



## Full wwPDB EM Validation Report ⓘ

Feb 10, 2025 – 12:25 PM JST

PDB ID : 8YBY  
EMDB ID : EMD-34563  
Title : State - I: Spike 2-up RBD with THSC20.HVTR26 (Fab26) - single Fab masked  
Authors : Rencilin, C.F.; Bhattacharya, J.; Dutta, S.  
Deposited on : 2024-02-16  
Resolution : 4.40 Å(reported)

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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.dev113  
MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.40

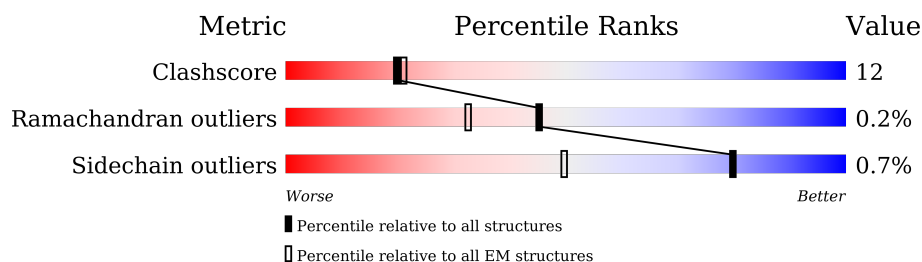
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 4.40 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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

Mol	Chain	Length	Quality of chain
1	B	1273	
1	D	1273	
1	E	1273	
2	A	231	
3	C	216	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 28024 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	B	1033	Total	C	N	O	S	0	0
			8031	5134	1334	1526	37		
1	D	1067	Total	C	N	O	S	0	0
			8350	5327	1395	1589	39		
1	E	1067	Total	C	N	O	S	0	0
			8350	5327	1395	1589	39		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	986	PRO	LYS	conflict	UNP P0DTC2
B	987	PRO	VAL	conflict	UNP P0DTC2
D	986	PRO	LYS	conflict	UNP P0DTC2
D	987	PRO	VAL	conflict	UNP P0DTC2
E	986	PRO	LYS	conflict	UNP P0DTC2
E	987	PRO	VAL	conflict	UNP P0DTC2

- Molecule 2 is a protein called THSC20.HVTR26 (Fab26) - Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	230	Total	C	N	O	S	0	0
			1693	1063	287	335	8		

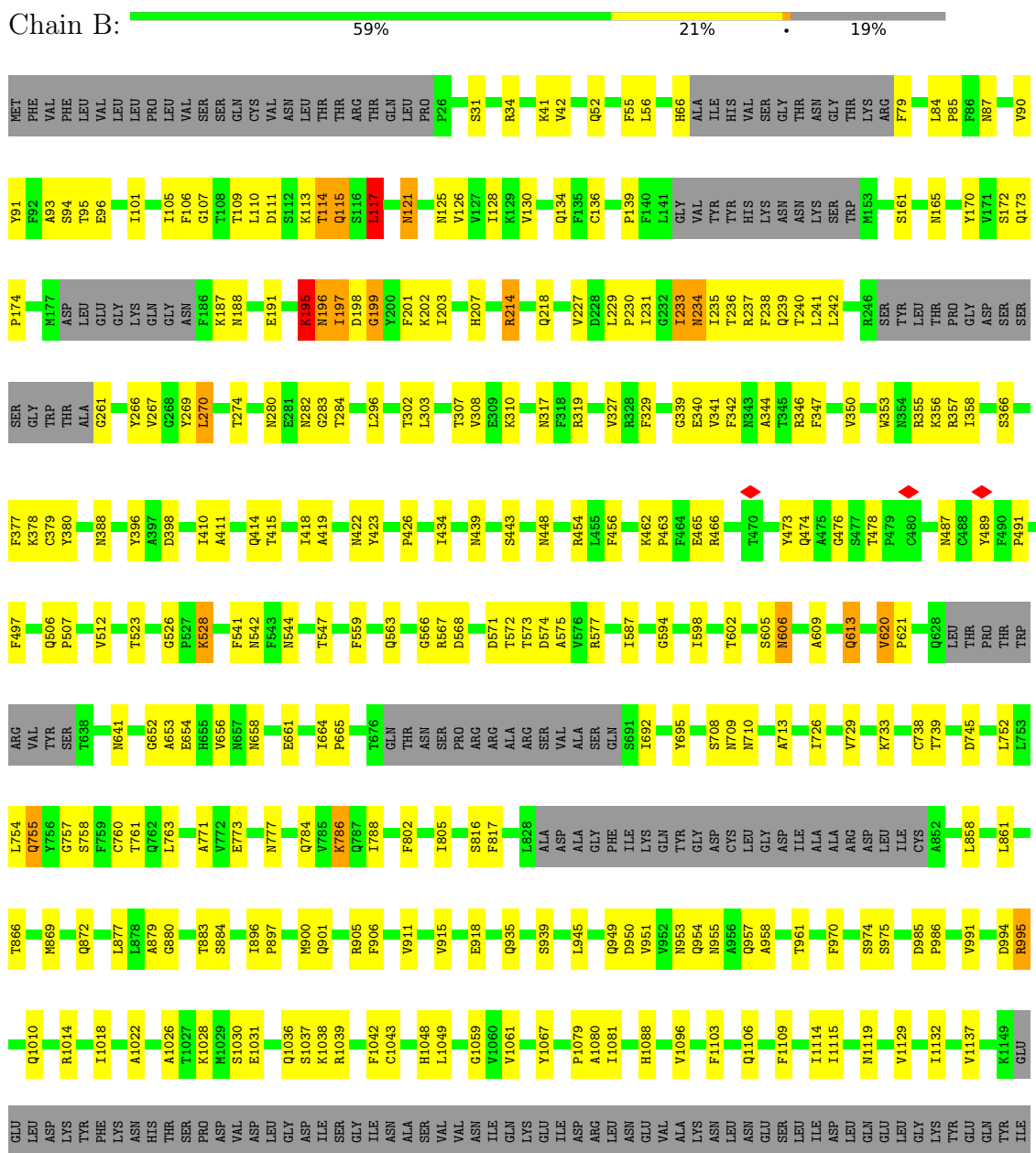
- Molecule 3 is a protein called THSC20.HVTR26 (Fab26) - Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	215	Total	C	N	O	S	0	0
			1600	999	263	331	7		

### 3 Residue-property plots [i](#)

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

#### • Molecule 1: Spike glycoprotein



LYS	TRP	PRO	TRP	TRP	TYR	THR	ILE	TRP	LEU	GLY	ILE	ALA	GLY	LEU	ILE	ALA	VAL	MET	VAL	THR	THR	ILE	MET	LEU	CYS	CYS	MET	THR	ASN	CYS	CYS	SER	CYS	LYS	GLY	CYS	CYS	LYS	PHE	ASP	GLU	ASP	SER	GLU	PRO	VAL	LEU	LYS	GLY	VAL	LYS	LEU
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• Molecule 1: Spike glycoprotein

Chain D: 

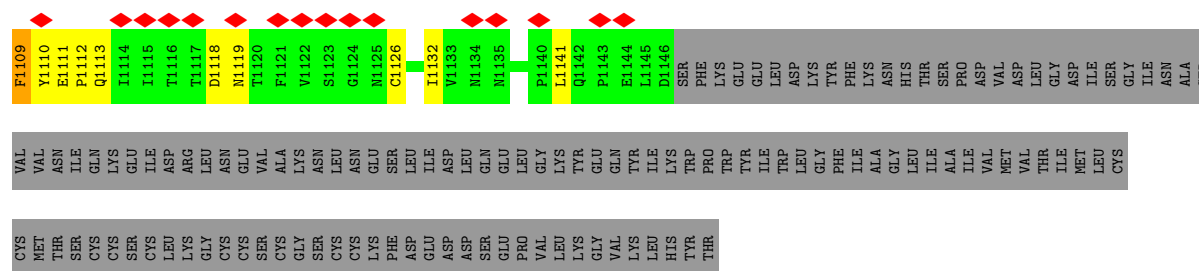
ASP	VAL	ASP	LEU	GLY	ASP	ILE	ASP	LEU	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
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• Molecule 1: Spike glycoprotein

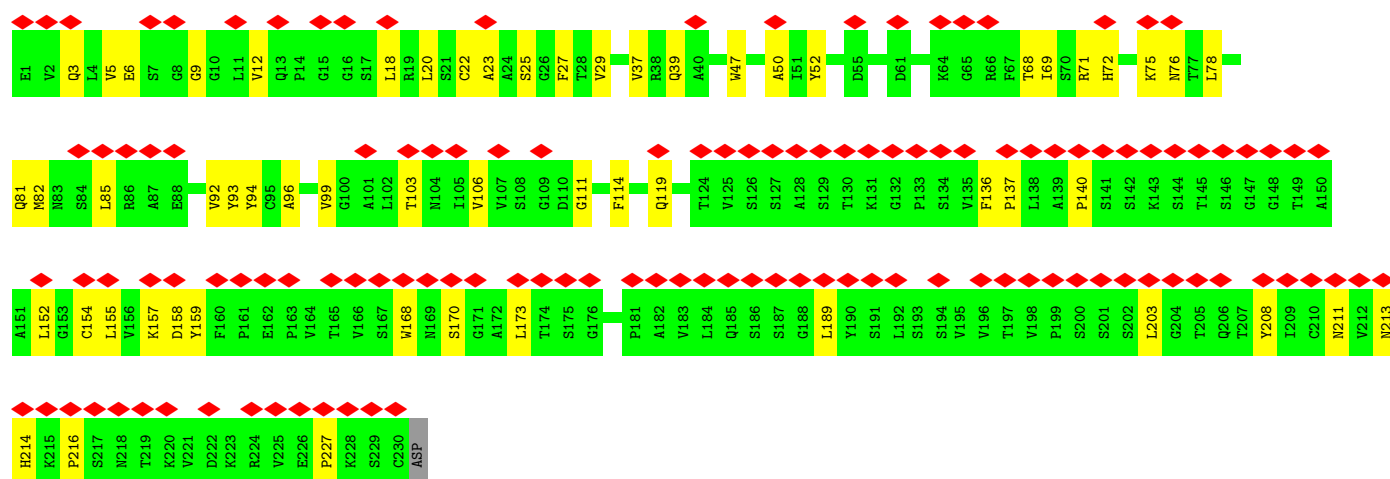
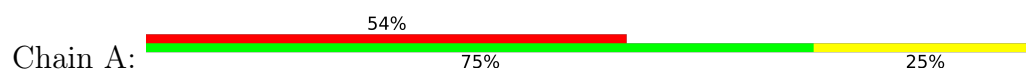
Chain E: 

MET	PHE	VAL	PHE	LEU	VAL	VAL	LEU	LEU	PRO	VAL	SER	Q14	M17	T22	T29	N30	R34	G35	V36	Y37	D40	R44	S45	S46	V47	L48	H49	Q52	D53	L54	F55	L56	P57	F58	V62	H66	A67	I68	H69	V70	SER	GLY	THR	THR	ASN	ASN	GLY	T76	K77	D80
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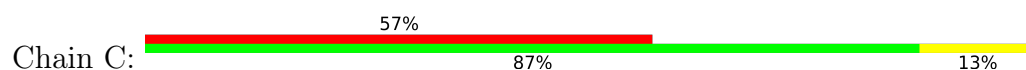




### • Molecule 2: THSC20.HVTR26 (Fab26) - Heavy Chain



### • Molecule 3: THSC20.HVTR26 (Fab26) - Light Chain



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	109596	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2750	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	27.086	Depositor
Minimum map value	-12.980	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	2.82	Depositor
Map size ( $\text{\AA}$ )	331.2, 331.2, 331.2	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.92, 0.92, 0.92	Depositor



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	B	0.47	2/8214 (0.0%)	0.70	13/11178 (0.1%)
1	D	0.51	1/8544 (0.0%)	0.65	5/11626 (0.0%)
1	E	0.46	0/8544	0.84	23/11626 (0.2%)
2	A	0.28	0/1731	0.54	0/2359
3	C	0.27	0/1639	0.51	0/2236
All	All	0.46	3/28672 (0.0%)	0.71	41/39025 (0.1%)

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	B	0	10
1	D	0	6
1	E	0	23
3	C	0	1
All	All	0	40

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	1000	ARG	CA-C	20.94	2.07	1.52
1	B	234	ASN	CG-ND2	10.10	1.58	1.32
1	B	234	ASN	CB-CG	5.36	1.63	1.51

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	425	LEU	CA-CB-CG	16.27	152.72	115.30
1	D	1000	ARG	O-C-N	-15.28	98.26	122.70
1	D	1000	ARG	CB-CA-C	12.79	135.97	110.40
1	D	1000	ARG	N-CA-CB	-11.59	89.73	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	1034	LEU	CA-CB-CG	10.39	139.19	115.30
1	B	117	LEU	CA-CB-CG	10.19	138.73	115.30
1	E	560	LEU	CA-CB-CG	8.76	135.45	115.30
1	E	422	ASN	CB-CA-C	-8.69	93.01	110.40
1	B	197	ILE	CG1-CB-CG2	-8.33	93.07	111.40
1	E	465	GLU	CA-CB-CG	7.62	130.15	113.40
1	E	461	LEU	CA-CB-CG	7.50	132.54	115.30
1	E	906	PHE	CB-CG-CD1	7.43	126.00	120.80
1	E	560	LEU	N-CA-CB	7.41	125.23	110.40
1	E	424	LYS	C-N-CA	7.21	139.74	121.70
1	E	241	LEU	CA-CB-CG	7.17	131.79	115.30
1	D	1000	ARG	CA-C-N	7.17	132.97	117.20
1	E	425	LEU	CB-CA-C	-6.72	97.42	110.20
1	E	900	MET	CA-CB-CG	6.67	124.64	113.30
1	E	117	LEU	CA-CB-CG	6.55	130.37	115.30
1	B	234	ASN	CB-CG-OD1	-6.37	108.86	121.60
1	E	421	TYR	C-N-CA	6.34	137.55	121.70
1	B	117	LEU	CB-CG-CD2	6.30	121.70	111.00
1	E	861	LEU	CA-CB-CG	6.16	129.48	115.30
1	D	984	LEU	C-N-CA	6.12	137.00	121.70
1	E	452	LEU	CA-CB-CG	6.10	129.33	115.30
1	B	195	LYS	CB-CA-C	6.02	122.44	110.40
1	B	199	GLY	N-CA-C	6.00	128.09	113.10
1	E	763	LEU	CA-CB-CG	5.94	128.95	115.30
1	B	559	PHE	C-N-CA	5.83	136.28	121.70
1	E	464	PHE	CB-CA-C	-5.64	99.11	110.40
1	B	197	ILE	C-N-CA	5.64	135.79	121.70
1	B	130	VAL	CG1-CB-CG2	-5.59	101.96	110.90
1	E	822	LEU	CA-CB-CG	-5.30	103.11	115.30
1	B	234	ASN	CB-CG-ND2	5.26	129.32	116.70
1	E	821	LEU	CB-CG-CD2	-5.23	102.10	111.00
1	B	745	ASP	CB-CG-OD1	5.12	122.91	118.30
1	E	296	LEU	CA-CB-CG	5.11	127.06	115.30
1	E	906	PHE	CB-CG-CD2	-5.06	117.26	120.80
1	E	466	ARG	CG-CD-NE	5.03	122.37	111.80
1	B	270	LEU	CA-CB-CG	5.02	126.85	115.30
1	B	571	ASP	CB-CG-OD1	5.00	122.80	118.30

There are no chirality outliers.

All (40) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	114	THR	Peptide
1	B	115	GLN	Peptide
1	B	117	LEU	Peptide
1	B	195	LYS	Mainchain
1	B	528	LYS	Peptide
1	B	613	GLN	Peptide
1	B	754	LEU	Peptide
1	B	755	GLN	Peptide
1	B	786	LYS	Peptide
1	B	945	LEU	Peptide
3	C	52	GLU	Peptide
1	D	1073	LYS	Peptide
1	D	1124	GLY	Peptide
1	D	215	ASP	Peptide
1	D	714	ILE	Peptide
1	D	756	TYR	Peptide
1	D	954	GLN	Peptide
1	E	1004	LEU	Peptide
1	E	1013	ILE	Peptide
1	E	1047	TYR	Peptide
1	E	1059	GLY	Peptide
1	E	1094	VAL	Peptide
1	E	1109	PHE	Peptide
1	E	348	ALA	Peptide
1	E	418	ILE	Peptide
1	E	422	ASN	Peptide
1	E	425	LEU	Mainchain
1	E	451	TYR	Mainchain,Peptide
1	E	463	PRO	Mainchain
1	E	468	ILE	Peptide
1	E	469	SER	Peptide
1	E	528	LYS	Peptide
1	E	599	THR	Peptide
1	E	719	THR	Peptide
1	E	738	CYS	Peptide
1	E	753	LEU	Peptide
1	E	789	TYR	Peptide
1	E	806	LEU	Peptide
1	E	886	TRP	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	B	8031	0	7820	211	0
1	D	8350	0	8140	195	0
1	E	8350	0	8141	351	0
2	A	1693	0	1664	34	0
3	C	1600	0	1544	19	0
All	All	28024	0	27309	677	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 12.

All (677) 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:D:1002:GLN:HG2	1:E:756:TYR:CE1	1.40	1.51
1:D:1000:ARG:CA	1:D:1000:ARG:C	2.07	1.22
1:D:1002:GLN:CG	1:E:756:TYR:CE1	2.31	1.12
1:E:753:LEU:HD12	1:E:756:TYR:CZ	1.85	1.11
1:D:1002:GLN:HG2	1:E:756:TYR:CD1	1.88	1.08
1:E:753:LEU:CD1	1:E:756:TYR:CZ	2.39	1.04
1:D:1002:GLN:HB3	1:E:756:TYR:CD1	1.94	1.02
1:D:1002:GLN:HB3	1:E:756:TYR:CG	1.96	1.00
1:D:1002:GLN:HG2	1:E:756:TYR:HE1	1.31	0.95
1:D:1002:GLN:CG	1:E:756:TYR:CD1	2.48	0.93
1:E:753:LEU:HD11	1:E:756:TYR:CE1	2.03	0.93
1:D:1002:GLN:CB	1:E:756:TYR:CD1	2.55	0.88
1:E:753:LEU:CD1	1:E:756:TYR:OH	2.22	0.88
1:E:753:LEU:CD1	1:E:756:TYR:CE1	2.57	0.87
1:D:1000:ARG:C	1:D:1000:ARG:HA	2.00	0.80
1:B:234:ASN:ND2	1:E:422:ASN:HD22	1.78	0.79
1:E:870:ILE:HA	1:E:873:TYR:HB2	1.67	0.77
1:B:230:PRO:HG3	1:E:398:ASP:HB3	1.68	0.75
1:B:195:LYS:H	1:E:465:GLU:HB2	1.52	0.74
1:E:882:ILE:HA	1:E:901:GLN:HE22	1.52	0.74
1:B:235:ILE:HG12	1:E:467:ASP:HB2	1.69	0.74
1:B:198:ASP:H	1:E:424:LYS:HG2	1.51	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:GLY:HA2	1:D:244:LEU:H	1.53	0.73
1:E:722:VAL:HG11	1:E:1063:LEU:HB3	1.72	0.71
1:E:1016:ALA:HA	1:E:1019:ARG:HB3	1.71	0.70
1:E:308:VAL:HG12	1:E:601:GLY:HA2	1.71	0.70
1:E:753:LEU:HD13	1:E:756:TYR:OH	1.89	0.70
1:B:93:ALA:HB3	1:B:266:TYR:HB2	1.73	0.70
1:B:418:ILE:HA	1:B:422:ASN:HB2	1.74	0.70
1:B:786:LYS:HG3	1:E:699:LEU:HA	1.74	0.70
1:E:443:SER:HB3	1:E:499:PRO:HD3	1.74	0.69
1:E:713:ALA:HA	1:E:1074:ASN:HA	1.74	0.69
1:D:712:ILE:HD12	1:E:901:GLN:HB2	1.74	0.69
1:D:1086:LYS:HA	1:E:1112:PRO:HG2	1.75	0.69
1:E:612:TYR:HB3	1:E:615:VAL:HB	1.75	0.68
1:B:109:THR:HB	1:E:471:GLU:HB2	1.75	0.67
1:D:968:SER:HB2	1:E:755:GLN:HG2	1.77	0.67
1:E:1081:ILE:HB	1:E:1088:HIS:HB2	1.76	0.67
1:D:1079:PRO:HB3	1:E:918:GLU:HG2	1.76	0.66
3:C:4:LEU:HB2	3:C:102:GLY:HA2	1.76	0.66
1:E:756:TYR:CD1	1:E:756:TYR:O	2.48	0.66
1:E:799:GLY:O	1:E:928:ASN:ND2	2.28	0.66
1:D:1009:THR:OG1	1:E:1005:GLN:NE2	2.29	0.66
1:E:396:TYR:HB2	1:E:514:SER:HB3	1.77	0.66
1:E:29:THR:HG23	1:E:62:VAL:HG23	1.77	0.66
1:B:113:LYS:HG2	1:E:472:ILE:HD13	1.77	0.65
1:E:142:GLY:HA3	1:E:156:GLU:HB2	1.76	0.65
1:B:113:LYS:HD3	1:E:492:LEU:HD12	1.76	0.65
1:E:90:VAL:HB	1:E:194:PHE:HB2	1.79	0.65
1:D:93:ALA:HB1	1:D:189:LEU:HD11	1.78	0.65
2:A:22:CYS:HB3	2:A:78:LEU:HB3	1.78	0.65
1:D:516:GLU:HB3	1:D:519:HIS:HD2	1.62	0.65
1:D:189:LEU:HB3	1:D:208:THR:HB	1.79	0.64
1:D:702:GLU:HA	1:E:791:THR:H	1.63	0.64
1:E:742:ILE:HG23	1:E:1000:ARG:HD2	1.77	0.64
1:D:703:ASN:H	1:E:790:LYS:HB3	1.63	0.64
2:A:106:VAL:HG21	2:A:111:GLY:HA2	1.80	0.64
1:B:991:VAL:HG12	1:B:995:ARG:HH21	1.62	0.64
1:E:395:VAL:HG13	1:E:515:PHE:HB3	1.78	0.63
1:E:379:CYS:HA	1:E:432:CYS:HA	1.79	0.63
1:B:188:ASN:HB3	1:B:207:HIS:HB3	1.81	0.63
1:D:1123:SER:N	1:E:1111:GLU:O	2.28	0.63
1:E:875:SER:O	1:E:879:ALA:N	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:949:GLN:OE1	1:E:953:ASN:ND2	2.31	0.63
1:B:113:LYS:N	1:E:470:THR:O	2.32	0.63
1:B:236:THR:HG22	1:E:457:ARG:HG2	1.81	0.62
2:A:158:ASP:HA	2:A:189:LEU:HB3	1.79	0.62
1:D:712:ILE:HA	1:E:897:PRO:HD2	1.81	0.62
1:E:1018:ILE:O	1:E:1022:ALA:N	2.30	0.62
1:B:729:VAL:H	1:B:1059:GLY:HA2	1.64	0.62
1:E:40:ASP:OD1	1:E:44:ARG:NH2	2.33	0.62
1:D:86:PHE:H	1:D:237:ARG:HA	1.64	0.62
1:B:462:LYS:HB2	1:B:465:GLU:HB2	1.82	0.62
1:E:205:SER:HB3	1:E:226:LEU:HD22	1.81	0.62
1:B:113:LYS:H	1:E:470:THR:HB	1.65	0.61
3:C:37:TRP:HB2	3:C:50:ILE:HB	1.81	0.61
1:E:139:PRO:HB2	1:E:241:LEU:HG	1.83	0.61
1:E:712:ILE:HG12	1:E:1107:ARG:HD2	1.80	0.61
1:B:115:GLN:O	1:E:454:ARG:NH2	2.34	0.60
1:D:1002:GLN:CB	1:E:756:TYR:CE1	2.82	0.60
1:E:736:VAL:O	1:E:764:ASN:ND2	2.29	0.60
1:E:711:SER:HA	1:E:1076:THR:HA	1.84	0.60
1:B:199:GLY:HA3	1:E:425:LEU:HB2	1.82	0.60
1:B:227:VAL:HG12	1:B:229:LEU:HD11	1.83	0.60
3:C:39:GLN:HB3	3:C:47:LYS:HB3	1.84	0.60
1:E:115:GLN:HB3	1:E:233:ILE:HD12	1.84	0.60
1:D:958:ALA:H	1:E:765:ARG:HD3	1.66	0.60
1:D:995:ARG:N	1:E:990:GLU:OE2	2.34	0.60
1:B:107:GLY:HA2	1:E:469:SER:H	1.66	0.60
1:D:123:ALA:HA	1:D:177:MET:HB3	1.84	0.60
1:E:1106:GLN:HG3	1:E:1108:ASN:H	1.67	0.59
1:B:117:LEU:H	1:E:468:ILE:HG23	1.65	0.59
1:D:704:SER:OG	1:D:705:VAL:N	2.34	0.59
1:D:1037:SER:OG	1:D:1038:LYS:N	2.35	0.59
1:B:196:ASN:N	1:E:462:LYS:O	2.22	0.59
1:D:598:ILE:HB	1:D:609:ALA:HB3	1.84	0.59
1:E:1030:SER:HB3	1:E:1034:LEU:HG	1.84	0.59
1:B:726:ILE:HG12	1:B:1061:VAL:HG22	1.83	0.59
2:A:136:PHE:HB2	2:A:155:LEU:HB3	1.83	0.59
1:E:762:GLN:OE1	1:E:765:ARG:NH2	2.36	0.59
1:B:957:GLN:OE1	1:D:765:ARG:NH1	2.35	0.59
2:A:3:GLN:HB3	2:A:25:SER:HB2	1.85	0.59
3:C:136:LEU:HB2	3:C:182:LEU:HB3	1.84	0.59
1:E:714:ILE:HG13	1:E:1075:PHE:HB2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1028:LYS:NZ	1:B:1042:PHE:O	2.35	0.59
1:D:434:ILE:HB	1:D:511:VAL:HB	1.84	0.58
1:D:703:ASN:N	1:E:790:LYS:HB3	2.17	0.58
1:D:963:VAL:HA	1:D:966:LEU:HB2	1.85	0.58
1:E:756:TYR:OH	1:E:997:ILE:HG22	2.03	0.58
1:E:923:ILE:O	1:E:927:PHE:N	2.34	0.58
1:D:1010:GLN:HB2	1:E:762:GLN:HG3	1.85	0.58
1:D:1121:PHE:O	1:E:1113:GLN:NE2	2.36	0.58
1:B:355:ARG:HD2	1:B:396:TYR:HB3	1.84	0.58
1:D:356:LYS:HB3	1:D:397:ALA:HB3	1.84	0.58
1:E:898:PHE:O	1:E:902:MET:N	2.35	0.58
1:E:1033:VAL:HG13	1:E:1051:SER:HB2	1.86	0.58
1:B:195:LYS:HB3	1:E:462:LYS:HB2	1.84	0.58
3:C:6:GLN:NE2	3:C:90:CYS:SG	2.75	0.58
1:D:596:SER:HB2	1:D:611:LEU:HB3	1.85	0.58
1:D:97:LYS:NZ	1:D:262:ALA:O	2.37	0.58
1:E:816:SER:O	1:E:820:ASP:N	2.37	0.58
1:E:104:TRP:N	1:E:119:ILE:O	2.36	0.57
1:D:703:ASN:H	1:E:791:THR:H	1.52	0.57
1:B:641:ASN:ND2	1:B:653:ALA:O	2.37	0.57
2:A:27:PHE:O	2:A:76:ASN:ND2	2.37	0.57
1:E:465:GLU:HG2	1:E:466:ARG:H	1.68	0.57
1:E:875:SER:HA	1:E:878:LEU:HB2	1.87	0.57
1:E:925:ASN:O	1:E:929:SER:N	2.36	0.57
1:E:911:VAL:HG13	1:E:1108:ASN:HB3	1.85	0.57
1:B:106:PHE:O	1:E:468:ILE:N	2.38	0.57
1:B:366:SER:H	1:B:388:ASN:HD21	1.53	0.57
1:E:885:GLY:O	1:E:1035:GLY:N	2.37	0.57
1:B:357:ARG:HH12	1:B:523:THR:HG21	1.69	0.57
1:B:456:PHE:HD2	1:B:491:PRO:HA	1.69	0.57
1:D:712:ILE:HG12	1:E:896:ILE:HG12	1.86	0.57
1:E:1021:SER:HA	1:E:1024:LEU:HB3	1.85	0.57
1:B:456:PHE:HB3	1:B:473:TYR:HB2	1.87	0.57
1:D:393:THR:HB	1:D:516:GLU:HB2	1.87	0.57
1:D:476:GLY:H	1:D:487:ASN:HB3	1.69	0.57
1:E:400:PHE:O	1:E:510:VAL:N	2.32	0.57
1:E:870:ILE:HG22	1:E:874:THR:HG23	1.86	0.57
1:B:199:GLY:N	1:E:424:LYS:HA	2.19	0.57
1:E:378:LYS:HB3	1:E:433:VAL:HB	1.86	0.57
1:D:231:ILE:HD12	1:D:233:ILE:HG12	1.87	0.56
1:D:1003:SER:HB3	1:E:759:PHE:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:911:VAL:HG12	1:E:915:VAL:HG21	1.85	0.56
1:D:273:ARG:NH2	1:D:292:ALA:O	2.39	0.56
1:B:915:VAL:HA	1:B:918:GLU:HB2	1.87	0.56
1:D:957:GLN:O	1:D:961:THR:N	2.37	0.56
1:D:994:ASP:HB3	1:E:990:GLU:HG2	1.87	0.56
1:B:231:ILE:HG21	1:E:466:ARG:HG3	1.87	0.56
1:B:901:GLN:O	1:B:905:ARG:N	2.39	0.56
1:E:882:ILE:HD13	1:E:1052:PHE:HE1	1.70	0.56
1:E:1079:PRO:HB2	1:E:1132:ILE:HG13	1.87	0.56
1:B:199:GLY:HA3	1:E:425:LEU:HD12	1.88	0.56
1:B:378:LYS:NZ	1:B:379:CYS:O	2.38	0.56
1:E:970:PHE:HD2	1:E:1000:ARG:HH21	1.54	0.56
1:B:954:GLN:O	1:B:957:GLN:NE2	2.38	0.56
2:A:29:VAL:O	2:A:71:ARG:NH1	2.39	0.56
1:E:676:THR:OG1	1:E:690:GLN:NE2	2.39	0.56
1:D:106:PHE:HB2	1:D:117:LEU:HB3	1.88	0.56
1:B:340:GLU:OE2	1:B:356:LYS:NZ	2.38	0.55
1:B:353:TRP:O	1:B:466:ARG:NH1	2.39	0.55
1:D:995:ARG:HG3	1:E:990:GLU:HG3	1.87	0.55
1:E:1012:LEU:O	1:E:1016:ALA:N	2.28	0.55
2:A:52:TYR:O	2:A:71:ARG:NH2	2.39	0.55
1:D:1086:LYS:HG2	1:E:1112:PRO:HB2	1.89	0.55
1:D:950:ASP:O	1:D:954:GLN:N	2.39	0.55
1:D:995:ARG:HA	1:D:998:THR:HG22	1.88	0.55
1:E:970:PHE:HE2	1:E:1000:ARG:HE	1.53	0.55
1:E:1044:GLY:HA3	1:E:1066:THR:HG21	1.88	0.55
1:B:426:PRO:HB3	1:B:463:PRO:HG3	1.89	0.55
1:B:90:VAL:O	1:E:462:LYS:NZ	2.39	0.55
1:D:994:ASP:O	1:D:998:THR:N	2.38	0.55
1:E:1096:VAL:N	1:E:1103:PHE:O	2.40	0.55
1:B:563:GLN:O	1:B:577:ARG:NH2	2.39	0.55
1:B:1081:ILE:HB	1:B:1088:HIS:HB2	1.89	0.55
1:D:756:TYR:OH	1:E:994:ASP:OD2	2.24	0.55
1:D:903:ALA:HB1	1:D:913:GLN:HG2	1.88	0.55
1:B:346:ARG:NH1	1:B:347:PHE:O	2.39	0.55
1:E:784:GLN:HG3	1:E:1034:LEU:HD21	1.89	0.55
1:B:418:ILE:O	1:B:423:TYR:N	2.37	0.55
1:B:733:LYS:HE3	1:B:771:ALA:HB1	1.88	0.55
1:D:1117:THR:O	1:E:1119:ASN:ND2	2.39	0.54
1:D:716:THR:N	1:D:1071:GLN:O	2.38	0.54
1:D:605:SER:OG	1:D:606:ASN:N	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1089:PHE:HB3	1:E:912:THR:HG21	1.88	0.54
1:E:69:HIS:CD2	1:E:77:LYS:H	2.25	0.54
1:B:96:GLU:HB2	1:B:187:LYS:HB3	1.90	0.54
1:D:816:SER:OG	1:D:1054:GLN:NE2	2.41	0.54
1:E:778:THR:HB	1:E:865:LEU:HD12	1.90	0.54
1:B:1037:SER:OG	1:B:1038:LYS:N	2.41	0.54
1:B:1039:ARG:NE	1:D:1031:GLU:OE2	2.36	0.54
1:E:596:SER:OG	1:E:613:GLN:NE2	2.41	0.54
1:E:978:ASN:HA	1:E:981:LEU:HD12	1.90	0.54
1:B:214:ARG:NH1	1:B:266:TYR:OH	2.41	0.54
1:E:280:ASN:OD1	1:E:284:THR:N	2.41	0.54
1:E:748:GLU:HB3	1:E:981:LEU:HD22	1.90	0.54
1:E:1048:HIS:HD2	1:E:1066:THR:HG22	1.72	0.54
1:E:454:ARG:NH2	1:E:467:ASP:O	2.41	0.54
1:E:881:THR:O	1:E:905:ARG:NH2	2.41	0.54
1:D:773:GLU:O	1:D:777:ASN:ND2	2.41	0.53
1:B:90:VAL:HG13	1:B:267:VAL:HG13	1.91	0.53
1:E:34:ARG:HH12	1:E:221:SER:HB3	1.72	0.53
1:E:98:SER:OG	1:E:182:LYS:O	2.26	0.53
1:E:141:LEU:N	1:E:242:LEU:O	2.41	0.53
1:E:393:THR:N	1:E:516:GLU:O	2.38	0.53
1:B:620:VAL:HG23	1:B:621:PRO:HD3	1.89	0.53
1:B:203:ILE:HB	1:B:227:VAL:HB	1.90	0.53
1:B:661:GLU:O	1:B:695:TYR:OH	2.26	0.53
1:D:437:ASN:ND2	1:D:506:GLN:OE1	2.41	0.53
1:E:949:GLN:O	1:E:953:ASN:ND2	2.41	0.53
1:B:236:THR:OG1	1:B:237:ARG:N	2.40	0.53
1:D:116:SER:N	1:D:131:CYS:O	2.34	0.53
1:D:1086:LYS:HB3	1:E:1113:GLN:HA	1.91	0.53
1:E:914:ASN:HB3	1:E:917:TYR:HD2	1.73	0.53
1:D:379:CYS:HB2	1:D:384:PRO:HB3	1.91	0.53
1:E:1081:ILE:O	1:E:1088:HIS:N	2.38	0.53
1:E:99:ASN:HD22	1:E:180:GLU:H	1.55	0.53
1:E:800:PHE:HA	1:E:928:ASN:HD21	1.72	0.53
1:B:1043:CYS:HB2	1:B:1048:HIS:HD2	1.73	0.53
1:D:654:GLU:HB3	1:D:693:ILE:HG22	1.90	0.53
1:D:1090:PRO:HD2	1:E:912:THR:HB	1.90	0.53
1:E:347:PHE:H	1:E:509:ARG:HH12	1.56	0.53
1:B:949:GLN:HE21	1:B:953:ASN:HD21	1.57	0.53
1:D:64:TRP:HE1	1:D:264:ALA:HB1	1.74	0.53
2:A:140:PRO:HD2	2:A:227:PRO:HA	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:128:ILE:HB	1:E:170:TYR:HB3	1.90	0.52
1:D:1087:ALA:HB3	1:E:1112:PRO:HD2	1.90	0.52
1:B:280:ASN:OD1	1:B:284:THR:N	2.42	0.52
1:E:110:LEU:HD23	1:E:116:SER:HB3	1.91	0.52
1:D:1003:SER:O	1:D:1007:TYR:N	2.37	0.52
1:D:1145:LEU:HB2	1:E:1141:LEU:HD23	1.90	0.52
1:E:594:GLY:H	1:E:614:ASP:HB2	1.74	0.52
1:E:153:MET:SD	1:E:155:SER:OG	2.68	0.52
2:A:68:THR:HB	2:A:81:GLN:HB3	1.91	0.52
1:E:457:ARG:HH11	1:E:458:LYS:H	1.57	0.52
1:D:905:ARG:NH1	1:D:1049:LEU:O	2.43	0.52
1:E:103:GLY:HA2	1:E:120:VAL:HA	1.92	0.52
2:A:155:LEU:HG	2:A:157:LYS:HB2	1.91	0.52
1:E:141:LEU:HB2	1:E:243:ALA:HA	1.91	0.52
1:B:234:ASN:ND2	1:E:422:ASN:H	2.08	0.52
1:B:879:ALA:O	1:B:883:THR:N	2.40	0.52
1:B:1010:GLN:OE1	1:B:1014:ARG:NH2	2.43	0.52
1:E:22:THR:OG1	1:E:77:LYS:O	2.28	0.52
1:E:273:ARG:NH2	1:E:290:ASP:OD2	2.43	0.52
1:E:525:CYS:SG	1:E:544:ASN:ND2	2.81	0.52
1:B:85:PRO:HG2	1:B:87:ASN:HD21	1.75	0.51
1:B:114:THR:H	1:E:470:THR:HG22	1.75	0.51
1:E:1018:ILE:HA	1:E:1021:SER:HB2	1.90	0.51
1:B:31:SER:HB2	1:B:56:LEU:HD21	1.92	0.51
1:B:126:VAL:HG13	1:B:174:PRO:HA	1.90	0.51
1:B:308:VAL:N	1:B:602:THR:OG1	2.40	0.51
1:D:705:VAL:HG13	1:E:897:PRO:HA	1.92	0.51
1:E:888:PHE:HZ	1:E:1034:LEU:HD22	1.75	0.51
1:B:422:ASN:OD1	1:B:454:ARG:N	2.40	0.51
2:A:72:HIS:HB3	2:A:75:LYS:HB2	1.92	0.51
1:E:37:TYR:OH	1:E:54:LEU:O	2.27	0.51
1:B:773:GLU:O	1:B:777:ASN:N	2.40	0.51
1:D:457:ARG:NH1	1:D:459:SER:OG	2.43	0.51
1:D:784:GLN:NE2	1:D:1030:SER:OG	2.43	0.51
1:D:1081:ILE:HG12	1:D:1135:ASN:HB3	1.92	0.51
1:E:81:ASN:OD1	1:E:81:ASN:N	2.44	0.51
1:E:756:TYR:OH	1:E:997:ILE:O	2.10	0.51
1:B:866:THR:H	1:B:869:MET:HE2	1.74	0.51
3:C:120:THR:HA	3:C:208:LYS:HG3	1.92	0.51
1:D:1087:ALA:H	1:E:1112:PRO:HG2	1.76	0.51
1:E:1021:SER:O	1:E:1025:ALA:N	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:SER:OG	1:B:173:GLN:N	2.44	0.50
1:B:439:ASN:ND2	1:B:506:GLN:OE1	2.44	0.50
1:B:566:GLY:N	1:B:575:ALA:O	2.44	0.50
1:E:654:GLU:HB3	1:E:693:ILE:HG22	1.92	0.50
1:E:819:GLU:OE2	1:E:1055:SER:N	2.33	0.50
1:B:541:PHE:O	1:B:547:THR:OG1	2.28	0.50
1:B:1106:GLN:HE21	1:B:1109:PHE:HB3	1.77	0.50
1:D:329:PHE:O	1:D:580:GLN:NE2	2.44	0.50
1:D:1002:GLN:NE2	1:E:997:ILE:O	2.44	0.50
1:E:36:VAL:HG22	1:E:288:ALA:HB3	1.92	0.50
1:D:316:SER:OG	1:D:317:ASN:N	2.43	0.50
1:E:80:ASP:O	1:E:265:TYR:OH	2.29	0.50
1:E:301:CYS:O	1:E:304:LYS:NZ	2.33	0.50
1:E:1079:PRO:HG2	1:E:1132:ILE:HA	1.93	0.50
1:E:172:SER:OG	1:E:173:GLN:N	2.44	0.50
1:B:128:ILE:HB	1:B:170:TYR:HB3	1.94	0.50
1:B:317:ASN:ND2	1:D:737:ASP:OD2	2.45	0.50
1:B:101:ILE:HA	1:B:242:LEU:HG	1.93	0.50
1:B:235:ILE:HB	1:E:461:LEU:HD11	1.93	0.50
1:B:598:ILE:HB	1:B:609:ALA:HB3	1.94	0.50
1:E:900:MET:O	1:E:904:TYR:N	2.42	0.50
1:B:234:ASN:HD22	1:E:422:ASN:HD22	1.59	0.50
1:D:612:TYR:O	1:D:649:CYS:N	2.42	0.50
1:D:998:THR:HG1	1:E:998:THR:HG1	1.60	0.50
1:D:1104:VAL:HB	1:D:1115:ILE:HG13	1.93	0.50
1:E:1030:SER:O	1:E:1034:LEU:N	2.43	0.50
1:B:905:ARG:O	1:B:1036:GLN:NE2	2.43	0.50
1:E:97:LYS:HB3	1:E:186:PHE:HA	1.94	0.50
1:E:980:ILE:HG23	1:E:984:LEU:HD12	1.93	0.50
1:B:641:ASN:HD21	1:B:654:GLU:HB3	1.76	0.49
1:E:29:THR:OG1	1:E:30:ASN:N	2.44	0.49
1:D:1033:VAL:HA	1:D:1051:SER:HB2	1.94	0.49
1:D:122:ASN:HB3	1:D:125:ASN:H	1.77	0.49
1:B:656:VAL:HG12	1:B:658:ASN:H	1.76	0.49
1:E:17:ASN:O	1:E:255:SER:OG	2.28	0.49
1:E:788:ILE:HG23	1:E:873:TYR:HE1	1.77	0.49
1:B:52:GLN:OE1	1:B:274:THR:OG1	2.23	0.49
1:B:105:ILE:HA	1:E:468:ILE:HD11	1.94	0.49
1:E:928:ASN:O	1:E:932:GLY:N	2.45	0.49
1:B:473:TYR:HB3	1:B:489:TYR:HB2	1.94	0.49
2:A:99:VAL:HG22	2:A:103:THR:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:LEU:HD12	1:D:217:PRO:HD2	1.94	0.49
1:D:1094:VAL:HG21	1:E:904:TYR:HE1	1.78	0.49
1:E:318:PHE:HD2	1:E:593:GLY:HA3	1.77	0.49
1:B:802:PHE:HD1	1:B:805:ILE:HD11	1.78	0.49
1:D:31:SER:HG	1:D:60:SER:H	1.55	0.49
1:D:236:THR:HG22	1:D:237:ARG:HG3	1.94	0.49
1:E:777:ASN:OD1	1:E:1019:ARG:NH1	2.38	0.49
2:A:20:LEU:HD11	2:A:93:TYR:HD2	1.78	0.49
1:D:949:GLN:HE21	1:D:953:ASN:HD21	1.61	0.49
1:E:126:VAL:HB	1:E:172:SER:HB3	1.94	0.49
1:E:216:LEU:HD12	1:E:217:PRO:HD2	1.95	0.49
1:E:740:MET:SD	1:E:856:ASN:ND2	2.85	0.49
1:B:788:ILE:HD11	1:B:872:GLN:HG3	1.95	0.49
1:D:1107:ARG:HH22	1:E:1036:GLN:HB3	1.78	0.49
1:B:411:ALA:HB3	1:B:414:GLN:HB3	1.95	0.49
1:E:550:GLY:HA2	1:E:590:CYS:H	1.77	0.49
1:B:1026:ALA:O	1:B:1030:SER:N	2.45	0.48
1:E:277:LEU:HD12	1:E:285:ILE:HG21	1.95	0.48
1:E:727:LEU:HD11	1:E:1028:LYS:HB2	1.95	0.48
1:E:991:VAL:HA	1:E:994:ASP:HB2	1.93	0.48
1:B:42:VAL:O	1:E:519:HIS:NE2	2.44	0.48
1:E:753:LEU:HD11	1:E:756:TYR:HE1	1.72	0.48
1:D:661:GLU:O	1:D:695:TYR:OH	2.30	0.48
1:D:1123:SER:HB3	1:E:1110:TYR:HA	1.94	0.48
1:B:1129:VAL:HB	1:B:1132:ILE:HD11	1.94	0.48
1:D:406:GLU:HB3	1:D:418:ILE:HG13	1.95	0.48
1:D:998:THR:HB	1:E:994:ASP:HA	1.95	0.48
1:E:1039:ARG:HD3	1:E:1042:PHE:HB2	1.95	0.48
2:A:39:GLN:O	2:A:92:VAL:N	2.41	0.48
1:D:1011:GLN:OE1	1:E:762:GLN:NE2	2.38	0.48
1:E:736:VAL:H	1:E:767:LEU:HD13	1.78	0.48
1:B:198:ASP:C	1:E:425:LEU:H	2.17	0.48
1:D:494:SER:OG	1:D:495:TYR:N	2.47	0.48
1:D:1045:LYS:NZ	1:E:784:GLN:OE1	2.41	0.48
1:E:727:LEU:HB3	1:E:1025:ALA:HA	1.94	0.48
1:B:41:LYS:HB2	1:E:516:GLU:HB3	1.96	0.48
1:B:195:LYS:HB3	1:E:462:LYS:HE3	1.96	0.48
1:E:107:GLY:HA3	1:E:116:SER:HA	1.96	0.48
1:E:349:SER:HB2	1:E:352:ALA:HB3	1.95	0.48
1:E:406:GLU:HG2	1:E:418:ILE:HD11	1.96	0.48
3:C:39:GLN:HG3	3:C:88:TYR:HE1	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:69:ILE:HD11	2:A:78:LEU:HD11	1.96	0.48
1:E:401:VAL:HA	1:E:509:ARG:HA	1.95	0.48
1:E:1039:ARG:HH11	1:E:1042:PHE:HB2	1.77	0.48
3:C:37:TRP:CE2	3:C:75:LEU:HB2	2.49	0.48
1:D:858:LEU:HD13	1:D:962:LEU:HD13	1.95	0.47
1:D:959:LEU:H	1:E:765:ARG:NH1	2.12	0.47
1:E:380:TYR:O	1:E:431:GLY:N	2.44	0.47
1:E:106:PHE:HB2	1:E:117:LEU:HB3	1.95	0.47
1:D:128:ILE:HB	1:D:170:TYR:HB3	1.97	0.47
1:E:193:VAL:HG21	1:E:223:LEU:HD13	1.96	0.47
1:E:594:GLY:O	1:E:614:ASP:N	2.47	0.47
1:B:880:GLY:O	1:B:884:SER:N	2.42	0.47
1:D:1089:PHE:N	1:E:1111:GLU:OE2	2.48	0.47
1:E:956:ALA:O	1:E:960:ASN:ND2	2.48	0.47
1:D:699:LEU:HG	1:E:869:MET:HA	1.97	0.47
1:E:950:ASP:OD1	1:E:950:ASP:N	2.48	0.47
1:B:1115:ILE:HG22	1:B:1137:VAL:HG13	1.95	0.47
1:D:401:VAL:HG22	1:D:509:ARG:HA	1.97	0.47
1:B:974:SER:OG	1:B:975:SER:N	2.47	0.47
2:A:203:LEU:HA	2:A:208:TYR:HE2	1.79	0.47
1:E:48:LEU:HB3	1:E:276:LEU:HD11	1.97	0.47
1:E:909:ILE:HG13	1:E:911:VAL:HG23	1.96	0.47
1:B:339:GLY:HA2	1:B:342:PHE:HD2	1.79	0.47
1:B:1043:CYS:HB2	1:B:1048:HIS:CD2	2.50	0.47
1:E:185:ASN:ND2	1:E:211:ASN:OD1	2.48	0.47
1:E:931:ILE:O	1:E:935:GLN:HB2	2.15	0.47
1:D:391:CYS:SG	1:D:525:CYS:N	2.88	0.47
1:D:1027:THR:O	1:D:1031:GLU:N	2.47	0.47
1:D:1141:LEU:N	1:E:1118:ASP:OD2	2.47	0.47
1:E:944:ALA:HA	1:E:949:GLN:HG3	1.96	0.47
1:B:1114:ILE:O	1:B:1119:ASN:ND2	2.48	0.46
1:D:122:ASN:ND2	1:D:124:THR:OG1	2.48	0.46
1:E:453:TYR:HE2	1:E:455:LEU:HD13	1.80	0.46
1:B:66:HIS:O	1:B:79:PHE:N	2.48	0.46
1:B:713:ALA:HB3	1:D:895:GLN:H	1.80	0.46
1:E:1024:LEU:O	1:E:1027:THR:OG1	2.23	0.46
1:B:958:ALA:HA	1:B:961:THR:HG22	1.96	0.46
1:E:253:ASP:OD1	1:E:253:ASP:N	2.49	0.46
3:C:40:GLN:HA	3:C:46:PRO:HA	1.97	0.46
1:D:1002:GLN:CB	1:E:756:TYR:CG	2.84	0.46
1:E:748:GLU:HA	1:E:752:LEU:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1050:MET:HE3	1:E:1050:MET:HB2	1.66	0.46
1:D:926:GLN:O	1:D:930:ALA:N	2.44	0.46
1:E:741:TYR:CZ	1:E:966:LEU:HD11	2.51	0.46
1:B:280:ASN:OD1	1:B:283:GLY:N	2.49	0.46
1:B:906:PHE:HB3	1:B:911:VAL:HB	1.98	0.46
3:C:22:CYS:HB3	3:C:73:ALA:HB3	1.98	0.46
1:D:819:GLU:OE2	1:D:1055:SER:N	2.37	0.46
1:B:574:ASP:HA	1:B:587:ILE:HB	1.98	0.46
2:A:140:PRO:HG3	2:A:152:LEU:HB3	1.97	0.46
1:E:95:THR:HG22	1:E:189:LEU:HD13	1.98	0.46
1:E:97:LYS:HZ2	1:E:186:PHE:HE1	1.64	0.46
1:E:729:VAL:HG11	1:E:782:PHE:CE2	2.51	0.46
1:B:350:VAL:HG21	1:B:418:ILE:HD12	1.98	0.46
1:B:935:GLN:O	1:B:939:SER:N	2.49	0.46
3:C:140:ILE:HB	3:C:178:ALA:HB3	1.98	0.46
1:D:115:GLN:HE21	1:D:130:VAL:HG12	1.80	0.46
1:D:731:MET:N	1:D:774:GLN:OE1	2.43	0.46
1:D:985:ASP:N	1:D:985:ASP:OD1	2.48	0.46
1:D:1003:SER:OG	1:E:756:TYR:O	2.27	0.46
1:B:568:ASP:OD1	1:B:572:THR:OG1	2.32	0.46
1:E:737:ASP:OD1	1:E:738:CYS:N	2.49	0.46
1:E:344:ALA:H	1:E:347:PHE:HE2	1.64	0.45
1:E:965:GLN:NE2	1:E:1003:SER:O	2.48	0.45
1:B:87:ASN:HB2	1:B:269:TYR:HE1	1.82	0.45
2:A:5:VAL:HB	2:A:23:ALA:HB3	1.99	0.45
1:D:676:THR:HA	1:D:690:GLN:HA	1.97	0.45
1:E:805:ILE:HB	1:E:878:LEU:HD11	1.98	0.45
1:B:201:PHE:HD1	1:E:465:GLU:HA	1.81	0.45
1:B:1031:GLU:HG2	1:E:1039:ARG:HH22	1.82	0.45
1:B:1079:PRO:HB3	1:D:900:MET:HG2	1.99	0.45
1:E:409:GLN:NE2	1:E:415:THR:O	2.50	0.45
1:E:457:ARG:NH1	1:E:459:SER:H	2.15	0.45
1:E:804:GLN:HB3	1:E:818:ILE:HG13	1.97	0.45
1:D:117:LEU:HD22	1:D:231:ILE:HD13	1.98	0.45
1:E:99:ASN:HB3	1:E:102:ARG:HH21	1.81	0.45
1:B:110:LEU:H	1:E:470:THR:N	2.15	0.45
1:D:91:TYR:N	1:D:268:GLY:O	2.50	0.45
1:D:986:PRO:HA	1:D:989:ALA:HB3	1.99	0.45
1:B:233:ILE:N	1:E:422:ASN:O	2.50	0.45
1:B:1018:ILE:O	1:B:1022:ALA:N	2.48	0.45
1:E:119:ILE:HG23	1:E:128:ILE:HG12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:115:GLN:NE2	1:D:167:THR:OG1	2.50	0.45
1:D:958:ALA:O	1:D:962:LEU:N	2.49	0.45
1:E:117:LEU:HD13	1:E:130:VAL:HG22	1.99	0.45
1:E:966:LEU:HD22	1:E:1000:ARG:HG3	1.99	0.45
1:B:398:ASP:HB2	1:B:512:VAL:HB	1.98	0.45
1:D:1049:LEU:HD11	1:D:1067:TYR:HB2	1.99	0.45
1:E:349:SER:HB3	1:E:351:TYR:H	1.83	0.45
1:E:1015:ALA:O	1:E:1019:ARG:N	2.41	0.45
1:E:1080:ALA:HB3	1:E:1132:ILE:HD11	1.98	0.45
1:B:710:ASN:OD1	1:B:710:ASN:N	2.50	0.44
1:B:1080:ALA:HB3	1:B:1132:ILE:HG12	1.98	0.44
1:D:1081:ILE:HG23	1:D:1135:ASN:HB3	1.99	0.44
1:B:415:THR:HA	1:B:419:ALA:HB3	1.98	0.44
1:B:1037:SER:HG	1:B:1038:LYS:H	1.65	0.44
1:B:1049:LEU:HD23	1:B:1049:LEU:HA	1.83	0.44
1:D:222:ALA:HB2	1:D:285:ILE:HB	1.99	0.44
1:E:914:ASN:HB3	1:E:917:TYR:CD2	2.51	0.44
1:B:757:GLY:HA3	1:E:968:SER:HA	1.99	0.44
2:A:6:GLU:OE1	2:A:119:GLN:N	2.45	0.44
1:D:418:ILE:HA	1:D:422:ASN:HD22	1.81	0.44
1:D:1097:SER:HB3	1:D:1102:TRP:CE3	2.53	0.44
1:E:402:ILE:HG21	1:E:410:ILE:HG13	1.98	0.44
1:E:782:PHE:HE1	1:E:1060:VAL:HB	1.82	0.44
1:B:94:SER:OG	1:B:95:THR:N	2.50	0.44
1:B:327:VAL:HG22	1:B:542:ASN:HB3	2.00	0.44
1:D:1007:TYR:HB2	1:E:759:PHE:CD2	2.53	0.44
1:D:1107:ARG:NH2	1:E:905:ARG:HA	2.32	0.44
1:E:313:TYR:CE2	1:E:599:THR:HG21	2.53	0.44
1:B:201:PHE:CD1	1:E:465:GLU:HA	2.53	0.44
1:B:708:SER:OG	1:B:709:ASN:N	2.49	0.44
1:D:346:ARG:HH22	1:D:450:ASN:HB3	1.82	0.44
1:D:956:ALA:O	1:D:960:ASN:N	2.50	0.44
1:D:1083:HIS:CE1	1:D:1137:VAL:H	2.35	0.44
1:E:777:ASN:HD21	1:E:1019:ARG:HD3	1.82	0.44
1:E:905:ARG:NH2	1:E:1035:GLY:H	2.16	0.44
1:B:310:LYS:HG2	1:B:664:ILE:HD11	2.00	0.44
1:B:526:GLY:O	1:B:528:LYS:NZ	2.40	0.44
1:B:708:SER:OG	1:B:710:ASN:OD1	2.35	0.44
1:B:758:SER:HB2	1:B:761:THR:HB	1.99	0.44
1:E:722:VAL:HG13	1:E:1065:VAL:HG22	1.99	0.44
2:A:137:PRO:HA	2:A:154:CYS:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:609:ALA:HA	1:B:652:GLY:HA3	1.99	0.44
1:B:985:ASP:HA	1:B:986:PRO:HD3	1.85	0.44
3:C:136:LEU:HD22	3:C:182:LEU:HD23	1.99	0.44
1:E:56:LEU:HD13	1:E:270:LEU:HD23	2.00	0.44
1:B:136:CYS:HB3	1:B:139:PRO:HB3	2.00	0.44
1:B:897:PRO:HG2	1:B:900:MET:HG2	1.99	0.44
1:E:44:ARG:HD2	1:E:279:TYR:CE1	2.53	0.44
1:E:402:ILE:HD12	1:E:407:VAL:HA	1.99	0.44
1:D:914:ASN:OD1	1:D:914:ASN:N	2.50	0.43
1:D:1011:GLN:O	1:D:1015:ALA:N	2.46	0.43
1:D:1083:HIS:N	1:D:1086:LYS:O	2.37	0.43
1:B:93:ALA:HA	1:B:191:GLU:HA	2.00	0.43
1:B:197:ILE:N	1:E:424:LYS:HB3	2.32	0.43
1:B:738:CYS:SG	1:B:739:THR:N	2.91	0.43
1:B:995:ARG:O	1:E:995:ARG:NH2	2.49	0.43
1:B:1106:GLN:NE2	1:B:1109:PHE:O	2.51	0.43
1:D:15:CYS:HB2	1:D:136:CYS:HB3	1.88	0.43
1:D:466:ARG:HH12	1:E:167:THR:HA	1.83	0.43
1:E:735:SER:HA	1:E:767:LEU:HB3	2.00	0.43
1:B:199:GLY:H	1:E:424:LYS:HA	1.83	0.43
3:C:41:HIS:HB2	3:C:44:LYS:HB2	1.99	0.43
1:E:551:VAL:HG23	1:E:590:CYS:HB3	1.98	0.43
1:E:1012:LEU:HA	1:E:1015:ALA:HB3	1.99	0.43
1:B:121:ASN:ND2	1:B:125:ASN:O	2.51	0.43
1:B:227:VAL:HG13	1:E:355:ARG:CZ	2.49	0.43
1:B:594:GLY:N	1:B:613:GLN:OE1	2.51	0.43
1:D:996:LEU:HA	1:D:996:LEU:HD23	1.77	0.43
1:D:1079:PRO:O	1:E:914:ASN:ND2	2.27	0.43
1:B:79:PHE:N	1:B:261:GLY:O	2.52	0.43
1:B:307:THR:HA	1:B:602:THR:HG21	2.01	0.43
1:B:378:LYS:H	1:B:434:ILE:HG13	1.83	0.43
1:B:884:SER:HA	1:B:896:ILE:HG22	1.99	0.43
3:C:155:ASP:OD2	3:C:193:ARG:N	2.52	0.43
1:D:210:ILE:H	1:D:210:ILE:HG13	1.69	0.43
1:E:89:GLY:O	1:E:270:LEU:N	2.34	0.43
1:E:613:GLN:HA	1:E:648:GLY:HA3	2.01	0.43
1:B:991:VAL:O	1:B:995:ARG:NE	2.48	0.43
2:A:170:SER:N	2:A:211:ASN:OD1	2.50	0.43
1:D:807:PRO:HA	1:D:816:SER:HA	2.00	0.43
1:E:418:ILE:HD12	1:E:418:ILE:HG23	1.85	0.43
1:E:462:LYS:HB2	1:E:462:LYS:HE3	1.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:734:THR:HG21	1:E:1011:GLN:HG3	2.01	0.43
1:E:736:VAL:N	1:E:767:LEU:HD13	2.34	0.43
1:B:55:PHE:HA	1:B:270:LEU:HD13	1.98	0.43
1:B:282:ASN:HB3	1:E:563:GLN:HB3	2.01	0.43
1:B:1096:VAL:O	1:B:1103:PHE:N	2.47	0.43
1:D:897:PRO:HD2	1:D:900:MET:HB2	2.00	0.43
1:E:784:GLN:HE21	1:E:1034:LEU:HD11	1.83	0.43
1:B:110:LEU:H	1:E:470:THR:H	1.67	0.43
1:B:195:LYS:HG3	1:E:464:PHE:HB2	1.98	0.43
1:B:950:ASP:HA	1:B:953:ASN:HB2	2.01	0.43
1:D:555:SER:OG	1:D:556:ASN:N	2.52	0.43
1:B:653:ALA:HB2	1:B:692:ILE:HG22	2.01	0.43
1:B:949:GLN:O	1:B:953:ASN:ND2	2.52	0.43
2:A:47:TRP:HZ2	2:A:50:ALA:HB2	1.84	0.43
1:E:401:VAL:HG22	1:E:509:ARG:HG2	2.01	0.43
1:B:114:THR:HG22	1:E:491:PRO:HG3	2.01	0.43
1:B:478:THR:O	1:B:487:ASN:ND2	2.50	0.43
1:B:951:VAL:O	1:B:955:ASN:ND2	2.52	0.43
1:D:103:GLY:O	1:D:241:LEU:N	2.44	0.43
1:D:965:GLN:HB3	1:E:754:LEU:O	2.18	0.43
1:B:90:VAL:H	1:E:462:LYS:NZ	2.17	0.42
1:B:229:LEU:HB3	1:E:353:TRP:HE1	1.83	0.42
1:B:111:ASP:O	1:E:470:THR:OG1	2.26	0.42
1:B:456:PHE:H	1:B:491:PRO:HB3	1.84	0.42
2:A:168:TRP:HB2	2:A:173:LEU:HB3	2.01	0.42
1:D:307:THR:HA	1:D:602:THR:HG21	2.00	0.42
1:D:1017:GLU:O	1:D:1021:SER:N	2.52	0.42
1:B:196:ASN:HB2	1:E:461:LEU:HG	2.02	0.42
1:B:357:ARG:NH2	1:B:358:ILE:O	2.52	0.42
2:A:6:GLU:O	2:A:119:GLN:NE2	2.51	0.42
1:D:1040:VAL:HB	1:E:1031:GLU:HG2	2.01	0.42
1:E:57:PRO:HB3	1:E:273:ARG:HH11	1.83	0.42
1:E:551:VAL:O	1:E:588:THR:N	2.47	0.42
1:B:105:ILE:HB	1:B:239:GLN:HB3	2.02	0.42
1:D:963:VAL:HA	1:D:966:LEU:HD12	2.00	0.42
1:D:1107:ARG:NH2	1:E:1036:GLN:OE1	2.50	0.42
1:B:443:SER:HB2	1:B:507:PRO:HG3	2.01	0.42
1:B:605:SER:OG	1:B:606:ASN:N	2.52	0.42
1:B:665:PRO:HB2	1:D:864:LEU:HD21	2.00	0.42
1:D:343:ASN:HB2	1:D:436:TRP:HZ2	1.83	0.42
1:E:48:LEU:HD22	1:E:305:SER:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:513:LEU:HB3	1:E:515:PHE:HE2	1.84	0.42
1:E:720:ILE:HG22	1:E:930:ALA:HB2	2.01	0.42
1:E:859:THR:OG1	1:E:860:VAL:N	2.52	0.42
1:B:84:LEU:O	1:B:238:PHE:N	2.48	0.42
1:B:994:ASP:HB3	1:E:992:GLN:NE2	2.34	0.42
1:D:1005:GLN:HB2	1:E:1002:GLN:HE22	1.85	0.42
1:D:1077:THR:OG1	1:D:1078:ALA:O	2.32	0.42
1:E:66:HIS:HB3	1:E:68:ILE:HG22	2.01	0.42
1:E:551:VAL:N	1:E:588:THR:O	2.42	0.42
1:E:605:SER:OG	1:E:606:ASN:N	2.53	0.42
1:B:165:ASN:HD21	1:E:452:LEU:HG	1.84	0.42
1:D:112:SER:N	1:D:133:PHE:O	2.52	0.42
1:D:703:ASN:H	1:E:791:THR:N	2.17	0.42
1:D:710:ASN:OD1	1:D:710:ASN:N	2.51	0.42
1:E:108:THR:HG22	1:E:109:THR:HG23	2.01	0.42
1:E:994:ASP:HB3	1:E:995:ARG:HE	1.85	0.42
1:B:202:LYS:HG3	1:E:464:PHE:CG	2.55	0.42
1:B:240:THR:OG1	1:B:241:LEU:N	2.53	0.42
1:B:950:ASP:OD1	1:B:950:ASP:N	2.52	0.42
2:A:12:VAL:HG21	2:A:18:LEU:HB2	2.01	0.42
3:C:20:ILE:O	3:C:75:LEU:N	2.52	0.42
1:D:94:SER:HA	1:D:265:TYR:HA	2.02	0.42
1:D:516:GLU:HB3	1:D:519:HIS:CD2	2.49	0.42
1:D:711:SER:O	1:E:897:PRO:HB2	2.20	0.42
1:D:1082:CYS:HB2	1:D:1132:ILE:HD13	2.02	0.42
1:E:552:LEU:HD13	1:E:585:LEU:HD13	2.01	0.42
1:B:474:GLN:HE21	1:B:476:GLY:H	1.67	0.42
1:B:760:CYS:HA	1:B:763:LEU:HB2	2.02	0.42
2:A:39:GLN:OE1	3:C:40:GLN:NE2	2.40	0.42
2:A:157:LYS:NZ	3:C:135:THR:OG1	2.38	0.42
1:D:290:ASP:OD1	1:D:291:CYS:N	2.53	0.42
1:D:572:THR:OG1	1:D:573:THR:N	2.52	0.42
1:E:903:ALA:HB1	1:E:913:GLN:HA	2.02	0.42
1:B:34:ARG:NH2	1:B:218:GLN:O	2.52	0.41
2:A:96:ALA:HB1	2:A:114:PHE:HB3	2.02	0.41
1:D:950:ASP:OD1	1:D:951:VAL:N	2.52	0.41
1:B:214:ARG:O	1:B:266:TYR:OH	2.25	0.41
1:B:816:SER:OG	1:B:817:PHE:N	2.51	0.41
1:D:280:ASN:OD1	1:D:283:GLY:N	2.52	0.41
1:D:1068:VAL:HG11	1:E:890:ALA:H	1.85	0.41
1:E:898:PHE:HA	1:E:901:GLN:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:TYR:OH	1:B:191:GLU:OE1	2.38	0.41
1:D:656:VAL:HG23	1:D:658:ASN:H	1.84	0.41
1:D:733:LYS:NZ	1:D:775:ASP:OD1	2.46	0.41
1:E:765:ARG:O	1:E:768:THR:OG1	2.27	0.41
1:E:804:GLN:HG3	1:E:935:GLN:NE2	2.35	0.41
2:A:37:VAL:N	2:A:94:TYR:O	2.44	0.41
1:D:438:SER:HB2	1:D:441:LEU:HD12	2.02	0.41
1:E:280:ASN:ND2	1:E:284:THR:OG1	2.53	0.41
1:E:924:ALA:O	1:E:928:ASN:N	2.50	0.41
1:E:926:GLN:HA	1:E:929:SER:HB3	2.02	0.41
1:E:984:LEU:HB3	1:E:989:ALA:HB2	2.02	0.41
1:B:380:TYR:OH	1:B:410:ILE:O	2.35	0.41
1:B:448:ASN:HB3	1:B:497:PHE:HB2	2.03	0.41
1:B:752:LEU:HA	1:B:752:LEU:HD23	1.88	0.41
2:A:214:HIS:CD2	2:A:216:PRO:HD2	2.56	0.41
1:D:110:LEU:HB3	1:D:135:PHE:CD2	2.55	0.41
1:E:117:LEU:HD11	1:E:128:ILE:HG23	2.02	0.41
1:E:364:ASP:OD1	1:E:388:ASN:ND2	2.44	0.41
1:E:978:ASN:O	1:E:982:SER:N	2.49	0.41
1:B:567:ARG:HB3	1:B:573:THR:HG22	2.02	0.41
1:D:342:PHE:HE1	1:D:434:ILE:HG21	1.85	0.41
1:E:609:ALA:HB2	1:E:692:ILE:HB	2.03	0.41
1:E:756:TYR:OH	1:E:997:ILE:CG2	2.68	0.41
1:B:377:PHE:HA	1:B:434:ILE:HG23	2.03	0.41
1:B:858:LEU:HD13	1:B:858:LEU:HA	1.85	0.41
3:C:135:THR:HA	3:C:183:SER:HA	2.03	0.41
1:D:426:PRO:HB2	1:D:429:PHE:HB2	2.01	0.41
1:D:1005:GLN:O	1:D:1009:THR:N	2.52	0.41
1:B:134:GLN:N	1:B:161:SER:OG	2.53	0.41
1:B:196:ASN:HA	1:E:464:PHE:H	1.86	0.41
1:B:233:ILE:HG23	1:E:351:TYR:HB2	2.02	0.41
1:B:302:THR:HG23	1:B:303:LEU:HD12	2.02	0.41
3:C:37:TRP:CD1	3:C:75:LEU:HD13	2.56	0.41
1:D:208:THR:HG22	1:D:210:ILE:HG23	2.03	0.41
1:D:454:ARG:HG3	1:D:491:PRO:HB2	2.02	0.41
1:D:560:LEU:O	1:D:562:PHE:N	2.46	0.41
1:D:645:THR:OG1	1:D:648:GLY:N	2.42	0.41
1:D:1082:CYS:HA	1:D:1087:ALA:HA	2.03	0.41
1:E:383:SER:HB2	1:E:386:LYS:HB2	2.03	0.41
1:E:452:LEU:HD22	1:E:492:LEU:HD13	2.03	0.41
1:E:513:LEU:HB3	1:E:515:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:703:ASN:OD1	1:E:703:ASN:N	2.53	0.41
1:B:1049:LEU:HD11	1:B:1067:TYR:HB2	2.02	0.41
1:D:247:SER:O	1:D:250:THR:OG1	2.28	0.41
1:D:336:CYS:HA	1:D:337:PRO:HD3	1.95	0.41
1:D:821:LEU:HD21	1:D:939:SER:HB3	2.03	0.41
1:E:563:GLN:O	1:E:577:ARG:NH2	2.54	0.41
1:E:804:GLN:HG3	1:E:935:GLN:HE21	1.86	0.41
1:B:197:ILE:HB	1:E:463:PRO:HB3	2.03	0.40
1:B:329:PHE:HE1	1:B:544:ASN:HA	1.86	0.40
1:B:784:GLN:HE22	1:B:1030:SER:HB2	1.86	0.40
1:B:877:LEU:HD23	1:B:877:LEU:HA	1.88	0.40
1:D:66:HIS:H	1:D:78:ARG:NH2	2.18	0.40
1:D:338:PHE:HA	1:D:341:VAL:HB	2.03	0.40
1:D:712:ILE:HG22	1:E:900:MET:HB2	2.03	0.40
1:E:119:ILE:HG12	1:E:128:ILE:HG12	2.02	0.40
1:E:321:GLN:HA	1:E:322:PRO:HD3	1.81	0.40
1:E:400:PHE:N	1:E:510:VAL:O	2.45	0.40
1:B:111:ASP:HB2	1:E:471:GLU:HG2	2.03	0.40
1:B:319:ARG:HE	1:D:740:MET:HE1	1.85	0.40
1:B:341:VAL:HA	1:B:344:ALA:HB2	2.02	0.40
1:E:906:PHE:CD1	1:E:1049:LEU:HD22	2.57	0.40
1:B:296:LEU:HD22	1:B:606:ASN:HD21	1.87	0.40
1:B:1119:ASN:N	1:B:1119:ASN:OD1	2.51	0.40
1:D:112:SER:HA	1:D:132:GLU:HB3	2.03	0.40
1:D:723:THR:OG1	1:D:724:THR:N	2.54	0.40
1:E:806:LEU:HA	1:E:807:PRO:HD3	1.80	0.40
1:B:105:ILE:HG13	1:B:241:LEU:HG	2.03	0.40
1:B:861:LEU:HD23	1:B:861:LEU:HA	1.91	0.40
2:A:9:GLY:HA2	2:A:18:LEU:HD21	2.03	0.40
2:A:159:TYR:HB2	2:A:214:HIS:CE1	2.56	0.40
1:E:85:PRO:HA	1:E:237:ARG:HA	2.03	0.40
1:E:888:PHE:HA	1:E:893:ALA:HB2	2.03	0.40
1:B:970:PHE:HE1	1:D:756:TYR:HA	1.87	0.40
2:A:82:MET:HB3	2:A:85:LEU:HD21	2.03	0.40
1:D:707:TYR:HD1	1:E:898:PHE:HB3	1.85	0.40
1:D:1125:ASN:HA	1:E:1112:PRO:HB3	2.03	0.40
1:E:70:VAL:N	1:E:261:GLY:O	2.43	0.40
1:E:728:PRO:HG2	1:E:1021:SER:HB3	2.04	0.40
1:E:1030:SER:HA	1:E:1033:VAL:HB	2.04	0.40
1:E:1126:CYS:HB3	1:E:1132:ILE:HD13	2.03	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	B	1017/1273 (80%)	882 (87%)	134 (13%)	1 (0%)	48	83
1	D	1055/1273 (83%)	956 (91%)	98 (9%)	1 (0%)	48	83
1	E	1055/1273 (83%)	873 (83%)	178 (17%)	4 (0%)	30	67
2	A	228/231 (99%)	209 (92%)	19 (8%)	0	100	100
3	C	213/216 (99%)	199 (93%)	14 (7%)	0	100	100
All	All	3568/4266 (84%)	3119 (87%)	443 (12%)	6 (0%)	45	78

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	464	PHE
1	B	196	ASN
1	D	1079	PRO
1	E	452	LEU
1	E	463	PRO
1	E	600	PRO

### 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	B	891/1112 (80%)	884 (99%)	7 (1%)	79	85
1	D	933/1112 (84%)	932 (100%)	1 (0%)	92	95
1	E	933/1112 (84%)	919 (98%)	14 (2%)	60	75

*Continued on next page...*

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	190/191 (100%)	189 (100%)	1 (0%)	86	90
3	C	183/184 (100%)	183 (100%)	0	100	100
All	All	3130/3711 (84%)	3107 (99%)	23 (1%)	80	87

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	121	ASN
1	B	214	ARG
1	B	233	ILE
1	B	606	ASN
1	B	620	VAL
1	B	755	GLN
1	B	995	ARG
2	A	213	ASN
1	D	1122	VAL
1	E	422	ASN
1	E	451	TYR
1	E	457	ARG
1	E	468	ILE
1	E	470	THR
1	E	778	THR
1	E	906	PHE
1	E	911	VAL
1	E	914	ASN
1	E	995	ARG
1	E	1042	PHE
1	E	1094	VAL
1	E	1107	ARG
1	E	1109	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	87	ASN
1	B	121	ASN
1	B	207	HIS
1	B	234	ASN
1	B	282	ASN
1	B	474	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	481	ASN
1	B	487	ASN
1	B	606	ASN
1	B	641	ASN
1	B	751	ASN
1	B	755	GLN
1	B	953	ASN
1	B	955	ASN
1	B	1011	GLN
2	A	178	HIS
2	A	213	ASN
3	C	41	HIS
1	D	17	ASN
1	D	115	GLN
1	D	360	ASN
1	D	519	HIS
1	D	580	GLN
1	D	607	GLN
1	D	777	ASN
1	D	784	GLN
1	D	901	GLN
1	D	907	ASN
1	D	953	ASN
1	D	1002	GLN
1	D	1054	GLN
1	E	69	HIS
1	E	99	ASN
1	E	115	GLN
1	E	207	HIS
1	E	314	GLN
1	E	613	GLN
1	E	690	GLN
1	E	853	GLN
1	E	901	GLN
1	E	907	ASN
1	E	960	ASN
1	E	992	GLN
1	E	1002	GLN
1	E	1005	GLN
1	E	1054	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

### 5.7 Other polymers ⓘ

There are no such residues in this entry.

### 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.



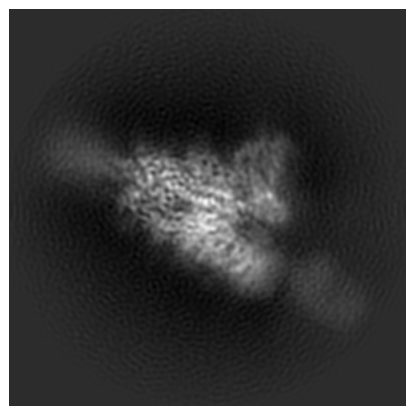
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34563. These allow visual inspection of the internal detail of the map and identification of artifacts.

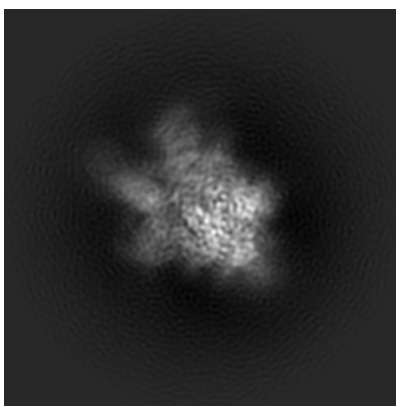
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

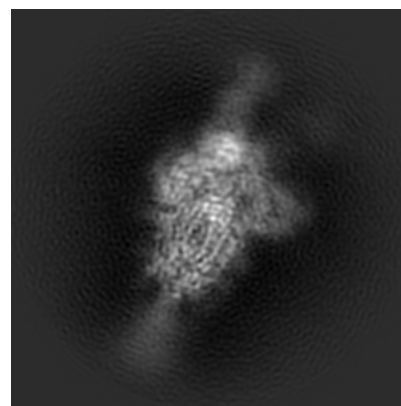
#### 6.1.1 Primary map



X

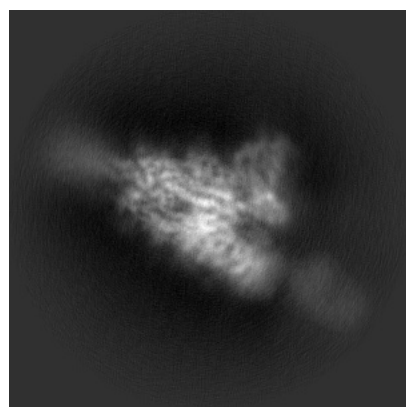


Y

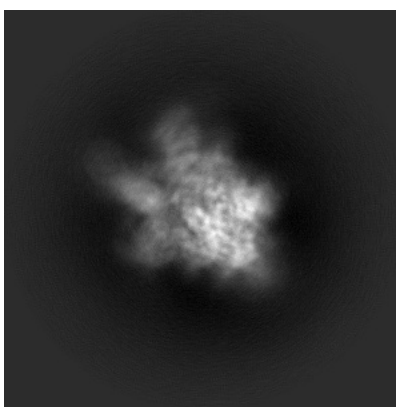


Z

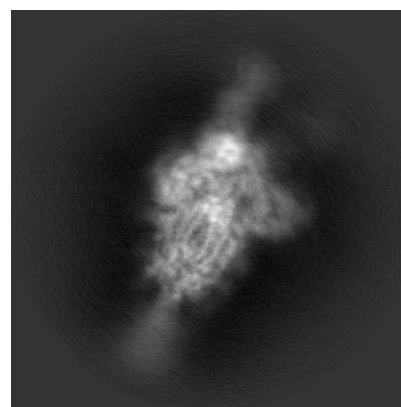
#### 6.1.2 Raw map



X



Y

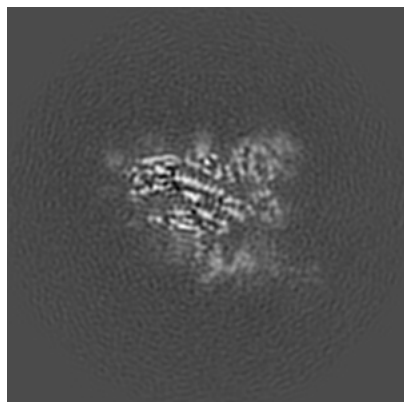


Z

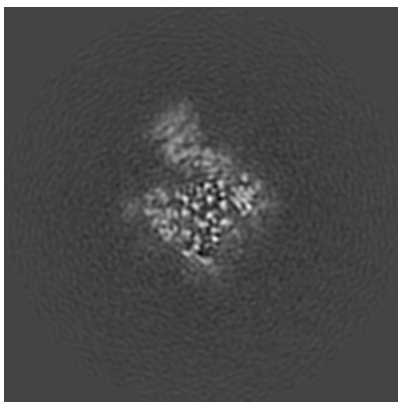
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

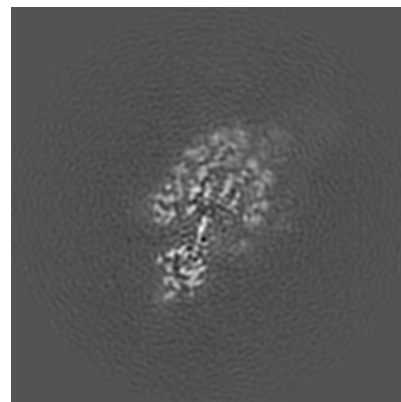
### 6.2.1 Primary map



X Index: 180

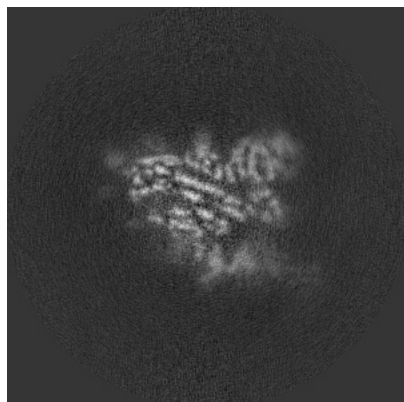


Y Index: 180

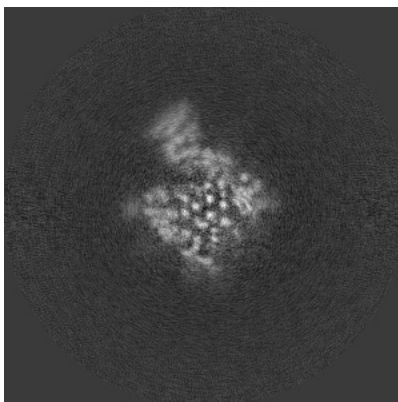


Z Index: 180

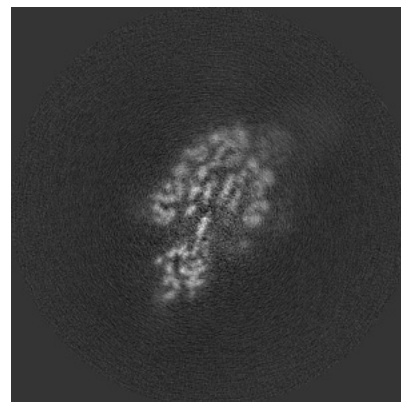
### 6.2.2 Raw map



X Index: 180



Y Index: 180

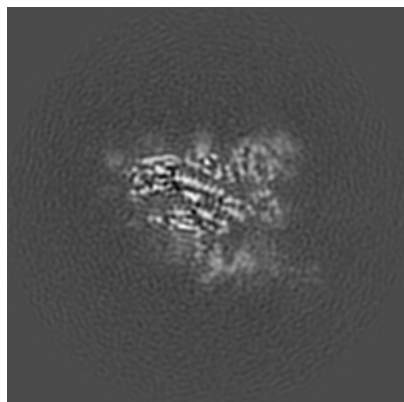


Z Index: 180

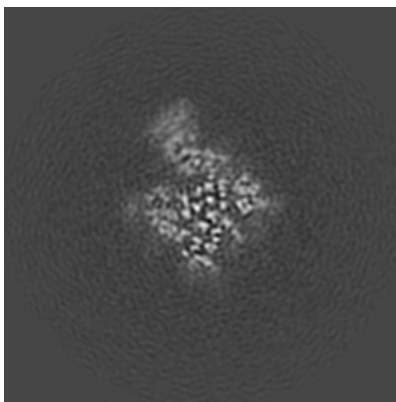
The images above show central slices of the map in three orthogonal directions.

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

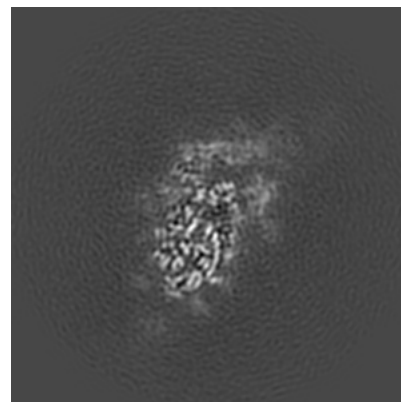
### 6.3.1 Primary map



X Index: 180

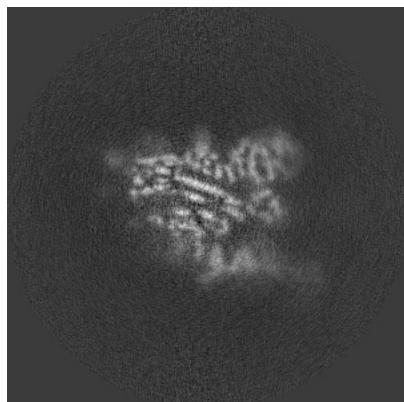


Y Index: 178

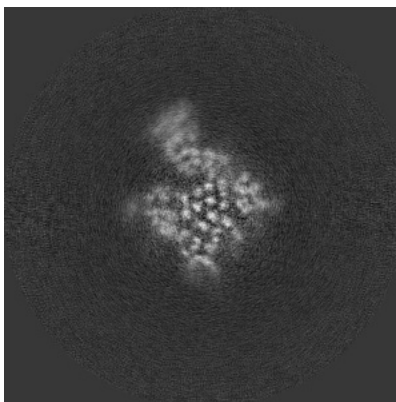


Z Index: 193

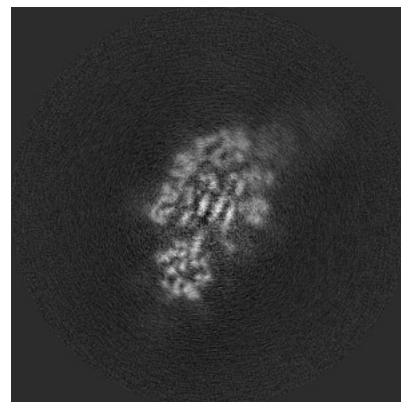
### 6.3.2 Raw map



X Index: 181



Y Index: 178

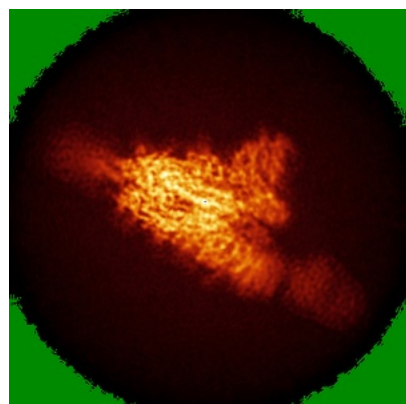


Z Index: 185

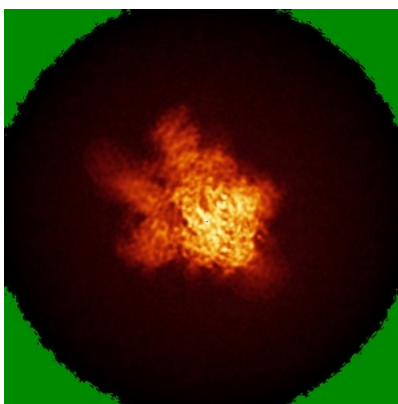
The images above show the largest variance slices of the map in three orthogonal directions.

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

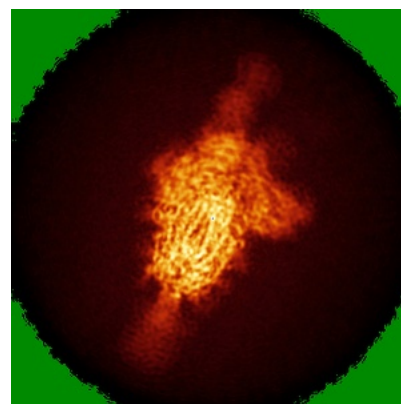
### 6.4.1 Primary map



X

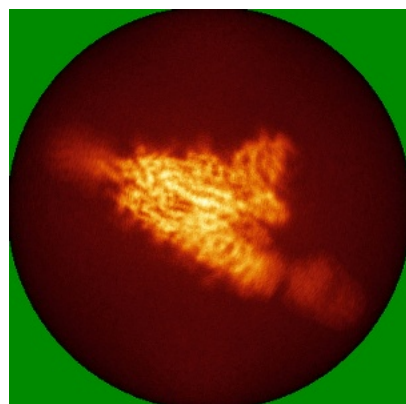


Y

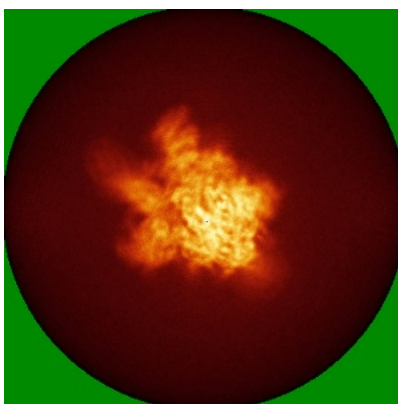


Z

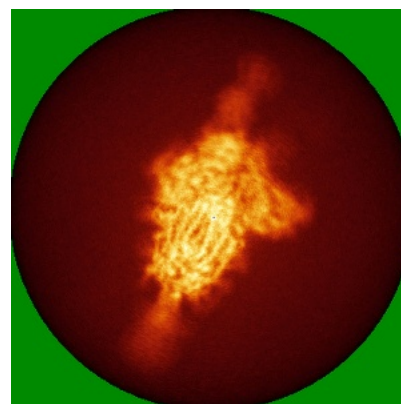
### 6.4.2 Raw map



X



Y

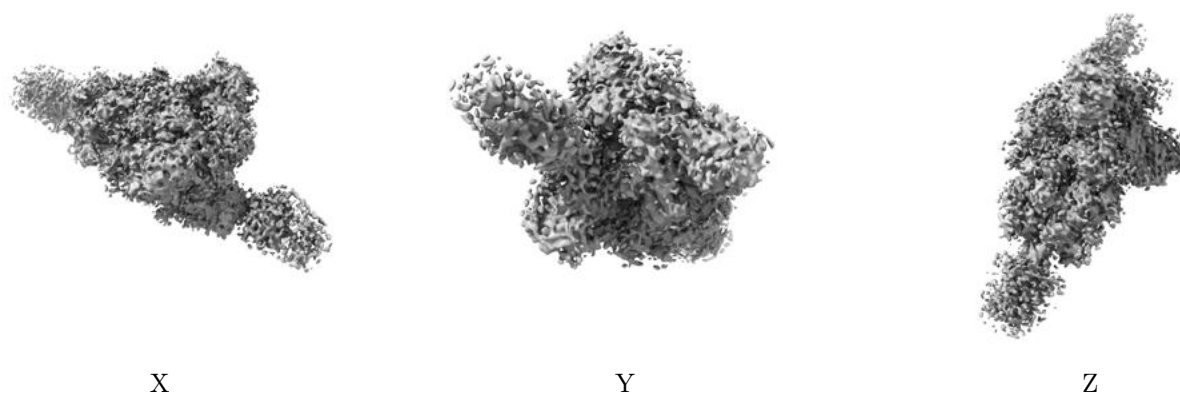


Z

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.

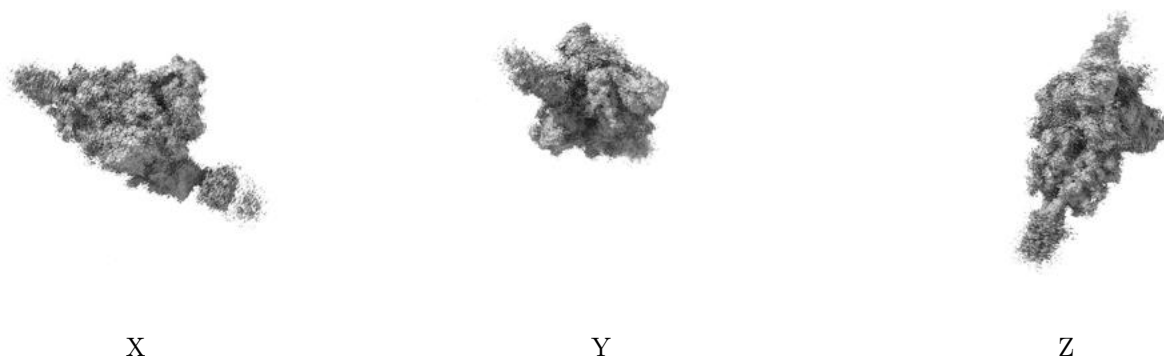
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 2.82. 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.

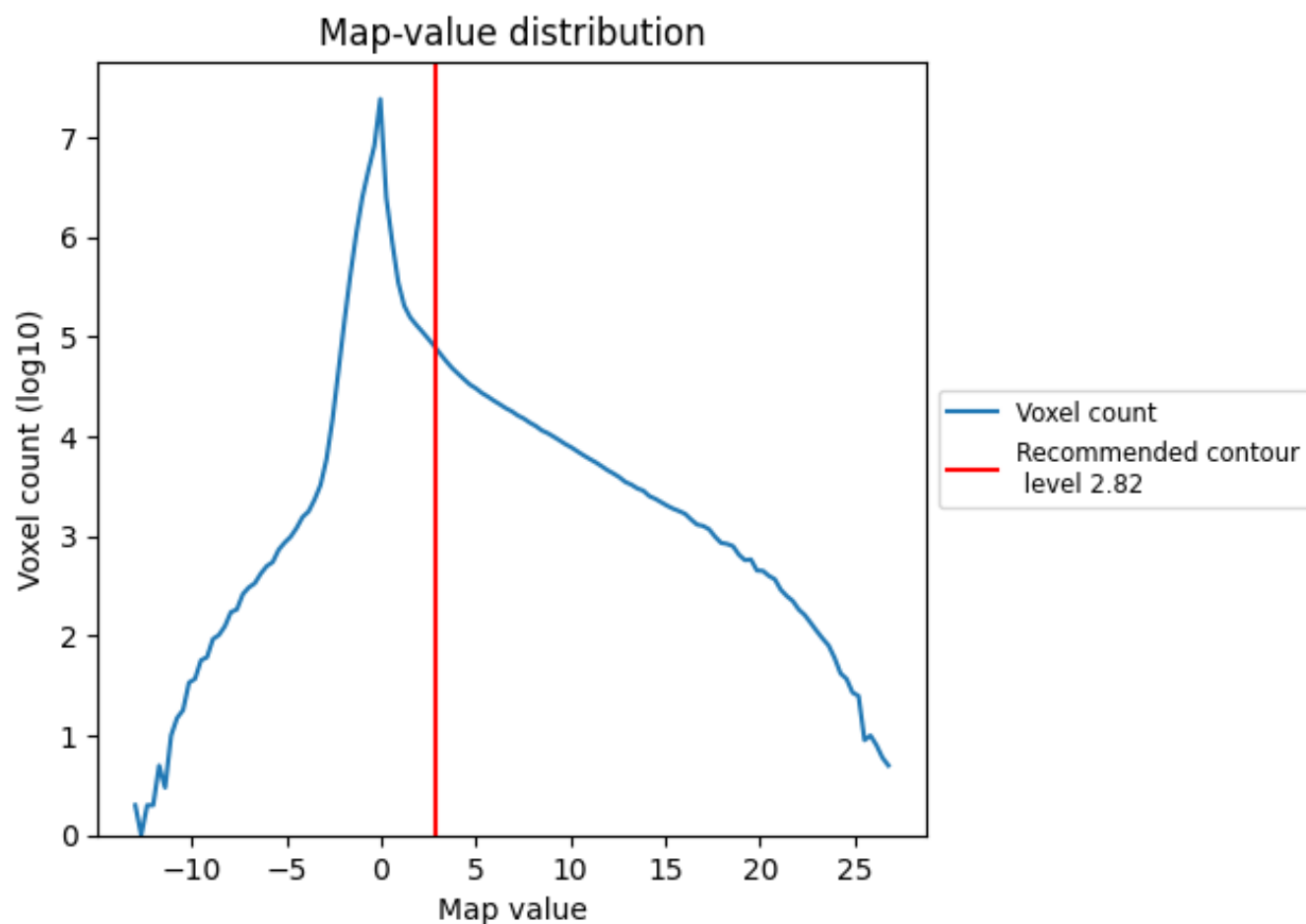
## 6.6 Mask visualisation [i](#)

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.

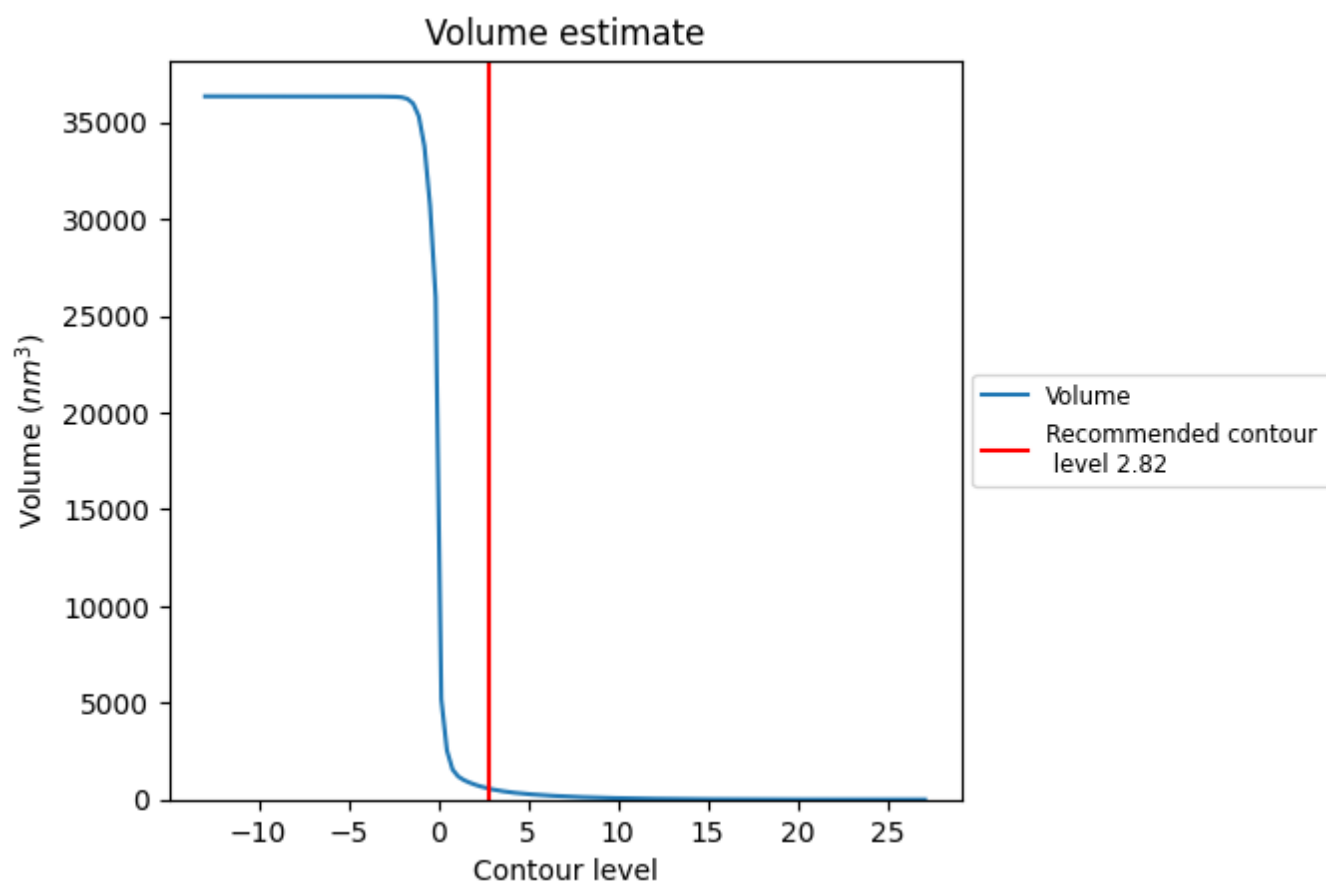
### 7.1 Map-value distribution [i](#)



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.



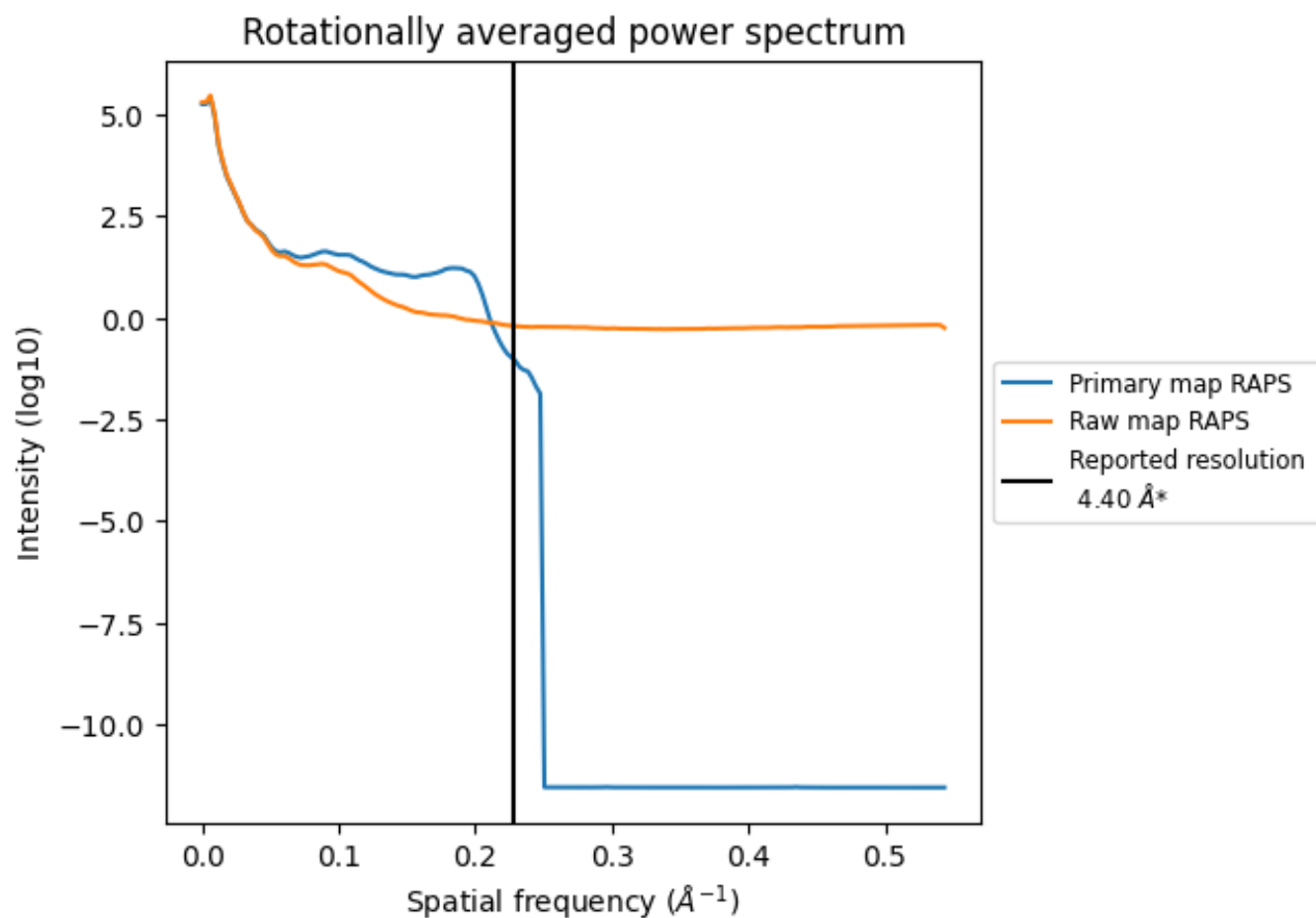
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 566 nm<sup>3</sup>; this corresponds to an approximate mass of 511 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 ⓘ



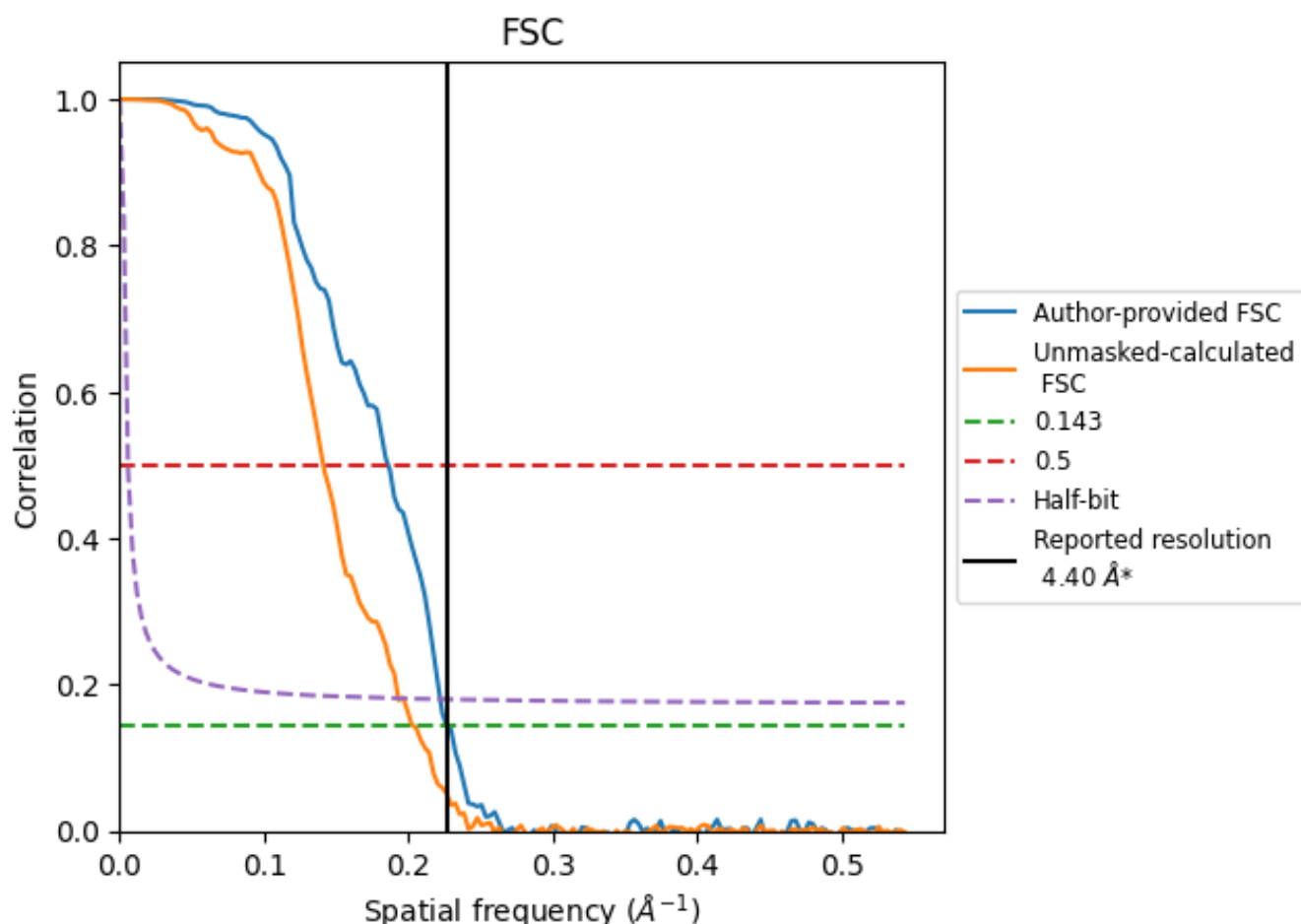
\*Reported resolution corresponds to spatial frequency of 0.227  $\text{\AA}^{-1}$



## 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 0.227 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

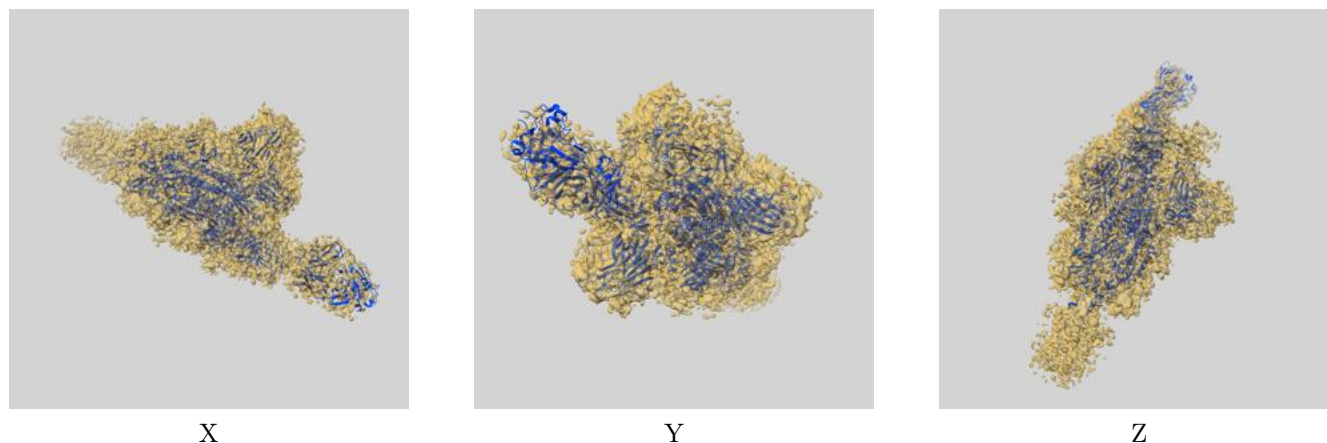
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.40	-	-
Author-provided FSC curve	4.41	5.38	4.50
Unmasked-calculated*	4.91	7.09	5.18

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.91 differs from the reported value 4.4 by more than 10 %

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

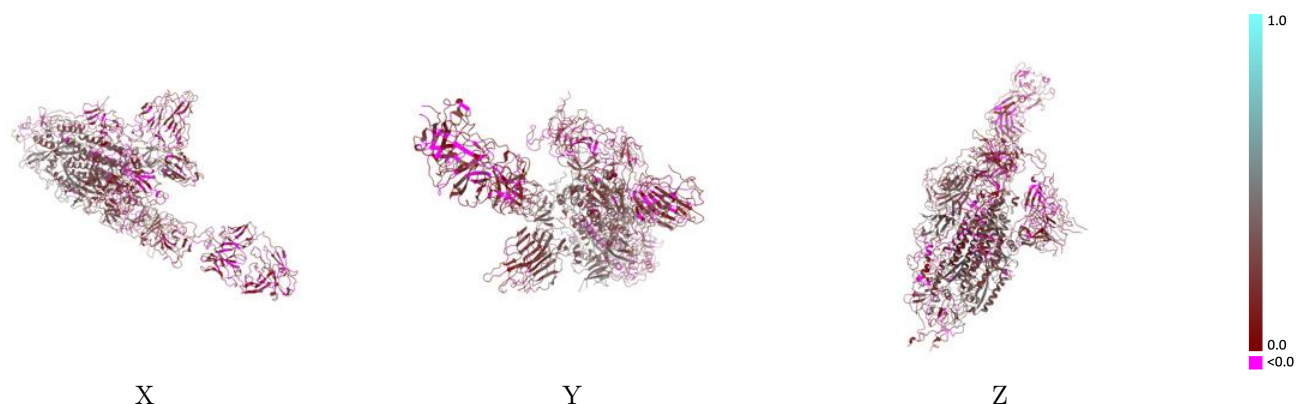
This section contains information regarding the fit between EMDB map EMD-34563 and PDB model 8YBY. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

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



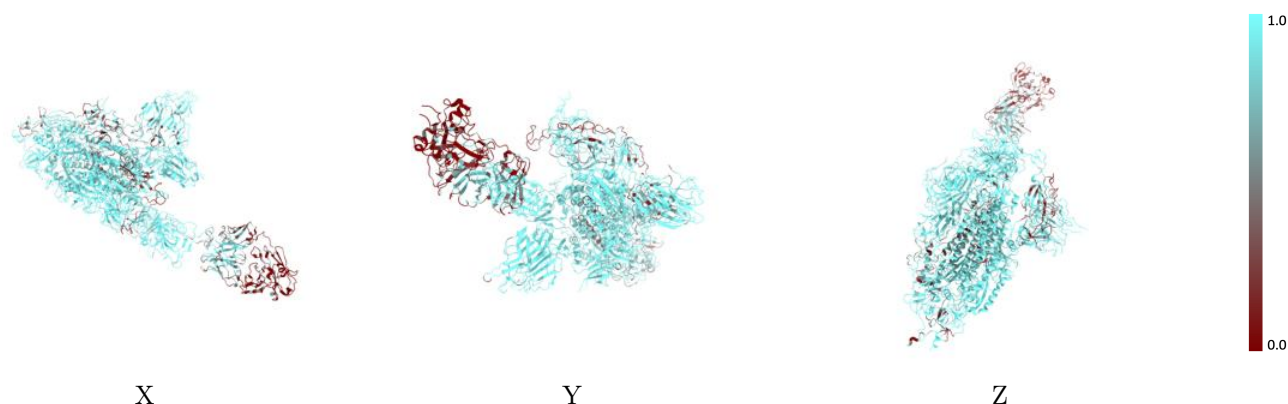
The images above show the 3D surface view of the map at the recommended contour level 2.82 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



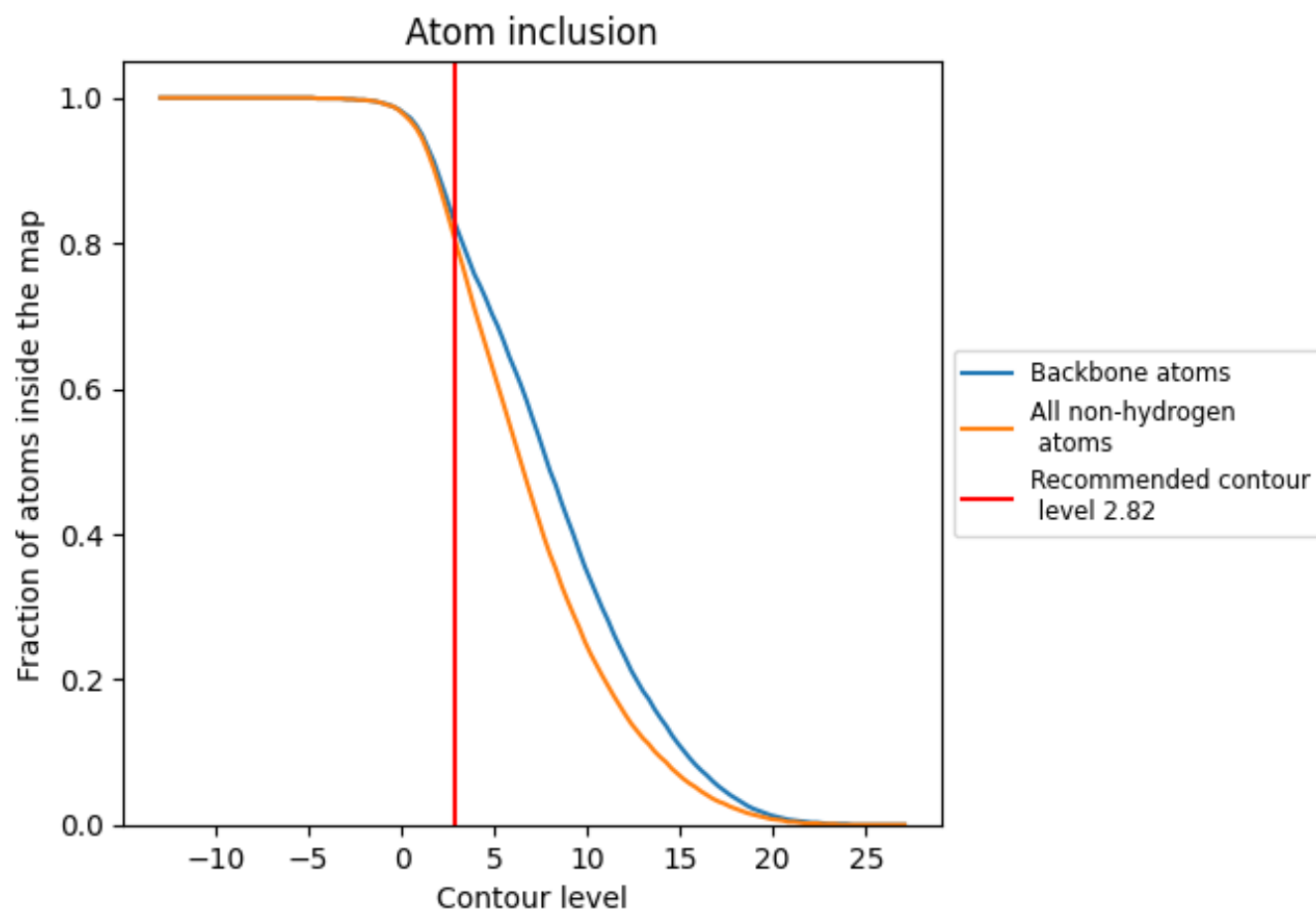
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.82).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 81% 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 (2.82) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8090	<div></div> 0.2290
A	<div></div> 0.4090	<div></div> 0.1310
B	<div></div> 0.9600	<div></div> 0.3040
C	<div></div> 0.3620	<div></div> 0.1290
D	<div></div> 0.9460	<div></div> 0.3030
E	<div></div> 0.6930	<div></div> 0.1210

