



# Full wwPDB EM Validation Report ⓘ

May 23, 2023 – 08:23 PM JST

PDB ID : 8H08  
EMDB ID : EMD-34411  
Title : SARS-CoV-2 BA.1 variants S ectodomain trimer in complex with neutralizing antibody 10-5B and 6-2C  
Authors : Wang, X.; Wang, Z.  
Deposited on : 2022-09-28  
Resolution : 3.20 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev50  
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 : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.33

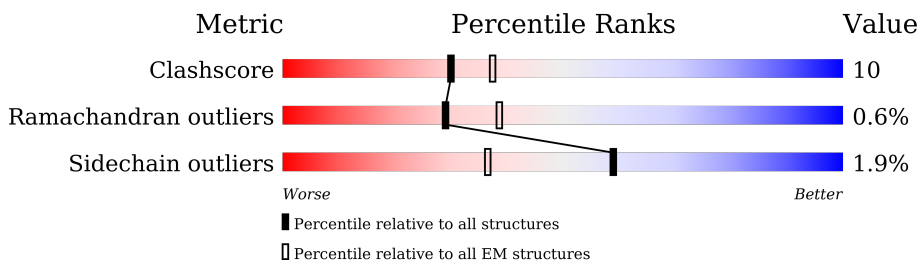
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.20 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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1300	
1	B	1300	
1	C	1300	
2	D	117	
2	F	117	
2	L	117	
3	H	107	
3	J	107	

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Mol	Chain	Length	Quality of chain
3	P	107	
4	I	122	
4	N	122	
4	Q	122	
5	K	107	
5	O	107	
5	R	107	

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 31277 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	A	948	7138	4564	1203	1344	27	0	0
1	B	933	7024	4476	1194	1329	25	0	0
1	C	948	7125	4544	1200	1355	26	0	0

There are 141 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	VAL	ALA	variant	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	95	ILE	THR	variant	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	145	ASP	TYR	variant	UNP P0DTC2
A	212	ILE	LEU	variant	UNP P0DTC2
A	214A	GLU	-	insertion	UNP P0DTC2
A	214B	PRO	-	insertion	UNP P0DTC2
A	214C	GLU	-	insertion	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	LEU	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2
A	493	ARG	GLN	variant	UNP P0DTC2
A	496	SER	GLY	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	547	LYS	THR	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	HIS	PRO	variant	UNP P0DTC2
A	682	GLY	ARG	engineered mutation	UNP P0DTC2
A	683	SER	ARG	engineered mutation	UNP P0DTC2
A	685	SER	ARG	engineered mutation	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	conflict	UNP P0DTC2
A	856	LYS	ASN	variant	UNP P0DTC2
A	892	PRO	ALA	conflict	UNP P0DTC2
A	899	PRO	ALA	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	981	PHE	LEU	variant	UNP P0DTC2
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
B	69	VAL	ALA	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	95	ILE	THR	variant	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	145	ASP	TYR	variant	UNP P0DTC2
B	212	ILE	LEU	variant	UNP P0DTC2
B	214A	GLU	-	insertion	UNP P0DTC2
B	214B	PRO	-	insertion	UNP P0DTC2
B	214C	GLU	-	insertion	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	LEU	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	446	SER	GLY	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	484	ALA	GLU	variant	UNP P0DTC2
B	493	ARG	GLN	variant	UNP P0DTC2
B	496	SER	GLY	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	547	LYS	THR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	HIS	PRO	variant	UNP P0DTC2
B	682	GLY	ARG	engineered mutation	UNP P0DTC2
B	683	SER	ARG	engineered mutation	UNP P0DTC2
B	685	SER	ARG	engineered mutation	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	conflict	UNP P0DTC2
B	856	LYS	ASN	variant	UNP P0DTC2
B	892	PRO	ALA	conflict	UNP P0DTC2
B	899	PRO	ALA	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	981	PHE	LEU	variant	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	69	VAL	ALA	variant	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	95	ILE	THR	variant	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	145	ASP	TYR	variant	UNP P0DTC2
C	212	ILE	LEU	variant	UNP P0DTC2
C	214A	GLU	-	insertion	UNP P0DTC2
C	214B	PRO	-	insertion	UNP P0DTC2
C	214C	GLU	-	insertion	UNP P0DTC2
C	339	ASP	GLY	variant	UNP P0DTC2
C	371	LEU	SER	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	446	SER	GLY	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2
C	484	ALA	GLU	variant	UNP P0DTC2
C	493	ARG	GLN	variant	UNP P0DTC2
C	496	SER	GLY	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	547	LYS	THR	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	681	HIS	PRO	variant	UNP P0DTC2
C	682	GLY	ARG	engineered mutation	UNP P0DTC2
C	683	SER	ARG	engineered mutation	UNP P0DTC2
C	685	SER	ARG	engineered mutation	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	817	PRO	PHE	conflict	UNP P0DTC2
C	856	LYS	ASN	variant	UNP P0DTC2
C	892	PRO	ALA	conflict	UNP P0DTC2
C	899	PRO	ALA	conflict	UNP P0DTC2
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C	981	PHE	LEU	variant	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2

- Molecule 2 is a protein called 10-5B H chain.

Mol	Chain	Residues	Atoms				AltConf	Trace	
2	D	116	Total	C	N	O	S	0	0
			839	519	150	166	4		
2	F	116	Total	C	N	O	S	0	0
			839	519	150	166	4		
2	L	116	Total	C	N	O	S	0	0
			839	519	150	166	4		

- Molecule 3 is a protein called 10-5B L chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	104	Total	C	N	O	S	0	0
			738	453	130	152	3		
3	H	104	Total	C	N	O	S	0	0
			738	453	130	152	3		
3	P	104	Total	C	N	O	S	0	0
			738	453	130	152	3		

- Molecule 4 is a protein called 6-2C H chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	N	122	Total	C	N	O	S	0	0
			945	597	165	181	2		
4	I	122	Total	C	N	O	S	0	0
			945	597	165	181	2		
4	Q	122	Total	C	N	O	S	0	0
			945	597	165	181	2		

- Molecule 5 is a protein called 6-2C L chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	O	106	Total	C	N	O	S	0	0
			794	506	133	153	2		
5	K	106	Total	C	N	O	S	0	0
			794	506	133	153	2		
5	R	106	Total	C	N	O	S	0	0
			794	506	133	153	2		

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>) (labeled as "Ligand of Interest" by depositor).



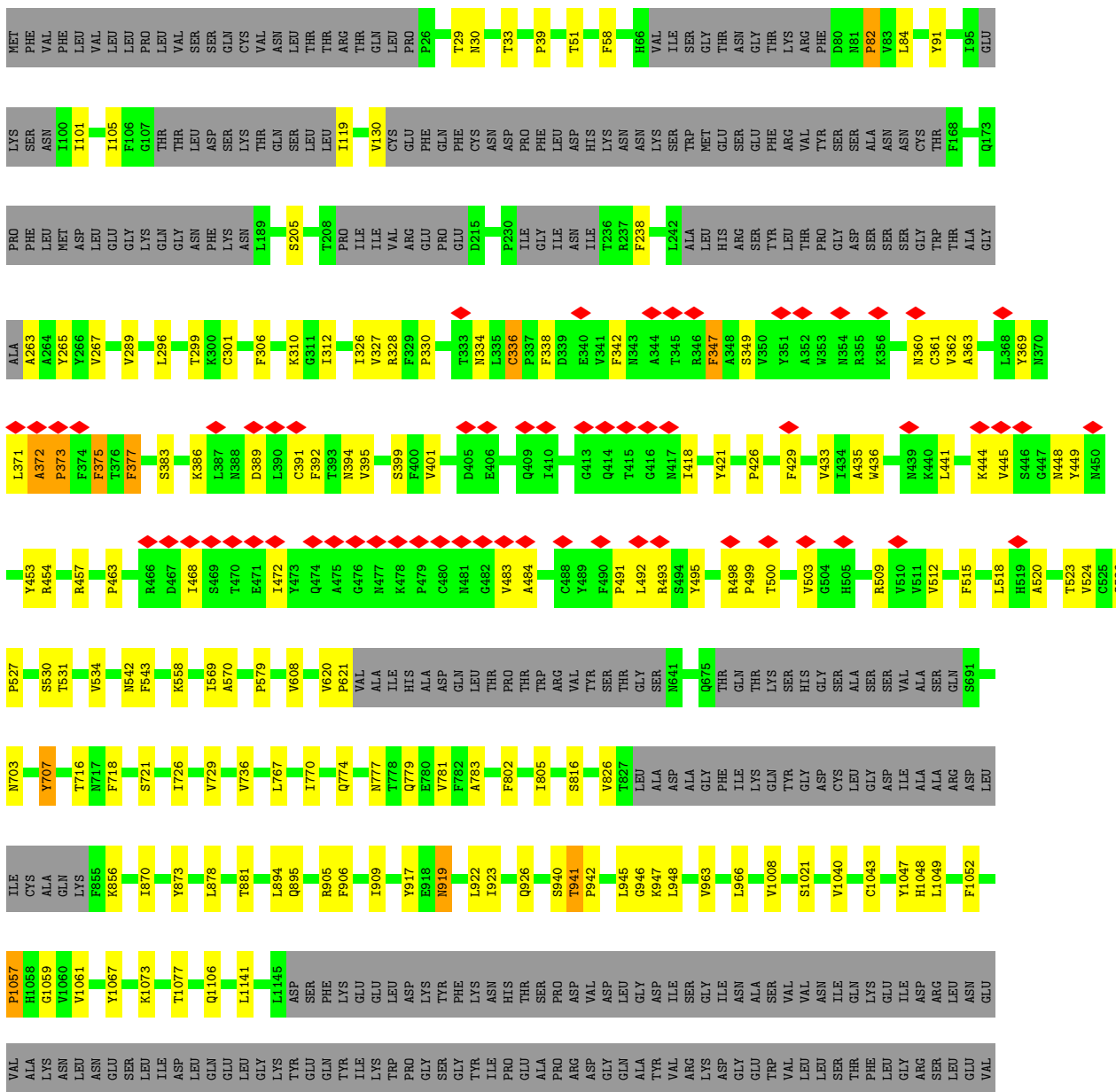


Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	

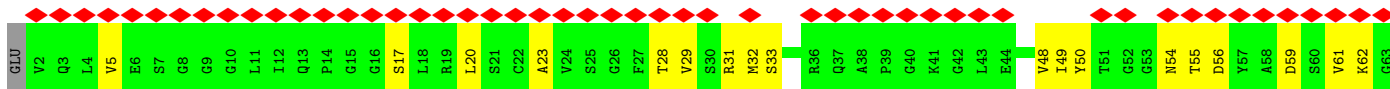
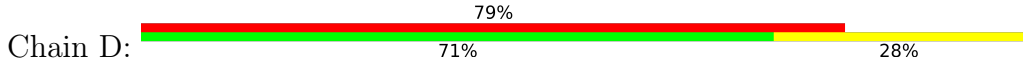


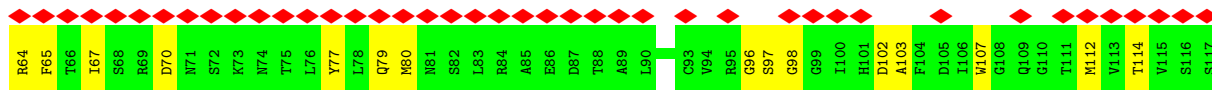


• Molecule 1: Spike glycoprotein

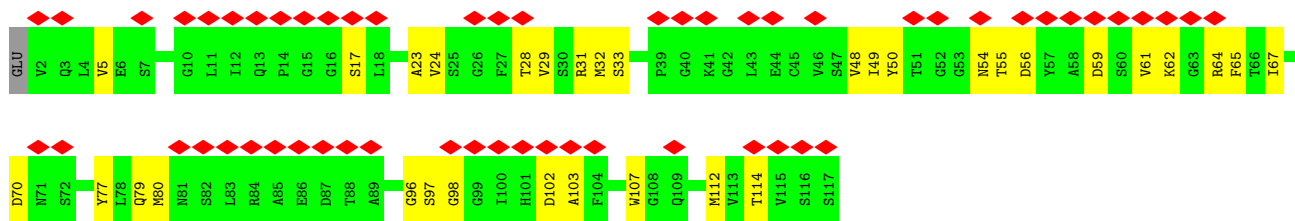


• Molecule 2: 10-5B H chain

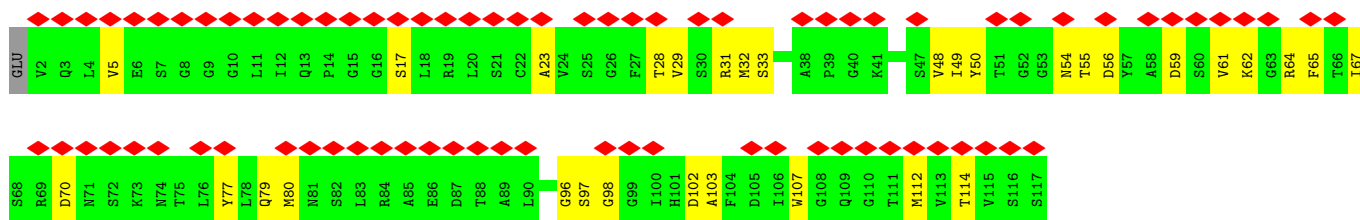




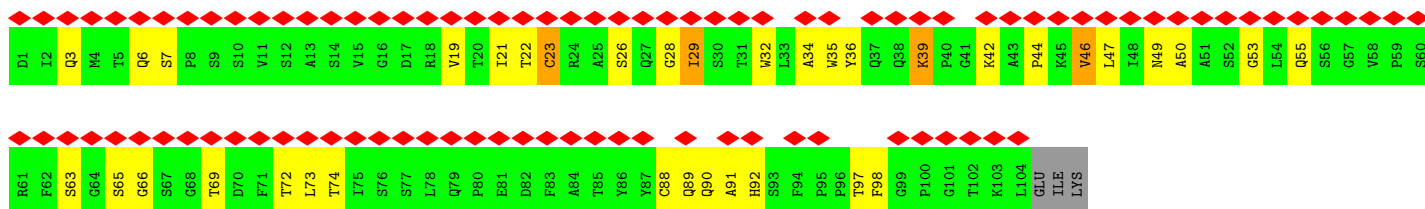
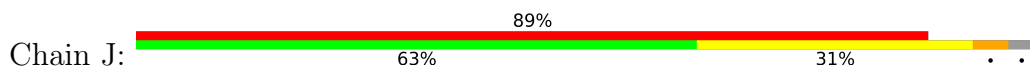
• Molecule 2: 10-5B H chain



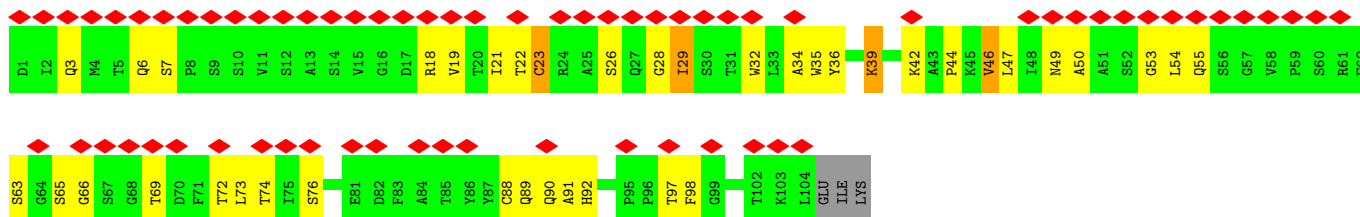
• Molecule 2: 10-5B H chain



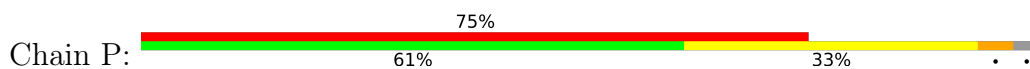
• Molecule 3: 10-5B L chain

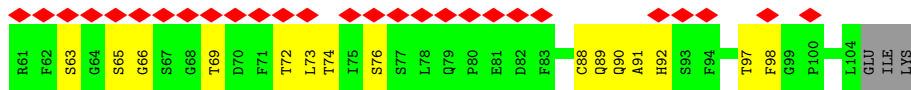
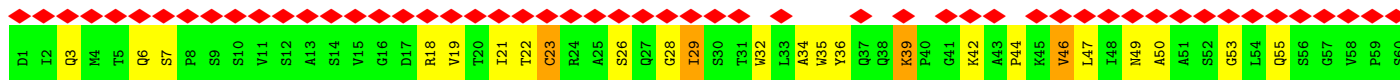


• Molecule 3: 10-5B L chain

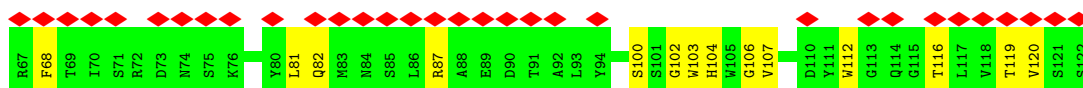
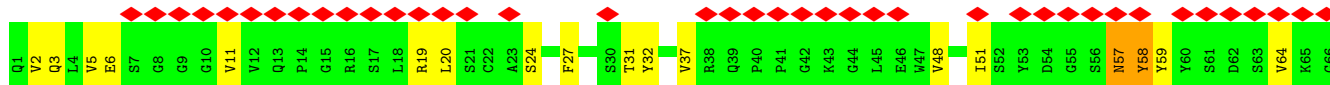
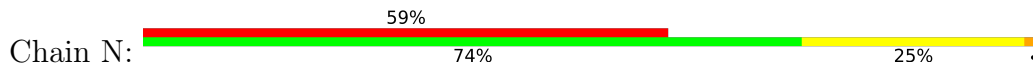


• Molecule 3: 10-5B L chain

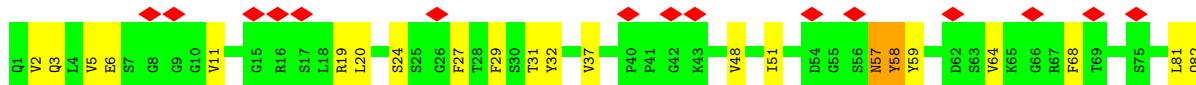




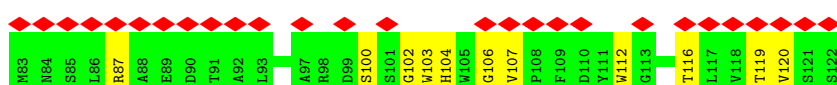
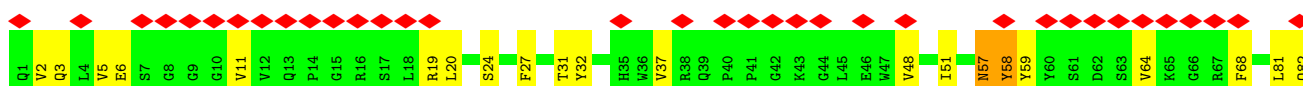
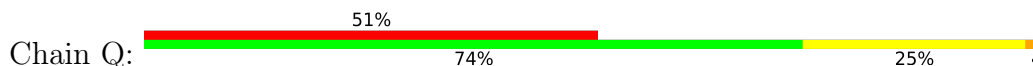
• Molecule 4: 6-2C H chain



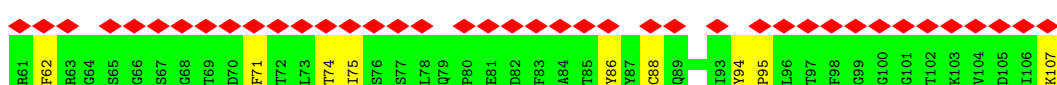
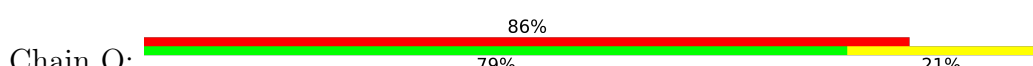
• Molecule 4: 6-2C H chain



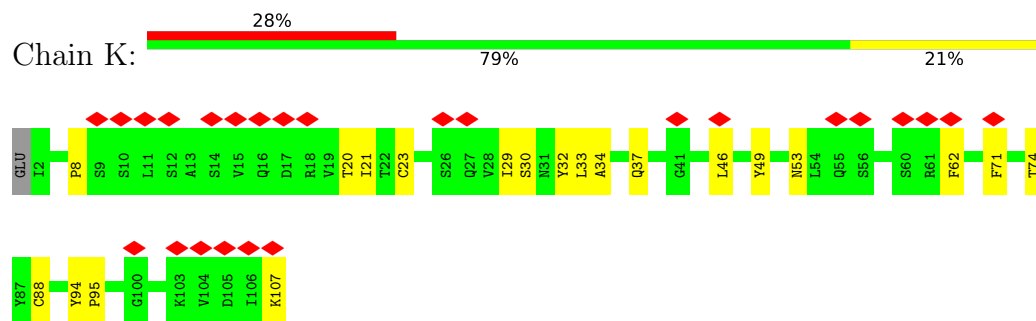
• Molecule 4: 6-2C H chain



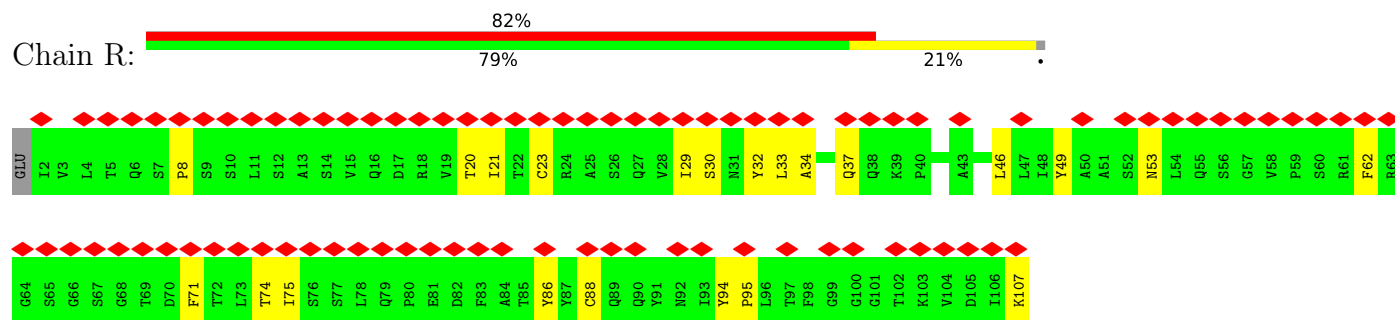
• Molecule 5: 6-2C L chain



• Molecule 5: 6-2C L chain



• Molecule 5: 6-2C L chain



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	372377	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.588	Depositor
Minimum map value	-1.650	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.054	Depositor
Recommended contour level	0.15	Depositor
Map size ( $\text{\AA}$ )	329.80002, 329.80002, 329.80002	wwPDB
Map dimensions	340, 340, 340	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.97, 0.97, 0.97	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: 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	0.66	1/7307 (0.0%)	0.90	5/9967 (0.1%)
1	B	0.66	2/7181 (0.0%)	0.90	10/9777 (0.1%)
1	C	0.67	4/7293 (0.1%)	0.91	7/9944 (0.1%)
2	D	0.61	0/853	0.72	0/1152
2	F	0.61	0/853	0.73	0/1152
2	L	0.61	0/853	0.73	0/1152
3	H	0.67	0/756	0.77	0/1032
3	J	0.67	0/756	0.77	0/1032
3	P	0.67	0/756	0.77	0/1032
4	I	0.68	2/975 (0.2%)	0.82	1/1326 (0.1%)
4	N	0.68	2/975 (0.2%)	0.82	1/1326 (0.1%)
4	Q	0.68	2/975 (0.2%)	0.82	1/1326 (0.1%)
5	K	0.63	0/812	0.89	1/1104 (0.1%)
5	O	0.63	0/812	0.89	1/1104 (0.1%)
5	R	0.63	0/812	0.89	1/1104 (0.1%)
All	All	0.66	13/31969 (0.0%)	0.87	28/43530 (0.1%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	57	ASN	C-N	6.50	1.49	1.34
4	N	57	ASN	C-N	6.49	1.49	1.34
4	Q	57	ASN	C-N	6.45	1.48	1.34
4	N	58	TYR	C-N	6.38	1.48	1.34
4	Q	58	TYR	C-N	6.38	1.48	1.34
4	I	58	TYR	C-N	6.34	1.48	1.34
1	A	875	SER	CA-CB	-5.46	1.44	1.52
1	B	875	SER	CA-CB	-5.46	1.44	1.52
1	B	1021	SER	CA-CB	-5.25	1.45	1.52
1	C	816	SER	CA-CB	-5.20	1.45	1.52
1	C	721	SER	CA-CB	-5.19	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1021	SER	CA-CB	-5.13	1.45	1.52
1	C	205	SER	CA-CB	-5.01	1.45	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	336	CYS	CB-CA-C	6.72	123.85	110.40
1	B	913	GLN	CB-CA-C	-6.66	97.08	110.40
1	B	812	PRO	N-CA-C	-6.57	95.01	112.10
1	C	336	CYS	CA-CB-SG	-6.41	102.47	114.00
1	B	336	CYS	CB-CA-C	-5.93	98.54	110.40
4	N	57	ASN	O-C-N	5.62	131.69	122.70
1	A	718	PHE	CB-CA-C	-5.61	99.19	110.40
4	Q	57	ASN	O-C-N	5.60	131.67	122.70
4	I	57	ASN	O-C-N	5.60	131.66	122.70
1	C	375	PHE	CB-CA-C	-5.54	99.32	110.40
1	B	375	PHE	CB-CA-C	-5.54	99.33	110.40
1	A	375	PHE	CB-CA-C	-5.53	99.34	110.40
1	B	1067	TYR	CB-CA-C	-5.53	99.34	110.40
1	A	906	PHE	CB-CA-C	-5.53	99.34	110.40
1	C	310	LYS	CB-CA-C	-5.53	99.35	110.40
1	B	814	LYS	CB-CA-C	5.42	121.25	110.40
1	C	1077	THR	CB-CA-C	-5.34	97.17	111.60
1	A	1067	TYR	CB-CA-C	-5.32	99.77	110.40
1	B	543	PHE	CB-CA-C	5.25	120.89	110.40
1	C	707	TYR	N-CA-CB	5.14	119.85	110.60
5	K	95	PRO	N-CA-C	-5.11	98.82	112.10
5	R	95	PRO	N-CA-C	-5.11	98.82	112.10
5	O	95	PRO	N-CA-C	-5.09	98.85	112.10
1	C	373	PRO	CB-CA-C	-5.08	99.30	112.00
1	B	373	PRO	CB-CA-C	-5.08	99.31	112.00
1	A	373	PRO	CB-CA-C	-5.07	99.34	112.00
1	B	901	GLN	CB-CA-C	-5.06	100.28	110.40
1	B	392	PHE	CB-CA-C	5.01	120.42	110.40

There are no chirality outliers.

There are no planarity outliers.

## 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	7138	0	6670	149	0
1	B	7024	0	6549	116	0
1	C	7125	0	6607	111	0
2	D	839	0	779	25	0
2	F	839	0	779	25	0
2	L	839	0	779	23	0
3	H	738	0	671	33	0
3	J	738	0	671	32	0
3	P	738	0	671	32	0
4	I	945	0	834	24	0
4	N	945	0	834	21	0
4	Q	945	0	834	23	0
5	K	794	0	768	15	0
5	O	794	0	768	15	0
5	R	794	0	768	15	0
6	A	14	0	13	0	0
6	B	14	0	13	0	0
6	C	14	0	13	0	0
All	All	31277	0	29021	602	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 10.

All (602) 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:722:VAL:CG2	1:A:934:ILE:HG13	1.67	1.24
1:A:722:VAL:HG21	1:A:934:ILE:HG13	1.37	1.05
1:A:722:VAL:CG2	1:A:934:ILE:CG1	2.38	1.01
1:C:483:VAL:CG1	2:D:54:ASN:OD1	2.11	0.99
1:B:483:VAL:CG1	2:F:54:ASN:OD1	2.11	0.99
1:A:483:VAL:CG1	2:L:54:ASN:OD1	2.11	0.98
1:C:483:VAL:HG11	2:D:54:ASN:OD1	1.77	0.85
1:A:722:VAL:HG23	1:A:934:ILE:HG13	1.58	0.85
1:B:822:LEU:HD22	1:B:945:LEU:HD21	1.60	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:VAL:HG11	2:L:54:ASN:OD1	1.77	0.82
1:B:483:VAL:HG11	2:F:54:ASN:OD1	1.77	0.82
1:C:330:PRO:HA	1:C:579:PRO:HB2	1.64	0.79
2:D:98:GLY:H	2:D:103:ALA:H	1.31	0.79
1:A:393:THR:HA	1:A:522:ALA:HA	1.64	0.77
3:J:29:ILE:HG23	3:J:92:HIS:HB2	1.68	0.76
3:P:29:ILE:HG23	3:P:92:HIS:HB2	1.68	0.76
2:F:98:GLY:H	2:F:103:ALA:H	1.31	0.76
2:L:98:GLY:H	2:L:103:ALA:H	1.31	0.76
2:F:112:MET:SD	2:F:114:THR:OG1	2.45	0.75
2:L:112:MET:SD	2:L:114:THR:OG1	2.45	0.74
2:D:61:VAL:HA	2:D:64:ARG:HE	1.52	0.74
3:H:29:ILE:HG23	3:H:92:HIS:HB2	1.67	0.74
4:N:100:SER:OG	4:N:106:GLY:HA2	1.88	0.74
2:F:61:VAL:HA	2:F:64:ARG:HE	1.52	0.74
4:I:100:SER:OG	4:I:106:GLY:HA2	1.88	0.74
2:L:61:VAL:HA	2:L:64:ARG:HE	1.52	0.73
4:Q:100:SER:OG	4:Q:106:GLY:HA2	1.88	0.73
2:D:112:MET:SD	2:D:114:THR:OG1	2.45	0.73
1:A:360:ASN:H	1:A:523:THR:HB	1.55	0.71
1:A:1030:SER:O	1:C:1040:VAL:HB	1.91	0.71
4:Q:37:VAL:HG11	4:Q:112:TRP:HZ3	1.57	0.70
4:I:37:VAL:HG11	4:I:112:TRP:HZ3	1.57	0.70
4:I:37:VAL:HG11	4:I:112:TRP:CZ3	2.28	0.69
1:A:499:PRO:HB2	5:R:94:TYR:HD2	1.58	0.68
2:F:31:ARG:HD3	2:F:97:SER:HB2	1.75	0.68
4:N:37:VAL:HG11	4:N:112:TRP:CZ3	2.28	0.68
3:P:63:SER:HB2	3:P:74:THR:H	1.59	0.68
3:H:3:GLN:HB2	3:H:26:SER:HB3	1.76	0.68
4:Q:37:VAL:HG11	4:Q:112:TRP:CZ3	2.28	0.68
3:H:63:SER:HB2	3:H:74:THR:H	1.59	0.68
2:L:31:ARG:HD3	2:L:97:SER:HB2	1.75	0.68
1:C:499:PRO:HB2	5:O:94:TYR:HD2	1.58	0.68
2:D:31:ARG:HD3	2:D:97:SER:HB2	1.75	0.68
3:J:35:TRP:CD1	3:J:88:CYS:SG	2.87	0.68
3:P:3:GLN:HB2	3:P:26:SER:HB3	1.76	0.68
3:J:3:GLN:HB2	3:J:26:SER:HB3	1.76	0.68
4:N:37:VAL:HG11	4:N:112:TRP:HZ3	1.57	0.67
1:B:499:PRO:HB2	5:K:94:TYR:HD2	1.58	0.67
3:H:35:TRP:CD1	3:H:88:CYS:SG	2.87	0.67
1:A:722:VAL:HG23	1:A:934:ILE:CD1	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:35:TRP:CD1	3:P:88:CYS:SG	2.87	0.67
3:J:63:SER:HB2	3:J:74:THR:H	1.59	0.67
2:F:49:ILE:HG13	2:F:55:THR:HG22	1.77	0.67
5:O:62:PHE:CE1	5:O:75:ILE:HD11	2.30	0.67
3:H:49:ASN:HB2	3:H:55:GLN:HA	1.77	0.66
3:P:49:ASN:HB2	3:P:55:GLN:HA	1.77	0.66
5:R:62:PHE:CE1	5:R:75:ILE:HD11	2.30	0.66
1:B:483:VAL:HG12	2:F:54:ASN:OD1	1.94	0.66
2:L:49:ILE:HG13	2:L:55:THR:HG22	1.77	0.66
5:K:62:PHE:CE1	5:K:75:ILE:HD11	2.30	0.66
3:J:49:ASN:HB2	3:J:55:GLN:HA	1.77	0.65
3:H:90:GLN:HB3	3:H:97:THR:H	1.61	0.65
1:B:371:LEU:C	1:B:373:PRO:HD2	2.17	0.65
3:J:90:GLN:HB3	3:J:97:THR:H	1.61	0.65
3:P:90:GLN:HB3	3:P:97:THR:H	1.61	0.65
1:C:483:VAL:HG12	2:D:54:ASN:OD1	1.94	0.65
1:A:722:VAL:CG2	1:A:934:ILE:CD1	2.75	0.64
1:A:483:VAL:HG12	2:L:54:ASN:OD1	1.94	0.64
1:C:906:PHE:O	1:C:909:ILE:HG12	1.97	0.64
2:D:49:ILE:HG13	2:D:55:THR:HG22	1.77	0.64
1:A:371:LEU:C	1:A:373:PRO:HD2	2.17	0.64
1:A:290:ASP:O	1:A:297:SER:HB3	1.98	0.64
1:C:327:VAL:HG12	1:C:542:ASN:HB3	1.80	0.64
1:C:371:LEU:C	1:C:373:PRO:HD2	2.17	0.64
1:B:804:GLN:HA	1:B:817:PRO:HG2	1.81	0.63
5:O:29:ILE:HD12	5:O:32:TYR:O	1.99	0.63
5:K:29:ILE:HD12	5:K:32:TYR:O	1.99	0.62
1:B:1089:PHE:HE2	1:C:917:TYR:HD2	1.47	0.62
1:A:930:ALA:O	1:A:934:ILE:HG12	2.00	0.62
5:R:29:ILE:HD12	5:R:32:TYR:O	1.99	0.62
1:B:62:VAL:HB	1:B:267:VAL:O	2.00	0.61
4:Q:48:VAL:HG13	4:Q:64:VAL:HG11	1.82	0.61
1:A:722:VAL:HG22	1:A:934:ILE:CG1	2.30	0.61
1:A:300:LYS:HG2	1:A:308:VAL:HG23	1.81	0.61
1:A:449:TYR:OH	2:L:28:THR:HG23	2.00	0.61
4:I:48:VAL:HG13	4:I:64:VAL:HG11	1.82	0.61
1:A:595:VAL:HG13	1:A:610:VAL:HG13	1.83	0.61
3:J:35:TRP:NE1	3:J:88:CYS:SG	2.74	0.61
3:J:89:GLN:HA	3:J:98:PHE:HA	1.82	0.61
4:N:48:VAL:HG13	4:N:64:VAL:HG11	1.82	0.61
3:P:89:GLN:HA	3:P:98:PHE:HA	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:821:LEU:HD13	1:B:935:GLN:HG3	1.82	0.61
3:H:89:GLN:HA	3:H:98:PHE:HA	1.82	0.61
1:B:449:TYR:OH	2:F:28:THR:HG23	2.00	0.61
1:A:454:ARG:HA	1:A:491:PRO:O	2.01	0.61
1:B:336:CYS:SG	1:B:362:VAL:O	2.58	0.60
1:C:449:TYR:OH	2:D:28:THR:HG23	2.00	0.60
1:B:671:CYS:SG	1:B:697:MET:HB3	2.41	0.60
1:C:454:ARG:HA	1:C:491:PRO:O	2.01	0.60
3:J:29:ILE:HG12	3:J:92:HIS:HB2	1.83	0.60
3:P:35:TRP:NE1	3:P:88:CYS:SG	2.74	0.60
1:A:65:PHE:CE1	1:A:82:PRO:HG2	2.36	0.60
1:A:453:TYR:O	1:A:492:LEU:HA	2.02	0.60
1:B:454:ARG:HA	1:B:491:PRO:O	2.01	0.60
1:C:453:TYR:O	1:C:492:LEU:HA	2.02	0.60
1:C:770:ILE:O	1:C:774:GLN:HG2	2.01	0.60
4:I:6:GLU:HG3	4:I:116:THR:HG22	1.84	0.60
1:B:383:SER:HB3	1:B:386:LYS:HB2	1.84	0.60
4:N:6:GLU:HG3	4:N:116:THR:HG22	1.84	0.60
1:A:383:SER:HB3	1:A:386:LYS:HB2	1.84	0.60
1:C:383:SER:HB3	1:C:386:LYS:HB2	1.84	0.60
1:A:722:VAL:HG23	1:A:934:ILE:CG1	2.21	0.59
3:H:35:TRP:NE1	3:H:88:CYS:SG	2.74	0.59
5:O:62:PHE:CD1	5:O:75:ILE:HD11	2.37	0.59
1:C:726:ILE:HD13	1:C:945:LEU:HD13	1.83	0.59
4:I:24:SER:OG	4:I:27:PHE:CE2	2.50	0.59
1:B:453:TYR:O	1:B:492:LEU:HA	2.02	0.59
1:B:1116:THR:HG22	1:B:1138:TYR:HB3	1.84	0.59
3:P:29:ILE:HG12	3:P:92:HIS:HB2	1.83	0.59
5:R:62:PHE:CD1	5:R:75:ILE:HD11	2.38	0.59
1:B:822:LEU:CD2	1:B:945:LEU:HD21	2.32	0.58
1:A:394:ASN:HD21	1:A:518:LEU:HD12	1.68	0.58
5:K:62:PHE:CD1	5:K:75:ILE:HD11	2.37	0.58
1:A:336:CYS:HB3	1:A:337:PRO:HD2	1.84	0.58
1:B:770:ILE:O	1:B:774:GLN:HG2	2.03	0.58
2:L:32:MET:HA	2:L:96:GLY:H	1.68	0.58
1:A:909:ILE:HD13	1:A:1049:LEU:HD21	1.85	0.58
3:H:21:ILE:HG22	3:H:73:LEU:HB3	1.86	0.58
3:H:29:ILE:HG12	3:H:92:HIS:HB2	1.84	0.58
1:A:598:ILE:HG23	1:A:664:ILE:HG21	1.86	0.58
1:C:726:ILE:HG12	1:C:1061:VAL:HG22	1.85	0.58
1:B:905:ARG:HD3	1:B:1049:LEU:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:32:MET:HA	2:F:96:GLY:H	1.69	0.57
3:P:21:ILE:HG22	3:P:73:LEU:HB3	1.86	0.57
4:Q:6:GLU:HG3	4:Q:116:THR:HG22	1.84	0.57
1:A:188:ASN:HB3	1:A:190:ARG:CZ	2.35	0.57
2:D:32:MET:HA	2:D:96:GLY:H	1.68	0.57
1:B:468:ILE:O	1:B:468:ILE:HG22	2.05	0.57
1:A:338:PHE:CE2	1:A:363:ALA:HB1	2.41	0.56
1:A:500:THR:CG2	4:Q:59:TYR:OH	2.53	0.56
3:P:7:SER:HB3	3:P:22:THR:OG1	2.05	0.56
1:B:338:PHE:CE2	1:B:363:ALA:HB1	2.41	0.56
1:C:500:THR:CG2	4:N:59:TYR:OH	2.53	0.56
3:J:21:ILE:HG22	3:J:73:LEU:HB3	1.86	0.56
1:B:500:THR:CG2	4:I:59:TYR:OH	2.53	0.56
3:J:7:SER:HB3	3:J:22:THR:OG1	2.05	0.56
1:A:468:ILE:HG22	1:A:468:ILE:O	2.05	0.56
3:H:7:SER:HB3	3:H:22:THR:OG1	2.05	0.56
1:B:620:VAL:HG21	1:B:651:ILE:HD11	1.87	0.56
1:C:338:PHE:CE2	1:C:363:ALA:HB1	2.41	0.56
1:A:498:ARG:HG2	1:A:499:PRO:HD2	1.87	0.56
1:B:276:LEU:HB3	1:B:289:VAL:HG22	1.88	0.56
1:B:503:VAL:HG11	5:K:30:SER:HB2	1.88	0.56
1:A:347:PHE:CE2	1:A:399:SER:HB2	2.41	0.56
1:B:347:PHE:CE2	1:B:399:SER:HB2	2.41	0.56
1:C:945:LEU:O	1:C:947:LYS:N	2.38	0.56
1:A:40:ASP:HB3	1:A:42:VAL:HG23	1.87	0.56
1:B:498:ARG:HG2	1:B:499:PRO:HD2	1.87	0.56
1:B:574:ASP:O	1:B:586:ASP:OD1	2.23	0.56
1:C:347:PHE:CE2	1:C:399:SER:HB2	2.41	0.56
5:R:20:THR:HG22	5:R:74:THR:HG22	1.88	0.56
1:C:326:ILE:HD13	1:C:534:VAL:H	1.71	0.55
3:P:34:ALA:HB3	3:P:89:GLN:HG2	1.89	0.55
2:L:17:SER:HA	2:L:80:MET:O	2.07	0.55
1:A:503:VAL:HG11	5:R:30:SER:HB2	1.88	0.55
1:C:500:THR:HG21	4:N:59:TYR:OH	2.07	0.55
5:O:20:THR:HG22	5:O:74:THR:HG22	1.88	0.55
5:K:20:THR:HG22	5:K:74:THR:HG22	1.88	0.55
1:A:472:ILE:CD1	1:A:484:ALA:N	2.70	0.55
1:B:325:SER:H	1:B:539:VAL:HG23	1.71	0.55
2:D:17:SER:HA	2:D:80:MET:O	2.06	0.55
1:C:472:ILE:CD1	1:C:484:ALA:N	2.70	0.55
3:J:34:ALA:HB3	3:J:89:GLN:HG2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:472:ILE:CD1	1:B:484:ALA:N	2.70	0.55
1:B:500:THR:HG21	4:I:59:TYR:OH	2.07	0.55
2:F:17:SER:HA	2:F:80:MET:O	2.07	0.55
3:H:29:ILE:CG2	3:H:92:HIS:HB2	2.36	0.55
1:C:498:ARG:HG2	1:C:499:PRO:HD2	1.87	0.54
1:A:500:THR:HG21	4:Q:59:TYR:OH	2.07	0.54
3:H:34:ALA:HB3	3:H:89:GLN:HG2	1.89	0.54
1:B:389:ASP:HA	1:B:527:PRO:HB3	1.90	0.54
1:C:941:THR:N	1:C:942:PRO:HD3	2.21	0.54
1:C:468:ILE:HG22	1:C:468:ILE:O	2.05	0.54
2:F:97:SER:HB3	2:F:102:ASP:HA	1.90	0.54
1:C:503:VAL:HG11	5:O:30:SER:HB2	1.88	0.53
5:K:37:GLN:HG3	5:K:86:TYR:CE2	2.44	0.53
2:L:65:PHE:HA	2:L:79:GLN:O	2.08	0.53
2:L:97:SER:HB3	2:L:102:ASP:HA	1.90	0.53
1:B:779:GLN:O	1:B:783:ALA:HB3	2.08	0.53
1:C:729:VAL:H	1:C:1059:GLY:HA2	1.74	0.53
2:D:65:PHE:HA	2:D:79:GLN:O	2.08	0.53
2:D:97:SER:HB3	2:D:102:ASP:HA	1.90	0.53
3:J:29:ILE:CG2	3:J:92:HIS:HB2	2.36	0.53
2:F:70:ASP:HB2	2:F:77:TYR:HE2	1.73	0.53
1:B:106:PHE:HB3	1:B:235:ILE:HD13	1.90	0.53
4:N:100:SER:HG	4:N:107:VAL:H	1.55	0.53
3:P:29:ILE:CG2	3:P:92:HIS:HB2	2.37	0.53
5:R:37:GLN:HG3	5:R:86:TYR:CE2	2.44	0.53
1:C:620:VAL:HG23	1:C:621:PRO:HD3	1.90	0.53
1:A:826:VAL:HB	1:A:1057:PRO:HG2	1.91	0.53
1:C:401:VAL:HG22	1:C:509:ARG:HG2	1.91	0.53
2:F:59:ASP:HA	2:F:62:LYS:HE3	1.91	0.53
1:B:911:VAL:HG13	1:B:1106:GLN:HE22	1.72	0.53
2:D:70:ASP:HB2	2:D:77:TYR:HE2	1.73	0.53
5:O:37:GLN:HG3	5:O:86:TYR:CE2	2.44	0.53
2:L:70:ASP:HB2	2:L:77:TYR:HE2	1.73	0.53
1:A:725:GLU:HG3	1:A:1064:HIS:CD2	2.44	0.53
2:L:59:ASP:HA	2:L:62:LYS:HE3	1.91	0.53
1:A:642:VAL:HG13	1:A:649:CYS:SG	2.50	0.52
1:B:656:VAL:HG12	1:B:658:ASN:H	1.74	0.52
1:A:722:VAL:CG2	1:A:934:ILE:HD11	2.39	0.52
1:A:401:VAL:HG22	1:A:509:ARG:HG2	1.91	0.52
2:D:59:ASP:HA	2:D:62:LYS:HE3	1.91	0.52
2:F:65:PHE:HA	2:F:79:GLN:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:963:VAL:HG11	1:C:570:ALA:HB1	1.91	0.52
1:B:401:VAL:HG22	1:B:509:ARG:HG2	1.91	0.52
1:A:391:CYS:SG	1:A:524:VAL:O	2.68	0.52
1:C:391:CYS:SG	1:C:524:VAL:O	2.68	0.52
1:A:1106:GLN:HE21	1:A:1109:PHE:HB3	1.74	0.52
1:C:444:LYS:HE2	1:C:448:ASN:HA	1.92	0.52
1:C:296:LEU:O	1:C:299:THR:HG22	2.09	0.52
1:B:106:PHE:CB	1:B:235:ILE:HD13	2.40	0.52
1:B:391:CYS:SG	1:B:524:VAL:O	2.68	0.52
1:B:1056:ALA:HB1	1:B:1057:PRO:HD2	1.92	0.52
2:F:33:SER:HB2	2:F:48:VAL:HG23	1.92	0.52
4:I:3:GLN:O	4:I:5:VAL:HG23	2.10	0.52
1:A:617:CYS:SG	1:A:644:GLN:HB2	2.50	0.51
2:D:33:SER:HB2	2:D:48:VAL:HG23	1.92	0.51
4:I:100:SER:OG	4:I:107:VAL:N	2.42	0.51
1:B:338:PHE:HE2	1:B:363:ALA:HB1	1.75	0.51
1:B:934:ILE:HG22	1:B:938:LEU:CD1	2.41	0.51
2:L:33:SER:HB2	2:L:48:VAL:HG23	1.92	0.51
4:Q:3:GLN:O	4:Q:5:VAL:HG23	2.10	0.51
4:N:3:GLN:O	4:N:5:VAL:HG23	2.10	0.51
1:C:736:VAL:HG22	1:C:767:LEU:CD1	2.40	0.51
1:C:767:LEU:HD21	1:C:1008:VAL:HG22	1.91	0.51
1:C:909:ILE:HD13	1:C:1049:LEU:HD21	1.91	0.51
1:C:360:ASN:H	1:C:523:THR:HB	1.76	0.51
1:A:338:PHE:HE2	1:A:363:ALA:HB1	1.75	0.51
1:A:906:PHE:O	1:A:909:ILE:HG12	2.11	0.51
1:C:289:VAL:HG23	1:C:306:PHE:CZ	2.46	0.51
1:C:338:PHE:HE2	1:C:363:ALA:HB1	1.75	0.51
4:I:51:ILE:HA	4:I:57:ASN:O	2.11	0.51
1:A:357:ARG:HH12	1:B:169:GLU:H	1.60	0.50
4:N:51:ILE:HA	4:N:57:ASN:O	2.11	0.50
1:A:444:LYS:HE2	1:A:448:ASN:HA	1.92	0.50
1:B:444:LYS:HE2	1:B:448:ASN:HA	1.92	0.50
1:C:779:GLN:O	1:C:783:ALA:HB3	2.11	0.50
4:I:100:SER:HG	4:I:107:VAL:H	1.59	0.50
2:L:98:GLY:N	2:L:103:ALA:H	2.06	0.50
4:Q:51:ILE:HA	4:Q:57:ASN:O	2.11	0.50
3:P:34:ALA:HB2	3:P:50:ALA:HA	1.94	0.50
1:A:811:LYS:CB	1:A:812:PRO:HD3	2.42	0.49
3:H:39:LYS:HE3	3:H:42:LYS:HB2	1.94	0.49
3:H:34:ALA:HB2	3:H:50:ALA:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LYS:HG2	1:A:308:VAL:CG2	2.41	0.49
1:B:618:THR:OG1	1:B:619:GLU:HG3	2.13	0.49
3:J:34:ALA:HB2	3:J:50:ALA:HA	1.93	0.49
2:L:50:TYR:HD2	2:L:54:ASN:HB2	1.78	0.49
3:P:32:TRP:O	3:P:91:ALA:HB3	2.13	0.49
1:A:98:SER:HB2	1:A:100:ILE:HG12	1.94	0.49
1:A:472:ILE:CD1	1:A:483:VAL:C	2.81	0.49
1:A:707:TYR:HE1	1:B:898:PHE:H	1.59	0.49
1:B:826:VAL:CB	1:B:1057:PRO:HG3	2.43	0.49
3:P:39:LYS:HE3	3:P:42:LYS:HB2	1.95	0.49
1:A:977:LEU:HD11	1:A:1000:ARG:HH12	1.77	0.49
1:B:472:ILE:HD13	1:B:484:ALA:N	2.28	0.49
1:C:878:LEU:HA	1:C:881:THR:HG22	1.95	0.49
2:D:67:ILE:HA	2:D:77:TYR:O	2.13	0.49
1:A:433:VAL:HG13	1:A:512:VAL:HG22	1.95	0.49
1:B:538:CYS:HA	1:B:551:VAL:HG22	1.95	0.49
1:B:911:VAL:HA	1:B:1106:GLN:NE2	2.27	0.49
1:C:312:ILE:HG23	1:C:312:ILE:O	2.13	0.49
2:D:50:TYR:HD2	2:D:54:ASN:HB2	1.78	0.49
1:A:81:ASN:N	1:A:82:PRO:HD3	2.28	0.49
1:C:472:ILE:HD11	1:C:484:ALA:N	2.28	0.49
2:F:50:TYR:HD2	2:F:54:ASN:HB2	1.78	0.49
1:A:472:ILE:HD13	1:A:484:ALA:N	2.28	0.48
1:B:718:PHE:HB3	1:B:1067:TYR:CE2	2.47	0.48
1:C:445:VAL:HA	1:C:499:PRO:HG3	1.95	0.48
1:C:472:ILE:HD13	1:C:483:VAL:C	2.33	0.48
3:J:32:TRP:O	3:J:91:ALA:HB3	2.13	0.48
1:A:870:ILE:O	1:A:874:THR:HG23	2.14	0.48
1:B:472:ILE:CD1	1:B:483:VAL:C	2.81	0.48
1:B:472:ILE:HD13	1:B:483:VAL:C	2.33	0.48
1:C:472:ILE:HD13	1:C:484:ALA:N	2.28	0.48
1:A:472:ILE:HD13	1:A:483:VAL:C	2.33	0.48
1:A:472:ILE:HD11	1:A:484:ALA:N	2.28	0.48
1:C:433:VAL:HG13	1:C:512:VAL:HG22	1.96	0.48
1:C:441:LEU:HD21	4:N:103:TRP:CD2	2.49	0.48
1:C:472:ILE:CD1	1:C:483:VAL:C	2.81	0.48
2:F:67:ILE:HA	2:F:77:TYR:O	2.13	0.48
4:Q:31:THR:OG1	4:Q:32:TYR:CD1	2.66	0.48
1:A:445:VAL:HA	1:A:499:PRO:HG3	1.95	0.48
1:A:1114:ILE:HG22	1:A:1115:ILE:N	2.28	0.48
1:C:826:VAL:CB	1:C:1057:PRO:HG3	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:441:LEU:HD13	4:N:102:GLY:HA3	1.95	0.48
4:I:31:THR:OG1	4:I:32:TYR:CD1	2.66	0.48
2:L:67:ILE:HA	2:L:77:TYR:O	2.13	0.48
3:J:39:LYS:HE3	3:J:42:LYS:HB2	1.94	0.48
4:N:31:THR:OG1	4:N:32:TYR:CD1	2.66	0.48
4:I:19:ARG:HG3	4:I:82:GLN:HG2	1.95	0.48
1:B:468:ILE:O	1:B:468:ILE:CG2	2.62	0.48
1:B:472:ILE:HD11	1:B:484:ALA:N	2.28	0.48
1:B:1095:PHE:CE1	1:B:1104:VAL:HG22	2.49	0.48
1:C:51:THR:O	1:C:51:THR:OG1	2.26	0.48
4:N:19:ARG:HG3	4:N:82:GLN:HG2	1.95	0.48
4:N:100:SER:OG	4:N:107:VAL:N	2.42	0.48
1:B:441:LEU:HD13	4:I:102:GLY:HA3	1.95	0.47
1:C:394:ASN:HD21	1:C:518:LEU:HD12	1.79	0.47
3:H:32:TRP:O	3:H:91:ALA:HB3	2.13	0.47
3:H:65:SER:HB3	3:H:72:THR:O	2.14	0.47
4:Q:19:ARG:HG3	4:Q:82:GLN:HG2	1.95	0.47
1:A:441:LEU:HD21	4:Q:103:TRP:CD2	2.49	0.47
1:A:741:TYR:CE1	1:A:966:LEU:HG	2.49	0.47
1:A:85:PRO:HD2	1:A:269:TYR:OH	2.14	0.47
1:A:66:HIS:HB2	1:A:263:ALA:O	2.15	0.47
1:B:433:VAL:HG13	1:B:512:VAL:HG22	1.95	0.47
3:H:35:TRP:HB3	3:H:47:LEU:HD23	1.97	0.47
1:A:441:LEU:HD13	4:Q:102:GLY:HA3	1.95	0.47
1:B:441:LEU:HD21	4:I:103:TRP:CD2	2.49	0.47
1:B:445:VAL:HA	1:B:499:PRO:HG3	1.95	0.47
1:C:468:ILE:O	1:C:468:ILE:CG2	2.62	0.47
5:R:62:PHE:CE1	5:R:75:ILE:CD1	2.98	0.47
3:J:65:SER:HB3	3:J:72:THR:O	2.14	0.47
1:A:468:ILE:O	1:A:468:ILE:CG2	2.62	0.47
1:B:1006:THR:O	1:B:1010:GLN:HG2	2.15	0.47
4:Q:100:SER:HG	4:Q:107:VAL:H	1.60	0.47
1:A:421:TYR:HD1	1:A:457:ARG:HB3	1.80	0.47
1:B:394:ASN:HD21	1:B:518:LEU:HD12	1.79	0.47
1:C:805:ILE:HD12	1:C:878:LEU:HD11	1.97	0.47
1:C:905:ARG:HD3	1:C:1049:LEU:O	2.14	0.47
1:A:789:TYR:HA	1:C:703:ASN:O	2.15	0.46
1:A:1031:GLU:HG2	1:C:1040:VAL:H	1.81	0.46
1:B:421:TYR:HD1	1:B:457:ARG:HB3	1.80	0.46
1:B:898:PHE:N	1:B:899:PRO:HD2	2.30	0.46
3:P:65:SER:HB3	3:P:72:THR:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:768:THR:O	1:A:772:VAL:HG23	2.15	0.46
2:D:29:VAL:HG13	2:D:32:MET:SD	2.56	0.46
1:B:1141:LEU:HD11	1:C:1141:LEU:HD13	1.97	0.46
1:C:84:LEU:HD13	1:C:267:VAL:HG21	1.97	0.46
1:C:426:PRO:HG2	1:C:429:PHE:HA	1.98	0.46
3:P:35:TRP:HB3	3:P:47:LEU:HD23	1.97	0.46
1:B:426:PRO:HB3	1:B:463:PRO:HB3	1.98	0.46
1:B:811:LYS:N	1:B:812:PRO:HD3	2.31	0.46
4:I:6:GLU:OE2	4:I:20:LEU:HB3	2.16	0.46
1:B:934:ILE:HG22	1:B:938:LEU:HD11	1.98	0.46
3:J:35:TRP:HB3	3:J:47:LEU:HD23	1.97	0.46
4:N:87:ARG:O	4:N:120:VAL:HG11	2.16	0.46
1:C:909:ILE:HG22	1:C:1047:TYR:HB3	1.98	0.46
2:D:107:TRP:CZ3	3:J:44:PRO:HB2	2.51	0.46
2:F:107:TRP:CZ3	3:H:44:PRO:HB2	2.51	0.46
1:A:64:TRP:HD1	1:A:65:PHE:N	2.13	0.46
4:Q:100:SER:OG	4:Q:107:VAL:N	2.42	0.46
1:B:426:PRO:HG2	1:B:429:PHE:HA	1.98	0.45
1:B:802:PHE:CD1	1:B:805:ILE:HD11	2.50	0.45
1:C:491:PRO:HG2	1:C:492:LEU:HG	1.98	0.45
1:C:856:LYS:HD3	1:C:966:LEU:HD12	1.98	0.45
2:D:97:SER:CB	2:D:102:ASP:HA	2.46	0.45
1:C:421:TYR:HD1	1:C:457:ARG:HB3	1.80	0.45
4:I:87:ARG:O	4:I:120:VAL:HG11	2.16	0.45
4:Q:24:SER:OG	4:Q:27:PHE:CE2	2.50	0.45
4:Q:87:ARG:O	4:Q:120:VAL:HG11	2.16	0.45
1:A:426:PRO:HG2	1:A:429:PHE:HA	1.98	0.45
1:A:541:PHE:CE2	1:A:587:ILE:HG21	2.51	0.45
1:A:558:LYS:HA	1:A:558:LYS:HD3	1.84	0.45
1:B:785:VAL:HG12	1:B:787:GLN:O	2.16	0.45
1:C:91:TYR:CG	1:C:91:TYR:O	2.70	0.45
1:C:296:LEU:HB2	1:C:608:VAL:HG21	1.97	0.45
2:L:29:VAL:HG13	2:L:32:MET:SD	2.56	0.45
1:A:62:VAL:HB	1:A:267:VAL:O	2.16	0.45
1:B:28:TYR:HE1	1:B:63:THR:HG22	1.81	0.45
1:B:491:PRO:HG2	1:B:492:LEU:HG	1.98	0.45
2:F:97:SER:CB	2:F:102:ASP:HA	2.46	0.45
1:A:101:ILE:HG13	1:A:242:LEU:HD12	1.99	0.45
1:A:426:PRO:HB3	1:A:463:PRO:HB3	1.98	0.45
2:F:29:VAL:HG13	2:F:32:MET:SD	2.56	0.45
5:K:29:ILE:HD11	5:K:71:PHE:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:107:TRP:CZ3	3:P:44:PRO:HB2	2.51	0.45
2:D:98:GLY:N	2:D:103:ALA:H	2.06	0.45
4:Q:11:VAL:HG22	4:Q:119:THR:HG21	1.99	0.45
1:A:296:LEU:CD1	1:A:599:THR:HG21	2.47	0.45
1:A:491:PRO:HG2	1:A:492:LEU:HG	1.98	0.45
1:A:722:VAL:HG21	1:A:934:ILE:CG1	2.24	0.45
1:A:1114:ILE:CG2	1:A:1115:ILE:N	2.79	0.45
4:N:6:GLU:OE2	4:N:20:LEU:HB3	2.16	0.45
2:F:98:GLY:N	2:F:103:ALA:H	2.06	0.45
1:A:1028:LYS:O	1:A:1032:CYS:HB2	2.17	0.45
1:C:426:PRO:HB3	1:C:463:PRO:HB3	1.98	0.45
1:C:922:LEU:HG	1:C:926:GLN:HE21	1.81	0.45
4:N:11:VAL:HG22	4:N:119:THR:HG21	1.99	0.45
5:K:62:PHE:CE1	5:K:75:ILE:CD1	2.98	0.45
4:Q:6:GLU:OE2	4:Q:20:LEU:HB3	2.16	0.45
5:R:46:LEU:HD21	5:R:49:TYR:HB3	1.99	0.45
1:A:928:ASN:O	1:A:931:ILE:HG13	2.17	0.44
3:J:36:TYR:CD1	3:J:46:VAL:HA	2.52	0.44
5:O:29:ILE:HD11	5:O:71:PHE:CE2	2.52	0.44
4:I:2:VAL:HG13	4:I:27:PHE:CD2	2.52	0.44
4:I:11:VAL:HG22	4:I:119:THR:HG21	1.99	0.44
2:L:97:SER:CB	2:L:102:ASP:HA	2.46	0.44
3:P:29:ILE:HA	3:P:92:HIS:CD2	2.52	0.44
3:P:36:TYR:CD1	3:P:46:VAL:HA	2.52	0.44
1:A:318:PHE:CZ	1:A:620:VAL:O	2.70	0.44
1:C:389:ASP:HA	1:C:527:PRO:HG3	2.00	0.44
1:C:518:LEU:C	1:C:520:ALA:H	2.20	0.44
3:J:29:ILE:HA	3:J:92:HIS:CD2	2.52	0.44
4:N:24:SER:OG	4:N:27:PHE:CD2	2.70	0.44
3:H:36:TYR:CD1	3:H:46:VAL:HA	2.52	0.44
5:K:46:LEU:HD21	5:K:49:TYR:HB3	1.99	0.44
1:A:39:PRO:HG2	1:A:40:ASP:H	1.82	0.44
1:C:395:VAL:HG21	1:C:524:VAL:HG11	1.99	0.44
5:O:62:PHE:CE1	5:O:75:ILE:CD1	2.97	0.44
1:A:979:ASP:O	1:A:983:ARG:HG2	2.16	0.44
1:B:395:VAL:HG21	1:B:524:VAL:HG11	1.99	0.44
4:N:2:VAL:HG13	4:N:27:PHE:CD2	2.52	0.44
3:H:29:ILE:HA	3:H:92:HIS:CD2	2.53	0.44
1:B:802:PHE:HD1	1:B:805:ILE:HD11	1.82	0.44
1:C:453:TYR:N	1:C:493:ARG:O	2.45	0.44
1:C:716:THR:HG21	1:C:1073:LYS:HE2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:3:GLN:HB2	3:H:26:SER:CB	2.46	0.44
3:H:6:GLN:HA	3:H:23:CYS:HA	1.99	0.44
3:P:29:ILE:HG12	3:P:92:HIS:CB	2.48	0.44
3:P:35:TRP:HH2	3:P:65:SER:O	2.01	0.44
5:R:29:ILE:HD11	5:R:71:PHE:CE2	2.52	0.44
1:A:29:THR:HG23	1:A:62:VAL:HG23	1.98	0.44
1:A:1081:ILE:HG13	1:A:1095:PHE:CD2	2.53	0.44
1:C:328:ARG:HG3	1:C:531:THR:O	2.18	0.44
3:J:29:ILE:CG1	3:J:92:HIS:HB2	2.48	0.44
3:J:35:TRP:HH2	3:J:65:SER:O	2.01	0.44
4:Q:24:SER:OG	4:Q:27:PHE:CD2	2.70	0.44
1:A:919:ASN:O	1:A:923:ILE:HG13	2.18	0.44
1:B:825:LYS:NZ	1:B:939:SER:HA	2.33	0.44
3:J:6:GLN:HA	3:J:23:CYS:HA	1.99	0.44
4:Q:2:VAL:HG13	4:Q:27:PHE:CD2	2.52	0.44
1:A:645:THR:OG1	1:A:646:ARG:N	2.50	0.43
1:A:1001:LEU:HD12	1:A:1001:LEU:HA	1.83	0.43
3:H:28:GLY:HA2	3:H:69:THR:HG22	2.00	0.43
3:H:35:TRP:HH2	3:H:65:SER:O	2.01	0.43
1:A:395:VAL:HG21	1:A:524:VAL:HG11	1.99	0.43
1:A:980:ILE:HD11	1:A:996:LEU:HD11	1.99	0.43
1:B:1089:PHE:CE2	1:C:917:TYR:HD2	2.31	0.43
3:H:29:ILE:CG1	3:H:92:HIS:HB2	2.48	0.43
3:P:6:GLN:HA	3:P:23:CYS:HA	1.99	0.43
1:B:329:PHE:HB3	1:B:330:PRO:HD2	2.00	0.43
1:B:362:VAL:HG13	1:B:526:GLY:HA2	1.98	0.43
3:P:28:GLY:HA2	3:P:69:THR:HG22	2.01	0.43
1:A:91:TYR:O	1:A:91:TYR:CG	2.71	0.43
1:A:360:ASN:HD21	1:B:170:TYR:HB3	1.84	0.43
1:A:736:VAL:HG11	1:A:1004:LEU:HD11	1.99	0.43
3:P:3:GLN:HB2	3:P:26:SER:CB	2.46	0.43
1:C:963:VAL:O	1:C:966:LEU:HB2	2.18	0.43
1:A:371:LEU:HD12	1:A:371:LEU:HA	1.89	0.43
1:A:543:PHE:CZ	1:A:585:LEU:HD12	2.54	0.43
1:A:802:PHE:HB3	1:A:805:ILE:HG12	2.01	0.43
1:B:205:SER:HB3	1:B:226:LEU:HD22	2.01	0.43
1:A:96:GLU:HG3	1:A:99:ASN:H	1.82	0.43
1:B:961:THR:O	1:B:965:GLN:HG3	2.18	0.43
5:O:46:LEU:HD21	5:O:49:TYR:HB3	2.00	0.43
5:K:23:CYS:SG	5:K:33:LEU:HD11	2.59	0.43
1:A:560:LEU:HB2	1:A:563:GLN:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:878:LEU:CD1	1:C:1052:PHE:CD1	3.02	0.43
1:C:940:SER:C	1:C:942:PRO:HD3	2.39	0.43
5:R:23:CYS:SG	5:R:33:LEU:HD11	2.59	0.43
1:A:713:ALA:O	1:B:894:LEU:HD22	2.19	0.43
1:A:718:PHE:HB3	1:A:1067:TYR:CE2	2.54	0.43
1:A:375:PHE:HA	1:A:435:ALA:O	2.19	0.43
1:B:499:PRO:CB	5:K:94:TYR:HD2	2.29	0.43
1:C:726:ILE:CG2	1:C:948:LEU:HG	2.48	0.43
2:D:5:VAL:HB	2:D:23:ALA:HB3	2.01	0.43
3:P:29:ILE:CG1	3:P:92:HIS:HB2	2.48	0.43
1:A:330:PRO:HA	1:A:579:PRO:HB2	2.00	0.42
1:C:375:PHE:HA	1:C:435:ALA:O	2.19	0.42
3:J:28:GLY:HA2	3:J:69:THR:HG22	2.01	0.42
4:Q:68:PHE:HD2	4:Q:81:LEU:HD21	1.85	0.42
1:A:372:ALA:N	1:A:373:PRO:HD2	2.35	0.42
1:B:990:GLU:HA	1:B:993:ILE:HD12	2.00	0.42
1:C:418:ILE:HD13	1:C:418:ILE:HA	1.88	0.42
3:J:3:GLN:HB2	3:J:26:SER:CB	2.46	0.42
4:I:24:SER:OG	4:I:27:PHE:CD2	2.70	0.42
1:A:121:ASN:HA	1:A:126:VAL:HG13	2.00	0.42
1:A:276:LEU:HD23	1:A:306:PHE:CE1	2.54	0.42
1:C:777:ASN:O	1:C:781:VAL:HG23	2.19	0.42
3:J:39:LYS:H	3:J:42:LYS:HE2	1.85	0.42
1:A:224:GLU:O	1:A:225:PRO:C	2.58	0.42
1:A:963:VAL:O	1:A:966:LEU:HB2	2.19	0.42
1:A:1089:PHE:CE2	1:B:917:TYR:HD2	2.38	0.42
1:B:372:ALA:N	1:B:373:PRO:HD2	2.34	0.42
1:C:29:THR:HG22	1:C:30:ASN:H	1.84	0.42
1:C:499:PRO:CB	5:O:94:TYR:HD2	2.29	0.42
1:C:966:LEU:HA	1:C:966:LEU:HD23	1.83	0.42
4:I:68:PHE:HD2	4:I:81:LEU:HD21	1.85	0.42
2:L:5:VAL:HB	2:L:23:ALA:HB3	2.01	0.42
3:P:39:LYS:H	3:P:42:LYS:HE2	1.85	0.42
1:A:120:VAL:O	1:A:126:VAL:HA	2.20	0.42
1:A:722:VAL:HG22	1:A:934:ILE:HD11	2.00	0.42
1:B:655:TYR:OH	1:B:696:THR:HG23	2.20	0.42
1:C:33:THR:HA	1:C:58:PHE:CE1	2.55	0.42
1:C:372:ALA:N	1:C:373:PRO:HD2	2.35	0.42
5:O:8:PRO:HG2	5:O:21:ILE:HA	2.01	0.42
1:A:296:LEU:HD13	1:A:599:THR:CG2	2.50	0.42
1:A:392:PHE:CG	1:A:515:PHE:HB3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:707:TYR:HB3	1:B:792:PRO:HG3	2.01	0.42
1:B:572:THR:HG22	1:C:856:LYS:HE2	2.00	0.42
1:B:612:TYR:O	1:B:648:GLY:HA3	2.20	0.42
1:C:802:PHE:CD1	1:C:805:ILE:HD11	2.54	0.42
1:C:870:ILE:HD13	1:C:870:ILE:HA	1.87	0.42
1:B:569:ILE:O	1:B:570:ALA:HB3	2.19	0.42
3:H:39:LYS:H	3:H:42:LYS:HE2	1.85	0.42
1:A:334:ASN:HD22	1:A:334:ASN:HA	1.70	0.42
1:A:517:LEU:HD23	1:A:517:LEU:HA	1.92	0.42
1:B:392:PHE:CG	1:B:515:PHE:HB3	2.55	0.42
1:C:718:PHE:HB3	1:C:1067:TYR:CE2	2.55	0.42
3:J:29:ILE:HG12	3:J:92:HIS:CB	2.48	0.42
3:H:29:ILE:HG12	3:H:92:HIS:CB	2.48	0.42
3:H:34:ALA:O	3:H:88:CYS:HA	2.20	0.42
5:K:8:PRO:HG2	5:K:21:ILE:HA	2.01	0.42
3:P:34:ALA:O	3:P:88:CYS:HA	2.20	0.42
5:R:8:PRO:HG2	5:R:21:ILE:HA	2.01	0.42
1:A:713:ALA:HB2	1:B:895:GLN:HG2	2.01	0.42
1:A:1094:VAL:HG12	1:A:1095:PHE:N	2.35	0.42
1:B:106:PHE:HB3	1:B:235:ILE:CG2	2.50	0.42
1:C:336:CYS:SG	1:C:362:VAL:O	2.78	0.42
5:O:23:CYS:SG	5:O:33:LEU:HD11	2.59	0.42
1:B:572:THR:CG2	1:C:856:LYS:HE3	2.49	0.42
2:F:5:VAL:HB	2:F:23:ALA:HB3	2.01	0.42
1:A:296:LEU:HD13	1:A:599:THR:HG21	2.02	0.41
1:A:777:ASN:O	1:A:781:VAL:HG23	2.20	0.41
1:A:705:VAL:CG1	1:B:895:GLN:HB3	2.50	0.41
1:B:375:PHE:HA	1:B:435:ALA:O	2.19	0.41
1:C:856:LYS:HD2	1:C:963:VAL:HG13	2.02	0.41
1:C:342:PHE:HD1	1:C:436:TRP:HH2	1.68	0.41
1:C:1043:CYS:HB3	1:C:1048:HIS:CD2	2.55	0.41
3:H:19:VAL:O	3:H:74:THR:HA	2.20	0.41
4:I:11:VAL:HA	4:I:119:THR:HG22	2.03	0.41
1:A:326:ILE:HG23	1:A:539:VAL:HG21	2.02	0.41
1:A:617:CYS:O	1:A:620:VAL:HG22	2.21	0.41
1:B:373:PRO:HB3	5:K:53:ASN:CG	2.41	0.41
1:B:538:CYS:CA	1:B:551:VAL:HG22	2.50	0.41
1:B:1117:THR:HA	1:B:1120:THR:OG1	2.20	0.41
1:C:569:ILE:O	1:C:570:ALA:HB3	2.19	0.41
1:A:600:PRO:HG3	1:A:605:SER:HB3	2.02	0.41
3:J:34:ALA:O	3:J:88:CYS:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:O:34:ALA:O	5:O:88:CYS:HA	2.20	0.41
3:H:32:TRP:HB3	3:H:91:ALA:HB3	2.03	0.41
1:A:39:PRO:HD2	1:A:285:ILE:HD11	2.01	0.41
1:A:342:PHE:HD1	1:A:436:TRP:HH2	1.68	0.41
1:A:595:VAL:HG13	1:A:610:VAL:CG1	2.49	0.41
1:C:373:PRO:HB3	5:O:53:ASN:CG	2.41	0.41
4:N:68:PHE:HD2	4:N:81:LEU:HD21	1.85	0.41
3:P:18:ARG:HA	3:P:76:SER:HA	2.03	0.41
1:A:373:PRO:HB3	5:R:53:ASN:CG	2.40	0.41
3:J:32:TRP:HB3	3:J:91:ALA:HB3	2.03	0.41
4:Q:11:VAL:HA	4:Q:119:THR:HG22	2.03	0.41
5:R:34:ALA:O	5:R:88:CYS:HA	2.20	0.41
1:A:104:TRP:HB3	1:A:106:PHE:CE1	2.56	0.41
1:B:948:LEU:HD21	1:B:1059:GLY:HA3	2.02	0.41
5:K:34:ALA:O	5:K:88:CYS:HA	2.21	0.41
1:A:360:ASN:ND2	1:B:170:TYR:HB3	2.36	0.41
1:B:276:LEU:O	1:B:288:ALA:HA	2.21	0.41
1:B:436:TRP:NE1	1:B:509:ARG:HB2	2.36	0.41
1:B:902:MET:O	1:B:916:LEU:HD22	2.20	0.41
1:C:919:ASN:O	1:C:923:ILE:HG13	2.20	0.41
2:D:5:VAL:O	2:D:23:ALA:N	2.54	0.41
3:P:19:VAL:O	3:P:74:THR:HA	2.20	0.41
1:A:436:TRP:NE1	1:A:509:ARG:HB2	2.36	0.41
1:A:569:ILE:O	1:A:570:ALA:HB3	2.21	0.41
1:A:961:THR:HA	1:A:964:LYS:HE3	2.02	0.41
1:B:369:TYR:HB2	1:B:377:PHE:CZ	2.56	0.41
1:C:39:PRO:HG3	1:C:51:THR:HG21	2.03	0.41
1:C:82:PRO:HG2	1:C:84:LEU:HD21	2.03	0.41
1:C:558:LYS:HE2	1:C:558:LYS:HB2	1.92	0.41
2:D:20:LEU:HD13	2:D:20:LEU:HA	1.95	0.41
3:J:19:VAL:O	3:J:74:THR:HA	2.20	0.41
1:A:395:VAL:HG22	1:A:515:PHE:CD1	2.56	0.40
1:B:453:TYR:N	1:B:493:ARG:O	2.45	0.40
1:B:865:LEU:HD23	1:B:865:LEU:HA	1.86	0.40
1:C:334:ASN:O	1:C:362:VAL:HB	2.21	0.40
1:C:392:PHE:CG	1:C:515:PHE:HB3	2.55	0.40
1:C:526:GLY:N	1:C:527:PRO:HD3	2.35	0.40
3:P:29:ILE:HG23	3:P:92:HIS:CB	2.46	0.40
1:A:722:VAL:HG23	1:A:934:ILE:HD11	1.99	0.40
1:A:878:LEU:HD21	1:A:1052:PHE:HB3	2.03	0.40
1:C:101:ILE:HD11	1:C:263:ALA:HB1	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:HIS:HA	1:A:264:ALA:HA	2.03	0.40
1:A:334:ASN:HB3	1:A:362:VAL:HG23	2.02	0.40
1:B:117:LEU:HA	1:B:130:VAL:HA	2.04	0.40
1:B:708:SER:HB3	1:B:711:SER:OG	2.20	0.40
1:C:369:TYR:HB2	1:C:377:PHE:CZ	2.56	0.40
2:F:5:VAL:O	2:F:23:ALA:N	2.54	0.40
1:B:106:PHE:HD1	1:B:238:PHE:HA	1.86	0.40
1:B:1004:LEU:HD12	1:B:1004:LEU:HA	1.86	0.40
1:C:105:ILE:HD13	1:C:119:ILE:O	2.21	0.40
2:F:24:VAL:HG21	2:F:29:VAL:HG22	2.04	0.40
3:H:18:ARG:HA	3:H:76:SER:HA	2.03	0.40
3:H:54:LEU:HD12	3:H:54:LEU:HA	1.94	0.40
1:A:369:TYR:HB2	1:A:377:PHE:CZ	2.56	0.40
1:A:499:PRO:CB	5:R:94:TYR:HD2	2.29	0.40
1:A:898:PHE:N	1:A:899:PRO:HD2	2.36	0.40
1:B:37:TYR:HA	1:B:223:LEU:H	1.86	0.40
1:B:395:VAL:HG22	1:B:515:PHE:CD1	2.56	0.40
3:J:28:GLY:O	3:J:29:ILE:C	2.59	0.40
4:I:24:SER:HB3	4:I:29:PHE:HD1	1.86	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	926/1300 (71%)	875 (94%)	47 (5%)	4 (0%)	34 69
1	B	903/1300 (70%)	852 (94%)	47 (5%)	4 (0%)	34 69
1	C	924/1300 (71%)	867 (94%)	49 (5%)	8 (1%)	17 56
2	D	114/117 (97%)	110 (96%)	4 (4%)	0	100 100
2	F	114/117 (97%)	110 (96%)	4 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	L	114/117 (97%)	110 (96%)	4 (4%)	0	100	100
3	H	102/107 (95%)	95 (93%)	4 (4%)	3 (3%)	4	28
3	J	102/107 (95%)	95 (93%)	4 (4%)	3 (3%)	4	28
3	P	102/107 (95%)	95 (93%)	4 (4%)	3 (3%)	4	28
4	I	120/122 (98%)	110 (92%)	10 (8%)	0	100	100
4	N	120/122 (98%)	110 (92%)	10 (8%)	0	100	100
4	Q	120/122 (98%)	110 (92%)	10 (8%)	0	100	100
5	K	104/107 (97%)	100 (96%)	4 (4%)	0	100	100
5	O	104/107 (97%)	100 (96%)	4 (4%)	0	100	100
5	R	104/107 (97%)	100 (96%)	4 (4%)	0	100	100
All	All	4073/5259 (77%)	3839 (94%)	209 (5%)	25 (1%)	29	64

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	811	LYS
1	C	946	GLY
1	C	1057	PRO
1	B	325	SER
1	C	530	SER
1	C	707	TYR
1	A	813	SER
1	C	543	PHE
1	A	903	ALA
1	B	543	PHE
1	B	813	SER
1	C	941	THR
3	J	53	GLY
3	H	53	GLY
3	P	53	GLY
1	A	372	ALA
1	B	372	ALA
1	C	372	ALA
3	H	66	GLY
3	P	66	GLY
3	J	29	ILE
3	J	66	GLY
3	H	29	ILE

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Mol	Chain	Res	Type
3	P	29	ILE
1	C	82	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	738/1130 (65%)	720 (98%)	18 (2%)	49	77
1	B	723/1130 (64%)	714 (99%)	9 (1%)	71	88
1	C	734/1130 (65%)	720 (98%)	14 (2%)	57	81
2	D	83/95 (87%)	82 (99%)	1 (1%)	71	88
2	F	83/95 (87%)	82 (99%)	1 (1%)	71	88
2	L	83/95 (87%)	82 (99%)	1 (1%)	71	88
3	H	76/89 (85%)	73 (96%)	3 (4%)	32	67
3	J	76/89 (85%)	73 (96%)	3 (4%)	32	67
3	P	76/89 (85%)	73 (96%)	3 (4%)	32	67
4	I	92/104 (88%)	90 (98%)	2 (2%)	52	79
4	N	92/104 (88%)	90 (98%)	2 (2%)	52	79
4	Q	92/104 (88%)	90 (98%)	2 (2%)	52	79
5	K	85/92 (92%)	84 (99%)	1 (1%)	71	88
5	O	85/92 (92%)	84 (99%)	1 (1%)	71	88
5	R	85/92 (92%)	84 (99%)	1 (1%)	71	88
All	All	3203/4530 (71%)	3141 (98%)	62 (2%)	59	81

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ASP
1	A	102	ARG
1	A	333	THR
1	A	334	ASN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	347	PHE
1	A	349	SER
1	A	361	CYS
1	A	377	PHE
1	A	495	TYR
1	A	518	LEU
1	A	574	ASP
1	A	591	SER
1	A	650	LEU
1	A	660	TYR
1	A	725	GLU
1	A	816	SER
1	A	855	PHE
1	A	906	PHE
1	B	347	PHE
1	B	349	SER
1	B	377	PHE
1	B	495	TYR
1	B	660	TYR
1	B	935	GLN
1	B	1017	GLU
1	B	1121	PHE
1	B	1144	GLU
1	C	130	VAL
1	C	238	PHE
1	C	265	TYR
1	C	301	CYS
1	C	347	PHE
1	C	349	SER
1	C	361	CYS
1	C	377	PHE
1	C	495	TYR
1	C	873	TYR
1	C	894	LEU
1	C	895	GLN
1	C	919	ASN
1	C	1106	GLN
2	D	56	ASP
3	J	23	CYS
3	J	39	LYS
3	J	46	VAL
4	N	58	TYR

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Mol	Chain	Res	Type
4	N	104	HIS
5	O	107	LYS
2	F	56	ASP
3	H	23	CYS
3	H	39	LYS
3	H	46	VAL
4	I	58	TYR
4	I	104	HIS
5	K	107	LYS
2	L	56	ASP
3	P	23	CYS
3	P	39	LYS
3	P	46	VAL
4	Q	58	TYR
4	Q	104	HIS
5	R	107	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	317	ASN
1	A	334	ASN
1	A	540	ASN
1	A	919	ASN
1	A	935	GLN
1	A	1005	GLN
1	A	1011	GLN
1	A	1054	GLN
1	B	49	HIS
1	B	52	GLN
1	B	536	ASN
1	B	751	ASN
1	B	787	GLN
1	B	965	GLN
1	B	1005	GLN
1	B	1083	HIS
1	B	1106	GLN
1	C	317	ASN
1	C	658	ASN
1	C	703	ASN
1	C	895	GLN
1	C	926	GLN

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Mol	Chain	Res	Type
1	C	1023	ASN
1	C	1106	GLN
2	D	79	GLN
3	J	92	HIS
2	F	79	GLN
3	H	92	HIS
2	L	79	GLN
3	P	92	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](#)

3 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	$\# Z  > 2$
6	NAG	B	1401	1	14,14,15	0.49	0	17,19,21	0.75	1 (5%)
6	NAG	C	1401	1	14,14,15	0.50	0	17,19,21	0.75	1 (5%)
6	NAG	A	1401	1	14,14,15	0.48	0	17,19,21	0.74	1 (5%)

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
6	NAG	B	1401	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1401	1	-	2/6/23/26	0/1/1/1
6	NAG	A	1401	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
6	C	1401	NAG	C1-O5-C5	2.31	115.33	112.19
6	A	1401	NAG	C1-O5-C5	2.28	115.28	112.19
6	B	1401	NAG	C1-O5-C5	2.27	115.27	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

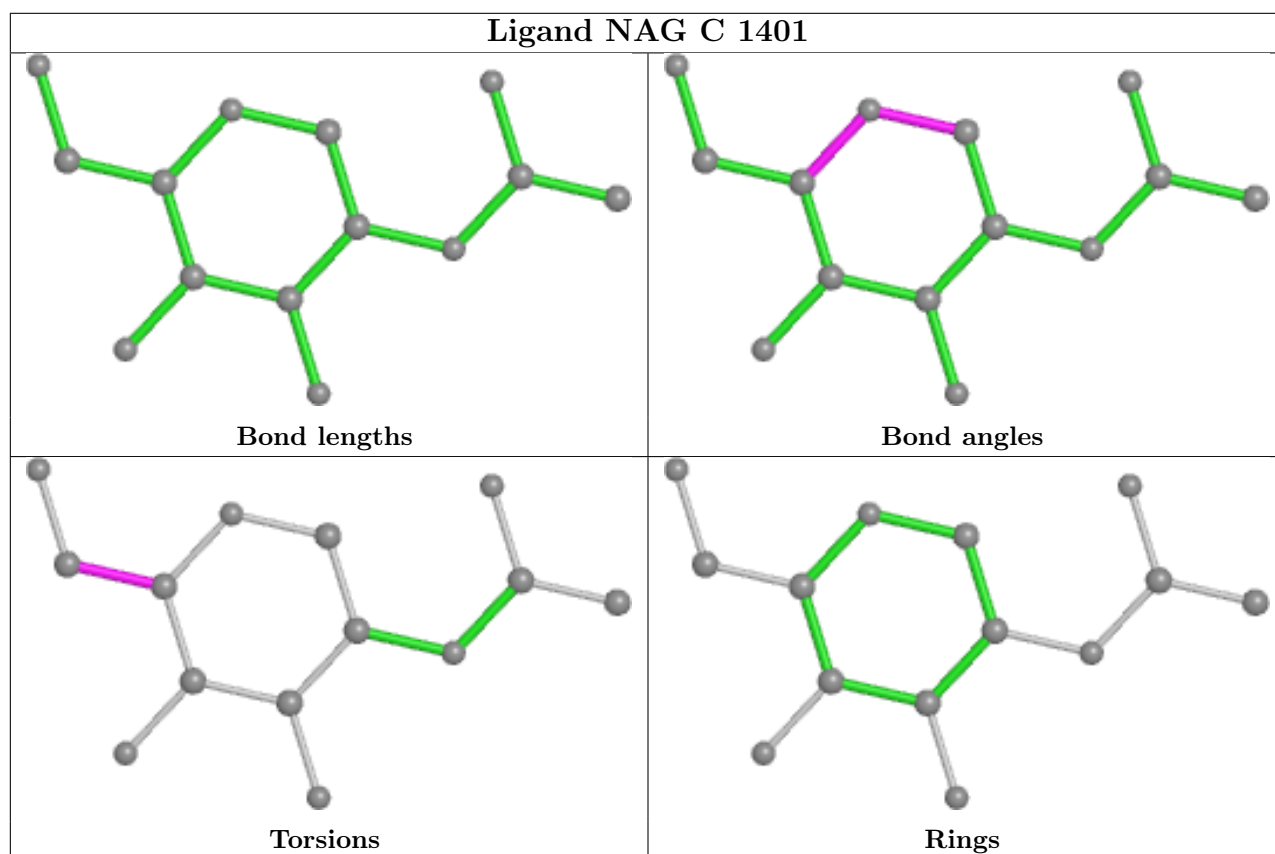
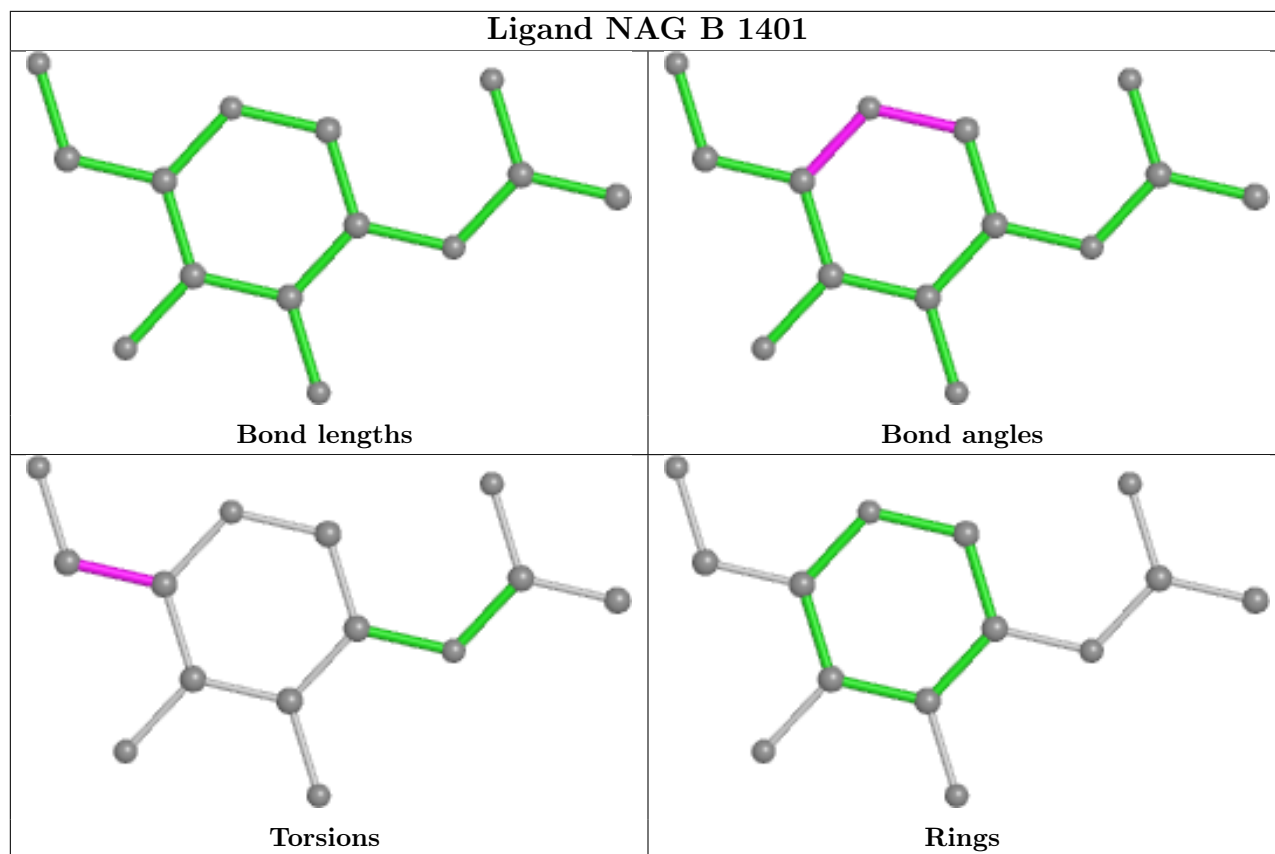
Mol	Chain	Res	Type	Atoms
6	A	1401	NAG	O5-C5-C6-O6
6	B	1401	NAG	O5-C5-C6-O6
6	C	1401	NAG	O5-C5-C6-O6
6	B	1401	NAG	C4-C5-C6-O6
6	C	1401	NAG	C4-C5-C6-O6
6	A	1401	NAG	C4-C5-C6-O6

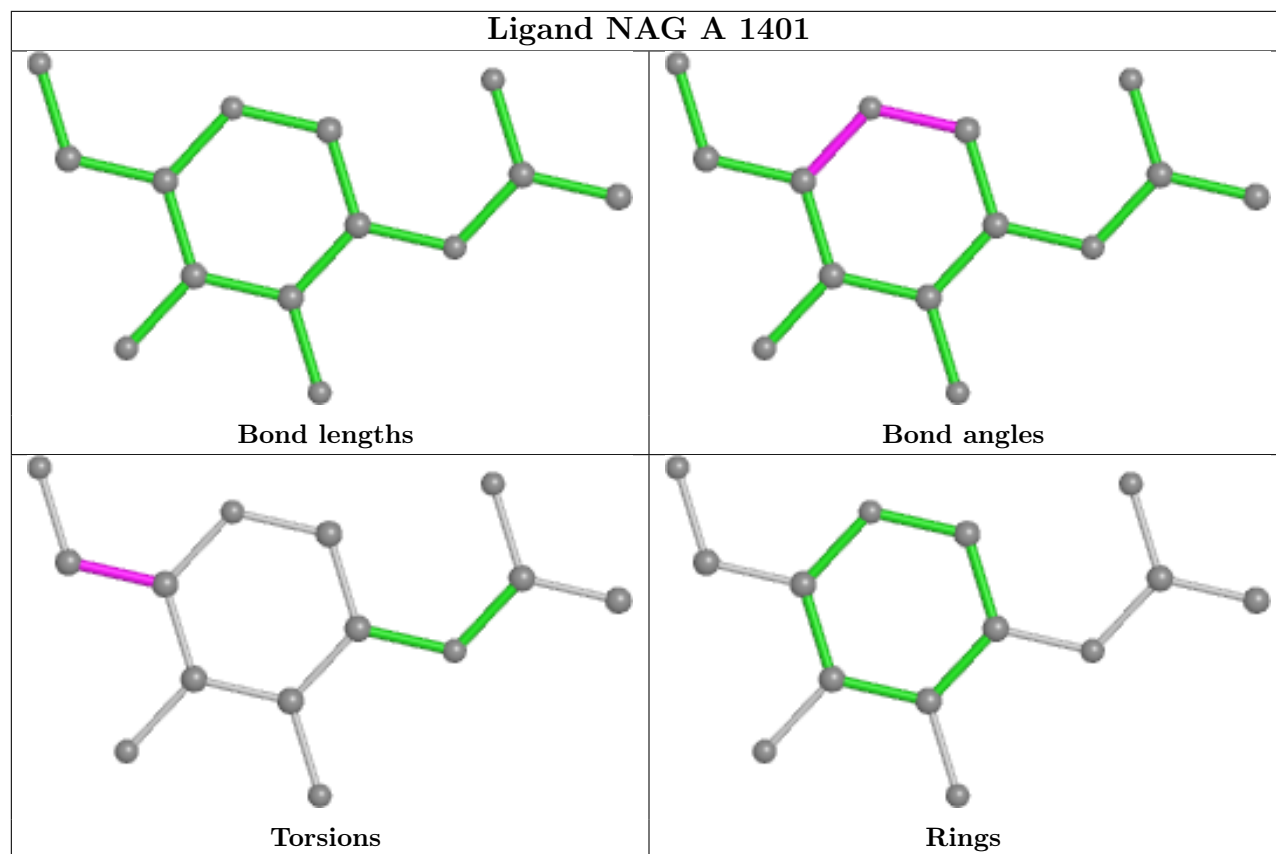
There are no ring outliers.

No monomer is involved in short contacts.

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.

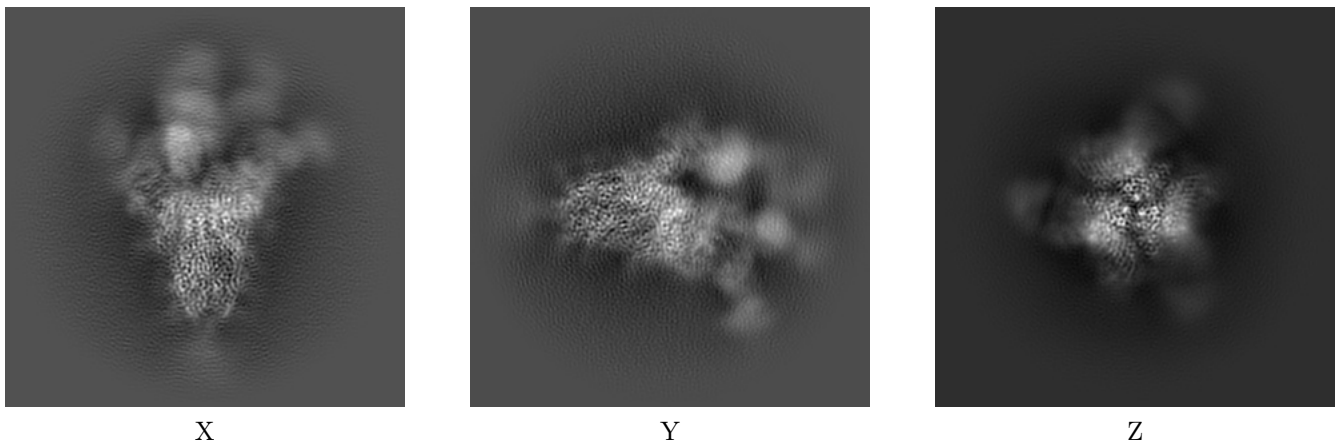
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34411. These allow visual inspection of the internal detail of the map and identification of artifacts.

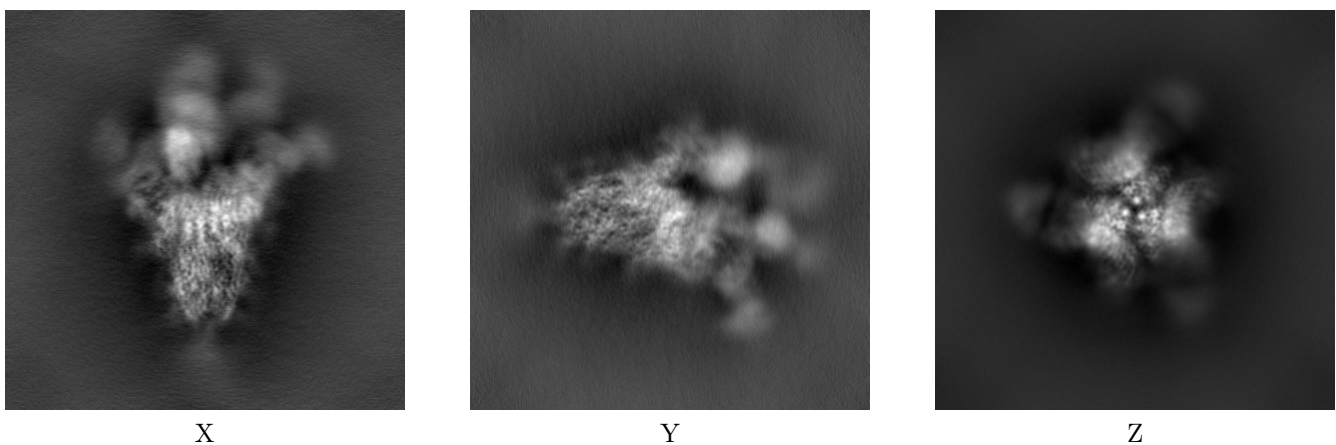
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

#### 6.1.1 Primary map



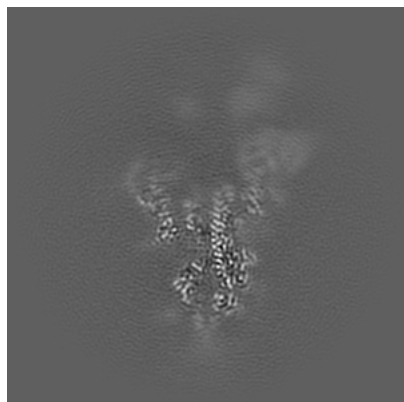
#### 6.1.2 Raw map



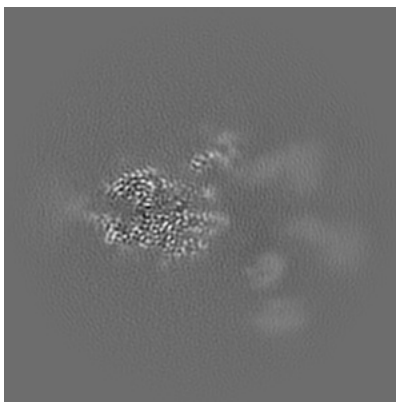
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

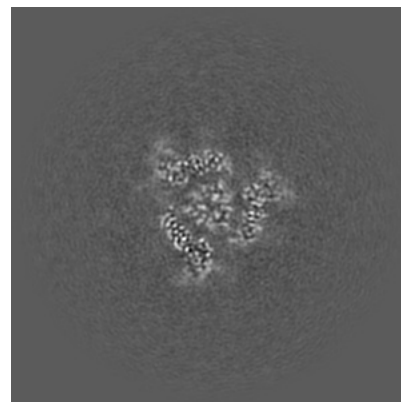
### 6.2.1 Primary map



X Index: 170

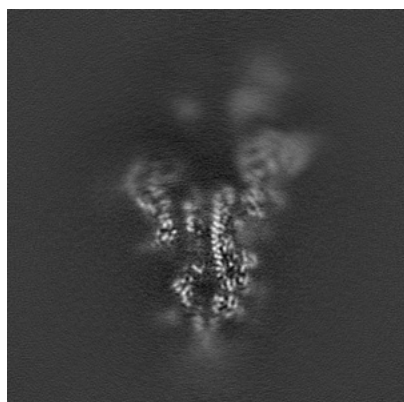


Y Index: 170

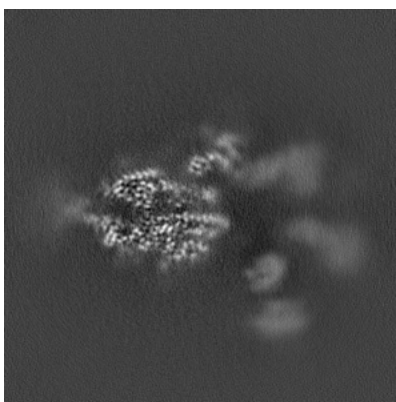


Z Index: 170

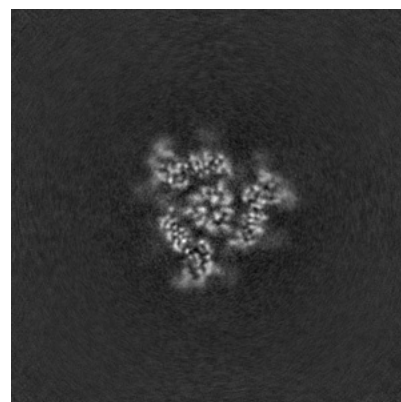
### 6.2.2 Raw map



X Index: 170



Y Index: 170

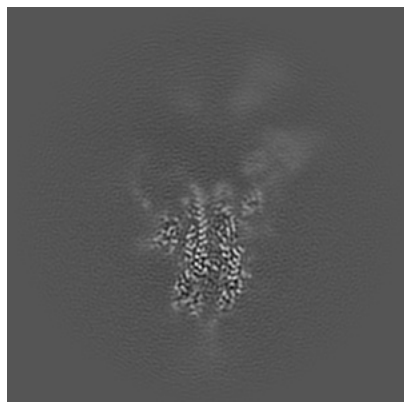


Z Index: 170

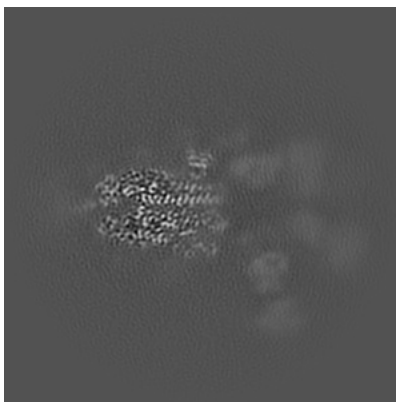
The images above show central slices of the map in three orthogonal directions.

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

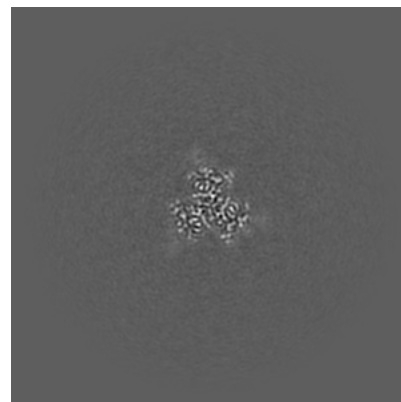
### 6.3.1 Primary map



X Index: 176

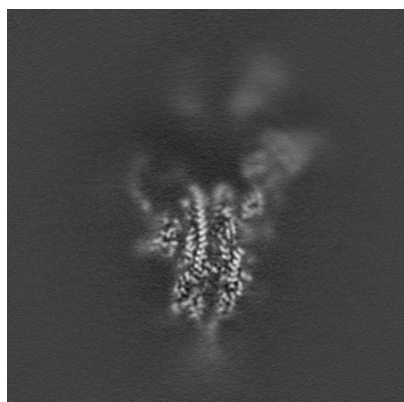


Y Index: 164

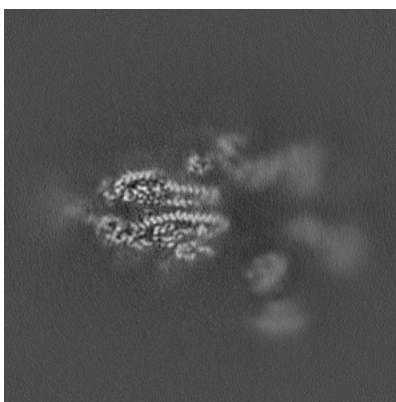


Z Index: 119

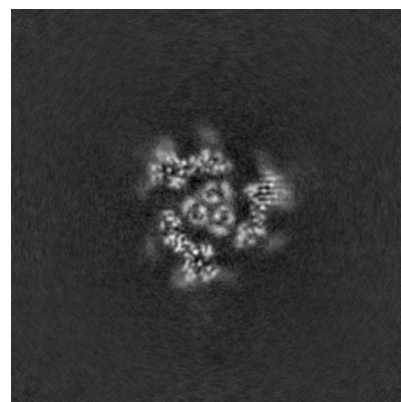
### 6.3.2 Raw map



X Index: 176



Y Index: 168

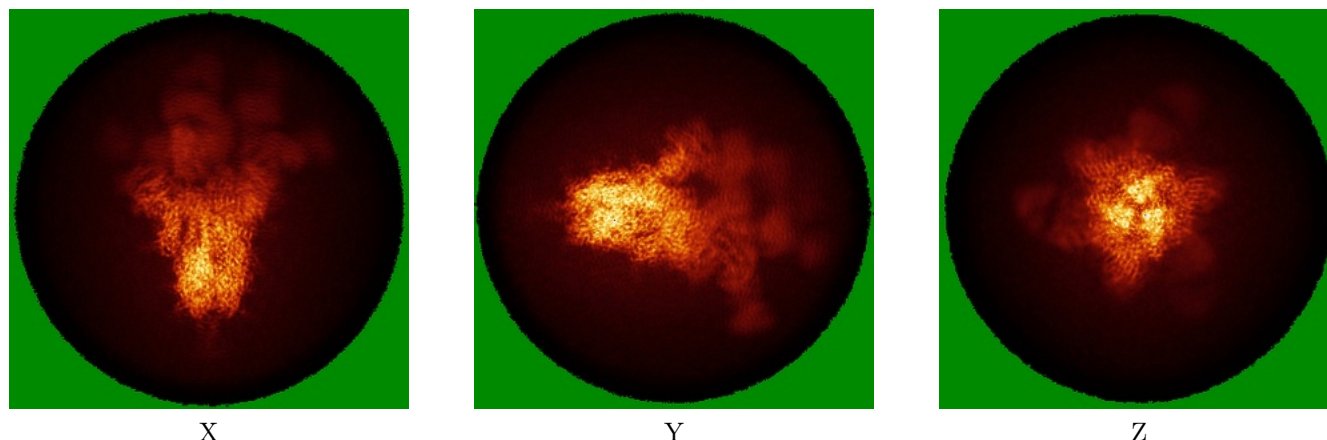


Z Index: 175

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

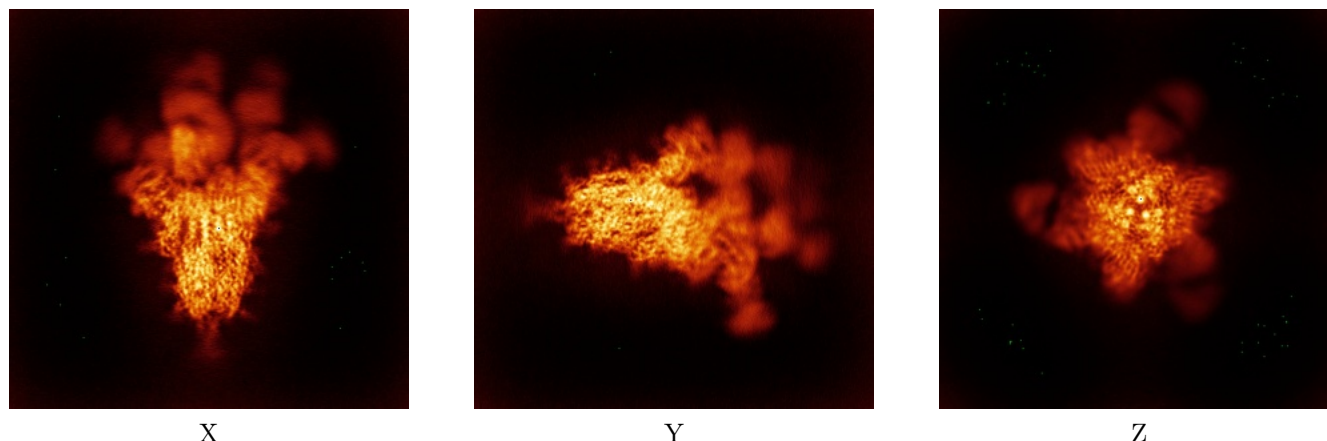


X

Y

Z

### 6.4.2 Raw map



X

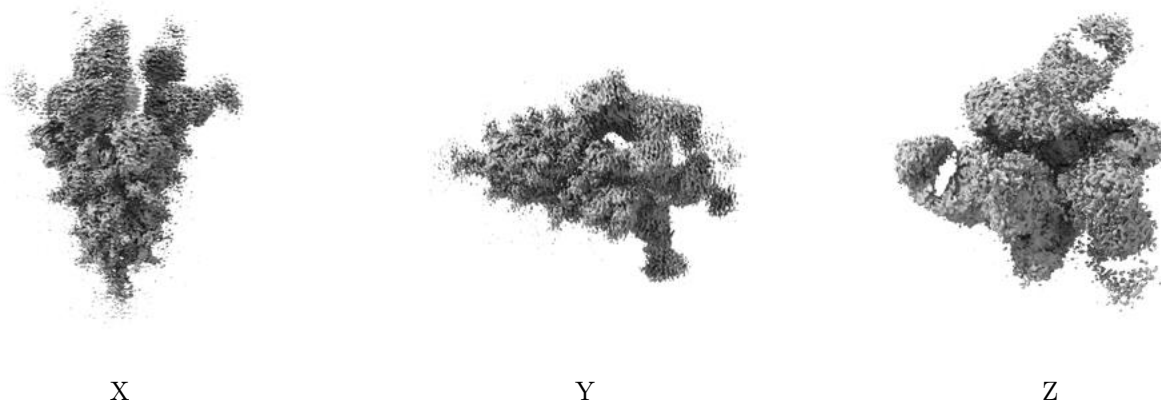
Y

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.

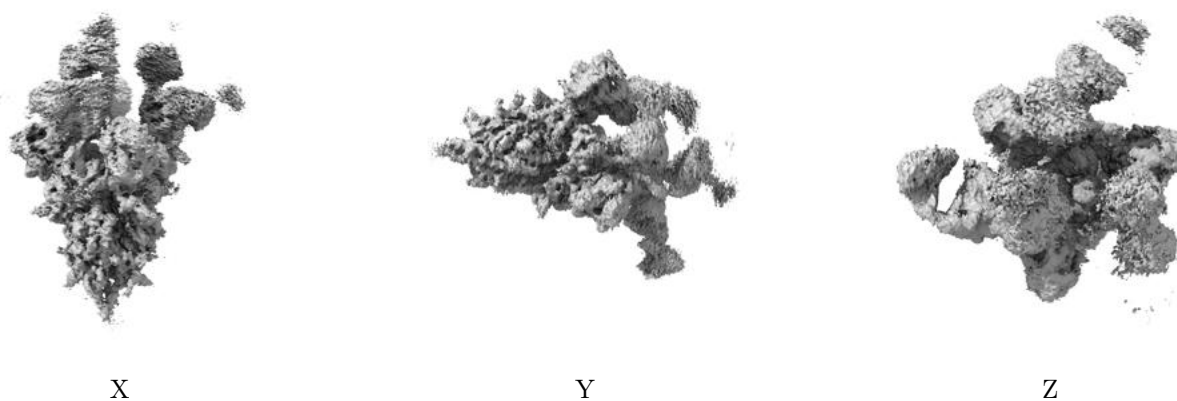
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.15. 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.

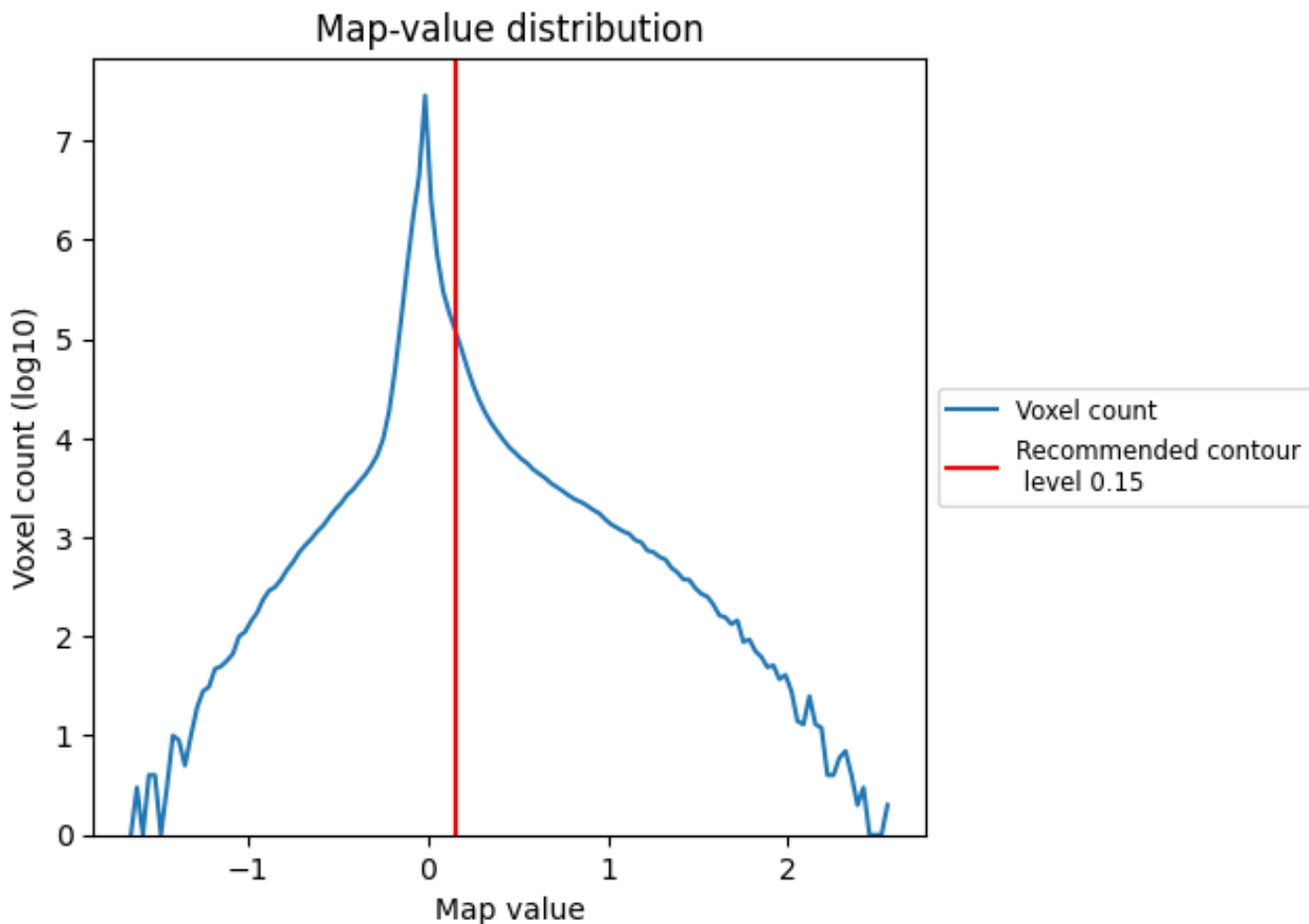
## 6.6 Mask visualisation [i](#)

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.

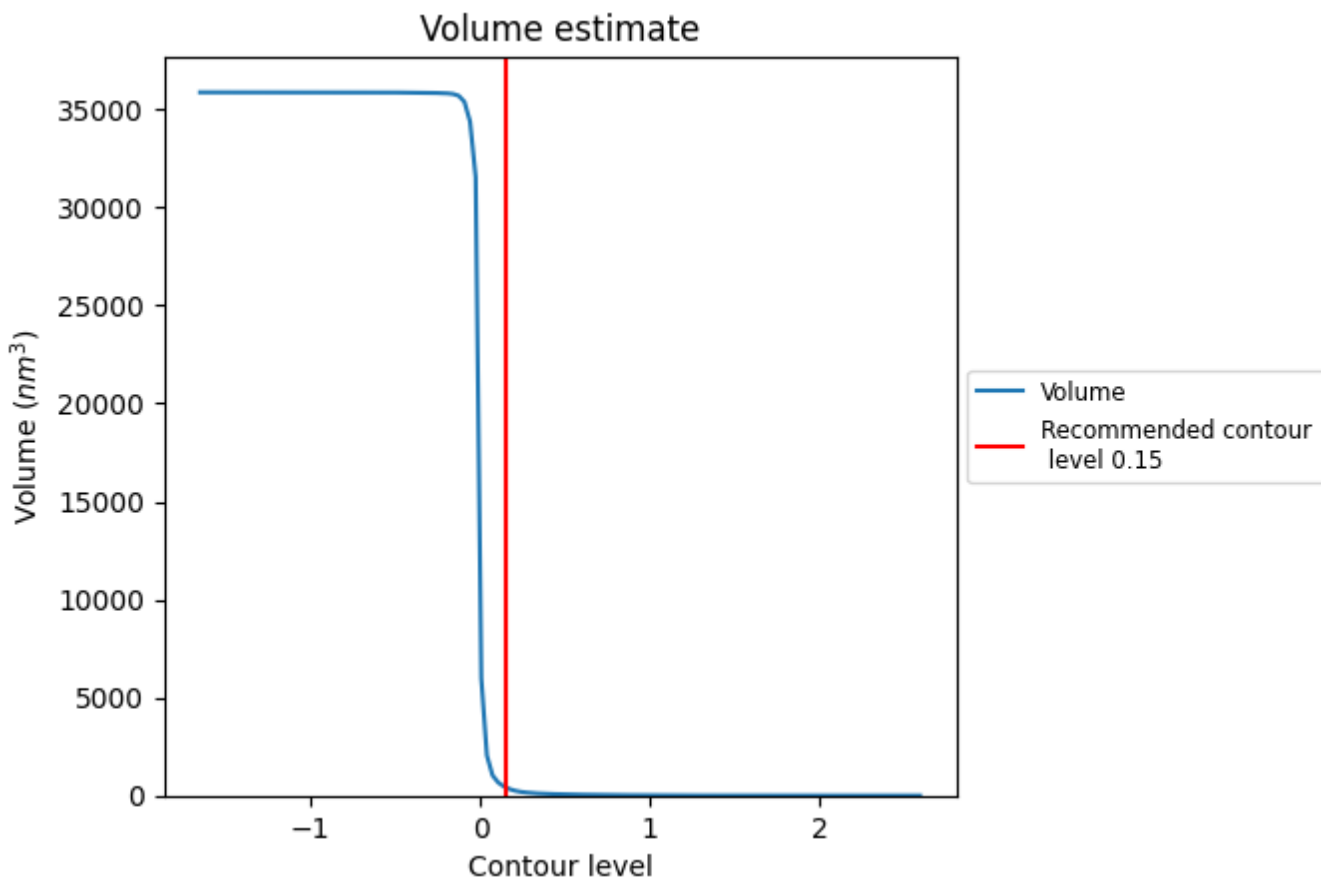
### 7.1 Map-value distribution [i](#)



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.



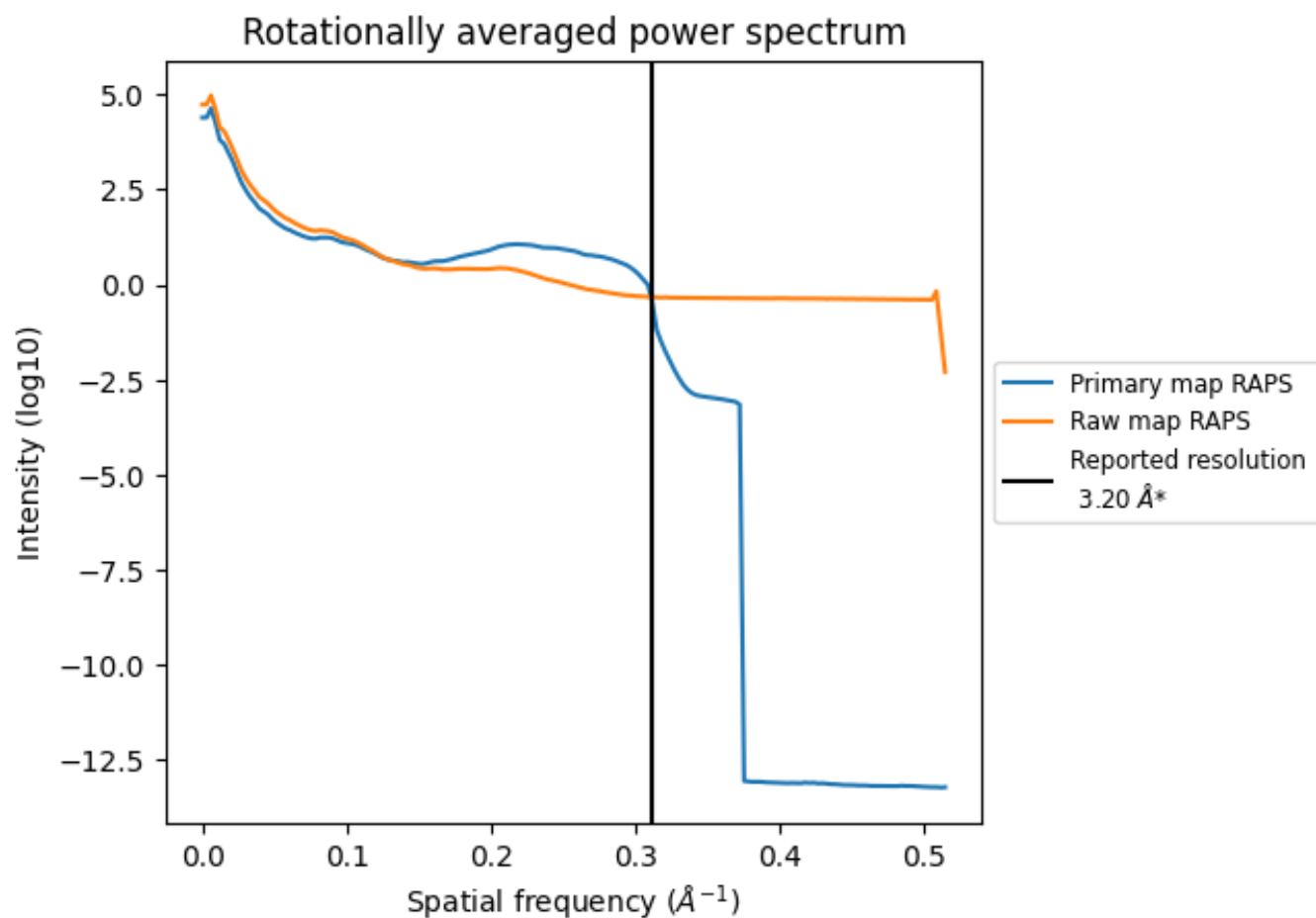
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 419 nm<sup>3</sup>; this corresponds to an approximate mass of 378 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](#)

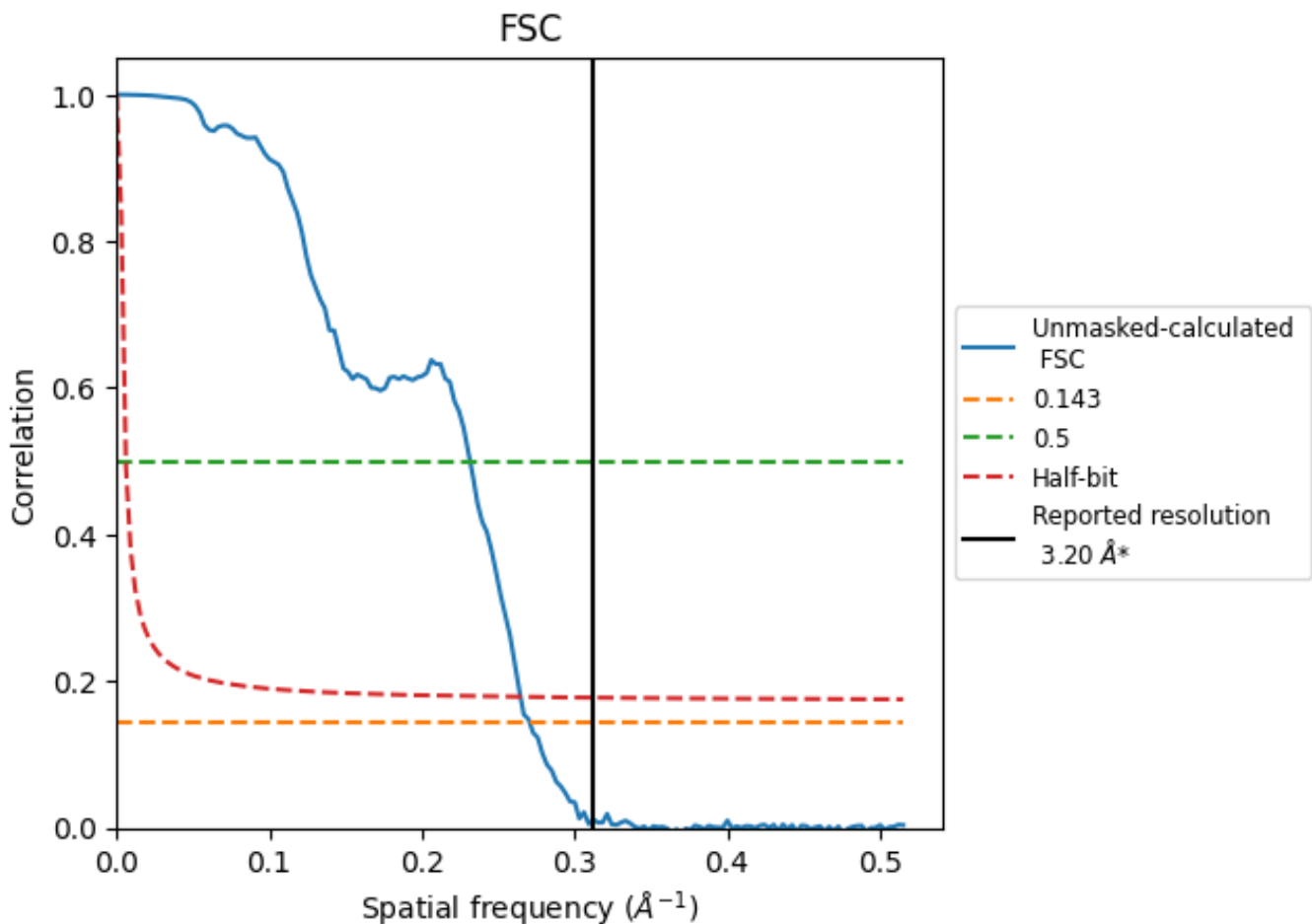


\*Reported resolution corresponds to spatial frequency of 0.312 Å<sup>-1</sup>

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

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

### 8.1 FSC [i](#)



\*Reported resolution corresponds to spatial frequency of 0.312  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

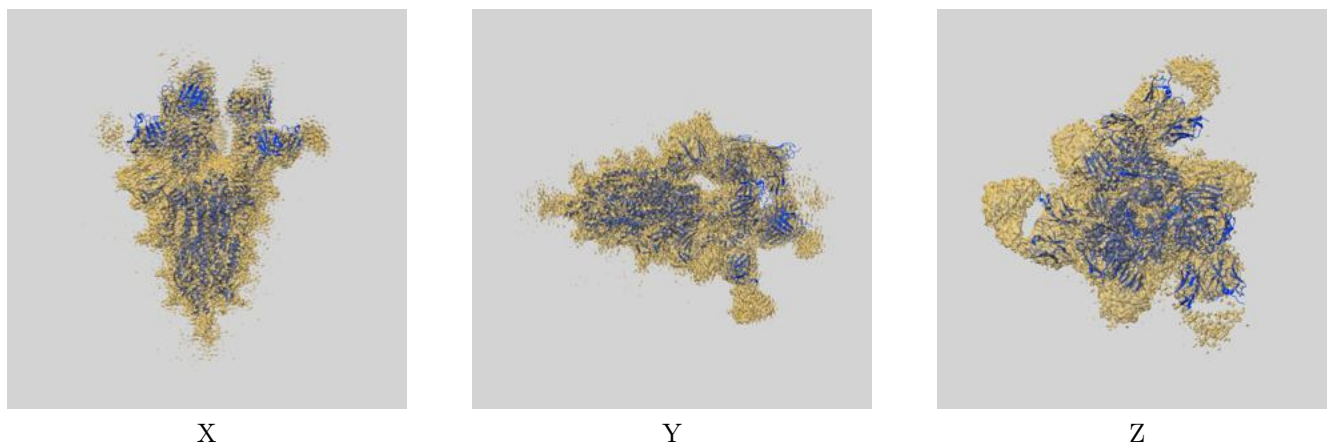
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.69	4.32	3.78

\*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 3.69 differs from the reported value 3.2 by more than 10 %

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

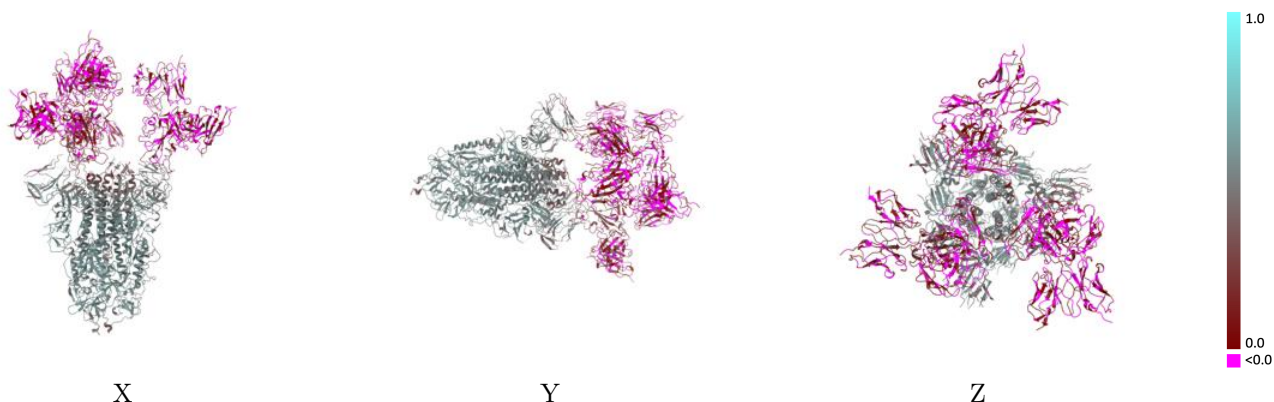
This section contains information regarding the fit between EMDB map EMD-34411 and PDB model 8H08. Per-residue inclusion information can be found in section 3 on page 10.

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



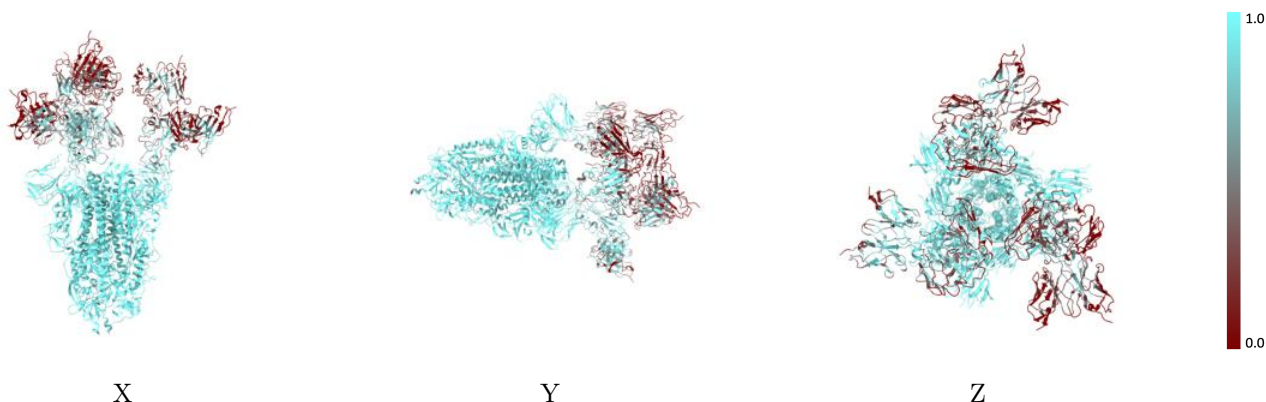
The images above show the 3D surface view of the map at the recommended contour level 0.15 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



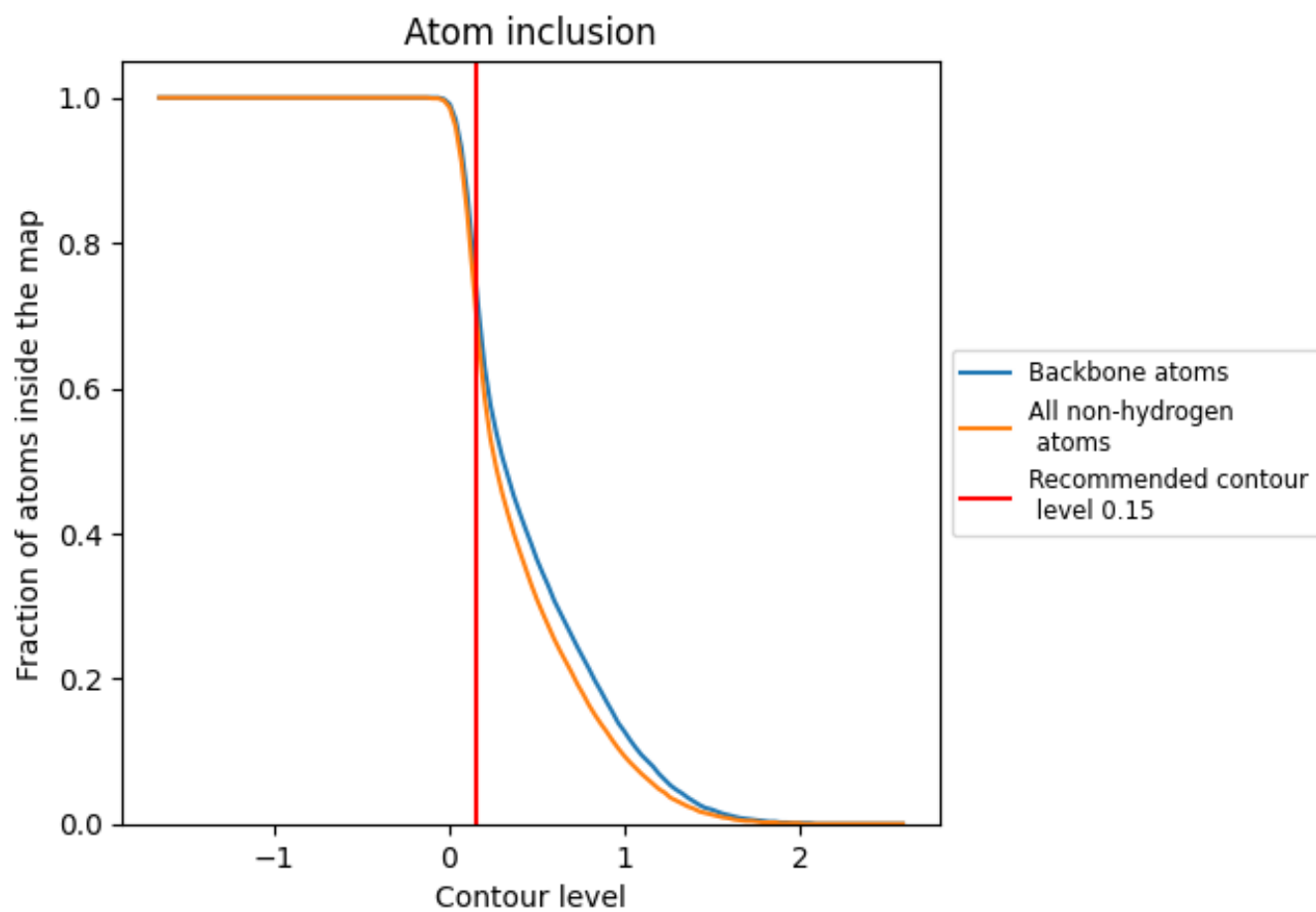
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.15).

































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 72% of all non-hydrogen atoms, are inside the map.

## 9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7190	 0.3000
A	 0.8760	 0.4160
B	 0.9170	 0.4410
C	 0.8730	 0.4210
D	 0.1820	 -0.0010
F	 0.4510	 0.0240
H	 0.3300	 0.0390
I	 0.6750	 0.1020
J	 0.1310	 0.0150
K	 0.5760	 0.0820
L	 0.3310	 0.0100
N	 0.3720	 0.0000
O	 0.1940	 0.0240
P	 0.2500	 0.0290
Q	 0.4400	 0.0180
R	 0.1940	 0.0280

