CA	5.29	1.53	1.46

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	PRO	CA-N-CD	-13.78	92.71	112.00
1	A	88	VAL	CA-C-N	13.25	146.84	121.54
1	A	88	VAL	C-N-CA	13.25	146.84	121.54
1	A	58	PHE	CA-C-N	10.80	142.17	121.54
1	A	58	PHE	C-N-CA	10.80	142.17	121.54
1	A	56	LEU	O-C-N	-10.00	109.82	121.32
1	A	92	PHE	N-CA-C	9.67	124.60	110.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1143	PRO	CA-N-CD	-9.22	99.09	112.00
1	B	1140	PRO	CA-N-CD	-9.15	99.19	112.00
1	A	39	PRO	N-CD-CG	-8.21	90.88	103.20
1	A	91	TYR	CA-C-N	8.11	132.38	120.87
1	A	91	TYR	C-N-CA	8.11	132.38	120.87
1	B	1138	TYR	CA-C-N	7.93	131.75	120.49
1	B	1138	TYR	C-N-CA	7.93	131.75	120.49
1	A	386	LYS	N-CA-C	7.55	123.94	112.54
1	A	58	PHE	N-CA-C	7.39	119.56	110.91
1	A	91	TYR	N-CA-C	6.95	125.60	110.80
1	A	34	ARG	N-CA-C	6.77	119.83	110.35
1	A	57	PRO	N-CA-C	6.69	126.25	112.47
1	C	979	ASP	N-CA-C	6.63	118.39	110.44
1	A	544	ASN	N-CA-C	6.59	121.15	113.18
1	B	91	TYR	N-CA-C	6.38	118.21	108.07
1	A	389	ASP	CA-C-N	6.34	133.22	121.62
1	A	389	ASP	C-N-CA	6.34	133.22	121.62
1	A	541	PHE	N-CA-C	6.27	119.03	107.99
1	A	59	PHE	N-CA-C	6.16	123.92	110.80
1	A	83	PRO	CA-C-N	5.89	132.40	122.15
1	A	83	PRO	C-N-CA	5.89	132.40	122.15
1	A	539	VAL	CA-C-N	5.81	132.64	121.54
1	A	539	VAL	C-N-CA	5.81	132.64	121.54
1	A	392	PHE	N-CA-C	-5.70	99.48	108.14
1	A	541	PHE	CA-C-N	5.61	129.19	121.33
1	A	541	PHE	C-N-CA	5.61	129.19	121.33
1	A	57	PRO	CA-C-N	5.52	131.26	122.93
1	A	57	PRO	C-N-CA	5.52	131.26	122.93
1	A	518	LEU	CA-C-N	5.51	132.06	121.54
1	A	518	LEU	C-N-CA	5.51	132.06	121.54
1	A	391	CYS	N-CA-CB	5.45	118.10	109.71
1	A	82	LEU	N-CA-C	5.44	114.56	108.25
1	C	230	PRO	CA-C-N	-5.39	117.44	123.16
1	C	230	PRO	C-N-CA	-5.39	117.44	123.16
1	A	88	VAL	N-CA-C	5.39	120.55	109.34
1	A	268	GLY	N-CA-C	5.35	125.86	113.18
1	A	389	ASP	N-CA-C	5.35	117.45	108.73
1	A	522	ALA	N-CA-C	5.33	122.16	110.80
1	A	548	GLY	N-CA-C	5.27	118.68	110.88
1	B	447	GLY	N-CA-C	5.10	117.24	112.04
1	A	32	PHE	CA-C-N	-5.01	115.51	123.12
1	A	32	PHE	C-N-CA	-5.01	115.51	123.12

There are no chirality outliers.

All (33) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	ILE	Peptide
1	A	222	ALA	Peptide
1	A	265	TYR	Peptide
1	A	266	TYR	Peptide
1	A	325	SER	Peptide
1	A	327	VAL	Peptide
1	A	378	LYS	Peptide
1	A	385	THR	Peptide
1	A	387	LEU	Peptide
1	A	388	ASN	Peptide
1	A	389	ASP	Peptide
1	A	391	CYS	Peptide
1	A	515	PHE	Peptide
1	A	526	GLY	Peptide
1	A	533	LEU	Peptide
1	A	538	CYS	Peptide
1	A	540	ASN	Peptide
1	A	56	LEU	Mainchain
1	A	88	VAL	Peptide
1	A	91	TYR	Peptide
1	A	92	PHE	Peptide
1	A	99	ASN	Peptide
1	B	379	CYS	Peptide
1	B	383	SER	Peptide
1	B	384	PRO	Peptide
1	B	464	PHE	Peptide
1	B	520	ALA	Peptide
1	B	524	VAL	Peptide
1	B	55	PHE	Peptide
1	C	133	PHE	Peptide
1	C	392	PHE	Peptide
1	C	419	ALA	Peptide
1	C	424	LYS	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	7810	0	7636	566	0
1	B	7810	0	7636	344	0
1	C	7810	0	7636	307	0
All	All	23430	0	22908	1113	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 24.

All (1113) 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:58:PHE:N	1:A:88:VAL:H	1.32	1.26
1:A:58:PHE:H	1:A:88:VAL:N	1.49	1.10
1:A:328:ARG:N	1:A:542:ASN:H	1.54	1.06
1:A:60:SER:HA	1:A:91:TYR:O	1.56	1.05
1:A:391:CYS:SG	1:A:545:GLY:N	2.34	1.01
1:A:60:SER:N	1:A:88:VAL:O	1.94	1.00
1:A:56:LEU:HG	1:A:88:VAL:HA	1.40	1.00
1:A:107:GLY:HA2	1:A:237:ARG:HB3	1.44	0.99
1:A:99:ASN:HD22	1:A:188:ASN:H	1.12	0.98
1:B:383:SER:HB3	1:C:977:LEU:HA	1.46	0.97
1:A:59:PHE:H	1:A:88:VAL:HB	1.28	0.96
1:A:391:CYS:SG	1:A:546:LEU:N	2.40	0.94
1:A:392:PHE:HB3	1:A:518:LEU:N	1.83	0.94
1:A:58:PHE:N	1:A:88:VAL:N	2.12	0.92
1:B:423:TYR:N	1:B:466:ARG:O	2.04	0.90
1:A:56:LEU:O	1:A:87:GLY:N	2.04	0.89
1:B:519:HIS:N	1:B:562:PHE:O	2.06	0.88
1:A:555:SER:HB3	1:A:586:ASP:HB2	1.56	0.88
1:A:104:TRP:H	1:A:191:GLU:HA	1.40	0.87
1:A:328:ARG:HE	1:A:548:GLY:HA3	1.41	0.86
1:A:58:PHE:O	1:A:92:PHE:N	2.08	0.85
1:B:468:ILE:HG22	1:C:234:ASN:HB3	1.56	0.85
1:B:56:LEU:HD22	1:B:91:TYR:HB2	1.57	0.85
1:B:468:ILE:HG23	1:C:108:THR:HG21	1.58	0.85
1:A:524:VAL:H	1:A:544:ASN:HA	1.42	0.84
1:A:325:SER:O	1:A:540:ASN:N	2.10	0.84
1:A:393:THR:H	1:A:517:LEU:C	1.84	0.84
1:A:337:PRO:HB3	1:A:363:ALA:HB3	1.59	0.84
1:A:525:CYS:N	1:A:543:PHE:O	2.10	0.84
1:A:525:CYS:SG	1:A:545:GLY:N	2.50	0.84
1:B:415:THR:HA	1:B:420:ASP:HB3	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:GLN:HG3	1:C:233:ILE:HD13	1.61	0.83
1:B:187:LYS:HA	1:B:209:PRO:HA	1.57	0.82
1:A:58:PHE:H	1:A:88:VAL:H	0.82	0.82
1:A:205:SER:N	1:A:224:GLU:OE1	2.13	0.82
1:A:55:PHE:O	1:A:87:GLY:N	2.13	0.81
1:C:659:SER:HB2	1:C:698:SER:HB2	1.62	0.81
1:B:424:LYS:HE2	1:C:195:LYS:HZ1	1.43	0.81
1:A:330:PRO:HD2	1:A:525:CYS:H	1.42	0.81
1:B:53:ASP:OD2	1:B:54:LEU:N	2.12	0.81
1:C:135:PHE:HD1	1:C:136:CYS:H	1.26	0.81
1:B:413:GLY:HA2	1:B:463:PRO:HA	1.62	0.81
1:A:390:LEU:O	1:A:392:PHE:HB2	1.81	0.80
1:A:391:CYS:HB2	1:A:521:PRO:HA	1.61	0.80
1:C:525:CYS:SG	1:C:526:GLY:N	2.47	0.80
1:A:215:ASP:HA	1:A:263:ALA:HA	1.63	0.79
1:A:536:ASN:HA	1:A:553:THR:HG22	1.62	0.79
1:A:563:GLN:O	1:A:577:ARG:NH2	2.16	0.79
1:B:55:PHE:HA	1:B:87:GLY:HA3	1.64	0.79
1:A:523:THR:OG1	1:A:544:ASN:ND2	2.15	0.79
1:A:96:GLU:H	1:A:218:GLN:HB2	1.47	0.79
1:A:328:ARG:HD3	1:A:549:THR:C	2.08	0.79
1:A:396:TYR:HD1	1:A:397:ALA:H	1.30	0.78
1:A:524:VAL:N	1:A:544:ASN:HA	1.98	0.78
1:A:99:ASN:N	1:A:210:ILE:O	2.17	0.78
1:A:466:ARG:HH21	1:B:130:VAL:HB	1.48	0.78
1:A:230:PRO:HB2	1:C:355:ARG:HH12	1.48	0.78
1:A:236:THR:HB	1:A:237:ARG:HH11	1.49	0.77
1:A:82:LEU:HD13	1:A:268:GLY:HA2	1.64	0.77
1:A:396:TYR:CD2	1:A:516:GLU:HB2	2.20	0.77
1:B:522:ALA:H	1:B:563:GLN:HB3	1.49	0.77
1:C:86:ASP:O	1:C:195:LYS:NZ	2.17	0.77
1:B:336:CYS:SG	1:B:361:CYS:N	2.58	0.76
1:B:976:VAL:HG22	1:B:978:ASN:H	1.50	0.76
1:C:482:GLY:HA2	1:C:490:PHE:HA	1.67	0.76
1:A:58:PHE:C	1:A:91:TYR:HA	2.10	0.76
1:A:392:PHE:HB3	1:A:518:LEU:H	1.49	0.76
1:C:442:ASP:H	1:C:507:PRO:HG3	1.51	0.75
1:C:334:ASN:ND2	1:C:359:SER:O	2.19	0.75
1:A:290:ASP:OD1	1:A:291:CYS:N	2.20	0.75
1:A:55:PHE:C	1:A:87:GLY:H	1.95	0.75
1:A:57:PRO:HB2	1:A:91:TYR:HB2	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:VAL:HG21	1:A:552:LEU:HB2	1.69	0.75
1:A:383:SER:O	1:B:984:LEU:N	2.20	0.75
1:A:385:THR:H	1:B:984:LEU:H	1.35	0.74
1:C:419:ALA:H	1:C:424:LYS:HD2	1.51	0.74
1:A:462:LYS:HE3	1:A:471:GLU:HG2	1.68	0.74
1:A:525:CYS:SG	1:A:543:PHE:C	2.71	0.74
1:B:85:ASN:OD1	1:B:86:ASP:N	2.20	0.74
1:B:384:PRO:HA	1:C:980:ILE:H	1.50	0.74
1:A:327:VAL:HA	1:A:542:ASN:N	2.02	0.74
1:A:519:HIS:HE1	1:A:546:LEU:HD12	1.50	0.74
1:C:127:VAL:HG22	1:C:170:TYR:H	1.53	0.74
1:A:386:LYS:N	1:B:983:ARG:HB2	2.02	0.74
1:B:520:ALA:O	1:B:562:PHE:N	2.21	0.73
1:A:86:ASP:HB2	1:A:273:ARG:H	1.53	0.73
1:B:381:GLY:HA3	1:C:974:SER:H	1.53	0.73
1:A:35:GLY:C	1:A:224:GLU:HB3	2.14	0.73
1:B:385:THR:OG1	1:C:977:LEU:O	2.07	0.73
1:C:402:ILE:H	1:C:508:TYR:HH	1.33	0.73
1:B:384:PRO:HB3	1:C:980:ILE:HB	1.70	0.72
1:B:518:LEU:HD22	1:B:562:PHE:HB2	1.70	0.72
1:A:328:ARG:H	1:A:541:PHE:HA	1.55	0.72
1:C:412:PRO:HA	1:C:425:LEU:HB3	1.70	0.72
1:A:55:PHE:N	1:A:85:ASN:O	2.22	0.72
1:A:99:ASN:HD22	1:A:188:ASN:N	1.87	0.72
1:A:411:ALA:H	1:C:988:GLU:HA	1.55	0.72
1:A:391:CYS:HB2	1:A:522:ALA:H	1.53	0.71
1:A:392:PHE:CD1	1:A:517:LEU:HA	2.26	0.71
1:C:465:GLU:HG3	1:C:466:ARG:HG2	1.71	0.71
1:A:55:PHE:O	1:A:86:ASP:N	2.23	0.71
1:A:462:LYS:HD2	1:B:108:THR:HG21	1.71	0.71
1:B:117:LEU:HB3	1:B:119:ILE:HD11	1.72	0.71
1:C:458:LYS:HE3	1:C:477:SER:HA	1.73	0.71
1:A:195:LYS:NZ	1:A:237:ARG:O	2.23	0.71
1:A:327:VAL:H	1:A:539:VAL:HB	1.56	0.71
1:A:414:GLN:NE2	1:C:986:PRO:O	2.24	0.71
1:B:443:SER:HB3	1:B:507:PRO:HG2	1.74	0.70
1:B:53:ASP:O	1:B:87:GLY:N	2.17	0.70
1:C:520:ALA:HB3	1:C:563:GLN:HA	1.73	0.70
1:A:574:ASP:HA	1:A:587:ILE:HB	1.74	0.70
1:B:1091:ARG:NH1	1:B:1118:ASP:O	2.25	0.70
1:B:376:THR:O	1:C:983:ARG:NH1	2.24	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:SER:OG	1:A:268:GLY:HA3	1.90	0.70
1:A:466:ARG:HB2	1:B:130:VAL:HG21	1.74	0.70
1:B:273:ARG:NH1	1:B:290:ASP:OD2	2.22	0.70
1:A:524:VAL:H	1:A:544:ASN:ND2	1.90	0.70
1:A:196:ASN:HB3	1:A:235:ILE:HB	1.73	0.69
1:A:328:ARG:H	1:A:542:ASN:H	1.38	0.69
1:A:391:CYS:HB2	1:A:522:ALA:N	2.07	0.69
1:A:31:SER:OG	1:A:32:PHE:N	2.18	0.69
1:C:452:LEU:HD21	1:C:493:GLN:H	1.56	0.69
1:A:599:THR:HB	1:A:608:VAL:HG23	1.75	0.69
1:C:413:GLY:HA3	1:C:429:PHE:HA	1.73	0.69
1:A:99:ASN:ND2	1:A:188:ASN:H	1.87	0.69
1:A:196:ASN:H	1:A:235:ILE:HD12	1.56	0.69
1:B:277:LEU:HD12	1:B:285:ILE:HD13	1.75	0.69
1:A:327:VAL:HG22	1:A:539:VAL:HG11	1.73	0.69
1:B:518:LEU:HB3	1:B:562:PHE:O	1.91	0.69
1:C:518:LEU:HD22	1:C:562:PHE:HB2	1.75	0.68
1:B:654:GLU:HB3	1:B:693:ILE:HG22	1.75	0.68
1:C:350:VAL:HA	1:C:400:PHE:HB2	1.76	0.68
1:A:346:ARG:NH2	1:A:352:ALA:O	2.25	0.68
1:A:98:SER:H	1:A:211:ASN:HA	1.56	0.68
1:C:992:GLN:OE1	1:C:992:GLN:N	2.25	0.68
1:A:644:GLN:NE2	1:A:645:THR:O	2.26	0.68
1:A:59:PHE:HZ	1:A:216:LEU:HD13	1.59	0.68
1:A:56:LEU:HG	1:A:88:VAL:CA	2.22	0.67
1:A:402:ILE:HD11	1:A:509:ARG:HA	1.77	0.67
1:B:187:LYS:HB2	1:B:210:ILE:HG23	1.75	0.67
1:A:64:TRP:HB2	1:A:78:ASP:N	2.08	0.67
1:A:95:THR:HA	1:A:218:GLN:HG2	1.75	0.67
1:A:58:PHE:CA	1:A:88:VAL:H	2.07	0.67
1:A:59:PHE:CE2	1:A:94:SER:HB3	2.30	0.67
1:A:329:PHE:H	1:A:541:PHE:CA	2.06	0.67
1:B:53:ASP:CG	1:B:54:LEU:H	2.03	0.67
1:A:331:ASN:HA	1:A:527:PRO:HA	1.75	0.67
1:A:391:CYS:HA	1:A:520:ALA:C	2.19	0.67
1:A:528:LYS:NZ	1:A:529:LYS:O	2.26	0.66
1:C:31:SER:HB3	1:C:62:VAL:HG11	1.78	0.66
1:A:194:PHE:HB2	1:A:201:PHE:HE1	1.59	0.66
1:A:391:CYS:CB	1:A:522:ALA:H	2.08	0.66
1:C:442:ASP:HB2	1:C:507:PRO:HB3	1.77	0.66
1:A:55:PHE:O	1:A:85:ASN:C	2.39	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:VAL:HG12	1:B:82:LEU:H	1.60	0.66
1:A:59:PHE:CZ	1:A:216:LEU:HD13	2.30	0.66
1:A:59:PHE:N	1:A:88:VAL:HB	2.07	0.66
1:A:98:SER:H	1:A:210:ILE:C	2.04	0.66
1:B:520:ALA:C	1:B:563:GLN:HB2	2.20	0.66
1:A:469:SER:HB2	1:B:114:THR:HB	1.78	0.66
1:C:119:ILE:HG23	1:C:126:VAL:HG13	1.78	0.66
1:A:389:ASP:HA	1:A:520:ALA:HB2	1.77	0.65
1:A:206:LYS:NZ	1:A:207:HIS:O	2.29	0.65
1:B:398:ASP:OD1	1:B:399:SER:N	2.30	0.65
1:A:135:PHE:HB3	1:A:139:PRO:HA	1.77	0.65
1:C:419:ALA:HB2	1:C:424:LYS:H	1.61	0.65
1:A:33:THR:HB	1:A:221:SER:H	1.62	0.65
1:B:30:ASN:OD1	1:B:31:SER:N	2.30	0.65
1:B:641:ASN:ND2	1:B:653:ALA:O	2.29	0.65
1:C:384:PRO:HA	1:C:387:LEU:HG	1.79	0.65
1:A:522:ALA:HB3	1:A:544:ASN:HB2	1.79	0.64
1:A:553:THR:HG21	1:A:588:THR:HB	1.79	0.64
1:B:62:VAL:HG11	1:B:266:TYR:HB3	1.78	0.64
1:C:418:ILE:HB	1:C:424:LYS:HA	1.78	0.64
1:C:866:THR:OG1	1:C:869:MET:SD	2.55	0.64
1:A:346:ARG:NH1	1:A:354:ASN:OD1	2.31	0.64
1:C:125:ASN:HB2	1:C:171:VAL:HA	1.79	0.64
1:C:697:MET:HE3	1:C:698:SER:H	1.63	0.64
1:A:56:LEU:C	1:A:87:GLY:C	2.65	0.64
1:B:99:ASN:O	1:B:245:HIS:N	2.30	0.64
1:B:644:GLN:NE2	1:B:645:THR:O	2.30	0.64
1:B:83:PRO:HG2	1:B:270:LEU:HD22	1.80	0.64
1:C:84:PHE:HB3	1:C:236:THR:HG22	1.78	0.64
1:B:29:THR:HG22	1:B:266:TYR:CD1	2.32	0.64
1:A:54:LEU:HB3	1:A:84:PHE:CE1	2.33	0.64
1:A:328:ARG:N	1:A:542:ASN:N	2.39	0.64
1:C:523:THR:HG21	1:C:582:LEU:HD13	1.79	0.64
1:C:402:ILE:N	1:C:508:TYR:HH	1.96	0.64
1:A:329:PHE:H	1:A:541:PHE:HA	1.61	0.64
1:B:118:LEU:HB3	1:B:129:LYS:HB2	1.79	0.64
1:B:290:ASP:OD1	1:B:291:CYS:N	2.30	0.64
1:B:328:ARG:NH1	1:B:530:SER:OG	2.31	0.64
1:A:85:ASN:HD22	1:A:270:LEU:HD11	1.63	0.64
1:A:59:PHE:O	1:A:92:PHE:HA	1.98	0.63
1:C:457:ARG:HH22	1:C:471:GLU:HB3	1.61	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:ILE:HD11	1:A:528:LYS:HZ1	1.62	0.63
1:A:516:GLU:HG2	1:A:517:LEU:H	1.62	0.63
1:A:44:ARG:NH1	1:A:279:TYR:OH	2.31	0.63
1:A:278:LYS:HB3	1:A:287:ASP:HB3	1.78	0.63
1:A:327:VAL:HG12	1:A:542:ASN:HB3	1.81	0.63
1:A:383:SER:HB3	1:A:384:PRO:HD2	1.79	0.63
1:A:195:LYS:HB3	1:A:235:ILE:HD11	1.80	0.63
1:B:351:TYR:CZ	1:C:112:SER:HA	2.33	0.63
1:A:384:PRO:HA	1:B:982:SER:HA	1.80	0.63
1:C:197:ILE:HG13	1:C:202:LYS:HE2	1.81	0.63
1:C:402:ILE:N	1:C:508:TYR:OH	2.25	0.63
1:A:54:LEU:HA	1:A:85:ASN:O	1.98	0.63
1:A:755:GLN:OE1	1:C:969:ASN:ND2	2.31	0.63
1:C:465:GLU:HG3	1:C:466:ARG:H	1.63	0.63
1:A:327:VAL:HG23	1:A:539:VAL:HG21	1.81	0.63
1:C:337:PRO:HG2	1:C:356:LYS:HE3	1.80	0.63
1:A:387:LEU:H	1:B:983:ARG:CB	2.11	0.63
1:B:420:ASP:OD1	1:B:421:TYR:N	2.32	0.63
1:A:237:ARG:O	1:A:238:PHE:HD1	1.80	0.62
1:A:392:PHE:HB3	1:A:518:LEU:C	2.25	0.62
1:A:513:LEU:HD22	1:A:518:LEU:HD13	1.81	0.62
1:A:330:PRO:CD	1:A:525:CYS:H	2.13	0.62
1:A:976:VAL:HG12	1:A:979:ASP:H	1.65	0.62
1:B:61:ASN:HB3	1:B:269:TYR:HD2	1.65	0.62
1:A:773:GLU:OE1	1:A:774:GLN:NE2	2.32	0.62
1:A:34:ARG:N	1:A:34:ARG:HE	1.97	0.62
1:A:57:PRO:HG3	1:A:271:GLN:H	1.65	0.62
1:B:616:ASN:HB3	1:B:619:GLU:HG2	1.82	0.62
1:A:392:PHE:HA	1:A:517:LEU:HG	1.81	0.62
1:A:41:LYS:HE3	1:C:517:LEU:HA	1.82	0.62
1:A:619:GLU:OE1	1:A:619:GLU:N	2.33	0.62
1:B:115:GLN:NE2	1:B:165:ASN:O	2.33	0.62
1:C:133:PHE:O	1:C:135:PHE:N	2.33	0.62
1:A:115:GLN:HA	1:A:118:LEU:HG	1.82	0.61
1:A:403:ARG:NH1	1:A:454:ARG:O	2.33	0.61
1:A:713:ALA:HB3	1:B:894:LEU:HB3	1.81	0.61
1:C:994:ASP:OD1	1:C:995:ARG:N	2.33	0.61
1:A:330:PRO:HD2	1:A:525:CYS:N	2.15	0.61
1:B:190:ARG:HH11	1:B:192:PHE:H	1.49	0.61
1:C:565:PHE:HD1	1:C:566:GLY:N	1.98	0.61
1:A:80:PRO:HG2	1:A:82:LEU:HD11	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:TYR:O	1:A:267:VAL:HG23	2.00	0.61
1:B:366:SER:O	1:B:370:ASN:HB2	2.01	0.61
1:C:335:LEU:HD12	1:C:336:CYS:H	1.63	0.61
1:A:59:PHE:O	1:A:59:PHE:CD2	2.54	0.61
1:A:127:VAL:HG13	1:A:129:LYS:HE3	1.83	0.61
1:A:82:LEU:HD13	1:A:268:GLY:CA	2.31	0.61
1:C:419:ALA:CB	1:C:424:LYS:H	2.13	0.61
1:C:33:THR:OG1	1:C:219:GLY:O	2.18	0.61
1:B:382:VAL:HB	1:C:979:ASP:OD1	2.01	0.61
1:A:324:GLU:O	1:A:540:ASN:ND2	2.34	0.61
1:B:200:TYR:CG	1:B:230:PRO:HB3	2.36	0.61
1:C:29:THR:H	1:C:64:TRP:HB3	1.66	0.61
1:A:35:GLY:O	1:A:224:GLU:HB3	2.01	0.61
1:A:66:HIS:CD2	1:A:68:ILE:HG12	2.35	0.61
1:B:520:ALA:HB1	1:B:559:PHE:HB3	1.83	0.61
1:C:136:CYS:O	1:C:165:ASN:N	2.34	0.61
1:A:391:CYS:HB2	1:A:521:PRO:CA	2.31	0.60
1:B:520:ALA:O	1:B:563:GLN:N	2.34	0.60
1:B:422:ASN:HB3	1:B:466:ARG:HD3	1.82	0.60
1:C:200:TYR:CD1	1:C:230:PRO:HB3	2.36	0.60
1:A:239:GLN:HA	1:A:267:VAL:HG11	1.83	0.60
1:C:644:GLN:NE2	1:C:645:THR:O	2.34	0.60
1:A:759:PHE:O	1:A:762:GLN:HG2	2.02	0.60
1:B:200:TYR:CD2	1:B:230:PRO:HB3	2.37	0.60
1:C:398:ASP:N	1:C:398:ASP:OD1	2.33	0.60
1:A:214:ARG:HB3	1:A:263:ALA:HB1	1.83	0.60
1:A:385:THR:N	1:B:983:ARG:H	1.99	0.60
1:B:78:ASP:N	1:B:269:TYR:HH	1.98	0.60
1:A:556:ASN:OD1	1:A:557:LYS:NZ	2.35	0.60
1:B:465:GLU:O	1:B:467:ASP:N	2.34	0.60
1:C:364:ASP:OD1	1:C:388:ASN:ND2	2.35	0.60
1:A:327:VAL:HA	1:A:542:ASN:HB3	1.84	0.60
1:A:391:CYS:O	1:A:392:PHE:CD2	2.53	0.60
1:B:692:ILE:HD12	1:B:692:ILE:H	1.67	0.60
1:A:552:LEU:HD23	1:A:587:ILE:HG12	1.84	0.59
1:B:383:SER:O	1:C:978:ASN:N	2.35	0.59
1:C:453:PHE:CE1	1:C:455:LEU:HD23	2.37	0.59
1:A:138:ASP:O	1:A:140:PHE:N	2.36	0.59
1:C:84:PHE:HB2	1:C:236:THR:HA	1.82	0.59
1:C:135:PHE:CD1	1:C:136:CYS:N	2.69	0.59
1:C:376:THR:OG1	1:C:435:ALA:O	2.20	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:578:ASP:HB3	1:C:583:GLU:H	1.65	0.59
1:B:377:PHE:HB3	1:C:984:LEU:HB2	1.83	0.59
1:C:130:VAL:HG22	1:C:132:GLU:HB3	1.84	0.59
1:C:231:ILE:HG13	1:C:232:GLY:N	2.16	0.59
1:C:200:TYR:HB3	1:C:228:ASP:HB3	1.83	0.59
1:A:56:LEU:HB2	1:A:57:PRO:HD3	1.85	0.59
1:A:59:PHE:HB3	1:A:92:PHE:HA	1.84	0.59
1:A:98:SER:H	1:A:211:ASN:N	2.00	0.59
1:A:108:THR:HG22	1:A:237:ARG:HA	1.85	0.59
1:A:129:LYS:NZ	1:A:169:GLU:OE1	2.22	0.59
1:A:517:LEU:HD11	1:A:544:ASN:O	2.03	0.59
1:A:523:THR:H	1:A:544:ASN:CG	2.10	0.59
1:A:553:THR:N	1:A:586:ASP:O	2.24	0.59
1:B:355:ARG:HB2	1:C:128:ILE:HG21	1.85	0.59
1:B:386:LYS:H	1:C:978:ASN:HA	1.68	0.59
1:C:104:TRP:HZ2	1:C:119:ILE:HB	1.67	0.59
1:C:521:PRO:O	1:C:563:GLN:NE2	2.35	0.59
1:C:968:SER:N	1:C:975:SER:OG	2.36	0.59
1:A:524:VAL:H	1:A:544:ASN:HD22	1.50	0.59
1:B:28:TYR:CG	1:B:62:VAL:HB	2.37	0.59
1:B:94:SER:OG	1:B:190:ARG:NE	2.36	0.59
1:C:697:MET:HE3	1:C:698:SER:N	2.17	0.59
1:A:383:SER:HB2	1:B:985:ASP:HA	1.84	0.59
1:A:390:LEU:C	1:A:520:ALA:HB3	2.28	0.59
1:A:98:SER:H	1:A:211:ASN:CA	2.16	0.58
1:A:329:PHE:N	1:A:542:ASN:N	2.50	0.58
1:A:315:THR:OG1	1:A:595:VAL:O	2.20	0.58
1:A:392:PHE:HD1	1:A:518:LEU:H	1.50	0.58
1:A:755:GLN:HA	1:C:969:ASN:HD22	1.68	0.58
1:A:328:ARG:O	1:A:543:PHE:HB2	2.04	0.58
1:B:381:GLY:HA3	1:C:974:SER:N	2.18	0.58
1:A:98:SER:N	1:A:211:ASN:HA	2.18	0.58
1:B:377:PHE:CD2	1:C:983:ARG:HB3	2.38	0.58
1:A:81:VAL:HA	1:A:239:GLN:CD	2.29	0.58
1:A:520:ALA:HB1	1:A:521:PRO:HD2	1.85	0.58
1:A:328:ARG:NH1	1:A:521:PRO:HG3	2.18	0.58
1:C:58:PHE:HB3	1:C:290:ASP:HB2	1.85	0.58
1:B:66:HIS:CE1	1:B:68:ILE:HG23	2.39	0.58
1:A:36:VAL:HG23	1:A:223:LEU:O	2.04	0.58
1:A:63:THR:HG22	1:A:64:TRP:H	1.67	0.58
1:A:107:GLY:N	1:A:239:GLN:H	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:PRO:CD	1:A:542:ASN:HA	2.33	0.58
1:A:553:THR:O	1:A:586:ASP:N	2.32	0.58
1:B:209:PRO:HG2	1:B:216:LEU:HD11	1.84	0.58
1:C:339:GLY:O	1:C:343:ASN:ND2	2.36	0.58
1:C:520:ALA:O	1:C:565:PHE:N	2.36	0.58
1:A:56:LEU:N	1:A:87:GLY:O	2.37	0.58
1:A:382:VAL:HG13	1:A:432:CYS:HB3	1.86	0.58
1:C:31:SER:H	1:C:62:VAL:HG21	1.68	0.58
1:A:403:ARG:HB2	1:A:418:ILE:HG23	1.84	0.58
1:A:392:PHE:O	1:A:519:HIS:HA	2.02	0.57
1:A:196:ASN:OD1	1:A:200:TYR:N	2.37	0.57
1:B:67:ALA:HB3	1:B:242:LEU:HD21	1.86	0.57
1:B:384:PRO:HA	1:C:980:ILE:N	2.19	0.57
1:B:412:PRO:HB3	1:B:464:PHE:CD2	2.39	0.57
1:C:131:CYS:SG	1:C:166:CYS:N	2.66	0.57
1:C:520:ALA:HB1	1:C:521:PRO:HD2	1.86	0.57
1:A:189:LEU:O	1:A:208:THR:N	2.26	0.57
1:B:271:GLN:HB3	1:B:273:ARG:HE	1.69	0.57
1:A:194:PHE:HB2	1:A:201:PHE:CE1	2.39	0.57
1:A:394:ASN:C	1:A:518:LEU:HD22	2.28	0.57
1:A:535:LYS:HB2	1:A:553:THR:HA	1.86	0.57
1:C:456:PHE:HE2	1:C:477:SER:HB3	1.70	0.57
1:A:327:VAL:CG2	1:A:539:VAL:HG21	2.35	0.57
1:B:330:PRO:HG3	1:B:577:ARG:HH22	1.70	0.57
1:A:330:PRO:HD3	1:A:543:PHE:H	1.68	0.57
1:A:524:VAL:C	1:A:544:ASN:HA	2.29	0.57
1:B:731:MET:HE1	1:B:1018:ILE:HG21	1.86	0.57
1:C:122:ASN:ND2	1:C:169:GLU:OE1	2.37	0.57
1:A:200:TYR:HB3	1:A:230:PRO:HA	1.87	0.57
1:A:329:PHE:H	1:A:541:PHE:C	2.12	0.57
1:B:274:THR:H	1:B:291:CYS:HB2	1.69	0.57
1:B:48:LEU:HD13	1:B:276:LEU:HD21	1.87	0.57
1:C:402:ILE:HG22	1:C:403:ARG:H	1.70	0.57
1:A:59:PHE:N	1:A:88:VAL:C	2.63	0.57
1:A:105:ILE:HD11	1:A:241:LEU:HB2	1.86	0.57
1:B:389:ASP:HB3	1:B:542:ASN:HD21	1.70	0.57
1:B:467:ASP:O	1:B:469:SER:N	2.38	0.56
1:C:933:LYS:NZ	1:C:933:LYS:O	2.38	0.56
1:A:59:PHE:HD2	1:A:93:ALA:H	1.52	0.56
1:A:329:PHE:CD2	1:A:547:THR:HA	2.40	0.56
1:A:365:TYR:HB3	1:A:377:PHE:HE1	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:400:PHE:CD1	1:B:402:ILE:HG12	2.41	0.56
1:B:420:ASP:O	1:B:466:ARG:HA	2.04	0.56
1:A:598:ILE:HD11	1:A:664:ILE:HG21	1.87	0.56
1:B:36:VAL:O	1:B:223:LEU:N	2.39	0.56
1:B:108:THR:HG23	1:B:109:THR:HG23	1.88	0.56
1:B:409:GLN:HB3	1:B:455:LEU:HD13	1.87	0.56
1:C:346:ARG:HH22	1:C:444:LYS:HB2	1.70	0.56
1:A:32:PHE:O	1:A:219:GLY:HA3	2.06	0.56
1:A:391:CYS:HA	1:A:520:ALA:O	2.06	0.56
1:B:387:LEU:HD13	1:C:979:ASP:HB2	1.88	0.56
1:C:498:GLN:H	1:C:505:TYR:HB3	1.70	0.56
1:C:984:LEU:HD21	1:C:988:GLU:HB3	1.86	0.56
1:A:194:PHE:O	1:A:195:LYS:HD2	2.06	0.56
1:C:104:TRP:CZ2	1:C:119:ILE:HB	2.40	0.56
1:C:129:LYS:HG2	1:C:130:VAL:H	1.71	0.56
1:C:358:ILE:HG12	1:C:360:ASN:H	1.70	0.56
1:C:793:PRO:O	1:C:795:LYS:NZ	2.39	0.56
1:B:37:TYR:H	1:B:88:VAL:HG11	1.70	0.56
1:C:301:CYS:O	1:C:304:LYS:NZ	2.29	0.56
1:A:56:LEU:CG	1:A:88:VAL:HA	2.26	0.56
1:A:58:PHE:O	1:A:91:TYR:HA	2.05	0.56
1:A:466:ARG:NH2	1:B:130:VAL:HB	2.20	0.56
1:B:518:LEU:HB3	1:B:562:PHE:C	2.30	0.56
1:B:812:PRO:O	1:B:814:LYS:NZ	2.39	0.56
1:A:524:VAL:HG21	1:A:578:ASP:HA	1.86	0.56
1:A:551:VAL:HG12	1:A:587:ILE:HG23	1.87	0.56
1:C:560:LEU:HB2	1:C:562:PHE:CE2	2.41	0.56
1:C:563:GLN:HG3	1:C:577:ARG:NH2	2.21	0.56
1:B:82:LEU:HD23	1:B:92:PHE:HA	1.88	0.56
1:B:1002:GLN:OE1	1:B:1002:GLN:N	2.27	0.56
1:A:106:PHE:HB3	1:A:238:PHE:HA	1.87	0.55
1:B:30:ASN:C	1:B:32:PHE:H	2.15	0.55
1:B:866:THR:OG1	1:B:869:MET:SD	2.60	0.55
1:A:32:PHE:HA	1:A:34:ARG:NE	2.21	0.55
1:A:1125:ASN:ND2	1:A:1127:ASP:OD1	2.38	0.55
1:B:57:PRO:HD2	1:B:270:LEU:HA	1.87	0.55
1:B:1118:ASP:N	1:B:1118:ASP:OD1	2.31	0.55
1:A:59:PHE:HB3	1:A:92:PHE:CD1	2.41	0.55
1:A:394:ASN:O	1:A:518:LEU:HD22	2.06	0.55
1:A:991:VAL:HG12	1:A:995:ARG:HH21	1.72	0.55
1:B:421:TYR:HD1	1:B:467:ASP:H	1.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:LEU:HD12	1:B:84:PHE:CD1	2.42	0.55
1:C:392:PHE:CE2	1:C:513:LEU:HB2	2.41	0.55
1:A:195:LYS:HB3	1:A:235:ILE:CD1	2.36	0.55
1:B:55:PHE:HZ	1:B:273:ARG:H	1.52	0.55
1:C:365:TYR:HA	1:C:368:LEU:HD12	1.89	0.55
1:B:53:ASP:C	1:B:87:GLY:H	2.13	0.55
1:B:1142:GLN:N	1:B:1142:GLN:OE1	2.40	0.55
1:C:565:PHE:HD1	1:C:566:GLY:H	1.54	0.55
1:A:326:ILE:HG12	1:A:530:SER:HA	1.87	0.55
1:A:466:ARG:HH22	1:B:128:ILE:HG22	1.72	0.55
1:A:519:HIS:CE1	1:A:546:LEU:HD12	2.38	0.55
1:A:59:PHE:HB3	1:A:92:PHE:CG	2.42	0.55
1:A:203:ILE:HD13	1:A:227:VAL:HB	1.89	0.55
1:A:369:TYR:CG	1:A:377:PHE:HB2	2.42	0.55
1:B:520:ALA:HB3	1:B:560:LEU:O	2.06	0.55
1:C:796:ASP:OD1	1:C:796:ASP:N	2.38	0.55
1:A:328:ARG:HG3	1:A:543:PHE:HD1	1.72	0.54
1:A:392:PHE:HA	1:A:517:LEU:CG	2.37	0.54
1:A:366:SER:HA	1:A:369:TYR:CD2	2.41	0.54
1:B:557:LYS:NZ	1:B:574:ASP:OD2	2.40	0.54
1:C:106:PHE:HB3	1:C:235:ILE:HG23	1.90	0.54
1:A:98:SER:OG	1:A:211:ASN:O	2.17	0.54
1:A:114:THR:N	1:C:466:ARG:HH12	2.06	0.54
1:A:348:ALA:HB3	1:A:352:ALA:HB3	1.89	0.54
1:C:378:LYS:HZ3	1:C:435:ALA:HB3	1.73	0.54
1:A:97:LYS:H	1:A:211:ASN:N	2.05	0.54
1:A:273:ARG:HD2	1:A:292:ALA:HB3	1.90	0.54
1:B:202:LYS:HD3	1:B:228:ASP:OD1	2.08	0.54
1:A:86:ASP:HB3	1:A:273:ARG:O	2.07	0.54
1:A:1091:ARG:NH1	1:A:1118:ASP:O	2.40	0.54
1:B:52:GLN:HA	1:B:86:ASP:OD1	2.07	0.54
1:C:204:TYR:HA	1:C:225:PRO:HA	1.88	0.54
1:C:337:PRO:HG3	1:C:357:ARG:NH1	2.23	0.54
1:A:391:CYS:O	1:A:392:PHE:HD2	1.91	0.54
1:A:964:LYS:O	1:A:964:LYS:NZ	2.28	0.54
1:B:527:PRO:HD2	1:B:528:LYS:HZ1	1.72	0.54
1:A:659:SER:HB2	1:A:698:SER:HB2	1.90	0.53
1:C:129:LYS:HE3	1:C:168:PHE:HB2	1.90	0.53
1:A:236:THR:HB	1:A:237:ARG:NH1	2.20	0.53
1:A:326:ILE:HD12	1:A:541:PHE:H	1.74	0.53
1:A:390:LEU:N	1:A:520:ALA:HB3	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:992:GLN:OE1	1:B:995:ARG:NH2	2.41	0.53
1:A:32:PHE:HA	1:A:34:ARG:CZ	2.38	0.53
1:A:60:SER:HA	1:A:91:TYR:C	2.33	0.53
1:B:115:GLN:HE21	1:B:165:ASN:HD22	1.56	0.53
1:B:578:ASP:OD1	1:B:581:THR:N	2.35	0.53
1:C:314:GLN:NE2	1:C:595:VAL:O	2.41	0.53
1:A:387:LEU:H	1:B:983:ARG:HB3	1.73	0.53
1:A:107:GLY:N	1:A:239:GLN:O	2.26	0.53
1:A:556:ASN:H	1:A:557:LYS:HZ2	1.57	0.53
1:A:1092:GLU:OE1	1:A:1107:ARG:NH2	2.42	0.53
1:B:54:LEU:HA	1:B:84:PHE:HA	1.90	0.53
1:B:200:TYR:HA	1:B:230:PRO:HA	1.91	0.53
1:B:239:GLN:HG3	1:B:240:THR:H	1.72	0.53
1:C:203:ILE:O	1:C:226:LEU:N	2.40	0.53
1:A:97:LYS:O	1:A:209:PRO:HB2	2.08	0.53
1:A:321:GLN:HG2	1:A:322:PRO:HD2	1.91	0.53
1:B:64:TRP:HZ3	1:B:265:TYR:O	1.91	0.53
1:C:759:PHE:O	1:C:763:LEU:HG	2.09	0.53
1:C:976:VAL:HG22	1:C:977:LEU:H	1.73	0.53
1:A:58:PHE:H	1:A:87:GLY:HA3	1.73	0.53
1:A:206:LYS:HG2	1:A:207:HIS:H	1.74	0.53
1:B:317:ASN:HA	1:B:594:GLY:HA2	1.90	0.53
1:B:979:ASP:HA	1:B:983:ARG:NH1	2.24	0.53
1:A:59:PHE:C	1:A:91:TYR:C	2.77	0.53
1:A:383:SER:CB	1:B:986:PRO:HD3	2.39	0.53
1:C:458:LYS:NZ	1:C:474:GLN:HB3	2.24	0.53
1:A:196:ASN:O	1:A:235:ILE:HD12	2.09	0.52
1:C:977:LEU:HD12	1:C:1000:ARG:HH12	1.73	0.52
1:C:445:VAL:HG23	1:C:498:GLN:HE21	1.74	0.52
1:C:525:CYS:HB2	1:C:544:ASN:HA	1.90	0.52
1:A:59:PHE:CZ	1:A:94:SER:HB3	2.43	0.52
1:A:539:VAL:O	1:A:549:THR:HG23	2.09	0.52
1:C:200:TYR:CE1	1:C:230:PRO:HB3	2.44	0.52
1:A:556:ASN:H	1:A:557:LYS:NZ	2.07	0.52
1:A:389:ASP:OD2	1:B:983:ARG:NH1	2.42	0.52
1:A:586:ASP:OD1	1:A:587:ILE:N	2.40	0.52
1:A:31:SER:HB3	1:A:286:THR:C	2.34	0.52
1:A:57:PRO:C	1:A:91:TYR:H	2.18	0.52
1:A:85:ASN:HD22	1:A:270:LEU:CD1	2.21	0.52
1:A:327:VAL:N	1:A:540:ASN:O	2.42	0.52
1:A:524:VAL:HG11	1:A:579:PRO:HD2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:563:GLN:HA	1:B:41:LYS:HB3	1.92	0.52
1:B:351:TYR:HB2	1:C:113:LYS:O	2.10	0.52
1:B:567:ARG:HH22	1:C:44:ARG:HE	1.58	0.52
1:A:57:PRO:CB	1:A:271:GLN:HG2	2.40	0.52
1:A:58:PHE:C	1:A:91:TYR:CA	2.80	0.52
1:A:196:ASN:HD21	1:A:199:GLY:N	2.07	0.52
1:A:117:LEU:HD13	1:A:194:PHE:HE1	1.75	0.52
1:A:468:ILE:HG13	1:B:114:THR:OG1	2.10	0.52
1:B:509:ARG:HG3	1:B:510:VAL:H	1.74	0.52
1:A:101:ILE:O	1:A:189:LEU:HG	2.11	0.51
1:A:327:VAL:H	1:A:539:VAL:CB	2.23	0.51
1:A:330:PRO:HG3	1:A:542:ASN:OD1	2.10	0.51
1:B:114:THR:OG1	1:B:115:GLN:N	2.42	0.51
1:C:950:ASP:N	1:C:950:ASP:OD1	2.42	0.51
1:A:58:PHE:H	1:A:87:GLY:C	2.15	0.51
1:A:195:LYS:O	1:A:202:LYS:N	2.33	0.51
1:A:391:CYS:HB3	1:A:544:ASN:N	2.26	0.51
1:B:204:TYR:CE1	1:B:225:PRO:HB3	2.45	0.51
1:C:419:ALA:HB3	1:C:424:LYS:HG2	1.92	0.51
1:A:391:CYS:HB3	1:A:544:ASN:CA	2.40	0.51
1:A:676:THR:HA	1:A:690:GLN:HA	1.92	0.51
1:B:30:ASN:CG	1:B:59:PHE:H	2.18	0.51
1:C:105:ILE:H	1:C:240:THR:HB	1.75	0.51
1:C:234:ASN:OD1	1:C:235:ILE:N	2.43	0.51
1:B:60:SER:HB2	1:B:269:TYR:N	2.26	0.51
1:B:190:ARG:C	1:B:190:ARG:HD2	2.36	0.51
1:B:703:ASN:OD1	1:B:704:SER:N	2.43	0.51
1:A:59:PHE:HA	1:A:88:VAL:HB	1.93	0.51
1:A:190:ARG:HA	1:A:207:HIS:HA	1.93	0.51
1:A:984:LEU:O	1:C:386:LYS:NZ	2.44	0.51
1:B:329:PHE:HB3	1:B:526:GLY:HA3	1.92	0.51
1:B:377:PHE:CE2	1:C:983:ARG:HB3	2.46	0.51
1:C:458:LYS:HZ3	1:C:474:GLN:HB3	1.75	0.51
1:C:498:GLN:HB2	1:C:501:ASN:HB2	1.93	0.51
1:A:57:PRO:C	1:A:91:TYR:N	2.69	0.51
1:A:59:PHE:H	1:A:88:VAL:CB	2.14	0.51
1:A:387:LEU:H	1:B:983:ARG:HB2	1.76	0.51
1:A:388:ASN:N	1:B:983:ARG:HE	2.09	0.51
1:B:421:TYR:HA	1:B:467:ASP:N	2.26	0.51
1:C:555:SER:OG	1:C:584:ILE:HG22	2.11	0.51
1:A:329:PHE:H	1:A:542:ASN:N	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:GLU:HB3	1:B:204:TYR:O	2.11	0.51
1:C:116:SER:HB3	1:C:130:VAL:HG23	1.93	0.51
1:C:974:SER:OG	1:C:975:SER:N	2.42	0.51
1:A:386:LYS:HE3	1:B:992:GLN:HB3	1.92	0.50
1:A:394:ASN:N	1:A:518:LEU:HA	2.26	0.50
1:B:374:PHE:HB3	1:C:983:ARG:CZ	2.40	0.50
1:C:102:ARG:H	1:C:242:LEU:HD23	1.76	0.50
1:C:419:ALA:HA	1:C:422:ASN:N	2.27	0.50
1:A:586:ASP:CG	1:A:587:ILE:H	2.19	0.50
1:B:758:SER:OG	1:B:762:GLN:NE2	2.44	0.50
1:C:1000:ARG:O	1:C:1003:SER:OG	2.29	0.50
1:A:425:LEU:HD13	1:A:426:PRO:HD2	1.91	0.50
1:B:387:LEU:N	1:C:978:ASN:O	2.45	0.50
1:A:80:PRO:HD2	1:A:82:LEU:HD11	1.93	0.50
1:B:866:THR:H	1:B:869:MET:HE1	1.76	0.50
1:C:906:PHE:O	1:C:909:ILE:HG22	2.11	0.50
1:A:62:VAL:HG12	1:A:269:TYR:CG	2.45	0.50
1:A:324:GLU:HB3	1:A:537:LYS:NZ	2.26	0.50
1:B:697:MET:HE3	1:B:697:MET:HA	1.94	0.50
1:C:78:ASP:N	1:C:78:ASP:OD1	2.44	0.50
1:C:417:LYS:HE3	1:C:457:ARG:HB2	1.94	0.50
1:C:433:VAL:HA	1:C:510:VAL:HG23	1.93	0.50
1:C:450:ASN:OD1	1:C:451:TYR:N	2.44	0.50
1:A:329:PHE:CD2	1:A:541:PHE:HB3	2.46	0.50
1:A:390:LEU:O	1:A:520:ALA:N	2.44	0.50
1:B:468:ILE:N	1:C:234:ASN:HD22	2.10	0.50
1:B:761:THR:OG1	1:B:765:ARG:NH2	2.45	0.50
1:B:1106:GLN:NE2	1:B:1111:GLU:OE2	2.37	0.50
1:A:322:PRO:HA	1:A:324:GLU:CD	2.37	0.50
1:A:282:ASN:OD1	1:C:558:LYS:NZ	2.33	0.49
1:A:330:PRO:N	1:A:542:ASN:HA	2.26	0.49
1:A:298:GLU:OE1	1:A:315:THR:HG22	2.12	0.49
1:A:525:CYS:CB	1:A:545:GLY:N	2.75	0.49
1:B:32:PHE:HB2	1:B:34:ARG:HG2	1.93	0.49
1:B:326:ILE:N	1:B:540:ASN:O	2.42	0.49
1:A:1107:ARG:HD3	1:B:904:TYR:CZ	2.47	0.49
1:C:33:THR:HA	1:C:58:PHE:CE1	2.47	0.49
1:C:234:ASN:HD21	1:C:236:THR:HG23	1.77	0.49
1:C:1050:MET:HA	1:C:1050:MET:HE2	1.94	0.49
1:A:57:PRO:HB3	1:A:271:GLN:HG2	1.93	0.49
1:A:84:PHE:O	1:A:85:ASN:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:TYR:HB2	1:B:225:PRO:HD3	1.94	0.49
1:C:335:LEU:HD12	1:C:336:CYS:N	2.27	0.49
1:A:329:PHE:HA	1:A:543:PHE:H	1.76	0.49
1:B:468:ILE:H	1:C:234:ASN:HB3	1.77	0.49
1:B:577:ARG:HG2	1:B:583:GLU:O	2.12	0.49
1:B:703:ASN:HB2	1:C:787:GLN:NE2	2.28	0.49
1:C:103:GLY:O	1:C:240:THR:OG1	2.30	0.49
1:A:81:VAL:C	1:A:82:LEU:HG	2.37	0.49
1:A:529:LYS:NZ	1:A:530:SER:H	2.11	0.49
1:A:618:THR:OG1	1:A:619:GLU:OE1	2.24	0.49
1:C:665:PRO:HA	1:C:671:CYS:SG	2.53	0.49
1:A:189:LEU:HD23	1:A:190:ARG:H	1.77	0.49
1:A:391:CYS:HB2	1:A:521:PRO:C	2.37	0.49
1:A:990:GLU:HA	1:A:993:ILE:HD12	1.94	0.49
1:B:210:ILE:HG13	1:B:212:LEU:HD23	1.95	0.49
1:B:423:TYR:CD1	1:C:232:GLY:HA2	2.47	0.49
1:A:64:TRP:HH2	1:A:214:ARG:HH21	1.61	0.49
1:A:330:PRO:HD3	1:A:543:PHE:N	2.27	0.49
1:A:405:ASP:OD1	1:A:406:GLU:N	2.41	0.49
1:B:350:VAL:O	1:C:115:GLN:HB3	2.13	0.49
1:B:412:PRO:HB3	1:B:464:PHE:CG	2.48	0.49
1:B:424:LYS:HD3	1:C:84:PHE:HE2	1.76	0.49
1:B:521:PRO:HD3	1:B:559:PHE:HB3	1.95	0.49
1:C:428:ASP:OD1	1:C:428:ASP:N	2.39	0.49
1:A:216:LEU:CD2	1:A:265:TYR:HA	2.43	0.49
1:A:392:PHE:HB3	1:A:518:LEU:CA	2.41	0.49
1:B:347:PHE:CE2	1:B:509:ARG:HD3	2.48	0.49
1:B:982:SER:C	1:B:983:ARG:HG3	2.37	0.49
1:A:37:TYR:CE2	1:A:224:GLU:HB2	2.48	0.49
1:B:394:ASN:HA	1:C:200:TYR:OH	2.13	0.49
1:C:453:PHE:HE1	1:C:455:LEU:HD23	1.76	0.49
1:A:101:ILE:HB	1:A:189:LEU:HD21	1.94	0.48
1:A:473:TYR:HE2	1:B:236:THR:HG21	1.76	0.48
1:A:523:THR:H	1:A:544:ASN:CB	2.26	0.48
1:B:318:PHE:N	1:B:593:GLY:O	2.45	0.48
1:B:448:ASN:OD1	1:B:450:ASN:ND2	2.45	0.48
1:C:418:ILE:HB	1:C:424:LYS:HD2	1.94	0.48
1:C:448:ASN:N	1:C:494:SER:OG	2.46	0.48
1:A:38:TYR:HA	1:A:285:ILE:HD11	1.94	0.48
1:A:42:VAL:HG13	1:C:567:ARG:HB2	1.95	0.48
1:A:56:LEU:H	1:A:56:LEU:HD23	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:GLN:NE2	1:A:130:VAL:O	2.47	0.48
1:A:525:CYS:HA	1:A:545:GLY:N	2.29	0.48
1:A:654:GLU:OE1	1:A:655:HIS:N	2.46	0.48
1:A:1116:THR:OG1	1:A:1118:ASP:OD1	2.20	0.48
1:B:128:ILE:HD12	1:B:170:TYR:HD2	1.77	0.48
1:B:204:TYR:CD1	1:B:225:PRO:HA	2.49	0.48
1:B:325:SER:HA	1:B:540:ASN:HB3	1.95	0.48
1:B:915:VAL:O	1:B:919:ASN:ND2	2.44	0.48
1:C:993:ILE:HG13	1:C:994:ASP:N	2.28	0.48
1:A:1146:ASP:OD1	1:A:1147:SER:N	2.46	0.48
1:B:54:LEU:HD12	1:B:84:PHE:HD1	1.78	0.48
1:B:122:ASN:OD1	1:B:124:THR:OG1	2.31	0.48
1:C:521:PRO:HB2	1:C:577:ARG:HB2	1.95	0.48
1:A:38:TYR:HE1	1:A:279:TYR:CG	2.30	0.48
1:A:111:ASP:OD2	1:A:237:ARG:NH2	2.40	0.48
1:A:330:PRO:HD3	1:A:543:PHE:O	2.13	0.48
1:B:115:GLN:HB3	1:B:130:VAL:O	2.13	0.48
1:A:58:PHE:H	1:A:87:GLY:CA	2.27	0.48
1:B:758:SER:HB2	1:B:761:THR:OG1	2.13	0.48
1:A:34:ARG:N	1:A:34:ARG:NE	2.61	0.48
1:A:388:ASN:H	1:B:983:ARG:HE	1.61	0.48
1:B:377:PHE:CZ	1:B:387:LEU:HD22	2.48	0.48
1:B:993:ILE:O	1:B:997:ILE:HG23	2.14	0.48
1:C:393:THR:HG21	1:C:516:GLU:HB3	1.96	0.48
1:A:516:GLU:CG	1:A:517:LEU:H	2.26	0.48
1:A:529:LYS:HZ3	1:A:530:SER:H	1.61	0.48
1:A:931:ILE:O	1:A:934:ILE:HG22	2.14	0.48
1:B:325:SER:OG	1:B:540:ASN:OD1	2.31	0.48
1:B:421:TYR:HB3	1:B:468:ILE:HA	1.95	0.48
1:B:518:LEU:H	1:B:564:GLN:HB3	1.79	0.48
1:B:598:ILE:HG23	1:B:609:ALA:HB3	1.95	0.48
1:B:974:SER:HB3	1:B:979:ASP:CG	2.38	0.48
1:A:108:THR:H	1:A:237:ARG:HD2	1.78	0.48
1:A:186:PHE:CG	1:A:187:LYS:N	2.81	0.48
1:A:212:LEU:HD23	1:A:212:LEU:H	1.79	0.48
1:A:216:LEU:HD23	1:A:216:LEU:HA	1.70	0.48
1:C:87:GLY:HA3	1:C:270:LEU:HB2	1.95	0.48
1:C:498:GLN:OE1	1:C:501:ASN:ND2	2.46	0.48
1:A:237:ARG:C	1:A:238:PHE:HD1	2.21	0.48
1:A:387:LEU:N	1:B:983:ARG:HE	2.11	0.48
1:B:56:LEU:HD22	1:B:91:TYR:CB	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:994:ASP:HA	1:B:997:ILE:HG12	1.94	0.48
1:B:1077:THR:OG1	1:B:1078:ALA:N	2.46	0.48
1:A:365:TYR:CD1	1:A:368:LEU:HD12	2.49	0.47
1:A:905:ARG:NH1	1:A:1049:LEU:O	2.47	0.47
1:C:454:ARG:O	1:C:455:LEU:HD22	2.14	0.47
1:C:555:SER:HB3	1:C:586:ASP:OD2	2.14	0.47
1:A:101:ILE:HG22	1:A:102:ARG:N	2.28	0.47
1:A:194:PHE:C	1:A:195:LYS:HD2	2.40	0.47
1:B:352:ALA:HB2	1:C:132:GLU:OE2	2.14	0.47
1:B:425:LEU:HD12	1:B:426:PRO:HD2	1.96	0.47
1:B:527:PRO:HD2	1:B:528:LYS:NZ	2.29	0.47
1:B:537:LYS:HA	1:B:537:LYS:HD3	1.67	0.47
1:A:36:VAL:HG22	1:A:37:TYR:O	2.14	0.47
1:A:220:PHE:O	1:A:220:PHE:CG	2.66	0.47
1:A:386:LYS:O	1:B:973:ILE:HD11	2.14	0.47
1:A:703:ASN:OD1	1:A:704:SER:N	2.47	0.47
1:B:62:VAL:HG21	1:B:266:TYR:HB3	1.96	0.47
1:B:1029:MET:HB2	1:B:1062:PHE:HZ	1.79	0.47
1:C:748:GLU:HA	1:C:751:ASN:HD21	1.80	0.47
1:C:989:ALA:O	1:C:993:ILE:HG12	2.15	0.47
1:A:78:ASP:O	1:A:266:TYR:HA	2.15	0.47
1:A:101:ILE:O	1:A:189:LEU:HA	2.14	0.47
1:A:391:CYS:O	1:A:544:ASN:C	2.57	0.47
1:A:905:ARG:NH1	1:A:1050:MET:HG2	2.29	0.47
1:B:203:ILE:HB	1:B:227:VAL:HB	1.97	0.47
1:B:421:TYR:CE2	1:B:462:LYS:HD2	2.50	0.47
1:B:468:ILE:HG22	1:C:234:ASN:CB	2.37	0.47
1:C:403:ARG:CZ	1:C:453:PHE:HB2	2.44	0.47
1:A:57:PRO:HD3	1:A:269:TYR:O	2.13	0.47
1:A:60:SER:HB3	1:A:88:VAL:O	2.13	0.47
1:A:325:SER:OG	1:A:539:VAL:HG12	2.15	0.47
1:A:536:ASN:H	1:A:553:THR:HA	1.78	0.47
1:A:37:TYR:CD2	1:A:224:GLU:HB2	2.49	0.47
1:A:331:ASN:HB3	1:A:528:LYS:O	2.14	0.47
1:A:535:LYS:HE2	1:A:585:LEU:HB3	1.96	0.47
1:C:346:ARG:NH2	1:C:444:LYS:HD2	2.29	0.47
1:A:96:GLU:HB3	1:A:211:ASN:H	1.78	0.47
1:A:129:LYS:HB2	1:A:133:PHE:CZ	2.50	0.47
1:A:189:LEU:HD12	1:A:210:ILE:HG12	1.96	0.47
1:A:410:ILE:HG23	1:C:987:PRO:O	2.14	0.47
1:B:421:TYR:OH	1:B:462:LYS:HD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:567:ARG:HG2	1:B:572:THR:O	2.15	0.47
1:C:969:ASN:HA	1:C:974:SER:O	2.14	0.47
1:A:35:GLY:HA3	1:A:224:GLU:OE2	2.15	0.47
1:A:271:GLN:OE1	1:A:273:ARG:NH2	2.46	0.47
1:A:383:SER:HB2	1:B:986:PRO:HD3	1.97	0.47
1:A:643:PHE:CE2	1:A:645:THR:HG22	2.49	0.47
1:A:813:SER:OG	1:A:814:LYS:N	2.47	0.47
1:B:448:ASN:HB2	1:B:507:PRO:HB3	1.97	0.47
1:B:578:ASP:N	1:B:583:GLU:O	2.32	0.47
1:B:821:LEU:O	1:B:825:LYS:HG2	2.15	0.47
1:B:1139:ASP:HB3	1:B:1142:GLN:HE22	1.79	0.47
1:C:436:TRP:CE3	1:C:438:SER:HB3	2.49	0.47
1:A:41:LYS:HG2	1:C:565:PHE:HB3	1.96	0.47
1:A:80:PRO:O	1:A:239:GLN:NE2	2.47	0.47
1:A:326:ILE:HD11	1:A:528:LYS:NZ	2.29	0.47
1:A:386:LYS:NZ	1:B:973:ILE:HG12	2.30	0.47
1:A:525:CYS:SG	1:A:544:ASN:N	2.87	0.47
1:C:409:GLN:O	1:C:410:ILE:HD13	2.15	0.47
1:A:759:PHE:O	1:A:763:LEU:HG	2.15	0.47
1:A:32:PHE:CD1	1:A:34:ARG:HD2	2.49	0.46
1:A:391:CYS:HB3	1:A:544:ASN:H	1.80	0.46
1:B:659:SER:HB2	1:B:698:SER:HB2	1.96	0.46
1:C:518:LEU:HB2	1:C:564:GLN:HB2	1.97	0.46
1:A:170:TYR:CE1	1:A:229:LEU:HD21	2.50	0.46
1:A:790:LYS:HE2	1:C:704:SER:HB3	1.97	0.46
1:B:228:ASP:C	1:B:229:LEU:HD22	2.40	0.46
1:B:397:ALA:C	1:B:510:VAL:HG11	2.40	0.46
1:A:204:TYR:CG	1:A:224:GLU:O	2.68	0.46
1:B:697:MET:HE3	1:B:698:SER:H	1.81	0.46
1:B:796:ASP:OD1	1:B:796:ASP:N	2.48	0.46
1:B:1004:LEU:O	1:B:1008:VAL:HG13	2.15	0.46
1:C:905:ARG:CZ	1:C:1050:MET:HE1	2.45	0.46
1:C:1097:SER:HB2	1:C:1102:TRP:CE3	2.50	0.46
1:A:56:LEU:O	1:A:86:ASP:C	2.56	0.46
1:A:196:ASN:HD21	1:A:199:GLY:H	1.62	0.46
1:A:328:ARG:HG3	1:A:521:PRO:HB3	1.98	0.46
1:B:82:LEU:HB2	1:B:238:PHE:HE1	1.79	0.46
1:C:392:PHE:CD2	1:C:513:LEU:HD22	2.51	0.46
1:C:462:LYS:HB3	1:C:465:GLU:HB3	1.97	0.46
1:A:390:LEU:CA	1:A:520:ALA:HB3	2.45	0.46
1:A:451:TYR:HB3	1:A:495:TYR:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1111:GLU:OE1	1:A:1113:GLN:NE2	2.49	0.46
1:B:281:GLU:O	1:B:282:ASN:HB3	2.15	0.46
1:B:428:ASP:OD1	1:B:428:ASP:N	2.44	0.46
1:B:555:SER:OG	1:B:584:ILE:O	2.30	0.46
1:C:616:ASN:N	1:C:619:GLU:OE2	2.36	0.46
1:A:56:LEU:O	1:A:87:GLY:CA	2.62	0.46
1:A:88:VAL:O	1:A:88:VAL:HG12	2.15	0.46
1:A:453:PHE:O	1:A:493:GLN:N	2.45	0.46
1:B:203:ILE:N	1:B:227:VAL:O	2.46	0.46
1:C:53:ASP:OD1	1:C:54:LEU:N	2.38	0.46
1:C:117:LEU:HD12	1:C:117:LEU:HA	1.79	0.46
1:A:51:THR:OG1	1:A:52:GLN:N	2.48	0.46
1:A:330:PRO:HG3	1:A:542:ASN:CG	2.40	0.46
1:A:391:CYS:O	1:A:544:ASN:HB3	2.16	0.46
1:B:280:ASN:ND2	1:B:284:THR:OG1	2.49	0.46
1:B:997:ILE:HG13	1:B:998:THR:N	2.30	0.46
1:B:1097:SER:HB2	1:B:1102:TRP:CE3	2.51	0.46
1:A:64:TRP:HZ2	1:A:242:LEU:HD21	1.80	0.46
1:A:391:CYS:CB	1:A:522:ALA:N	2.74	0.46
1:C:117:LEU:HD22	1:C:231:ILE:HD11	1.98	0.46
1:A:396:TYR:HD1	1:A:397:ALA:N	2.07	0.46
1:A:996:LEU:HD23	1:A:996:LEU:HA	1.81	0.46
1:C:355:ARG:HA	1:C:396:TYR:HB2	1.98	0.46
1:C:977:LEU:O	1:C:977:LEU:HD23	2.16	0.46
1:A:993:ILE:O	1:A:997:ILE:HG12	2.16	0.46
1:A:1097:SER:HB2	1:A:1102:TRP:CE3	2.51	0.46
1:B:127:VAL:HA	1:B:171:VAL:HG22	1.98	0.46
1:B:383:SER:N	1:C:976:VAL:O	2.37	0.46
1:B:578:ASP:HB2	1:B:583:GLU:HB3	1.97	0.46
1:C:417:LYS:O	1:C:417:LYS:HG2	2.16	0.46
1:A:312:ILE:HB	1:A:598:ILE:HD13	1.98	0.45
1:A:327:VAL:HA	1:A:542:ASN:H	1.76	0.45
1:A:385:THR:OG1	1:B:984:LEU:HB3	2.16	0.45
1:A:755:GLN:O	1:C:970:PHE:N	2.48	0.45
1:C:104:TRP:HA	1:C:240:THR:OG1	2.16	0.45
1:A:394:ASN:H	1:A:518:LEU:HA	1.82	0.45
1:A:523:THR:HB	1:A:576:VAL:HB	1.98	0.45
1:B:199:GLY:HA3	1:B:232:GLY:HA2	1.98	0.45
1:C:419:ALA:H	1:C:424:LYS:CD	2.25	0.45
1:C:474:GLN:HE22	1:C:491:PRO:HG2	1.81	0.45
1:C:773:GLU:OE1	1:C:774:GLN:NE2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1048:HIS:NE2	1:C:1050:MET:O	2.49	0.45
1:C:1097:SER:HB2	1:C:1102:TRP:CD2	2.51	0.45
1:A:56:LEU:CD1	1:A:268:GLY:H	2.30	0.45
1:A:57:PRO:HB2	1:A:91:TYR:CB	2.43	0.45
1:A:190:ARG:HG2	1:A:207:HIS:HB3	1.98	0.45
1:C:474:GLN:OE1	1:C:475:ALA:N	2.50	0.45
1:A:91:TYR:O	1:A:91:TYR:CG	2.69	0.45
1:A:412:PRO:O	1:C:990:GLU:HB2	2.16	0.45
1:A:461:LEU:HD23	1:A:461:LEU:HA	1.80	0.45
1:A:529:LYS:HA	1:A:529:LYS:HD2	1.70	0.45
1:A:1001:LEU:HD12	1:A:1001:LEU:HA	1.82	0.45
1:C:714:ILE:HD12	1:C:715:PRO:HD2	1.99	0.45
1:A:900:MET:HE3	1:A:900:MET:HB2	1.83	0.45
1:B:119:ILE:HA	1:B:127:VAL:O	2.16	0.45
1:B:167:THR:O	1:B:167:THR:OG1	2.30	0.45
1:B:387:LEU:O	1:B:389:ASP:N	2.50	0.45
1:B:509:ARG:HG3	1:B:510:VAL:N	2.31	0.45
1:C:197:ILE:O	1:C:200:TYR:HD2	1.99	0.45
1:A:96:GLU:HG3	1:A:215:ASP:OD1	2.16	0.45
1:A:241:LEU:HD23	1:A:241:LEU:HA	1.81	0.45
1:A:386:LYS:H	1:B:983:ARG:HB2	1.79	0.45
1:A:517:LEU:C	1:A:518:LEU:HD23	2.42	0.45
1:A:31:SER:O	1:A:220:PHE:N	2.49	0.45
1:A:42:VAL:CG1	1:C:567:ARG:HB2	2.47	0.45
1:A:102:ARG:HA	1:A:102:ARG:HD3	1.68	0.45
1:A:662:CYS:HB2	1:A:697:MET:HE3	1.99	0.45
1:A:766:ALA:O	1:A:770:ILE:HG23	2.17	0.45
1:B:658:ASN:OD1	1:B:659:SER:N	2.49	0.45
1:C:106:PHE:CD1	1:C:235:ILE:HD13	2.52	0.45
1:C:241:LEU:HD12	1:C:242:LEU:O	2.16	0.45
1:C:417:LYS:HA	1:C:420:ASP:HB3	1.99	0.45
1:A:327:VAL:C	1:A:542:ASN:H	2.20	0.45
1:A:56:LEU:HD21	1:A:88:VAL:HG13	1.98	0.45
1:A:950:ASP:O	1:A:954:GLN:HG2	2.17	0.45
1:C:30:ASN:HD21	1:C:59:PHE:C	2.25	0.45
1:C:328:ARG:NH1	1:C:580:GLN:OE1	2.48	0.45
1:C:333:THR:HG22	1:C:527:PRO:CB	2.47	0.45
1:A:522:ALA:N	1:A:543:PHE:HB3	2.31	0.45
1:A:929:SER:O	1:A:933:LYS:HG2	2.16	0.45
1:B:521:PRO:HA	1:B:563:GLN:HB2	1.98	0.45
1:B:912:THR:HG23	1:B:915:VAL:HG23	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:405:ASP:OD1	1:C:504:GLY:HA3	2.17	0.45
1:A:545:GLY:O	1:A:546:LEU:HG	2.17	0.44
1:B:347:PHE:CD2	1:B:509:ARG:HD3	2.52	0.44
1:B:380:TYR:CE2	1:C:992:GLN:HG3	2.52	0.44
1:B:740:MET:SD	1:B:741:TYR:N	2.90	0.44
1:C:189:LEU:HD23	1:C:189:LEU:HA	1.80	0.44
1:A:57:PRO:HG3	1:A:271:GLN:N	2.30	0.44
1:A:60:SER:HG	1:A:268:GLY:HA3	1.79	0.44
1:A:383:SER:CB	1:B:985:ASP:HA	2.48	0.44
1:B:272:PRO:O	1:B:292:ALA:HB3	2.17	0.44
1:B:383:SER:O	1:C:979:ASP:N	2.51	0.44
1:B:780:GLU:O	1:B:784:GLN:NE2	2.51	0.44
1:A:528:LYS:NZ	1:A:541:PHE:HB2	2.32	0.44
1:B:422:ASN:HB3	1:B:466:ARG:CD	2.47	0.44
1:B:521:PRO:HB3	1:B:577:ARG:HG3	1.99	0.44
1:B:541:PHE:HE2	1:B:543:PHE:HB2	1.83	0.44
1:C:290:ASP:OD1	1:C:291:CYS:N	2.51	0.44
1:C:518:LEU:HD23	1:C:518:LEU:HA	1.82	0.44
1:C:772:VAL:O	1:C:776:LYS:HG3	2.17	0.44
1:A:58:PHE:O	1:A:92:PHE:CD2	2.71	0.44
1:A:119:ILE:HG23	1:A:127:VAL:O	2.17	0.44
1:A:328:ARG:H	1:A:542:ASN:N	2.08	0.44
1:B:563:GLN:OE1	1:B:565:PHE:N	2.50	0.44
1:C:106:PHE:HD1	1:C:235:ILE:HD13	1.83	0.44
1:A:104:TRP:N	1:A:191:GLU:HA	2.21	0.44
1:A:271:GLN:OE1	1:A:273:ARG:NE	2.46	0.44
1:B:999:GLY:HA2	1:B:1002:GLN:HE22	1.82	0.44
1:A:34:ARG:HB3	1:A:206:LYS:HE3	2.00	0.44
1:A:189:LEU:HD23	1:A:190:ARG:N	2.33	0.44
1:A:270:LEU:HD12	1:A:271:GLN:N	2.33	0.44
1:A:723:THR:OG1	1:A:724:THR:N	2.49	0.44
1:B:822:LEU:O	1:B:826:VAL:HG12	2.18	0.44
1:C:55:PHE:N	1:C:271:GLN:O	2.42	0.44
1:B:36:VAL:HA	1:B:88:VAL:CG1	2.48	0.44
1:B:231:ILE:HD13	1:B:231:ILE:HA	1.85	0.44
1:B:770:ILE:HD11	1:B:1012:LEU:HD22	2.00	0.44
1:C:101:ILE:HD11	1:C:240:THR:HG23	2.00	0.44
1:C:104:TRP:CE3	1:C:104:TRP:C	2.96	0.44
1:A:92:PHE:HZ	1:A:219:GLY:O	2.01	0.44
1:A:295:PRO:O	1:A:299:THR:HG23	2.18	0.44
1:A:1118:ASP:OD1	1:A:1118:ASP:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ASN:ND2	1:B:58:PHE:HB3	2.33	0.44
1:B:61:ASN:HB3	1:B:269:TYR:CD2	2.49	0.44
1:B:129:LYS:HG2	1:B:169:GLU:OE1	2.18	0.44
1:B:200:TYR:HA	1:B:229:LEU:O	2.18	0.44
1:A:206:LYS:HG2	1:A:207:HIS:N	2.31	0.44
1:A:290:ASP:O	1:A:297:SER:HB3	2.18	0.44
1:C:417:LYS:NZ	1:C:473:TYR:HD1	2.16	0.44
1:A:330:PRO:HG3	1:A:542:ASN:CB	2.48	0.43
1:A:738:CYS:HB3	1:A:760:CYS:HB2	1.84	0.43
1:B:327:VAL:HA	1:B:542:ASN:HB3	2.00	0.43
1:B:738:CYS:SG	1:B:739:THR:N	2.91	0.43
1:C:968:SER:HB3	1:C:970:PHE:CD1	2.53	0.43
1:A:66:HIS:CG	1:A:67:ALA:N	2.86	0.43
1:A:96:GLU:HB2	1:A:216:LEU:O	2.18	0.43
1:A:289:VAL:HG12	1:A:290:ASP:O	2.18	0.43
1:A:339:GLY:HA3	1:A:368:LEU:HD22	2.00	0.43
1:A:877:LEU:O	1:A:881:THR:HG22	2.17	0.43
1:B:29:THR:HG22	1:B:266:TYR:HD1	1.80	0.43
1:B:392:PHE:CD2	1:B:393:THR:N	2.84	0.43
1:B:568:ASP:CG	1:B:574:ASP:HB2	2.43	0.43
1:B:1002:GLN:H	1:B:1002:GLN:CD	2.21	0.43
1:C:106:PHE:HB3	1:C:235:ILE:CG2	2.48	0.43
1:C:128:ILE:H	1:C:170:TYR:HD2	1.66	0.43
1:C:394:ASN:ND2	1:C:514:SER:HB2	2.33	0.43
1:C:866:THR:H	1:C:869:MET:HE1	1.83	0.43
1:A:59:PHE:O	1:A:59:PHE:HD2	2.00	0.43
1:A:96:GLU:HB3	1:A:210:ILE:HB	2.00	0.43
1:A:195:LYS:HE2	1:A:238:PHE:HE1	1.82	0.43
1:B:62:VAL:HG12	1:B:63:THR:N	2.34	0.43
1:A:385:THR:O	1:B:983:ARG:HD3	2.19	0.43
1:B:54:LEU:O	1:B:88:VAL:N	2.52	0.43
1:C:338:PHE:N	1:C:338:PHE:CD1	2.86	0.43
1:A:54:LEU:C	1:A:238:PHE:HE2	2.26	0.43
1:A:55:PHE:O	1:A:56:LEU:O	2.36	0.43
1:A:327:VAL:CG2	1:A:552:LEU:HB2	2.45	0.43
1:A:536:ASN:OD1	1:A:553:THR:HB	2.19	0.43
1:C:738:CYS:HB3	1:C:760:CYS:HB2	1.96	0.43
1:C:753:LEU:HD12	1:C:997:ILE:HD11	2.00	0.43
1:A:57:PRO:HD2	1:A:91:TYR:H	1.83	0.43
1:A:99:ASN:ND2	1:A:188:ASN:O	2.52	0.43
1:A:107:GLY:O	1:A:239:GLN:HB3	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASN:OD1	1:A:137:ASN:N	2.51	0.43
1:B:48:LEU:HA	1:B:277:LEU:O	2.19	0.43
1:C:703:ASN:OD1	1:C:704:SER:N	2.50	0.43
1:C:1029:MET:HB2	1:C:1062:PHE:HZ	1.82	0.43
1:A:523:THR:N	1:A:544:ASN:N	2.67	0.43
1:B:113:LYS:O	1:B:113:LYS:HG3	2.18	0.43
1:A:97:LYS:HE3	1:A:98:SER:HB3	2.01	0.43
1:A:411:ALA:N	1:C:988:GLU:HA	2.30	0.43
1:A:543:PHE:CE2	1:A:552:LEU:HG	2.53	0.43
1:A:550:GLY:O	1:A:551:VAL:HG23	2.19	0.43
1:B:454:ARG:HG2	1:B:456:PHE:H	1.84	0.43
1:C:498:GLN:H	1:C:505:TYR:HD2	1.67	0.43
1:C:522:ALA:HA	1:C:563:GLN:NE2	2.34	0.43
1:C:559:PHE:O	1:C:560:LEU:HD13	2.18	0.43
1:A:55:PHE:O	1:A:86:ASP:CA	2.67	0.43
1:A:140:PHE:CZ	1:A:143:VAL:HB	2.54	0.43
1:A:423:TYR:HA	1:A:463:PRO:HB3	2.00	0.43
1:C:92:PHE:CE1	1:C:94:SER:HB3	2.54	0.43
1:C:346:ARG:HD2	1:C:346:ARG:HA	1.64	0.43
1:A:54:LEU:HD13	1:A:84:PHE:HE1	1.83	0.43
1:A:99:ASN:ND2	1:A:186:PHE:HB3	2.34	0.43
1:A:126:VAL:HB	1:A:172:SER:O	2.18	0.43
1:A:328:ARG:CD	1:A:521:PRO:HB3	2.49	0.43
1:A:458:LYS:HA	1:A:458:LYS:HD2	1.85	0.43
1:C:644:GLN:NE2	1:C:648:GLY:O	2.52	0.43
1:C:725:GLU:N	1:C:725:GLU:OE1	2.52	0.43
1:A:786:LYS:HA	1:A:786:LYS:HD3	1.86	0.42
1:B:85:ASN:CG	1:B:86:ASP:H	2.27	0.42
1:B:92:PHE:HB3	1:B:191:GLU:OE1	2.19	0.42
1:B:731:MET:SD	1:B:731:MET:N	2.91	0.42
1:B:784:GLN:N	1:B:784:GLN:OE1	2.52	0.42
1:C:448:ASN:O	1:C:494:SER:HA	2.18	0.42
1:A:100:ILE:HG13	1:A:101:ILE:HG12	2.00	0.42
1:A:116:SER:H	1:A:118:LEU:HG	1.84	0.42
1:A:366:SER:HA	1:A:369:TYR:CE2	2.54	0.42
1:A:384:PRO:HA	1:B:982:SER:CA	2.49	0.42
1:A:392:PHE:HA	1:A:517:LEU:CD1	2.49	0.42
1:A:869:MET:H	1:A:869:MET:HG3	1.63	0.42
1:A:1005:GLN:O	1:A:1009:THR:OG1	2.34	0.42
1:B:54:LEU:HA	1:B:84:PHE:O	2.18	0.42
1:B:105:ILE:HG22	1:B:110:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:541:PHE:CE2	1:B:543:PHE:HB2	2.54	0.42
1:B:729:VAL:HG23	1:B:1059:GLY:HA2	2.01	0.42
1:B:866:THR:H	1:B:869:MET:CE	2.31	0.42
1:C:190:ARG:HD2	1:C:192:PHE:HZ	1.84	0.42
1:A:328:ARG:CZ	1:A:521:PRO:HG3	2.49	0.42
1:C:58:PHE:HB2	1:C:275:PHE:CZ	2.54	0.42
1:C:877:LEU:O	1:C:881:THR:HG22	2.20	0.42
1:A:129:LYS:HE2	1:A:170:TYR:O	2.19	0.42
1:A:392:PHE:CB	1:A:518:LEU:N	2.70	0.42
1:A:546:LEU:HD23	1:A:546:LEU:HA	1.65	0.42
1:A:553:THR:OG1	1:A:586:ASP:HB3	2.20	0.42
1:A:741:TYR:CE1	1:A:966:LEU:HD21	2.54	0.42
1:B:556:ASN:N	1:B:556:ASN:OD1	2.51	0.42
1:B:1139:ASP:OD1	1:B:1140:PRO:HD3	2.19	0.42
1:C:107:GLY:HA2	1:C:116:SER:OG	2.18	0.42
1:C:141:LEU:HD21	1:C:143:VAL:HB	2.01	0.42
1:C:190:ARG:HB3	1:C:192:PHE:CE1	2.55	0.42
1:A:62:VAL:HG12	1:A:269:TYR:CD1	2.55	0.42
1:A:88:VAL:HG11	1:A:266:TYR:OH	2.20	0.42
1:A:189:LEU:CD1	1:A:210:ILE:HG12	2.49	0.42
1:A:191:GLU:C	1:A:192:PHE:HD1	2.28	0.42
1:A:272:PRO:O	1:A:273:ARG:HD3	2.18	0.42
1:A:319:ARG:HG2	1:A:320:VAL:N	2.34	0.42
1:A:523:THR:C	1:A:543:PHE:C	2.88	0.42
1:B:130:VAL:HG12	1:B:166:CYS:HB2	2.01	0.42
1:B:200:TYR:CE2	1:B:230:PRO:HB3	2.55	0.42
1:C:466:ARG:HG2	1:C:466:ARG:H	1.57	0.42
1:A:58:PHE:N	1:A:87:GLY:HA3	2.34	0.42
1:A:229:LEU:HD23	1:A:229:LEU:HA	1.90	0.42
1:A:524:VAL:H	1:A:544:ASN:CA	2.22	0.42
1:A:524:VAL:C	1:A:543:PHE:O	2.61	0.42
1:B:126:VAL:H	1:B:172:SER:C	2.27	0.42
1:B:382:VAL:HG12	1:C:979:ASP:H	1.84	0.42
1:C:333:THR:HG22	1:C:527:PRO:HB3	2.01	0.42
1:A:64:TRP:CZ2	1:A:242:LEU:HD21	2.54	0.42
1:A:271:GLN:HA	1:A:272:PRO:HD3	1.80	0.42
1:A:576:VAL:HG22	1:A:585:LEU:HB2	2.01	0.42
1:B:62:VAL:HG22	1:B:267:VAL:C	2.45	0.42
1:A:40:ASP:O	1:C:519:HIS:ND1	2.52	0.42
1:A:290:ASP:OD2	1:A:293:LEU:HB2	2.19	0.42
1:A:516:GLU:O	1:A:518:LEU:HG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:716:THR:N	1:A:1071:GLN:O	2.48	0.42
1:A:759:PHE:HA	1:A:762:GLN:NE2	2.35	0.42
1:B:96:GLU:OE1	1:B:242:LEU:HD22	2.20	0.42
1:B:415:THR:HG23	1:B:461:LEU:HD23	2.02	0.42
1:B:577:ARG:HG2	1:B:578:ASP:H	1.85	0.42
1:C:135:PHE:HD1	1:C:136:CYS:N	2.02	0.42
1:C:556:ASN:OD1	1:C:556:ASN:N	2.52	0.42
1:A:66:HIS:HD2	1:A:68:ILE:HG12	1.83	0.42
1:A:359:SER:N	1:A:361:CYS:SG	2.93	0.42
1:B:577:ARG:NE	1:B:578:ASP:O	2.44	0.42
1:C:57:PRO:HB2	1:C:60:SER:OG	2.20	0.42
1:C:406:GLU:OE2	1:C:455:LEU:HD21	2.20	0.42
1:C:418:ILE:HG22	1:C:423:TYR:O	2.19	0.42
1:C:458:LYS:HD3	1:C:458:LYS:HA	1.56	0.42
1:C:773:GLU:OE2	1:C:777:ASN:ND2	2.53	0.42
1:C:1125:ASN:OD1	1:C:1125:ASN:N	2.53	0.42
1:A:140:PHE:CD1	1:A:141:LEU:N	2.88	0.42
1:B:743:CYS:O	1:B:980:ILE:HD11	2.19	0.42
1:C:126:VAL:O	1:C:172:SER:HB2	2.20	0.42
1:C:204:TYR:HB3	1:C:223:LEU:HD11	2.02	0.42
1:A:41:LYS:HD2	1:C:517:LEU:HD23	2.02	0.41
1:A:366:SER:HA	1:A:369:TYR:HD2	1.81	0.41
1:A:429:PHE:O	1:A:430:THR:OG1	2.33	0.41
1:B:62:VAL:HG11	1:B:266:TYR:CB	2.49	0.41
1:B:377:PHE:CB	1:C:984:LEU:HB2	2.49	0.41
1:C:328:ARG:HB3	1:C:579:PRO:HG3	2.02	0.41
1:C:520:ALA:HB3	1:C:563:GLN:CA	2.47	0.41
1:C:969:ASN:HB3	1:C:972:ALA:O	2.20	0.41
1:A:364:ASP:OD1	1:A:365:TYR:N	2.53	0.41
1:A:383:SER:OG	1:B:985:ASP:OD1	2.37	0.41
1:A:391:CYS:HB3	1:A:544:ASN:HB2	2.02	0.41
1:B:450:ASN:HB3	1:B:451:TYR:CE1	2.54	0.41
1:B:738:CYS:HB3	1:B:760:CYS:HB2	1.83	0.41
1:C:234:ASN:CG	1:C:235:ILE:N	2.79	0.41
1:C:323:THR:HB	1:C:539:VAL:HG12	2.03	0.41
1:C:418:ILE:HG22	1:C:423:TYR:HD1	1.85	0.41
1:A:129:LYS:HD3	1:A:169:GLU:HG2	2.01	0.41
1:B:81:VAL:C	1:B:82:LEU:HD12	2.46	0.41
1:B:423:TYR:HD1	1:C:233:ILE:H	1.67	0.41
1:B:914:ASN:C	1:B:914:ASN:OD1	2.63	0.41
1:A:973:ILE:HD11	1:A:980:ILE:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1004:LEU:HD23	1:A:1004:LEU:HA	1.76	0.41
1:B:30:ASN:HB2	1:B:59:PHE:CG	2.55	0.41
1:B:36:VAL:HA	1:B:88:VAL:HG11	2.03	0.41
1:B:112:SER:O	1:B:112:SER:OG	2.35	0.41
1:B:131:CYS:HB3	1:B:166:CYS:HB3	1.50	0.41
1:B:377:PHE:HB3	1:C:984:LEU:HD13	2.01	0.41
1:B:468:ILE:HG23	1:C:108:THR:CG2	2.40	0.41
1:B:969:ASN:HB2	1:C:755:GLN:CD	2.44	0.41
1:A:59:PHE:CA	1:A:88:VAL:HB	2.50	0.41
1:A:168:PHE:O	1:A:170:TYR:HD1	2.04	0.41
1:B:414:GLN:HA	1:B:461:LEU:HD22	2.02	0.41
1:B:947:LYS:HB2	1:B:947:LYS:HE2	1.76	0.41
1:B:969:ASN:HB3	1:B:972:ALA:O	2.21	0.41
1:C:377:PHE:HE1	1:C:434:ILE:HG13	1.86	0.41
1:A:200:TYR:CD1	1:A:230:PRO:HA	2.56	0.41
1:A:216:LEU:HD21	1:A:265:TYR:HA	2.01	0.41
1:A:391:CYS:SG	1:A:545:GLY:CA	3.07	0.41
1:A:394:ASN:HB3	1:A:395:VAL:H	1.74	0.41
1:A:1107:ARG:HD3	1:B:904:TYR:CE2	2.55	0.41
1:B:577:ARG:HG2	1:B:578:ASP:N	2.36	0.41
1:B:773:GLU:OE1	1:B:773:GLU:C	2.64	0.41
1:A:105:ILE:HA	1:A:192:PHE:O	2.21	0.41
1:A:116:SER:N	1:A:118:LEU:HG	2.36	0.41
1:A:197:ILE:C	1:A:199:GLY:H	2.28	0.41
1:A:273:ARG:HG3	1:A:290:ASP:OD2	2.21	0.41
1:A:329:PHE:HB2	1:A:525:CYS:CB	2.50	0.41
1:A:523:THR:H	1:A:544:ASN:N	2.18	0.41
1:A:745:ASP:OD1	1:C:319:ARG:NH2	2.45	0.41
1:B:388:ASN:ND2	1:C:982:SER:HB2	2.36	0.41
1:B:389:ASP:HB3	1:B:542:ASN:ND2	2.34	0.41
1:B:959:LEU:HA	1:B:959:LEU:HD23	1.87	0.41
1:C:413:GLY:HA2	1:C:428:ASP:O	2.21	0.41
1:C:523:THR:HG23	1:C:577:ARG:HB3	2.02	0.41
1:A:98:SER:OG	1:A:99:ASN:N	2.53	0.41
1:B:37:TYR:HA	1:B:223:LEU:O	2.21	0.41
1:C:51:THR:OG1	1:C:52:GLN:N	2.53	0.41
1:C:867:ASP:N	1:C:867:ASP:OD1	2.52	0.41
1:C:1049:LEU:O	1:C:1050:MET:HE2	2.20	0.41
1:A:141:LEU:O	1:A:143:VAL:N	2.52	0.41
1:A:463:PRO:O	1:B:234:ASN:HB2	2.21	0.41
1:A:516:GLU:HG2	1:A:517:LEU:N	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:539:VAL:HG22	1:A:550:GLY:O	2.20	0.41
1:A:731:MET:SD	1:A:731:MET:N	2.94	0.41
1:A:1001:LEU:O	1:A:1005:GLN:HG2	2.21	0.41
1:B:426:PRO:HG2	1:B:428:ASP:CG	2.45	0.41
1:B:521:PRO:HA	1:B:563:GLN:CB	2.51	0.41
1:B:521:PRO:CD	1:B:559:PHE:HB3	2.50	0.41
1:B:900:MET:HE2	1:B:900:MET:HA	2.02	0.41
1:B:1014:ARG:O	1:B:1017:GLU:HG3	2.21	0.41
1:C:130:VAL:CG2	1:C:132:GLU:HB3	2.51	0.41
1:C:354:ASN:O	1:C:398:ASP:HA	2.21	0.41
1:C:421:TYR:CG	1:C:457:ARG:HD2	2.55	0.41
1:C:984:LEU:HD21	1:C:988:GLU:OE1	2.21	0.41
1:A:64:TRP:CE2	1:A:79:ASN:HB3	2.56	0.41
1:A:114:THR:OG1	1:A:115:GLN:N	2.52	0.41
1:A:525:CYS:SG	1:A:543:PHE:O	2.79	0.41
1:A:63:THR:O	1:A:80:PRO:HD3	2.21	0.40
1:A:520:ALA:HB1	1:A:521:PRO:CD	2.49	0.40
1:A:661:GLU:O	1:A:695:TYR:OH	2.37	0.40
1:B:85:ASN:CG	1:B:86:ASP:N	2.79	0.40
1:B:741:TYR:CE2	1:B:1004:LEU:HD21	2.55	0.40
1:B:1005:GLN:O	1:B:1008:VAL:HG22	2.21	0.40
1:C:317:ASN:OD1	1:C:318:PHE:N	2.53	0.40
1:C:453:PHE:CD1	1:C:455:LEU:HD23	2.56	0.40
1:C:1077:THR:OG1	1:C:1078:ALA:N	2.55	0.40
1:C:1116:THR:HB	1:C:1140:PRO:HG3	2.03	0.40
1:A:192:PHE:HA	1:A:204:TYR:O	2.21	0.40
1:A:337:PRO:HD2	1:A:361:CYS:HB2	2.03	0.40
1:B:382:VAL:HG13	1:C:976:VAL:O	2.22	0.40
1:B:406:GLU:HB3	1:B:409:GLN:CD	2.46	0.40
1:B:458:LYS:NZ	1:B:489:TYR:HB2	2.35	0.40
1:C:100:ILE:O	1:C:242:LEU:HD21	2.22	0.40
1:C:113:LYS:HD2	1:C:113:LYS:HA	1.81	0.40
1:C:196:ASN:C	1:C:197:ILE:HD13	2.46	0.40
1:C:396:TYR:CE1	1:C:512:VAL:HB	2.57	0.40
1:A:62:VAL:HG11	1:A:268:GLY:CA	2.52	0.40
1:A:63:THR:HG22	1:A:64:TRP:N	2.36	0.40
1:A:478:THR:O	1:A:480:CYS:N	2.48	0.40
1:A:796:ASP:OD1	1:A:796:ASP:N	2.41	0.40
1:A:915:VAL:O	1:A:919:ASN:ND2	2.45	0.40
1:B:128:ILE:N	1:B:170:TYR:O	2.52	0.40
1:A:97:LYS:H	1:A:211:ASN:H	1.67	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:LYS:HB2	1:A:133:PHE:CE2	2.57	0.40
1:A:271:GLN:OE1	1:A:293:LEU:HA	2.21	0.40
1:A:330:PRO:CG	1:A:542:ASN:HA	2.51	0.40
1:B:55:PHE:HE2	1:B:85:ASN:OD1	2.04	0.40
1:B:334:ASN:OD1	1:B:360:ASN:HA	2.22	0.40
1:B:385:THR:OG1	1:B:386:LYS:HG3	2.21	0.40
1:B:490:PHE:CD1	1:B:491:PRO:HD2	2.57	0.40
1:C:897:PRO:HB2	1:C:900:MET:HG2	2.03	0.40
1:A:329:PHE:HA	1:A:525:CYS:SG	2.62	0.40
1:B:200:TYR:CD1	1:B:230:PRO:HB3	2.56	0.40
1:B:374:PHE:HB3	1:C:983:ARG:NE	2.36	0.40
1:B:421:TYR:CZ	1:B:462:LYS:HD2	2.56	0.40
1:B:426:PRO:HD3	1:C:196:ASN:O	2.20	0.40
1:B:558:LYS:HE2	1:C:282:ASN:ND2	2.37	0.40
1:C:54:LEU:HB3	1:C:270:LEU:HB3	2.04	0.40
1:C:1127:ASP:OD1	1:C:1127:ASP:N	2.49	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	985/1269 (78%)	836 (85%)	140 (14%)	9 (1%)	14	46
1	B	985/1269 (78%)	883 (90%)	101 (10%)	1 (0%)	48	79
1	C	985/1269 (78%)	882 (90%)	103 (10%)	0	100	100
All	All	2955/3807 (78%)	2601 (88%)	344 (12%)	10 (0%)	37	65

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	LEU

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Mol	Chain	Res	Type
1	A	85	ASN
1	A	520	ALA
1	A	535	LYS
1	B	282	ASN
1	A	57	PRO
1	A	223	LEU
1	A	91	TYR
1	A	325	SER
1	A	545	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	877/1101 (80%)	872 (99%)	5 (1%)	78	79
1	B	877/1101 (80%)	876 (100%)	1 (0%)	88	87
1	C	877/1101 (80%)	873 (100%)	4 (0%)	81	80
All	All	2631/3303 (80%)	2621 (100%)	10 (0%)	81	81

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	PHE
1	A	56	LEU
1	A	207	HIS
1	A	740	MET
1	A	966	LEU
1	B	49	HIS
1	C	189	LEU
1	C	312	ILE
1	C	429	PHE
1	C	973	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	99	ASN
1	A	334	ASN
1	A	360	ASN
1	A	519	HIS
1	A	542	ASN
1	A	544	ASN
1	A	603	ASN
1	A	607	GLN
1	A	913	GLN
1	A	954	GLN
1	A	1002	GLN
1	A	1113	GLN
1	B	165	ASN
1	B	343	ASN
1	B	440	ASN
1	B	450	ASN
1	B	542	ASN
1	B	675	GLN
1	B	762	GLN
1	B	774	GLN
1	B	804	GLN
1	B	954	GLN
1	B	1010	GLN
1	B	1048	HIS
1	C	115	GLN
1	C	165	ASN
1	C	234	ASN
1	C	343	ASN
1	C	498	GLN
1	C	501	ASN
1	C	606	ASN
1	C	872	GLN
1	C	913	GLN
1	C	969	ASN
1	C	1011	GLN
1	C	1058	HIS
1	C	1071	GLN
1	C	1074	ASN
1	C	1108	ASN
1	C	1142	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

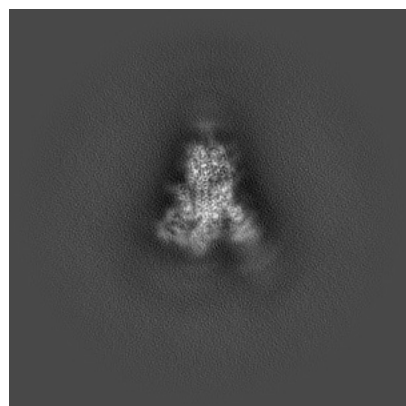
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40977. These allow visual inspection of the internal detail of the map and identification of artifacts.

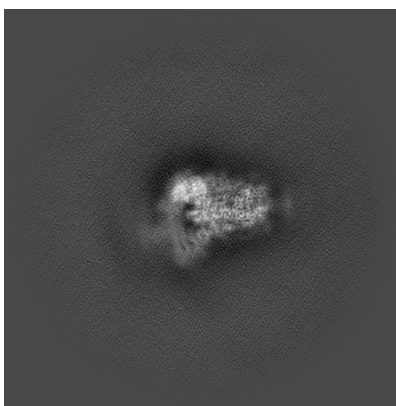
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

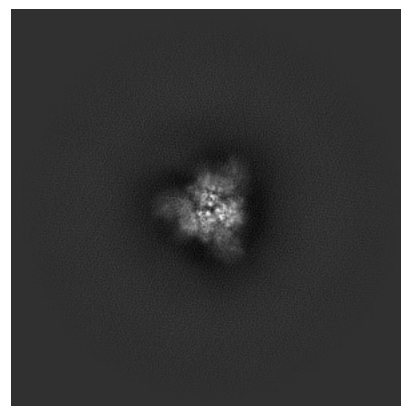
6.1.1 Primary map



X

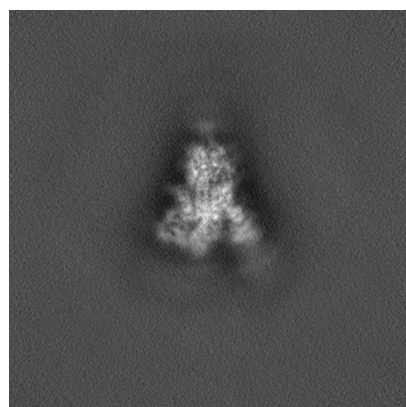


Y

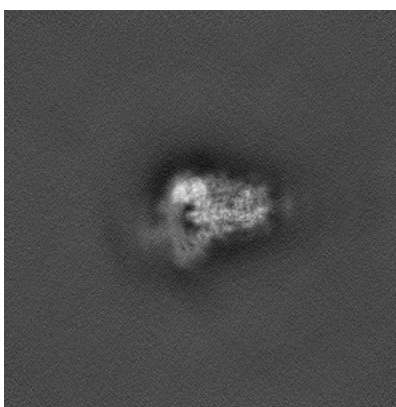


Z

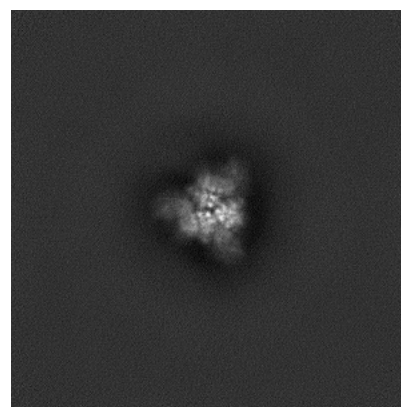
6.1.2 Raw map



X



Y



Z

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

6.2 Central slices [i](#)

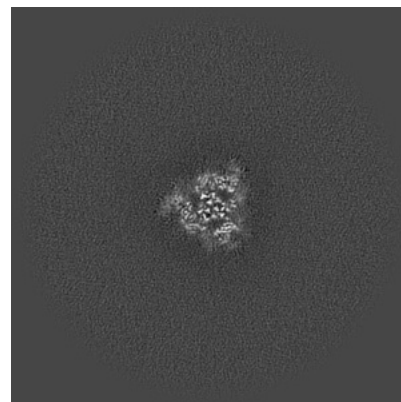
6.2.1 Primary map



X Index: 256



Y Index: 256



Z Index: 256

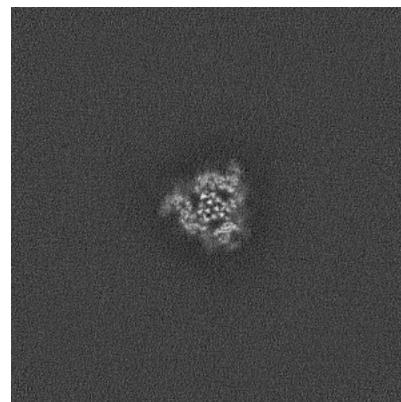
6.2.2 Raw map



X Index: 256



Y Index: 256

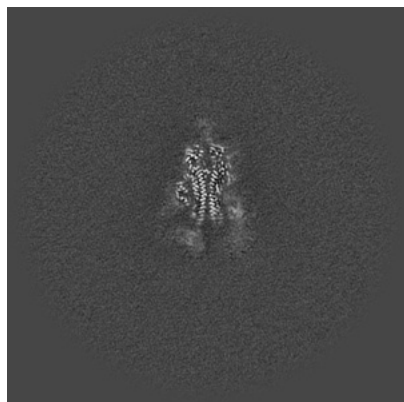


Z Index: 256

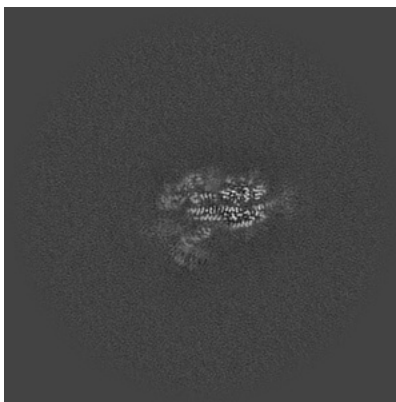
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

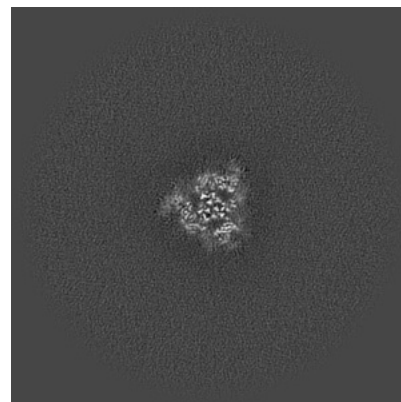
6.3.1 Primary map



X Index: 253



Y Index: 248



Z Index: 256

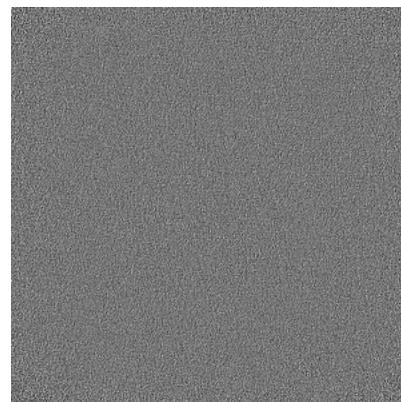
6.3.2 Raw map



X Index: 253



Y Index: 248

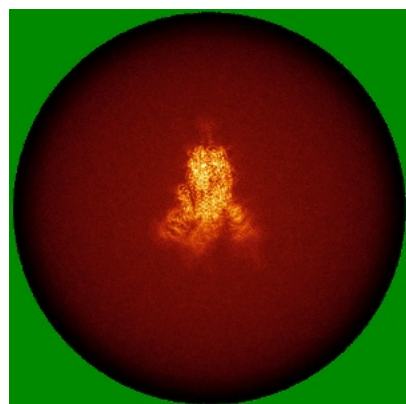


Z Index: 0

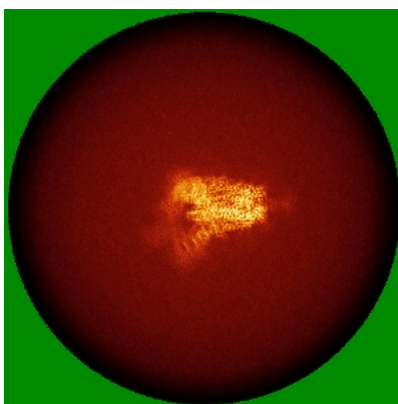
The images above show the largest variance slices of the map in three orthogonal directions.

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

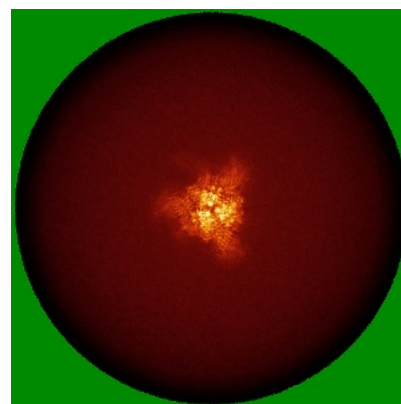
6.4.1 Primary map



X

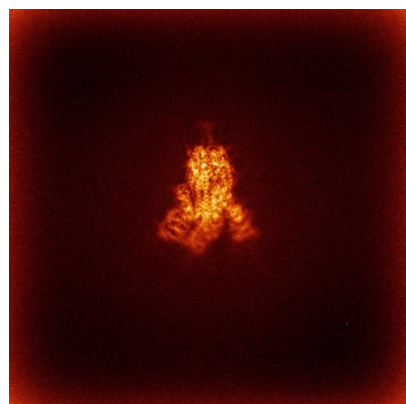


Y

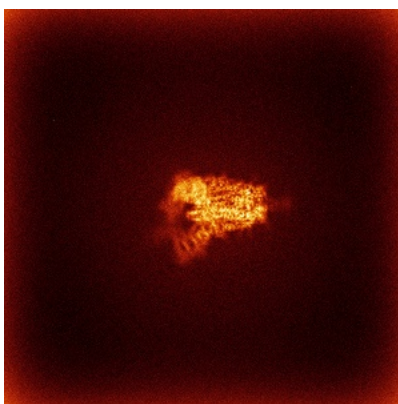


Z

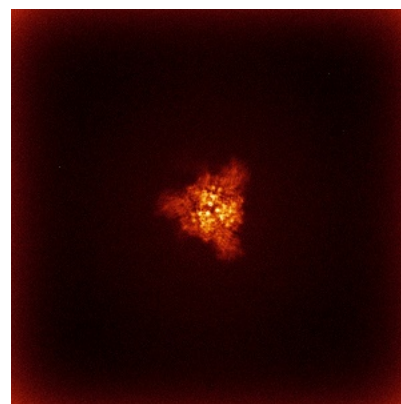
6.4.2 Raw map



X



Y

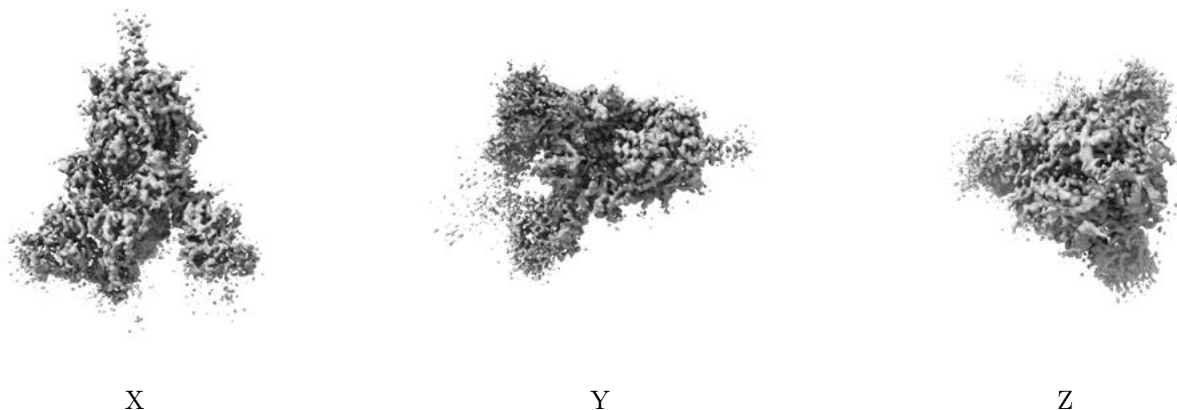


Z

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

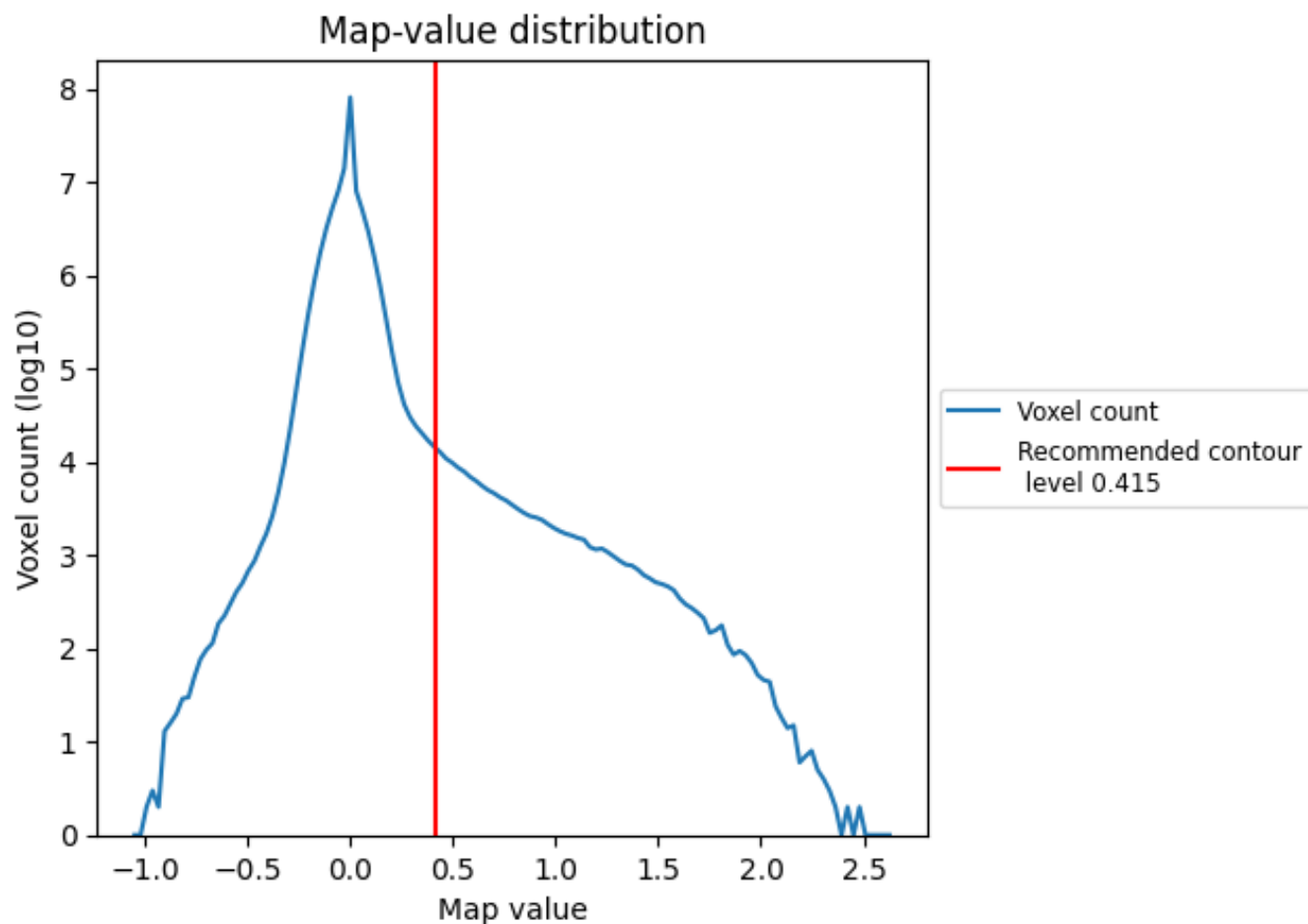
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

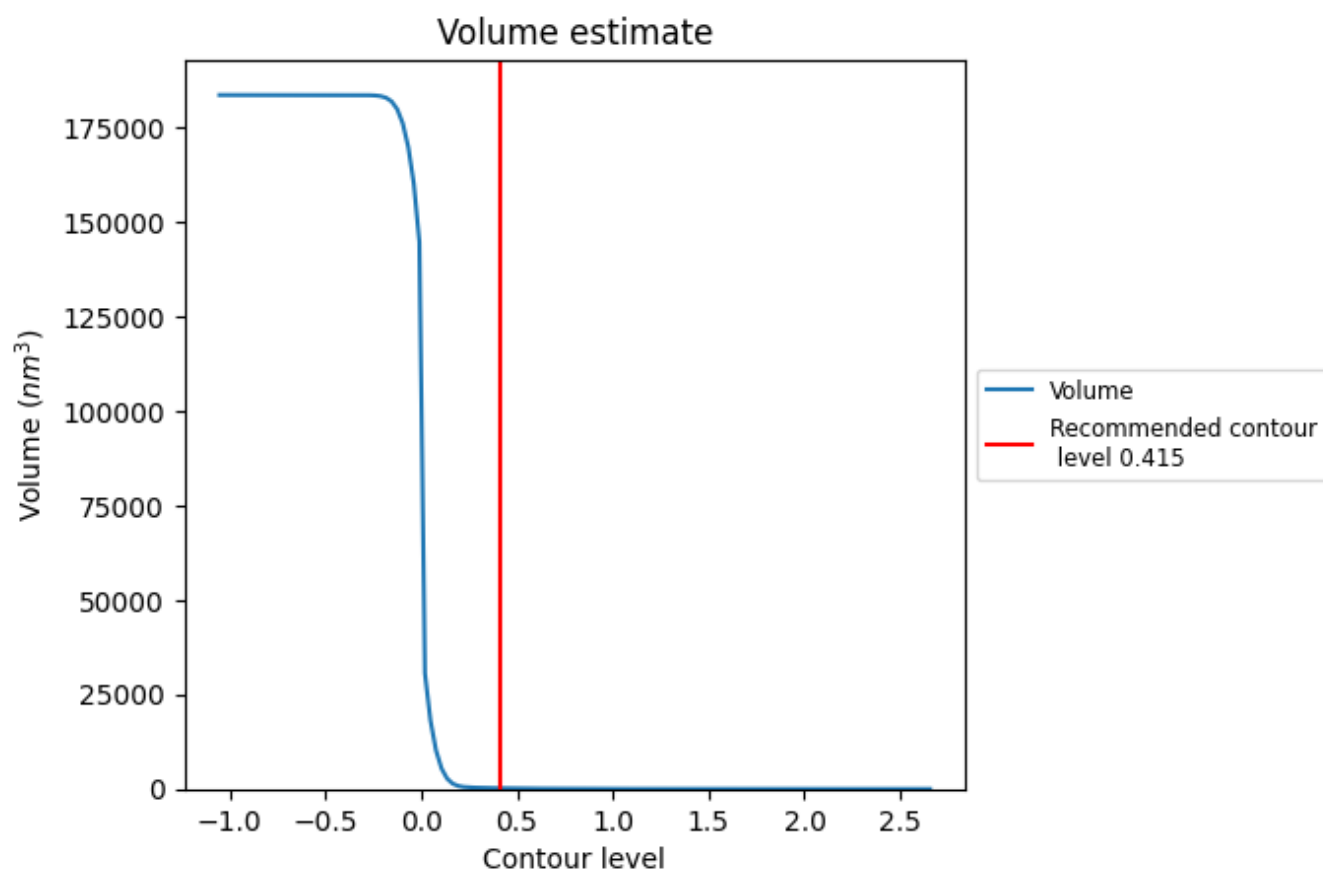
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

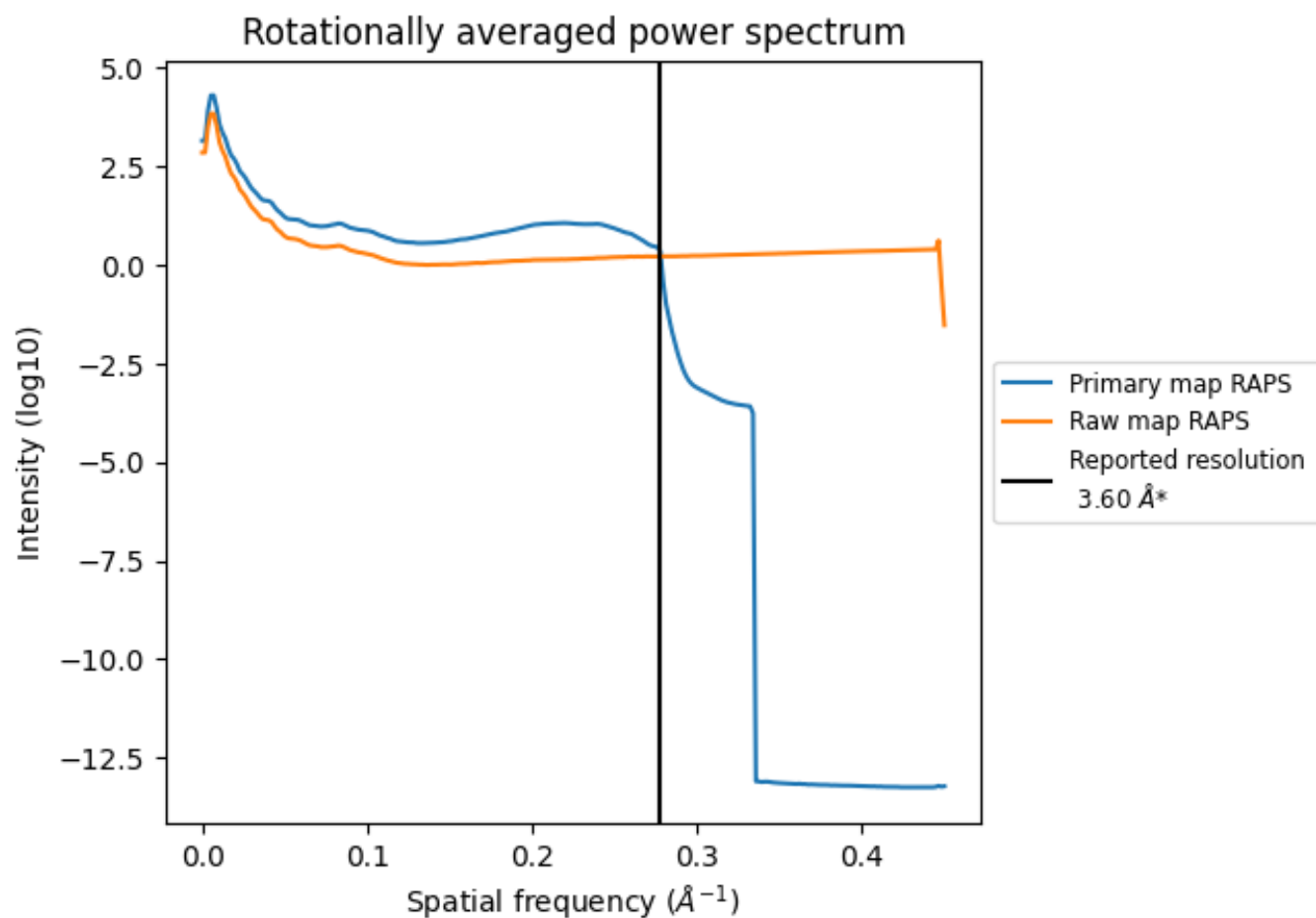
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 193 nm^3 ; this corresponds to an approximate mass of 174 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 ⓘ

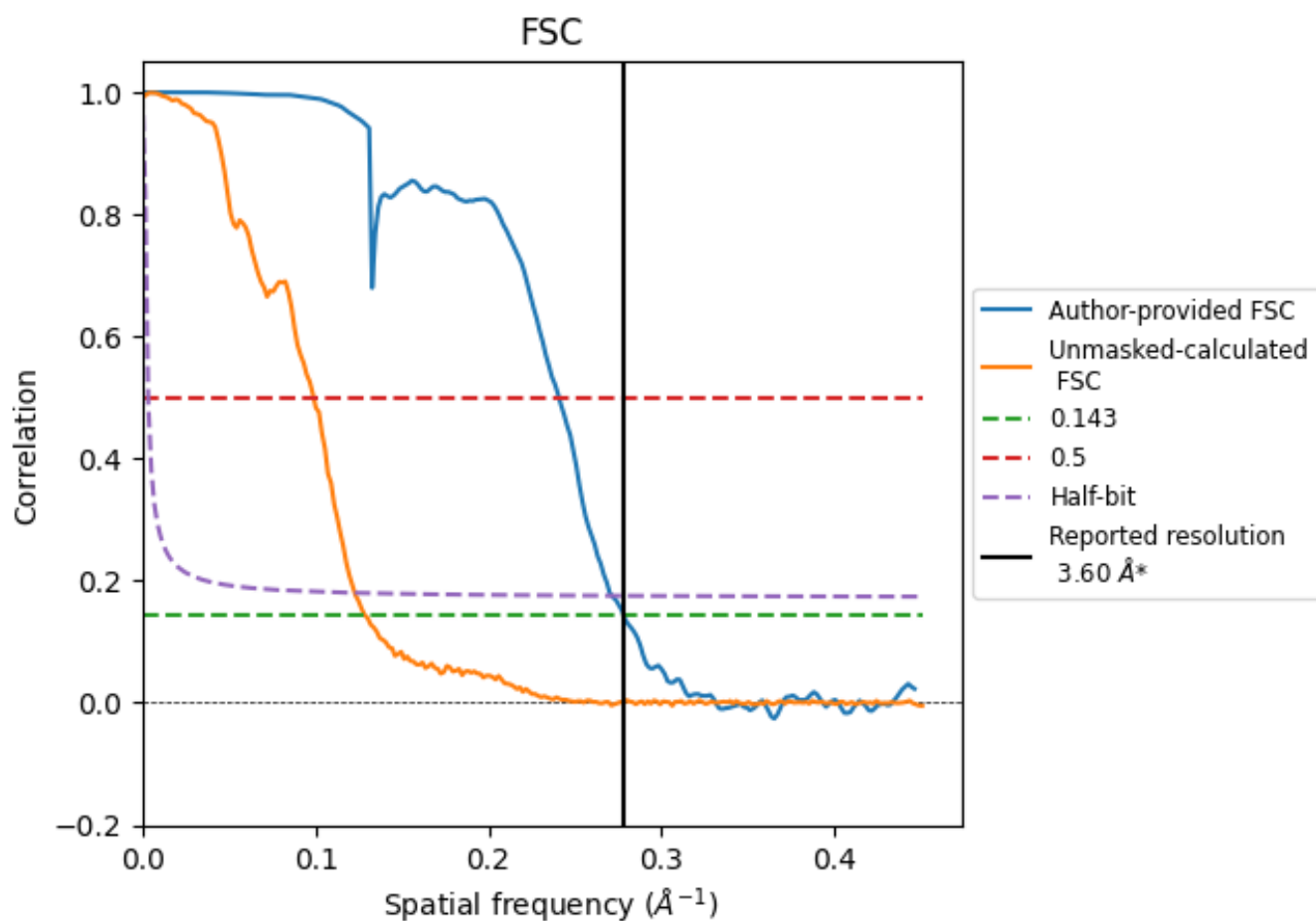


*Reported resolution corresponds to spatial frequency of 0.278 Å⁻¹

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 0.278 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

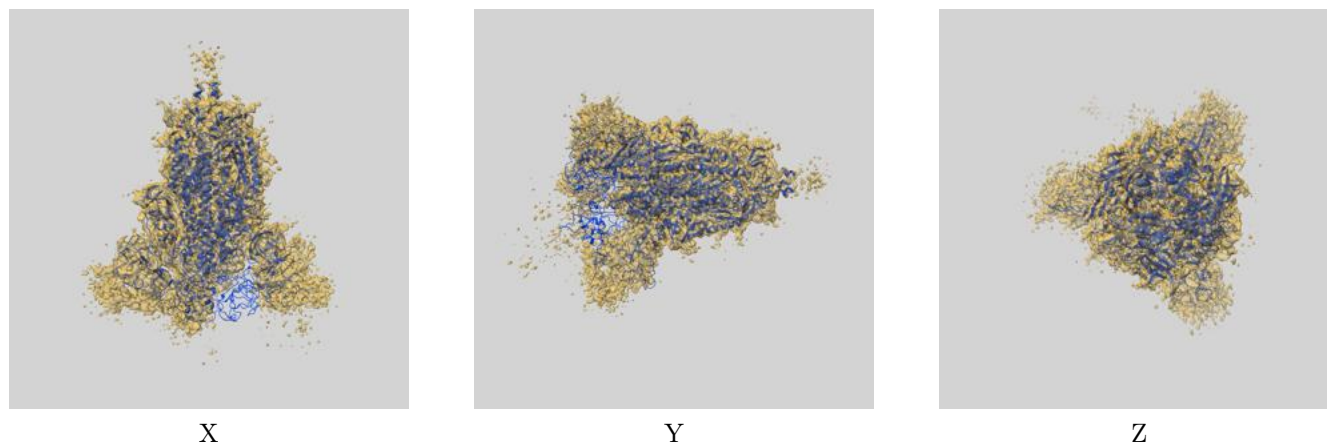
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.60	-	-
Author-provided FSC curve	3.60	4.15	3.69
Unmasked-calculated*	7.76	10.10	8.14

*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 7.76 differs from the reported value 3.6 by more than 10 %

9 Map-model fit [i](#)

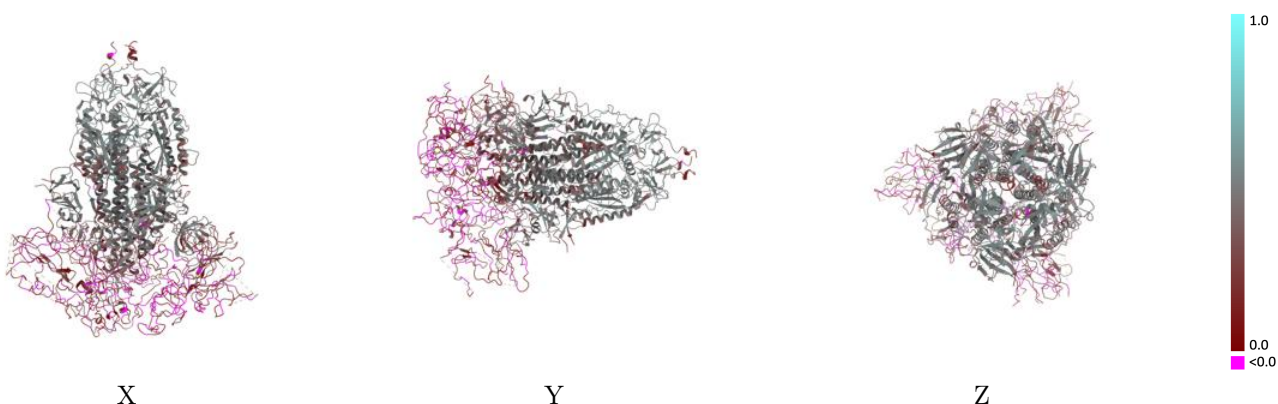
This section contains information regarding the fit between EMDB map EMD-40977 and PDB model 8T21. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



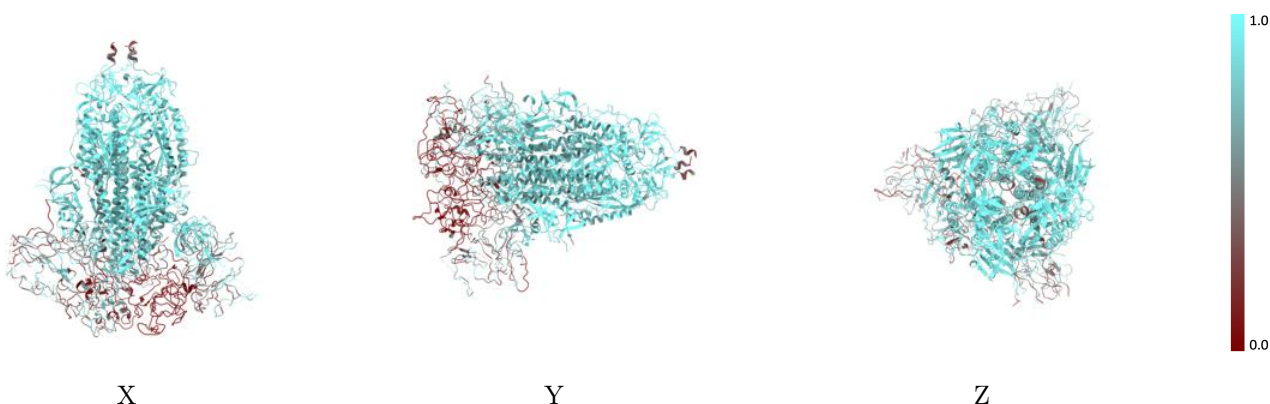
The images above show the 3D surface view of the map at the recommended contour level 0.415 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



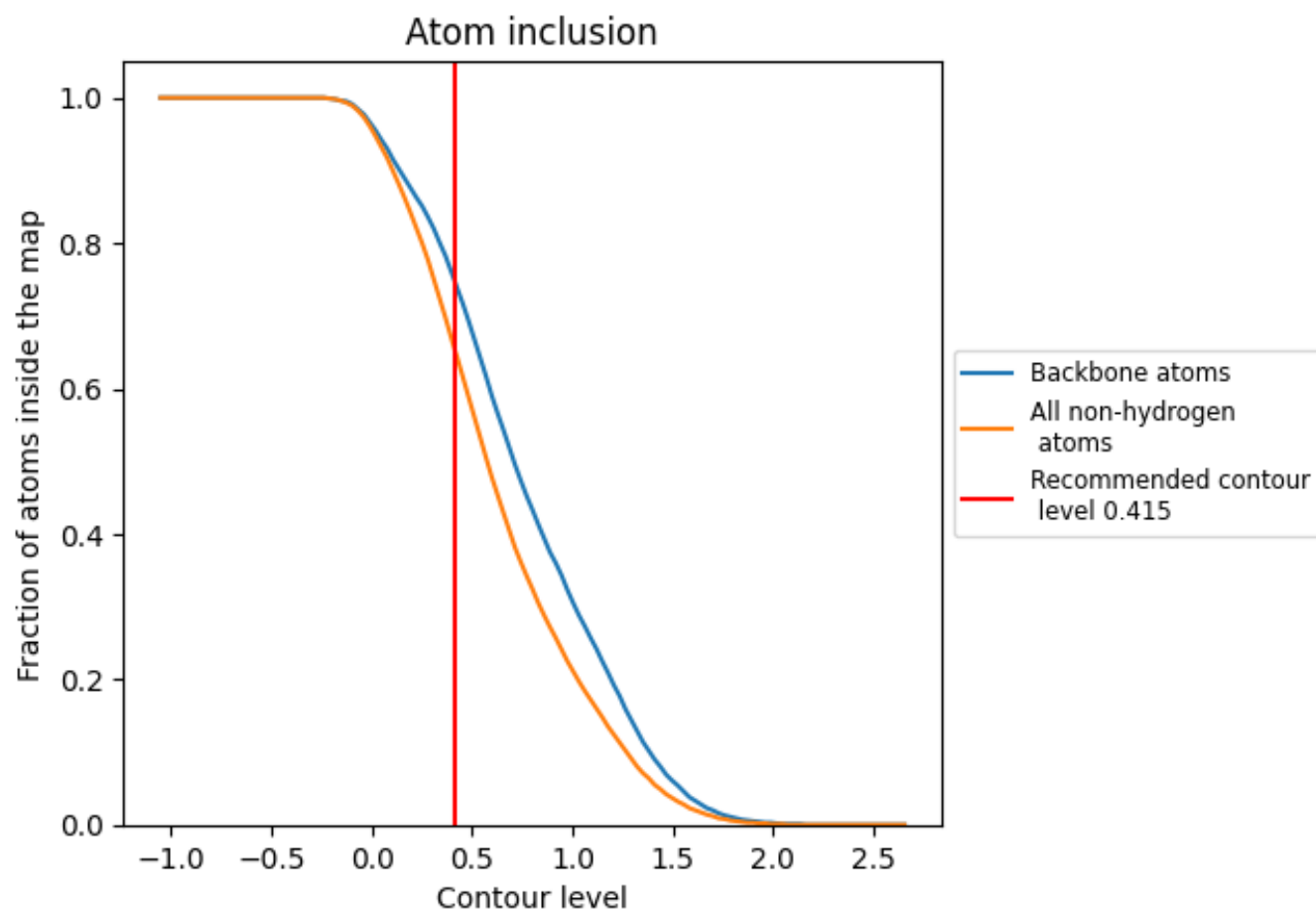
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.415).

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 65% 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 (0.415) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6520	0.3070
A	0.6170	0.2910
B	0.6120	0.2910
C	0.7280	0.3400

