



Full wwPDB EM Validation Report ⓘ

Oct 1, 2024 – 11:48 AM JST

PDB ID : 7V6O
EMDB ID : EMD-31744
Title : MERS S ectodomain trimer in complex with neutralizing antibody 111 (state 2)
Authors : Wang, X.; Zhao, J.; Wang, Z.; Zeng, J.; Zhang, S.; Wang, Y.
Deposited on : 2021-08-20
Resolution : 4.56 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.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) : 2.39

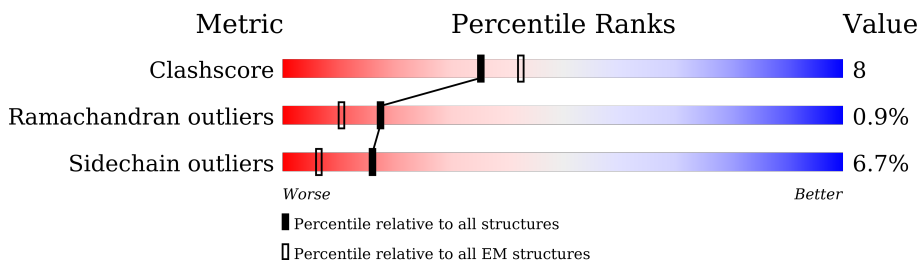
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.56 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 ≥ 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	1189	
1	B	1189	
1	C	1189	
2	D	216	
2	F	216	
2	H	216	
3	E	227	
3	G	227	

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Mol	Chain	Length	Quality of chain
3	I	227	<div><div></div><div>11%</div><div>78%</div><div>19%</div><div>..</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 34816 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1137	Total	C	N	O	S	0	0
			8325	5252	1396	1634	43		
1	B	1141	Total	C	N	O	S	1	0
			8446	5337	1401	1661	47		
1	C	1137	Total	C	N	O	S	1	0
			8401	5312	1400	1641	48		

- Molecule 2 is a protein called 111 L.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	216	Total	C	N	O	S	0	0
			1574	969	268	329	8		
2	F	216	Total	C	N	O	S	0	0
			1602	987	275	332	8		
2	H	216	Total	C	N	O	S	0	0
			1600	985	277	330	8		

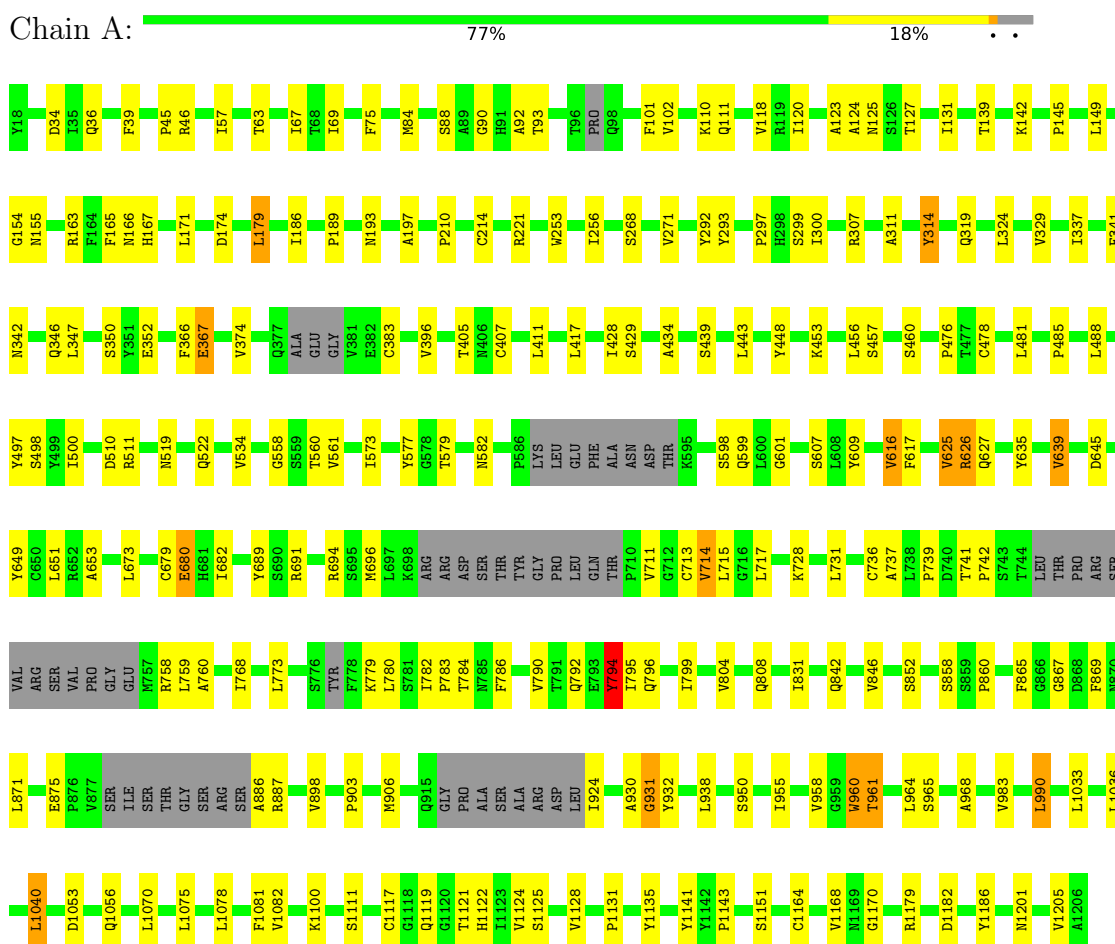
- Molecule 3 is a protein called 111 H.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	226	Total	C	N	O	S	0	0
			1615	999	278	332	6		
3	G	225	Total	C	N	O	S	0	0
			1644	1031	280	327	6		
3	I	225	Total	C	N	O	S	0	0
			1609	1001	277	326	5		

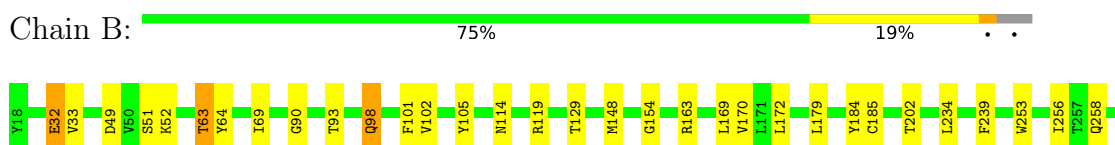
3 Residue-property plots

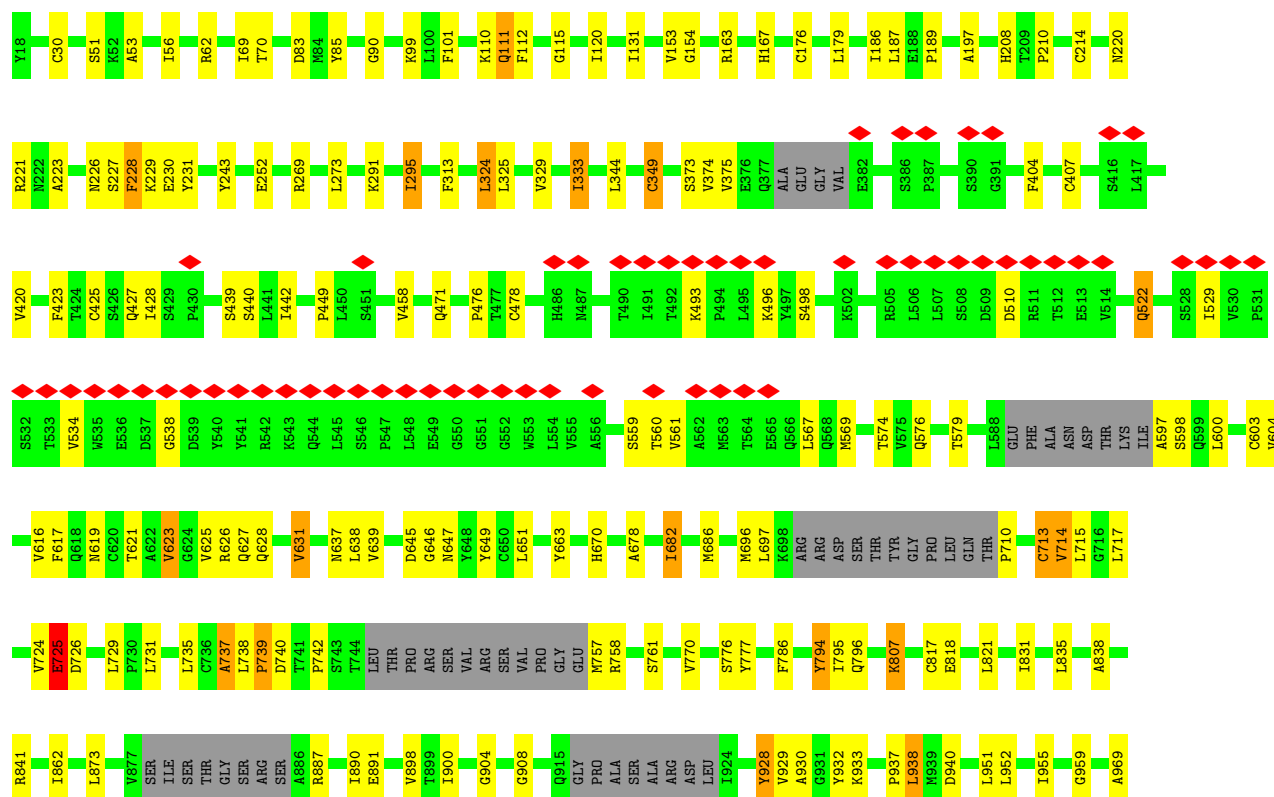
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

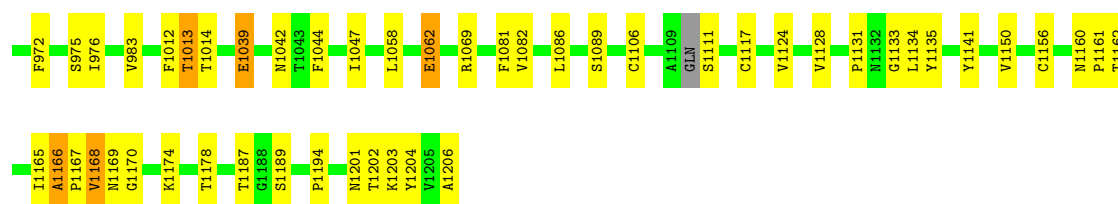
• Molecule 1: Spike glycoprotein



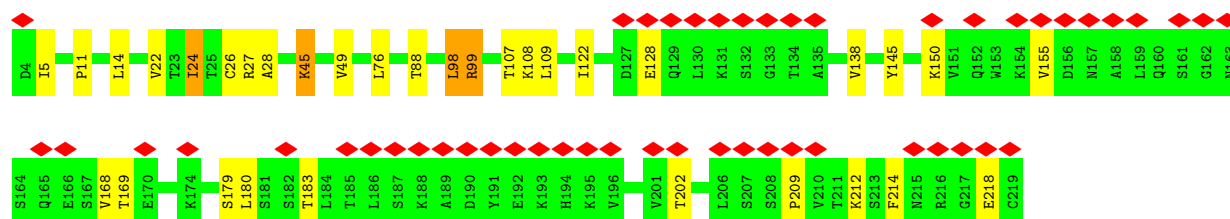
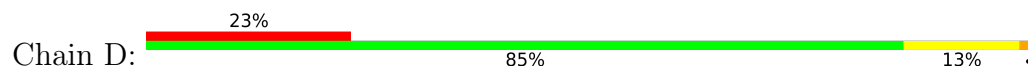
• Molecule 1: Spike glycoprotein



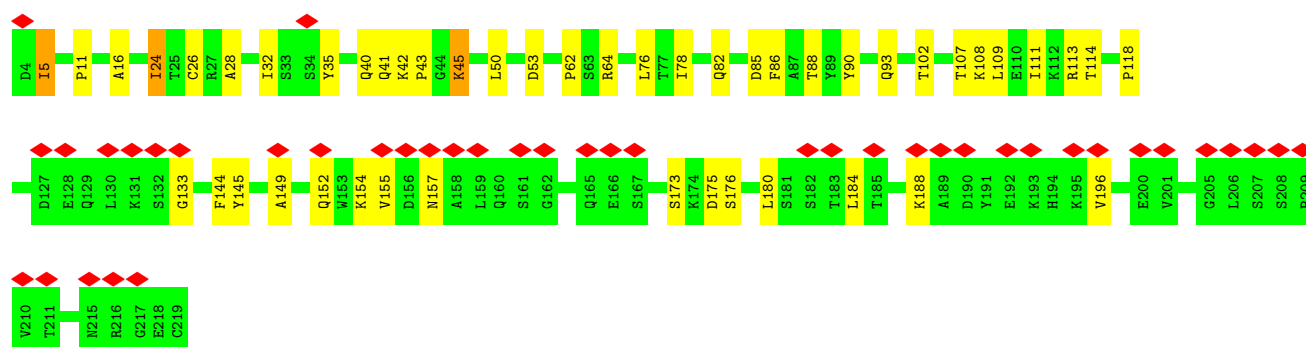
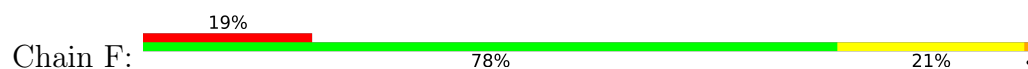




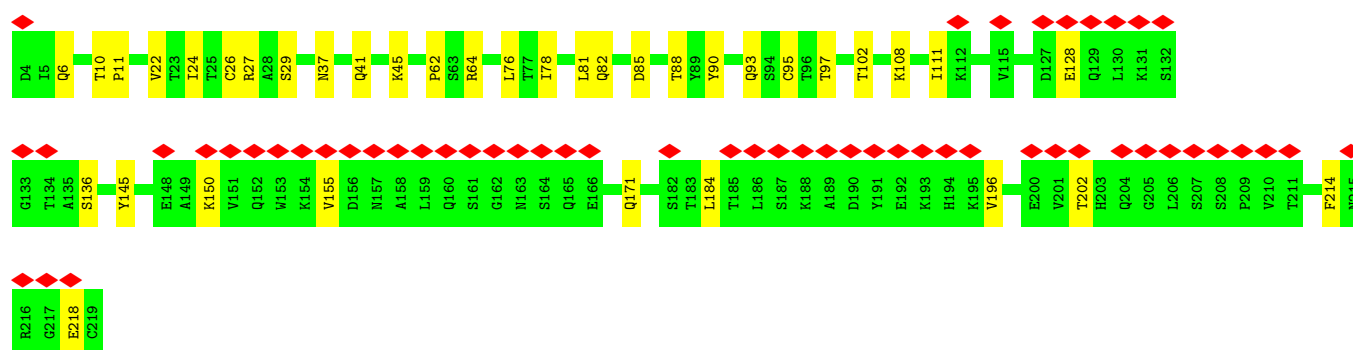
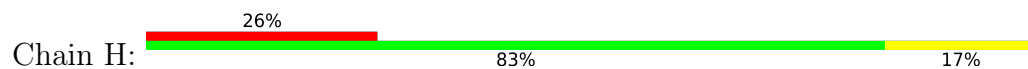
- Molecule 2: 111 L



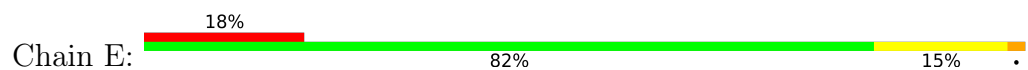
- Molecule 2: 111 L

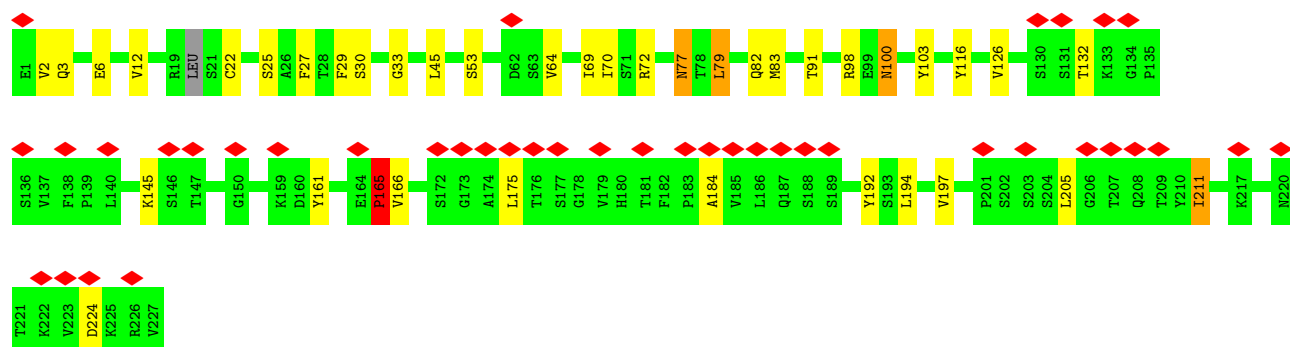


- Molecule 2: 111 L

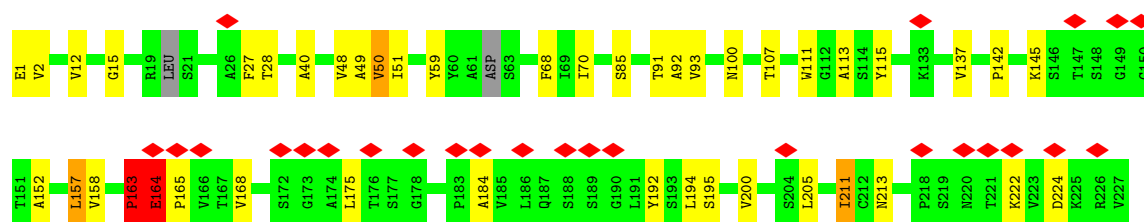
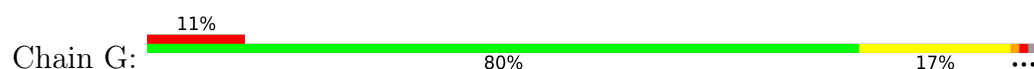


- Molecule 3: 111 H

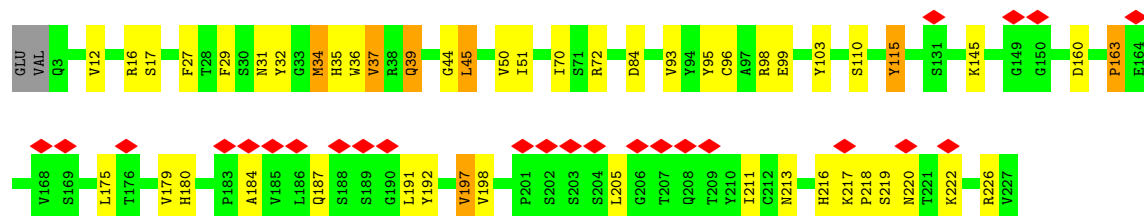
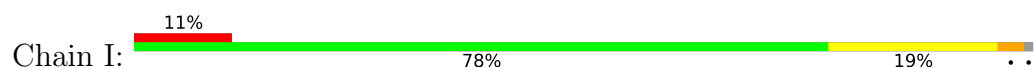




• Molecule 3: 111 H



• Molecule 3: 111 H



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1463548	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)	1500	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.048	Depositor
Minimum map value	-0.012	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.004	Depositor
Map size (\AA)	351.328, 351.328, 351.328	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.0979, 1.0979, 1.0979	Depositor

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.66	0/8490	0.64	1/11580 (0.0%)
1	B	0.66	0/8621	0.65	1/11751 (0.0%)
1	C	0.66	0/8572	0.66	1/11681 (0.0%)
2	D	0.66	0/1598	0.66	0/2171
2	F	0.65	0/1630	0.65	0/2214
2	H	0.65	0/1628	0.64	0/2210
3	E	0.67	0/1647	0.65	0/2244
3	G	0.66	0/1682	0.69	2/2293 (0.1%)
3	I	0.68	0/1643	0.67	0/2242
All	All	0.66	0/35511	0.65	5/48386 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2
1	C	0	3
All	All	0	5

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	163	PRO	CA-N-CD	-6.87	101.89	111.50
3	G	164	GLU	CB-CA-C	6.11	122.62	110.40
1	B	1166	ALA	CB-CA-C	6.07	119.20	110.10
1	C	725	GLU	CB-CA-C	5.47	121.35	110.40
1	A	794	TYR	CB-CA-C	5.02	120.44	110.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	51[A]	SER	Mainchain
1	B	51[B]	SER	Mainchain
1	C	51[A]	SER	Mainchain
1	C	51[B]	SER	Mainchain
1	C	928	TYR	Mainchain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8325	0	7768	129	0
1	B	8446	0	7904	161	0
1	C	8401	0	7876	113	0
2	D	1574	0	1493	13	0
2	F	1602	0	1510	24	0
2	H	1600	0	1509	15	0
3	E	1615	0	1498	16	0
3	G	1644	0	1542	19	0
3	I	1609	0	1502	46	0
All	All	34816	0	32602	516	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 8.

All (516) 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:B:640:GLY:HA3	1:B:651:LEU:CD2	1.58	1.34
1:A:519:ASN:HB3	1:A:522:GLN:CG	1.59	1.32
1:A:519:ASN:CB	1:A:522:GLN:HG3	1.69	1.21
3:I:216:HIS:CE1	3:I:218:PRO:CG	2.25	1.18
1:B:667:THR:HG22	1:B:744:THR:OG1	1.40	1.17
1:B:640:GLY:HA3	1:B:651:LEU:HD21	1.13	1.12
3:I:216:HIS:HE1	3:I:218:PRO:CB	1.64	1.10
1:B:664:ASP:CG	1:B:667:THR:OG1	1.91	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:217:LYS:O	3:I:220:ASN:OD1	1.71	1.07
3:I:216:HIS:CE1	3:I:218:PRO:HB2	1.90	1.07
3:I:216:HIS:HE1	3:I:218:PRO:HB2	1.12	1.06
3:I:216:HIS:CE1	3:I:218:PRO:CD	2.40	1.05
3:I:216:HIS:HB3	3:I:219:SER:HB3	1.33	1.05
3:I:216:HIS:CE1	3:I:218:PRO:CB	2.40	1.05
3:I:216:HIS:CE1	3:I:218:PRO:HD2	1.91	1.05
1:B:640:GLY:CA	1:B:651:LEU:HD21	1.89	1.03
1:B:640:GLY:CA	1:B:651:LEU:CD2	2.37	1.02
3:I:216:HIS:CE1	3:I:218:PRO:HG2	1.93	1.02
1:B:664:ASP:OD1	1:B:667:THR:OG1	1.76	1.00
1:B:359:GLY:O	1:B:361:TYR:CE2	2.14	1.00
1:B:332:TYR:HB2	1:B:334:ARG:HH12	1.29	0.94
1:B:359:GLY:O	1:B:361:TYR:CD2	2.23	0.91
3:I:216:HIS:ND1	3:I:218:PRO:HD2	1.84	0.90
3:I:216:HIS:ND1	3:I:219:SER:N	2.22	0.88
2:F:53:ASP:OD2	3:G:111:TRP:CH2	2.28	0.86
1:B:639:VAL:O	1:B:651:LEU:HD23	1.76	0.86
3:I:35:HIS:HB3	3:I:50:VAL:HA	1.57	0.86
1:B:640:GLY:HA3	1:B:651:LEU:HD22	1.56	0.83
1:B:667:THR:CG2	1:B:744:THR:OG1	2.28	0.78
3:I:36:TRP:HA	3:I:96:CYS:HA	1.64	0.78
1:B:332:TYR:HB2	1:B:334:ARG:NH1	1.98	0.77
1:B:332:TYR:CB	1:B:334:ARG:HH12	1.97	0.77
1:B:667:THR:O	1:B:668:LYS:HB2	1.84	0.77
1:B:330:ASP:OD2	1:B:334:ARG:NH2	2.17	0.76
1:A:383:CYS:H	1:A:407:CYS:HB2	1.48	0.76
1:C:70:THR:HA	1:C:324:LEU:HA	1.67	0.75
1:A:522:GLN:HE21	1:A:522:GLN:HA	1.51	0.75
1:C:189:PRO:HB2	1:C:197:ALA:HB2	1.69	0.73
1:B:347:LEU:HD13	1:B:361:TYR:CZ	2.24	0.73
1:B:651:LEU:HD23	1:B:651:LEU:H	1.54	0.72
1:C:220:ASN:HB3	1:C:223:ALA:HB2	1.72	0.71
1:A:1151:SER:HB2	1:A:1170:GLY:HA2	1.72	0.71
3:I:216:HIS:O	3:I:220:ASN:N	2.22	0.70
3:G:213:ASN:HB3	3:G:222:LYS:HE2	1.73	0.70
1:B:253:TRP:H	1:B:268:SER:HB3	1.56	0.70
1:A:167:HIS:H	1:A:186:ILE:HG12	1.57	0.69
1:A:34:ASP:OD2	1:A:36:GLN:NE2	2.24	0.69
2:H:95:CYS:SG	2:H:97:THR:OG1	2.49	0.69
1:C:1166:ALA:HB3	1:C:1167:PRO:HD3	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:781:SER:HB3	1:B:1146:HIS:HB3	1.75	0.69
1:B:358:SER:HB3	1:B:665:LYS:H	1.58	0.68
1:A:737:ALA:HA	1:B:938:LEU:HA	1.75	0.68
1:C:498:SER:HB3	1:C:534:VAL:HG23	1.75	0.68
1:A:898:VAL:HG21	1:A:1131:PRO:HG2	1.73	0.68
1:C:631:VAL:HG13	1:C:639:VAL:HB	1.74	0.68
1:A:519:ASN:HB3	1:A:522:GLN:HG3	0.74	0.67
1:B:1053:ASP:HB2	1:B:1063:GLN:HA	1.77	0.67
1:A:522:GLN:HA	1:A:522:GLN:NE2	2.09	0.67
1:A:968:ALA:HB1	1:C:770:VAL:HG11	1.76	0.67
3:I:216:HIS:HB3	3:I:219:SER:CB	2.20	0.67
1:B:347:LEU:HB2	1:B:361:TYR:CD1	2.29	0.66
3:I:35:HIS:CD2	3:I:99:GLU:HG2	2.30	0.66
2:H:6:GLN:H	2:H:29:SER:HB2	1.59	0.66
1:B:783:PRO:HB3	1:B:1143:PRO:HB3	1.78	0.66
1:A:405:THR:HB	1:A:582:ASN:HD21	1.60	0.66
1:A:741:THR:HA	1:A:759:LEU:HA	1.78	0.65
1:B:620:CYS:HB2	1:B:648:TYR:HE2	1.62	0.64
3:E:72:ARG:HB3	3:E:79:LEU:HD12	1.79	0.64
1:B:328:SER:HB2	1:B:330:ASP:OD1	1.96	0.64
1:A:253:TRP:H	1:A:268:SER:HB3	1.61	0.64
1:C:738:LEU:N	1:C:739:PRO:HD3	2.12	0.64
1:B:667:THR:HG22	1:B:744:THR:HG1	1.63	0.63
1:A:498:SER:HB3	1:A:534:VAL:HG23	1.81	0.63
3:G:211:ILE:HD11	3:G:224:ASP:HB2	1.80	0.63
1:C:1165:ILE:HG23	1:C:1169:ASN:HA	1.79	0.63
1:B:1197:ILE:HD13	1:C:1189:SER:HB3	1.80	0.63
2:D:5:ILE:HD11	2:D:28:ALA:HB1	1.80	0.63
1:C:731:LEU:HD22	1:C:742:PRO:HA	1.80	0.63
3:I:216:HIS:HE1	3:I:218:PRO:CG	1.81	0.63
1:B:639:VAL:O	1:B:651:LEU:CD2	2.47	0.62
3:G:142:PRO:HG2	3:G:145:LYS:HB3	1.79	0.62
1:A:34:ASP:CG	1:A:36:GLN:HE21	2.03	0.62
1:B:49:ASP:HB3	1:B:52:LYS:HG2	1.82	0.62
1:A:46:ARG:HD2	1:A:314:TYR:CD1	2.33	0.62
3:I:216:HIS:CG	3:I:218:PRO:HD2	2.34	0.62
1:C:904:GLY:HA3	1:C:908:GLY:HA3	1.81	0.62
1:A:102:VAL:HG13	1:A:299:SER:HB2	1.82	0.61
1:C:831:ILE:HG23	1:C:1082:VAL:HG21	1.81	0.61
1:A:903:PRO:HB3	1:A:924:ILE:HG12	1.83	0.61
1:B:258:GLN:HG2	1:B:263:VAL:HG22	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:35:TYR:HA	2:F:53:ASP:HA	1.83	0.61
1:C:807:LYS:HG2	1:C:821:LEU:HB3	1.83	0.60
3:I:27:PHE:HZ	3:I:32:TYR:HB2	1.65	0.60
1:A:626:ARG:HH22	1:A:645:ASP:HA	1.65	0.60
1:C:1058:LEU:HD12	1:C:1062:GLU:HB3	1.82	0.60
1:A:731:LEU:H	1:A:742:PRO:HB3	1.67	0.60
1:C:898:VAL:HG21	1:C:1131:PRO:HG2	1.83	0.60
1:A:189:PRO:HB2	1:A:197:ALA:HB2	1.82	0.60
1:C:439:SER:HB3	1:C:579:THR:OG1	2.01	0.60
1:A:210:PRO:HA	1:A:214:CYS:HB2	1.83	0.60
3:I:35:HIS:HE2	3:I:99:GLU:H	1.48	0.60
1:B:664:ASP:OD2	1:B:667:THR:OG1	2.20	0.59
1:B:601:GLY:C	1:B:603:CYS:H	2.04	0.59
3:G:184:ALA:HB1	3:G:192:TYR:HB3	1.83	0.59
1:A:341:PHE:HE1	1:A:696:MET:HB2	1.67	0.59
1:B:351:TYR:HE2	1:B:361:TYR:HE1	1.50	0.59
1:A:63:THR:HG23	1:C:631:VAL:HB	1.85	0.59
1:A:443:LEU:HD22	1:A:573:ILE:HG13	1.85	0.59
1:A:804:VAL:HG11	1:A:1078:LEU:HD11	1.85	0.58
1:A:127:THR:HA	1:A:139:THR:HA	1.84	0.58
1:C:807:LYS:HA	1:C:821:LEU:HD13	1.86	0.58
1:A:36:GLN:HB3	1:A:39:PHE:HD2	1.69	0.58
2:F:24:ILE:HG23	2:F:76:LEU:HB3	1.86	0.58
1:B:731:LEU:H	1:B:742:PRO:HB3	1.69	0.58
1:B:1053:ASP:HB3	1:B:1057:ARG:HB2	1.85	0.58
2:H:93:GLN:H	2:H:102:THR:HB	1.68	0.58
3:I:213:ASN:HB3	3:I:222:LYS:HE2	1.85	0.58
1:B:347:LEU:HD13	1:B:361:TYR:CE1	2.39	0.57
1:A:57:ILE:H	1:A:75:PHE:HE1	1.51	0.57
1:B:985:ILE:HG23	1:B:1180:ILE:HG21	1.86	0.57
1:C:449:PRO:HG3	1:C:561:VAL:HG21	1.86	0.57
3:I:34:MET:HG2	3:I:98:ARG:HB3	1.86	0.57
1:A:955:ILE:HA	1:A:958:VAL:HG12	1.87	0.57
1:B:664:ASP:OD1	1:B:667:THR:CB	2.51	0.57
1:B:692:SER:HB2	1:B:697:LEU:HB2	1.86	0.57
1:B:797:THR:HG21	1:B:1134:LEU:HB2	1.87	0.57
1:B:981:ASN:HD21	1:B:988:GLN:HB2	1.69	0.57
1:B:1152:ALA:H	1:B:1172:PHE:HA	1.69	0.57
1:C:428:ILE:HB	1:C:476:PRO:HB3	1.86	0.57
1:C:891:GLU:HG2	1:C:1128:VAL:HB	1.85	0.57
2:H:24:ILE:HG23	2:H:76:LEU:HB3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:664:ASP:CG	1:B:667:THR:HG1	1.94	0.57
1:B:1201:ASN:HD22	1:B:1204:TYR:HB2	1.70	0.56
1:A:983:VAL:HG22	1:A:1111:SER:HB3	1.86	0.56
1:B:90:GLY:HA2	1:B:101:PHE:HB2	1.86	0.56
1:B:1166:ALA:CB	1:B:1167:PRO:HD3	2.36	0.56
3:G:164:GLU:CB	3:G:165:PRO:HD3	2.36	0.56
1:B:93:THR:HG23	1:B:98:GLN:HG2	1.87	0.56
1:B:1200:LEU:O	1:B:1201:ASN:HB2	2.06	0.56
2:F:40:GLN:HE21	2:F:50:LEU:HD21	1.71	0.56
3:I:51:ILE:HB	3:I:70:ILE:HD12	1.87	0.56
1:A:329:VAL:HG12	1:C:623:VAL:HG11	1.87	0.56
1:A:790:VAL:HG21	1:A:1119:GLN:HB3	1.87	0.55
1:B:401:ARG:HD3	1:B:521:ASN:HB2	1.88	0.55
1:C:167:HIS:H	1:C:186:ILE:HA	1.71	0.55
3:I:216:HIS:ND1	3:I:218:PRO:CD	2.56	0.55
1:B:351:TYR:CE2	1:B:361:TYR:HE1	2.24	0.55
1:B:640:GLY:N	1:B:651:LEU:HD21	2.22	0.55
3:I:37:VAL:HG12	3:I:95:TYR:HB2	1.87	0.55
1:A:783:PRO:HB3	1:A:1143:PRO:HB3	1.86	0.55
3:G:137:VAL:HA	3:G:158:VAL:HA	1.89	0.55
1:A:84:MET:HG2	1:A:314:TYR:CE1	2.41	0.55
1:C:30:CYS:SG	1:C:231:TYR:HA	2.46	0.55
1:C:538:GLY:H	1:C:559:SER:HA	1.71	0.55
1:A:120:ILE:HG21	1:A:145:PRO:HD3	1.88	0.55
1:B:853:VAL:HG13	1:B:951:LEU:HG	1.89	0.55
1:B:1057:ARG:O	1:B:1059:ASP:N	2.40	0.55
2:F:145:TYR:CD1	2:F:176:SER:HB2	2.42	0.55
1:B:772:GLN:HA	1:C:969:ALA:HB1	1.89	0.55
1:A:118:VAL:HG11	1:A:145:PRO:HG3	1.88	0.54
1:C:325:LEU:HB3	1:C:333:ILE:HD13	1.88	0.54
2:F:16:ALA:O	2:F:111:ILE:HA	2.06	0.54
1:B:1130:ALA:HB2	1:B:1135:TYR:HB2	1.88	0.54
1:B:957:GLY:HA3	1:B:968:ALA:HA	1.89	0.54
1:B:524:SER:HB3	1:B:527:VAL:HG23	1.90	0.54
1:A:653:ALA:HB3	1:B:928:TYR:HB3	1.90	0.54
1:B:693:THR:HG22	1:B:694:ARG:H	1.73	0.54
1:A:906:MET:HG3	1:C:678:ALA:HA	1.88	0.54
1:C:619:ASN:HB3	1:C:647:ASN:HD22	1.72	0.53
1:A:84:MET:HG2	1:A:314:TYR:HE1	1.72	0.53
1:A:478:CYS:HB2	1:A:573:ILE:HB	1.90	0.53
1:C:440:SER:HB3	1:C:576:GLN:HB2	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:15:GLY:HA2	3:G:85:SER:HA	1.89	0.53
3:G:164:GLU:HB3	3:G:165:PRO:HD3	1.91	0.53
1:A:635:TYR:HD1	1:B:1042:ASN:HD21	1.55	0.53
1:B:263:VAL:HB	1:B:284:LEU:HB3	1.90	0.53
1:A:90:GLY:HA2	1:A:101:PHE:H	1.73	0.53
1:C:617:PHE:CD1	1:C:649:TYR:HB3	2.44	0.53
3:E:3:GLN:HB2	3:E:25:SER:HB2	1.90	0.53
1:B:1153:TYR:HA	1:B:1165:ILE:HG23	1.91	0.53
1:C:742:PRO:O	1:C:758:ARG:HB2	2.08	0.53
1:A:46:ARG:HD2	1:A:314:TYR:CE1	2.44	0.52
1:B:664:ASP:OD1	1:B:667:THR:CG2	2.56	0.52
3:E:33:GLY:HA2	3:E:72:ARG:HH22	1.74	0.52
2:F:111:ILE:HG22	2:F:145:TYR:OH	2.09	0.52
1:A:742:PRO:O	1:A:758:ARG:HB3	2.09	0.52
2:F:152:GLN:HE22	2:F:154:LYS:HB3	1.73	0.52
1:A:453:LYS:HG2	1:A:481:LEU:HD23	1.92	0.52
1:B:776:SER:HB2	1:B:1206:ALA:HB1	1.91	0.52
1:A:831:ILE:HG23	1:A:1082:VAL:HG21	1.91	0.52
1:B:33:VAL:H	1:B:202:THR:HG21	1.75	0.52
1:B:169:LEU:HA	1:B:184:TYR:HA	1.91	0.52
1:B:498:SER:HB3	1:B:534:VAL:HG23	1.92	0.52
1:B:983:VAL:HG22	1:B:1121:THR:HB	1.91	0.52
1:B:105:TYR:HB3	1:B:295:ILE:HG23	1.92	0.52
1:B:601:GLY:C	1:B:603:CYS:N	2.63	0.52
1:C:937:PRO:O	1:C:938:LEU:HB3	2.10	0.52
1:A:869:PHE:O	1:A:871:LEU:HG	2.10	0.51
1:C:838:ALA:HB2	1:C:1086:LEU:HD23	1.92	0.51
2:F:118:PRO:HB3	2:F:144:PHE:HB3	1.93	0.51
1:B:428:ILE:HB	1:B:476:PRO:HB3	1.92	0.51
1:B:794:TYR:HE1	1:B:1133:GLY:HA3	1.74	0.51
1:C:1166:ALA:HB2	1:C:1194:PRO:HD3	1.93	0.51
3:E:145:LYS:HB3	3:E:205:LEU:HD13	1.91	0.51
1:A:36:GLN:HB3	1:A:39:PHE:CD2	2.45	0.51
1:C:1165:ILE:HD13	1:C:1170:GLY:H	1.75	0.51
1:B:725:GLU:HB2	1:B:761:SER:HB2	1.93	0.51
1:B:664:ASP:OD1	1:B:667:THR:HG23	2.11	0.50
2:F:111:ILE:HD13	2:F:173:SER:O	2.11	0.50
3:I:217:LYS:HB3	3:I:218:PRO:HD3	1.93	0.50
1:A:366:PHE:HB3	1:A:689:TYR:HB3	1.94	0.50
2:D:24:ILE:HG23	2:D:76:LEU:HB3	1.94	0.50
3:G:51:ILE:HB	3:G:70:ILE:HD12	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LEU:HA	1:A:293:TYR:HA	1.93	0.50
1:A:625:VAL:HG23	1:B:63:THR:HB	1.94	0.50
1:C:420:VAL:HG11	1:C:423:PHE:HD2	1.77	0.50
1:C:458:VAL:HG13	1:C:471:GLN:HB3	1.94	0.50
1:C:493:LYS:HG2	1:C:567:LEU:HB2	1.93	0.50
1:A:990:LEU:HB2	1:A:1179:ARG:HD3	1.93	0.49
1:B:266:PHE:HA	1:B:280:GLN:HA	1.94	0.49
1:B:634:ALA:HB2	1:C:69:ILE:HG21	1.94	0.49
1:A:63:THR:HB	1:C:628:GLN:HE21	1.76	0.49
1:B:374:VAL:H	1:B:606:TYR:HA	1.77	0.49
1:C:442:ILE:HB	1:C:574:THR:HB	1.94	0.49
1:B:330:ASP:OD1	1:B:331:GLY:N	2.46	0.49
2:H:22:VAL:HG22	2:H:81:LEU:HD11	1.94	0.49
2:H:62:PRO:HB2	2:H:64:ARG:HG2	1.94	0.49
1:B:484:VAL:HG11	1:B:491:ILE:HG21	1.94	0.49
1:B:1164:CYS:HB2	1:B:1196:PRO:HD3	1.94	0.49
1:C:631:VAL:CG1	1:C:639:VAL:HB	2.41	0.49
1:B:500:ILE:HA	1:B:558:GLY:HA2	1.94	0.49
1:A:869:PHE:HB3	1:A:871:LEU:HD21	1.94	0.49
1:B:632:TYR:CD1	1:C:62:ARG:HG2	2.48	0.49
1:C:710:PRO:HB3	1:C:717:LEU:HD11	1.93	0.49
1:A:429:SER:HB2	1:B:1058:LEU:HA	1.92	0.49
1:A:714:VAL:O	1:A:715:LEU:C	2.51	0.49
1:B:332:TYR:CB	1:B:334:ARG:NH1	2.66	0.49
2:F:5:ILE:HD11	2:F:28:ALA:HB1	1.95	0.49
2:F:113:ARG:HG2	2:F:114:THR:H	1.77	0.49
1:A:500:ILE:HA	1:A:558:GLY:HA2	1.94	0.49
1:B:794:TYR:CZ	1:B:796:GLN:HA	2.48	0.48
2:D:202:THR:HG23	2:D:209:PRO:HG3	1.95	0.48
2:F:88:THR:HA	2:F:107:THR:O	2.13	0.48
1:B:738:LEU:HD22	1:C:940:ASP:HB2	1.94	0.48
1:C:930:ALA:C	1:C:932:TYR:H	2.16	0.48
1:A:118:VAL:HA	1:A:314:TYR:O	2.13	0.48
1:A:411:LEU:HD12	1:A:434:ALA:HB2	1.95	0.48
1:A:617:PHE:HB3	1:A:649:TYR:HB3	1.96	0.48
1:C:625:VAL:HG12	1:C:627:GLN:H	1.79	0.48
2:D:150:LYS:HB2	2:D:202:THR:HB	1.95	0.48
2:F:62:PRO:HB2	2:F:64:ARG:HG2	1.95	0.48
1:B:518:VAL:HG12	1:B:524:SER:HB2	1.96	0.48
1:A:88:SER:HB2	1:A:131:ILE:HB	1.96	0.48
1:B:448:TYR:CZ	1:B:456:LEU:HB2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:794:TYR:HE2	1:C:1133:GLY:HA3	1.79	0.48
3:G:152:ALA:HB3	3:G:205:LEU:HD12	1.95	0.48
1:A:694:ARG:HH21	1:A:696:MET:HG2	1.79	0.48
1:B:1036:LEU:O	1:B:1040:LEU:HG	2.14	0.48
1:C:959:GLY:HA3	1:C:975:SER:HB3	1.94	0.48
3:I:160:ASP:HA	3:I:191:LEU:HB3	1.95	0.48
3:I:184:ALA:HB1	3:I:192:TYR:HB3	1.94	0.48
1:A:193:ASN:HB3	3:E:103:TYR:CE1	2.48	0.48
1:A:497:TYR:HB2	1:A:561:VAL:HB	1.96	0.48
1:B:358:SER:HA	1:B:663:TYR:HB3	1.94	0.48
1:C:375:VAL:H	1:C:597:ALA:HA	1.78	0.48
3:I:39:GLN:HG3	3:I:45:LEU:HD22	1.95	0.48
1:A:741:THR:HG23	1:A:759:LEU:HB3	1.95	0.48
1:B:114:ASN:HB2	1:B:318:LEU:O	2.13	0.48
1:C:425:CYS:HA	1:C:478:CYS:HA	1.96	0.48
1:A:679:CYS:HB3	1:A:717:LEU:HD13	1.96	0.47
1:A:930:ALA:O	1:A:931:GLY:C	2.52	0.47
1:C:735:LEU:HD23	1:C:739:PRO:HG2	1.96	0.47
2:H:41:GLN:HB3	2:H:90:TYR:HE1	1.79	0.47
1:A:405:THR:HB	1:A:582:ASN:ND2	2.28	0.47
1:C:714:VAL:HG13	1:C:715:LEU:H	1.78	0.47
1:C:731:LEU:H	1:C:740:ASP:HB3	1.79	0.47
1:B:786:PHE:HE2	1:B:994:GLN:HG2	1.80	0.47
1:B:1152:ALA:HB2	1:B:1173:ILE:HB	1.96	0.47
1:C:794:TYR:CZ	1:C:796:GLN:HA	2.49	0.47
1:C:1168:VAL:HG23	1:C:1178:THR:HG22	1.95	0.47
1:A:179:LEU:HD21	1:A:221:ARG:HA	1.97	0.47
1:A:350:SER:C	1:A:352:GLU:H	2.18	0.47
1:A:1179:ARG:HB3	1:A:1186:TYR:CB	2.45	0.47
1:B:651:LEU:CD2	1:B:651:LEU:H	2.26	0.47
2:D:88:THR:HA	2:D:107:THR:O	2.15	0.47
3:I:115:TYR:CD1	3:I:115:TYR:C	2.87	0.47
1:C:110:LYS:NZ	1:C:295:ILE:HG12	2.29	0.47
3:G:49:ALA:HB1	3:G:70:ILE:HD13	1.97	0.47
1:B:1166:ALA:HB1	1:B:1167:PRO:HD3	1.97	0.47
1:C:1156:CYS:H	1:C:1204:TYR:HA	1.80	0.46
1:A:142:LYS:HA	1:A:311:ALA:HB3	1.96	0.46
1:A:457:SER:HB3	1:A:460:SER:HB3	1.96	0.46
1:B:786:PHE:HA	1:B:1141:TYR:CE2	2.50	0.46
1:C:324:LEU:HD22	1:C:349:CYS:HA	1.96	0.46
1:A:773:LEU:HB2	1:A:779:LYS:HE2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:TYR:HE2	1:B:361:TYR:CE1	2.32	0.46
1:B:450:LEU:HA	1:B:481:LEU:HD21	1.96	0.46
1:B:605:GLU:HA	1:B:614:ARG:HA	1.96	0.46
1:C:374:VAL:HA	1:C:597:ALA:HA	1.97	0.46
1:C:725:GLU:O	1:C:725:GLU:HG2	2.16	0.46
1:C:1156:CYS:HA	1:C:1162:THR:HB	1.98	0.46
3:I:35:HIS:ND1	3:I:50:VAL:HG13	2.31	0.46
1:B:742:PRO:HB2	1:B:758:ARG:HH21	1.80	0.46
1:B:794:TYR:CE1	1:B:1133:GLY:HA3	2.50	0.46
2:F:45:LYS:HE3	2:F:45:LYS:HB3	1.82	0.46
1:C:522:GLN:HE21	1:C:522:GLN:HB3	1.55	0.46
1:C:1166:ALA:CB	1:C:1167:PRO:HD3	2.44	0.46
1:A:519:ASN:CB	1:A:522:GLN:CG	2.55	0.46
1:C:269:ARG:HA	1:C:273:LEU:HB2	1.98	0.46
1:B:347:LEU:HD22	1:B:361:TYR:CD2	2.50	0.46
2:F:93:GLN:H	2:F:102:THR:HB	1.80	0.46
1:A:439:SER:HB2	1:A:577:TYR:HA	1.98	0.46
1:A:485:PRO:HG2	1:A:488:LEU:HB2	1.98	0.46
1:B:678:ALA:HB3	1:B:681:HIS:HB3	1.98	0.46
1:B:616:VAL:H	1:B:652:ARG:H	1.64	0.46
3:I:34:MET:HA	3:I:98:ARG:HA	1.98	0.46
1:A:366:PHE:HB3	1:A:689:TYR:CB	2.45	0.45
1:A:792:GLN:HA	1:A:792:GLN:OE1	2.16	0.45
2:D:122:ILE:HB	2:D:212:LYS:HE2	1.99	0.45
3:G:157:LEU:HD12	3:G:195:SER:HB2	1.98	0.45
1:A:731:LEU:N	1:A:742:PRO:HB3	2.30	0.45
1:A:680:GLU:H	1:A:680:GLU:HG3	1.48	0.45
3:G:100:ASN:HB3	3:G:115:TYR:HA	1.98	0.45
3:I:35:HIS:HD2	3:I:99:GLU:HG2	1.79	0.45
1:B:32:GLU:H	1:B:32:GLU:HG3	1.57	0.45
1:B:616:VAL:HB	1:B:652:ARG:HB2	1.97	0.45
2:F:149:ALA:HB3	2:F:180:LEU:HD22	1.99	0.45
1:B:373:SER:HA	1:B:605:GLU:O	2.17	0.45
2:F:86:PHE:HB3	2:F:109:LEU:HB2	1.99	0.45
1:B:666:GLU:OE2	1:B:666:GLU:HA	2.16	0.45
1:B:970:ILE:HG12	1:B:971:PRO:HD2	1.98	0.45
1:C:344:LEU:HD21	1:C:663:TYR:HA	1.98	0.45
1:A:367:GLU:H	1:A:367:GLU:HG3	1.61	0.45
1:A:858:SER:HB3	1:A:950:SER:HA	1.97	0.45
1:B:720:SER:HB2	1:B:759:LEU:HD21	1.99	0.45
2:F:41:GLN:HB3	2:F:90:TYR:HE2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:960:TRP:O	1:A:961:THR:C	2.56	0.45
1:B:319:GLN:HE21	1:B:321:LEU:HD12	1.82	0.45
3:E:161:TYR:CE1	3:E:192:TYR:HB2	2.52	0.45
1:B:148:MET:HG3	1:B:170:VAL:HG22	1.99	0.45
1:C:1160:ASN:N	1:C:1161:PRO:HD2	2.31	0.45
1:A:428:ILE:HB	1:A:476:PRO:HB3	2.00	0.44
1:B:636:GLN:HB2	1:C:1042:ASN:HA	1.99	0.44
1:C:930:ALA:O	1:C:932:TYR:N	2.48	0.44
2:D:168:VAL:HG22	2:D:180:LEU:HD13	1.99	0.44
3:I:35:HIS:CG	3:I:50:VAL:HG13	2.52	0.44
3:I:115:TYR:CD1	3:I:115:TYR:O	2.70	0.44
3:I:216:HIS:NE2	3:I:218:PRO:HG2	2.30	0.44
1:A:154:GLY:O	1:A:163:ARG:HB3	2.17	0.44
1:A:448:TYR:CZ	1:A:456:LEU:HB2	2.52	0.44
1:C:972:PHE:O	1:C:976:ILE:HG12	2.17	0.44
2:F:42:LYS:HD3	2:F:43:PRO:HD2	2.00	0.44
1:A:742:PRO:HB2	1:A:758:ARG:HH21	1.83	0.44
1:B:951:LEU:O	1:B:955:ILE:HG13	2.18	0.44
1:C:928:TYR:C	1:C:930:ALA:H	2.20	0.44
2:H:136:SER:HA	2:H:184:LEU:O	2.18	0.44
1:A:842:GLN:O	1:A:846:VAL:HG23	2.18	0.44
1:B:478:CYS:HB2	1:B:573:ILE:HB	1.99	0.44
1:B:623:VAL:HG21	1:C:329:VAL:HG12	1.99	0.44
3:G:40:ALA:HB2	3:G:92:ALA:HB2	1.99	0.44
1:A:625:VAL:HB	1:A:627:GLN:HG2	1.99	0.44
1:A:1053:ASP:HA	1:A:1056:GLN:HB2	1.99	0.44
3:E:100:ASN:HB2	3:E:116:TYR:H	1.83	0.44
2:H:88:THR:HB	2:H:108:LYS:HD2	2.00	0.44
1:B:664:ASP:OD1	1:B:664:ASP:O	2.35	0.44
1:A:616:VAL:O	1:A:651:LEU:HA	2.17	0.44
1:B:664:ASP:OD2	1:B:667:THR:HG21	2.18	0.44
2:F:175:ASP:O	2:F:176:SER:OG	2.26	0.44
2:H:111:ILE:HD12	2:H:111:ILE:H	1.82	0.44
3:E:161:TYR:CE1	3:E:165:PRO:HA	2.53	0.44
1:A:131:ILE:HD12	1:A:131:ILE:H	1.82	0.43
3:E:184:ALA:HB1	3:E:192:TYR:HB3	2.00	0.43
1:A:534:VAL:HB	1:A:560:THR:HG22	2.00	0.43
1:A:794:TYR:HB2	1:A:1135:TYR:HA	2.00	0.43
1:C:952:LEU:O	1:C:955:ILE:HG12	2.18	0.43
1:A:102:VAL:HB	1:A:297:PRO:HB2	1.99	0.43
1:B:357:GLU:O	1:B:361:TYR:OH	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:958:VAL:O	1:B:961:THR:HB	2.19	0.43
2:H:90:TYR:HE2	3:I:44:GLY:HA2	1.82	0.43
2:D:14:LEU:O	2:D:109:LEU:HA	2.18	0.43
1:B:64:TYR:HD2	1:B:69:ILE:HD13	1.82	0.43
1:A:366:PHE:HA	1:A:691:ARG:HB3	2.00	0.43
1:A:1121:THR:OG1	1:A:1141:TYR:HB3	2.19	0.43
3:E:211:ILE:HD11	3:E:224:ASP:HB2	2.01	0.43
3:I:37:VAL:N	3:I:95:TYR:O	2.52	0.43
1:A:300:ILE:H	1:A:300:ILE:HD12	1.83	0.43
1:A:794:TYR:CZ	1:A:796:GLN:HA	2.53	0.43
1:B:804:VAL:H	1:B:835:LEU:HD13	1.82	0.43
1:C:841:ARG:HH21	1:C:1089:SER:HB3	1.84	0.43
2:F:154:LYS:HD2	2:F:157:ASN:HA	2.01	0.43
1:B:725:GLU:CD	1:B:758:ARG:HG2	2.39	0.43
1:C:163:ARG:HH11	1:C:187:LEU:HD22	1.83	0.43
1:C:726:ASP:HA	1:C:761:SER:HA	1.99	0.43
3:E:2:VAL:HG13	3:E:27:PHE:HB2	2.01	0.43
1:B:381:VAL:O	1:B:382:GLU:C	2.57	0.43
1:C:404:PHE:HB3	1:C:407:CYS:SG	2.59	0.43
1:C:1013:THR:HB	1:C:1014:THR:H	1.53	0.43
1:B:831:ILE:HG23	1:B:1082:VAL:HG21	2.01	0.42
1:C:112:PHE:CE1	1:C:115:GLY:HA2	2.54	0.42
1:C:786:PHE:HA	1:C:1141:TYR:CE2	2.54	0.42
3:I:175:LEU:HD23	3:I:198:VAL:HG21	2.01	0.42
1:A:123:ALA:HB3	1:A:311:ALA:HA	2.01	0.42
1:A:324:LEU:HD22	1:A:337:ILE:HD12	2.00	0.42
1:A:739:PRO:HB3	1:A:760:ALA:HA	2.01	0.42
1:A:804:VAL:HA	1:A:932:TYR:HA	2.02	0.42
1:C:131:ILE:HD12	1:C:131:ILE:H	1.84	0.42
1:C:226:ASN:O	1:C:230:GLU:HG3	2.19	0.42
2:D:214:PHE:HB2	2:D:218:GLU:OE1	2.19	0.42
1:A:166:ASN:HA	1:A:186:ILE:HG23	2.00	0.42
1:B:269:ARG:HA	1:B:273:LEU:HB2	2.00	0.42
1:B:457:SER:HB3	1:B:460:SER:HB3	2.01	0.42
1:C:83:ASP:HB3	1:C:85:TYR:CE2	2.54	0.42
3:E:30:SER:O	3:E:53:SER:HB2	2.19	0.42
1:A:865:PHE:HB3	1:A:871:LEU:HD12	2.01	0.42
1:B:102:VAL:HB	1:B:297:PRO:HB2	2.01	0.42
1:C:625:VAL:O	1:C:626:ARG:HB2	2.20	0.42
1:B:185:CYS:HB2	1:B:234:LEU:HD12	2.01	0.42
1:B:357:GLU:HA	1:B:665:LYS:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:ILE:HB	1:C:252:GLU:HG3	2.01	0.42
1:C:715:LEU:HB2	1:C:737:ALA:HB2	2.00	0.42
1:C:777:TYR:HB3	1:C:1150:VAL:HG23	2.01	0.42
1:C:1203:LYS:HA	1:C:1203:LYS:HD2	1.91	0.42
1:A:111:GLN:HA	1:A:292:TYR:HA	2.00	0.42
1:A:439:SER:HB3	1:A:579:THR:OG1	2.19	0.42
1:B:328:SER:OG	1:B:334:ARG:HG2	2.18	0.42
1:B:408:ASN:HB3	1:B:587:LYS:HD3	2.02	0.42
1:C:210:PRO:HA	1:C:214:CYS:HB2	2.01	0.42
1:C:226:ASN:HA	1:C:229:LYS:HE3	2.01	0.42
2:D:138:VAL:HG22	2:D:183:THR:HG22	2.02	0.42
2:H:45:LYS:HE3	2:H:45:LYS:HB3	1.87	0.42
1:A:783:PRO:O	1:A:1182:ASP:HA	2.19	0.42
1:B:526:CYS:HB3	1:B:556:ALA:HB2	2.01	0.42
1:B:1160:ASN:N	1:B:1161:PRO:HD2	2.34	0.42
1:C:534:VAL:HB	1:C:560:THR:HG22	2.02	0.42
1:A:324:LEU:HB3	1:A:337:ILE:HB	2.02	0.42
2:D:45:LYS:HE3	2:D:45:LYS:HB3	1.88	0.42
3:I:16:ARG:HE	3:I:17:SER:H	1.66	0.42
1:A:34:ASP:CG	1:A:36:GLN:NE2	2.70	0.42
1:A:102:VAL:CG1	1:A:299:SER:HB2	2.50	0.42
1:A:110:LYS:HE2	1:A:110:LYS:HB2	1.89	0.42
1:A:124:ALA:O	1:A:125:ASN:HB2	2.20	0.42
1:A:875:GLU:OE1	1:A:886:ALA:N	2.52	0.42
1:C:496:LYS:HE2	1:C:560:THR:HG21	2.01	0.42
3:G:50:VAL:HG13	3:G:59:TYR:HB2	2.01	0.42
1:A:639:VAL:O	1:A:651:LEU:HD21	2.20	0.42
1:A:673:LEU:HD21	1:A:715:LEU:HD23	2.02	0.42
1:C:111:GLN:HE21	1:C:111:GLN:HB3	1.61	0.42
3:E:91:THR:HG23	3:E:126:VAL:O	2.20	0.42
2:F:82:GLN:HB2	2:F:85:ASP:HB2	2.02	0.42
2:H:214:PHE:HB2	2:H:218:GLU:OE1	2.19	0.42
1:B:328:SER:O	1:B:331:GLY:N	2.52	0.41
1:B:394:PRO:HG2	1:B:445:TYR:O	2.19	0.41
1:B:509:ASP:O	1:B:510:ASP:HB3	2.20	0.41
1:B:724:VAL:HG12	1:B:760:ALA:HB3	2.02	0.41
1:C:616:VAL:O	1:C:651:LEU:HA	2.19	0.41
1:C:776:SER:HB2	1:C:1206:ALA:HB1	2.01	0.41
3:I:34:MET:H	3:I:72:ARG:HH22	1.67	0.41
3:I:180:HIS:HB2	3:I:197:VAL:HG13	2.02	0.41
1:A:1164:CYS:HB2	1:A:1205:VAL:HG21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:GLN:H	1:B:98:GLN:HG3	1.71	0.41
1:B:442:ILE:HB	1:B:574:THR:HB	2.02	0.41
1:C:154:GLY:O	1:C:163:ARG:HB3	2.19	0.41
2:D:169:THR:HG22	2:D:179:SER:H	1.85	0.41
3:G:205:LEU:HD23	3:G:205:LEU:HA	1.83	0.41
1:B:408:ASN:HA	1:B:585:CYS:O	2.20	0.41
1:B:735:LEU:HD22	1:B:741:THR:HG22	2.02	0.41
1:B:738:LEU:HD12	1:B:738:LEU:HA	1.82	0.41
1:B:988:GLN:HE21	1:B:988:GLN:HB3	1.60	0.41
1:C:1128:VAL:HG23	1:C:1135:TYR:HB3	2.03	0.41
1:C:1201:ASN:HB3	1:C:1206:ALA:C	2.40	0.41
3:E:6:GLU:HA	3:E:22:CYS:HA	2.02	0.41
3:G:2:VAL:HG13	3:G:27:PHE:CG	2.55	0.41
1:B:850:PHE:HD1	1:B:850:PHE:HA	1.79	0.41
3:I:145:LYS:HB2	3:I:205:LEU:HD13	2.02	0.41
3:I:179:VAL:HG22	3:I:198:VAL:HG12	2.01	0.41
1:B:739:PRO:HB3	1:B:760:ALA:HA	2.02	0.41
1:C:873:LEU:HD13	1:C:890:ILE:HD12	2.03	0.41
1:B:344:LEU:O	1:B:347:LEU:HG	2.21	0.41
1:C:53:ALA:HA	1:C:56:ILE:HD12	2.02	0.41
1:C:738:LEU:N	1:C:739:PRO:CD	2.82	0.41
1:A:522:GLN:HE21	1:A:522:GLN:CA	2.18	0.41
1:B:545:LEU:HD21	1:B:554:LEU:HB2	2.02	0.41
1:B:154:GLY:O	1:B:163:ARG:HB3	2.20	0.41
1:B:722:LEU:HD23	1:B:722:LEU:HA	1.81	0.41
1:C:604:VAL:H	1:C:604:VAL:HG22	1.64	0.41
1:A:374:VAL:O	1:A:607:SER:N	2.54	0.41
1:A:1036:LEU:HD13	1:A:1081:PHE:HD1	1.86	0.41
1:B:1052:GLY:HA3	1:B:1066:GLN:HB3	2.03	0.41
3:E:69:ILE:HB	3:E:82:GLN:HB3	2.02	0.41
3:I:217:LYS:N	3:I:218:PRO:CD	2.84	0.41
1:A:786:PHE:HA	1:A:1141:TYR:CE2	2.55	0.41
1:B:119:ARG:HB3	1:B:314:TYR:HB2	2.03	0.41
1:B:357:GLU:HA	1:B:357:GLU:OE1	2.21	0.41
1:B:511:ARG:HD2	1:B:511:ARG:HA	1.82	0.41
1:A:852:SER:HB3	1:A:1100:LYS:HE3	2.03	0.40
1:C:228:PHE:HD1	1:C:228:PHE:HA	1.75	0.40
1:C:1106:CYS:HA	1:C:1111:SER:HB2	2.03	0.40
3:G:113:ALA:HB1	3:G:115:TYR:CD1	2.56	0.40
2:H:82:GLN:HB2	2:H:85:ASP:HB2	2.03	0.40
1:C:90:GLY:HA2	1:C:101:PHE:H	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:682:ILE:HG13	1:C:717:LEU:HD23	2.04	0.40
1:C:835:LEU:HD12	1:C:835:LEU:HA	1.89	0.40
1:C:1039:GLU:O	1:C:1044:PHE:HB2	2.21	0.40
3:E:29:PHE:HB2	3:E:77:ASN:HA	2.03	0.40
2:H:150:LYS:HB2	2:H:202:THR:HB	2.02	0.40
1:A:964:LEU:HB3	1:A:965:SER:H	1.59	0.40
1:A:983:VAL:HG21	1:A:1122:HIS:O	2.20	0.40
1:C:344:LEU:HD23	1:C:344:LEU:HA	1.87	0.40
2:D:98:LEU:HB3	2:D:99:ARG:H	1.73	0.40
2:F:133:GLY:HA2	2:F:188:LYS:HG3	2.02	0.40
1:C:818:GLU:HA	1:C:821:LEU:HB2	2.03	0.40
1:A:92:ALA:H	1:A:307:ARG:HH11	1.69	0.40
1:B:450:LEU:HD12	1:B:568:GLN:HG2	2.03	0.40
1:B:725:GLU:HB3	1:B:728:LYS:HB2	2.03	0.40
1:B:1054:ILE:H	1:B:1054:ILE:HG13	1.60	0.40
1:C:1202:THR:C	1:C:1204:TYR:H	2.24	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1119/1189 (94%)	1052 (94%)	52 (5%)	15 (1%)	10	42
1	B	1128/1189 (95%)	1070 (95%)	48 (4%)	10 (1%)	14	51
1	C	1122/1189 (94%)	1051 (94%)	62 (6%)	9 (1%)	16	54
2	D	214/216 (99%)	209 (98%)	3 (1%)	2 (1%)	14	51
2	F	214/216 (99%)	211 (99%)	2 (1%)	1 (0%)	25	64
2	H	214/216 (99%)	211 (99%)	2 (1%)	1 (0%)	25	64
3	E	222/227 (98%)	218 (98%)	3 (1%)	1 (0%)	25	64

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	G	219/227 (96%)	213 (97%)	4 (2%)	2 (1%)	14	51
3	I	223/227 (98%)	209 (94%)	13 (6%)	1 (0%)	30	68
All	All	4675/4896 (96%)	4444 (95%)	189 (4%)	42 (1%)	17	51

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	599	GLN
1	A	931	GLY
1	A	961	THR
1	B	604	VAL
1	B	1058	LEU
1	B	1166	ALA
1	B	1167	PRO
1	B	1201	ASN
1	C	646	GLY
1	C	737	ALA
1	C	1166	ALA
2	D	98	LEU
3	E	165	PRO
3	G	163	PRO
3	G	164	GLU
3	I	163	PRO
1	A	713	CYS
1	A	867	GLY
1	A	887	ARG
1	A	1040	LEU
1	A	1201	ASN
1	C	600	LEU
1	A	609	TYR
1	B	622	ALA
1	C	713	CYS
1	C	739	PRO
1	B	767	PRO
1	B	1181	VAL
1	C	724	VAL
1	C	887	ARG
1	A	601	GLY
1	A	1168	VAL
1	B	743	SER
2	F	11	PRO

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Mol	Chain	Res	Type
2	H	11	PRO
1	A	45	PRO
1	A	342	ASN
1	C	697	LEU
1	A	860	PRO
2	D	11	PRO
1	A	271	VAL
1	B	931	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	866/1022 (85%)	818 (94%)	48 (6%)	18	40
1	B	892/1022 (87%)	837 (94%)	55 (6%)	15	37
1	C	881/1022 (86%)	818 (93%)	63 (7%)	12	32
2	D	173/192 (90%)	162 (94%)	11 (6%)	14	36
2	F	175/192 (91%)	165 (94%)	10 (6%)	17	39
2	H	175/192 (91%)	165 (94%)	10 (6%)	17	39
3	E	170/191 (89%)	154 (91%)	16 (9%)	7	23
3	G	174/191 (91%)	158 (91%)	16 (9%)	7	24
3	I	169/191 (88%)	152 (90%)	17 (10%)	6	21
All	All	3675/4215 (87%)	3429 (93%)	246 (7%)	16	34

All (246) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ILE
1	A	69	ILE
1	A	93	THR
1	A	155	ASN
1	A	165	PHE
1	A	171	LEU

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Mol	Chain	Res	Type
1	A	174	ASP
1	A	179	LEU
1	A	256	ILE
1	A	314	TYR
1	A	319	GLN
1	A	346	GLN
1	A	347	LEU
1	A	367	GLU
1	A	396	VAL
1	A	417	LEU
1	A	510	ASP
1	A	511	ARG
1	A	598	SER
1	A	616	VAL
1	A	625	VAL
1	A	626	ARG
1	A	639	VAL
1	A	680	GLU
1	A	682	ILE
1	A	711	VAL
1	A	714	VAL
1	A	728	LYS
1	A	736	CYS
1	A	768	ILE
1	A	780	LEU
1	A	782	ILE
1	A	784	THR
1	A	794	TYR
1	A	795	ILE
1	A	799	ILE
1	A	808	GLN
1	A	938	LEU
1	A	960	TRP
1	A	990	LEU
1	A	1033	LEU
1	A	1040	LEU
1	A	1070	LEU
1	A	1075	LEU
1	A	1117	CYS
1	A	1124	VAL
1	A	1125	SER
1	A	1128	VAL

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Mol	Chain	Res	Type
1	B	32	GLU
1	B	63	THR
1	B	98	GLN
1	B	129	THR
1	B	172	LEU
1	B	179	LEU
1	B	239	PHE
1	B	256	ILE
1	B	313	PHE
1	B	321	LEU
1	B	324	LEU
1	B	333	ILE
1	B	411	LEU
1	B	414	LEU
1	B	450	LEU
1	B	502	LYS
1	B	503	CYS
1	B	517	LEU
1	B	529	ILE
1	B	581	THR
1	B	604	VAL
1	B	620	CYS
1	B	631	VAL
1	B	633	ASP
1	B	638	LEU
1	B	642	TYR
1	B	655	VAL
1	B	680	GLU
1	B	681	HIS
1	B	682	ILE
1	B	691	ARG
1	B	711	VAL
1	B	743	SER
1	B	744	THR
1	B	759	LEU
1	B	794	TYR
1	B	795	ILE
1	B	804	VAL
1	B	817	CYS
1	B	821	LEU
1	B	900	ILE
1	B	929	VAL

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Mol	Chain	Res	Type
1	B	938	LEU
1	B	967	PHE
1	B	988	GLN
1	B	1005	LEU
1	B	1013	THR
1	B	1054	ILE
1	B	1066	GLN
1	B	1078	LEU
1	B	1111	SER
1	B	1124	VAL
1	B	1127	VAL
1	B	1134	LEU
1	B	1162	THR
1	C	99	LYS
1	C	111	GLN
1	C	153	VAL
1	C	176	CYS
1	C	179	LEU
1	C	208	HIS
1	C	221	ARG
1	C	227	SER
1	C	228	PHE
1	C	243	TYR
1	C	291	LYS
1	C	295	ILE
1	C	313	PHE
1	C	324	LEU
1	C	333	ILE
1	C	349	CYS
1	C	373	SER
1	C	427	GLN
1	C	510	ASP
1	C	522	GLN
1	C	529	ILE
1	C	569	MET
1	C	598	SER
1	C	603	CYS
1	C	621	THR
1	C	623	VAL
1	C	631	VAL
1	C	637	ASN
1	C	638	LEU

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Mol	Chain	Res	Type
1	C	645	ASP
1	C	670	HIS
1	C	682	ILE
1	C	686	MET
1	C	696	MET
1	C	713	CYS
1	C	714	VAL
1	C	725	GLU
1	C	729	LEU
1	C	757	MET
1	C	794	TYR
1	C	795	ILE
1	C	807	LYS
1	C	817	CYS
1	C	862	ILE
1	C	900	ILE
1	C	929	VAL
1	C	933	LYS
1	C	938	LEU
1	C	951	LEU
1	C	983	VAL
1	C	1012	PHE
1	C	1013	THR
1	C	1039	GLU
1	C	1047	ILE
1	C	1062	GLU
1	C	1069	ARG
1	C	1081	PHE
1	C	1117	CYS
1	C	1124	VAL
1	C	1134	LEU
1	C	1168	VAL
1	C	1174	LYS
1	C	1187	THR
2	D	22	VAL
2	D	24	ILE
2	D	26	CYS
2	D	27	ARG
2	D	45	LYS
2	D	49	VAL
2	D	99	ARG
2	D	108	LYS

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Mol	Chain	Res	Type
2	D	128	GLU
2	D	145	TYR
2	D	155	VAL
3	E	12	VAL
3	E	45	LEU
3	E	64	VAL
3	E	70	ILE
3	E	77	ASN
3	E	79	LEU
3	E	83	MET
3	E	98	ARG
3	E	100	ASN
3	E	132	THR
3	E	165	PRO
3	E	166	VAL
3	E	175	LEU
3	E	194	LEU
3	E	197	VAL
3	E	211	ILE
2	F	5	ILE
2	F	24	ILE
2	F	26	CYS
2	F	32	ILE
2	F	45	LYS
2	F	78	ILE
2	F	108	LYS
2	F	155	VAL
2	F	184	LEU
2	F	196	VAL
3	G	1	GLU
3	G	12	VAL
3	G	28	THR
3	G	48	VAL
3	G	50	VAL
3	G	68	PHE
3	G	91	THR
3	G	93	VAL
3	G	107	THR
3	G	157	LEU
3	G	163	PRO
3	G	168	VAL
3	G	175	LEU

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Mol	Chain	Res	Type
3	G	194	LEU
3	G	200	VAL
3	G	211	ILE
2	H	10	THR
2	H	26	CYS
2	H	27	ARG
2	H	37	ASN
2	H	78	ILE
2	H	128	GLU
2	H	145	TYR
2	H	155	VAL
2	H	171	GLN
2	H	196	VAL
3	I	12	VAL
3	I	29	PHE
3	I	31	ASN
3	I	34	MET
3	I	37	VAL
3	I	39	GLN
3	I	45	LEU
3	I	84	ASP
3	I	93	VAL
3	I	103	TYR
3	I	110	SER
3	I	115	TYR
3	I	163	PRO
3	I	187	GLN
3	I	197	VAL
3	I	211	ILE
3	I	226	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	277	ASN
1	A	280	GLN
1	A	346	GLN
1	A	348	HIS
1	A	398	ASN
1	A	522	GLN
1	A	765	ASN

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Mol	Chain	Res	Type
1	A	785	ASN
1	A	839	ASN
1	A	993	ASN
1	A	999	ASN
1	A	1029	ASN
1	A	1042	ASN
1	A	1079	ASN
1	A	1132	ASN
1	A	1201	ASN
1	B	98	GLN
1	B	261	GLN
1	B	280	GLN
1	B	319	GLN
1	B	981	ASN
1	B	988	GLN
1	B	999	ASN
1	B	1201	ASN
1	C	111	GLN
1	C	114	ASN
1	C	226	ASN
1	C	421	ASN
1	C	522	GLN
1	C	808	GLN
1	C	812	ASN
1	C	1042	ASN
1	C	1169	ASN
2	D	157	ASN
2	D	160	GLN
3	E	100	ASN
2	F	37	ASN
2	F	40	GLN
2	F	105	GLN
2	F	152	GLN
2	F	157	ASN
2	F	160	GLN
3	G	187	GLN
2	H	40	GLN
2	H	105	GLN
2	H	157	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

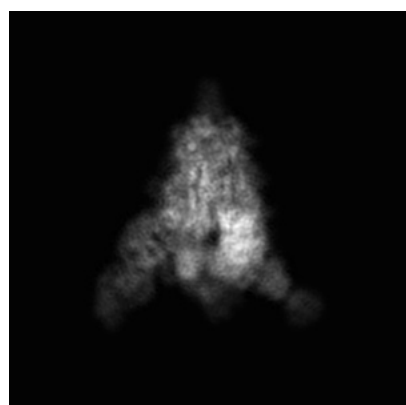
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31744. These allow visual inspection of the internal detail of the map and identification of artifacts.

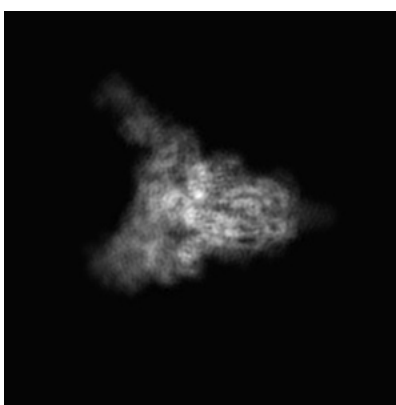
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

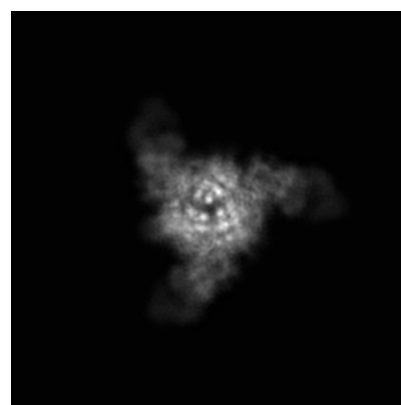
6.1.1 Primary map



X



Y

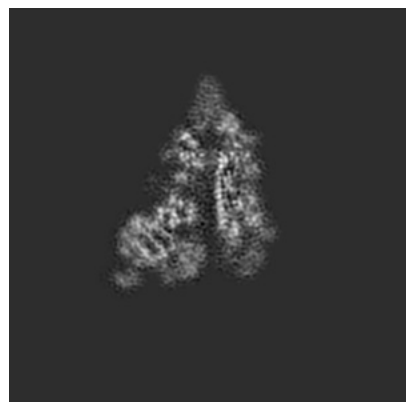


Z

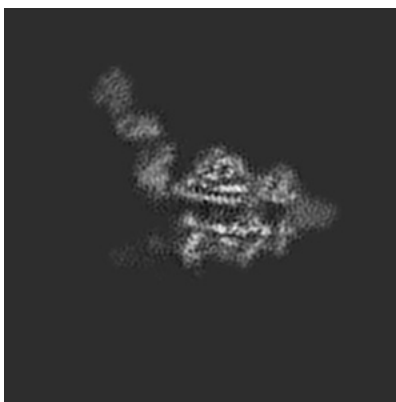
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

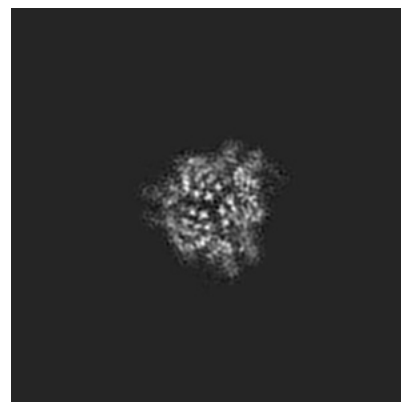
6.2.1 Primary map



X Index: 160



Y Index: 160



Z Index: 160

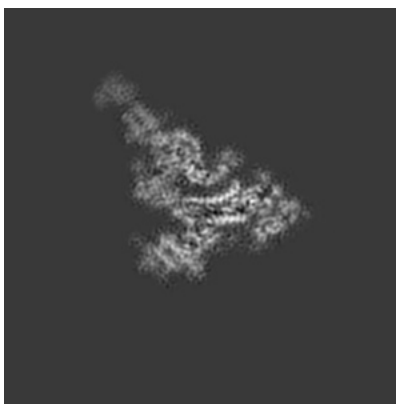
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

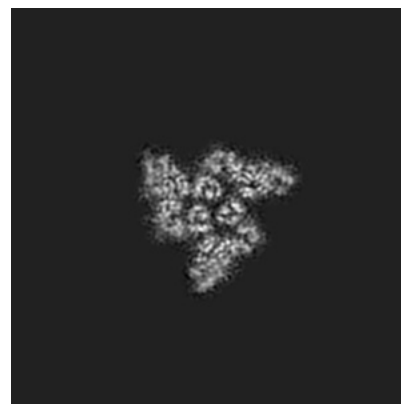
6.3.1 Primary map



X Index: 167



Y Index: 174

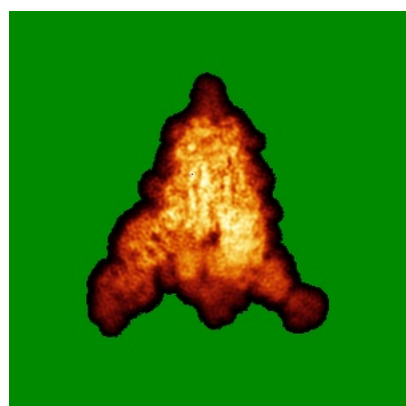


Z Index: 148

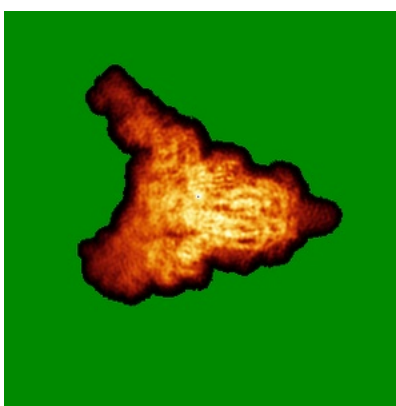
The images above show the largest variance slices of the map in three orthogonal directions.

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

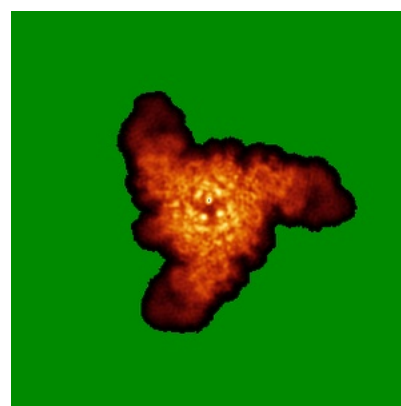
6.4.1 Primary map



X



Y

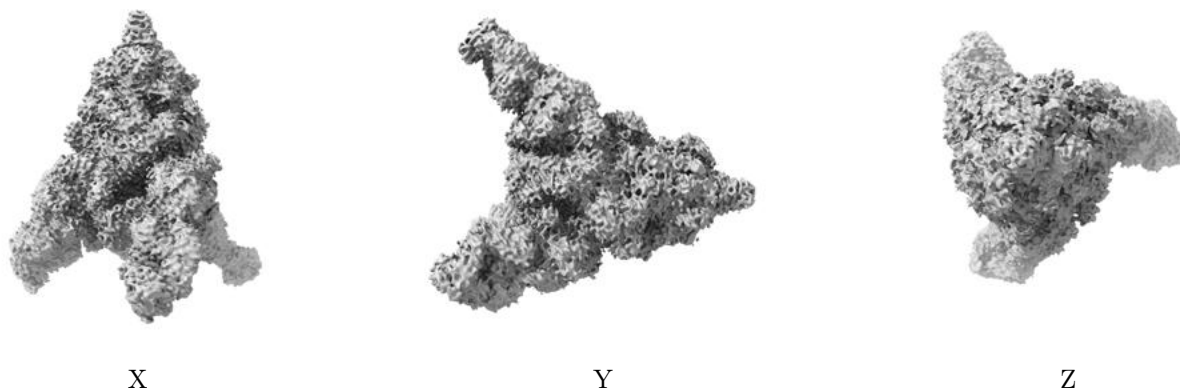


Z

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

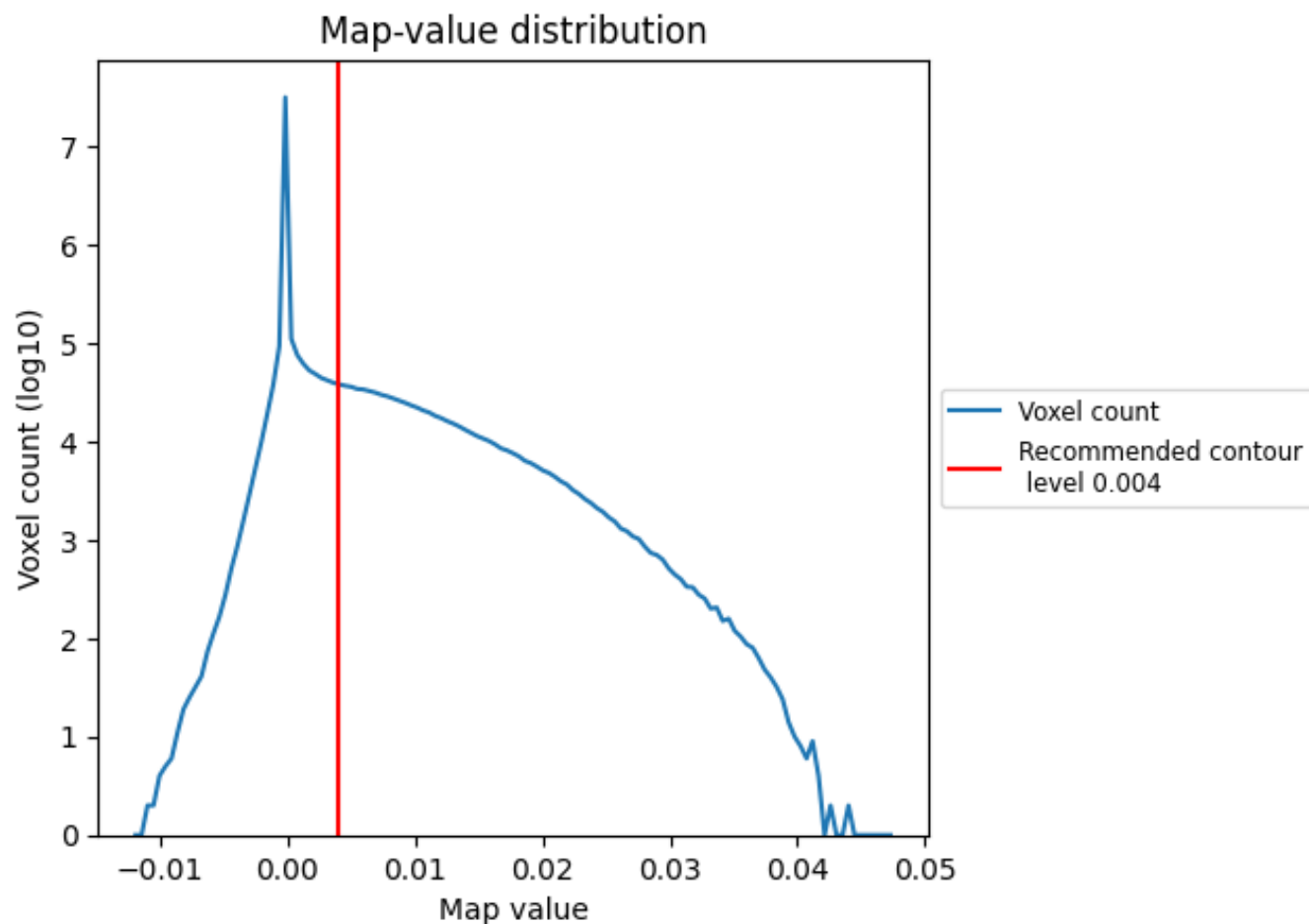
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

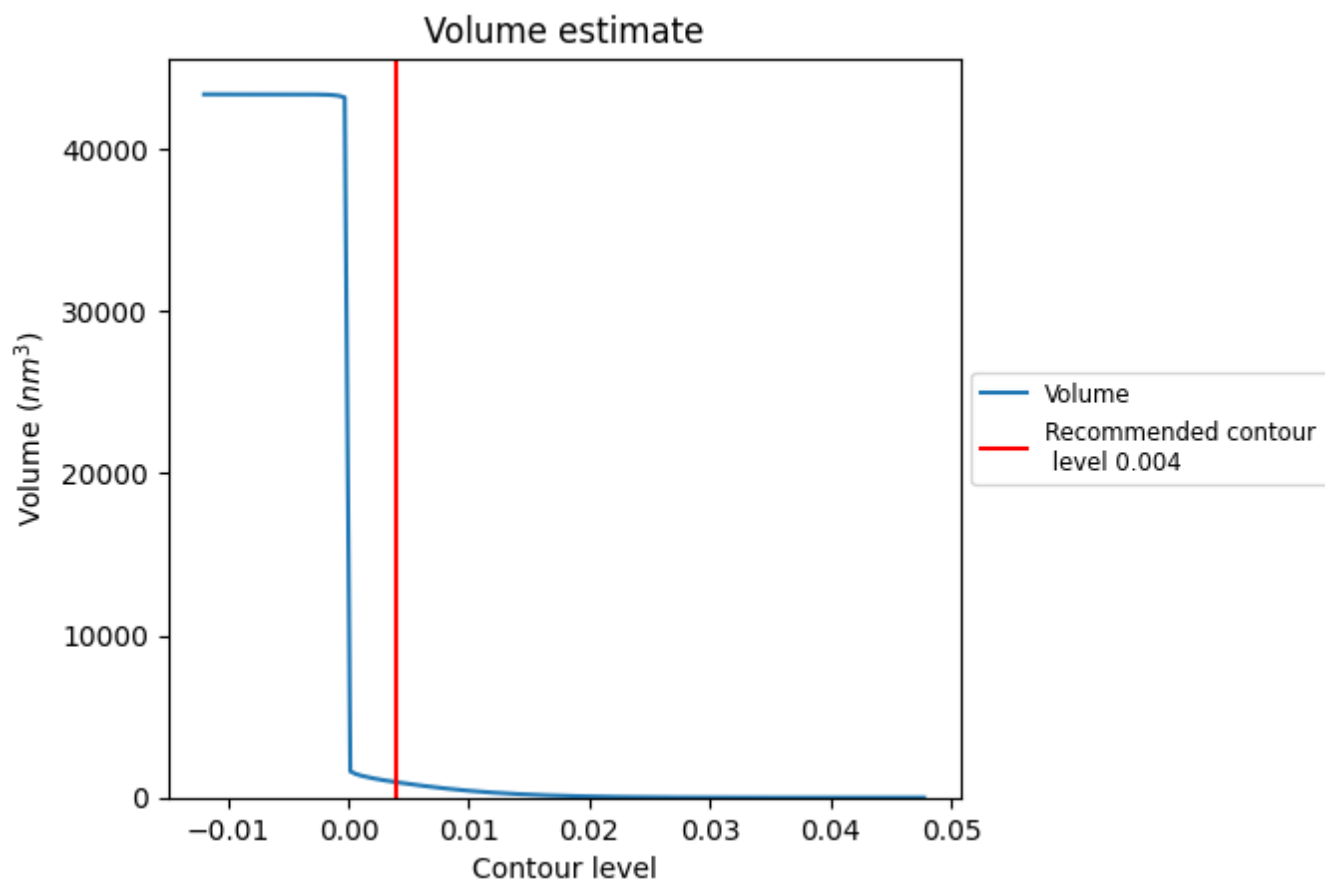
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

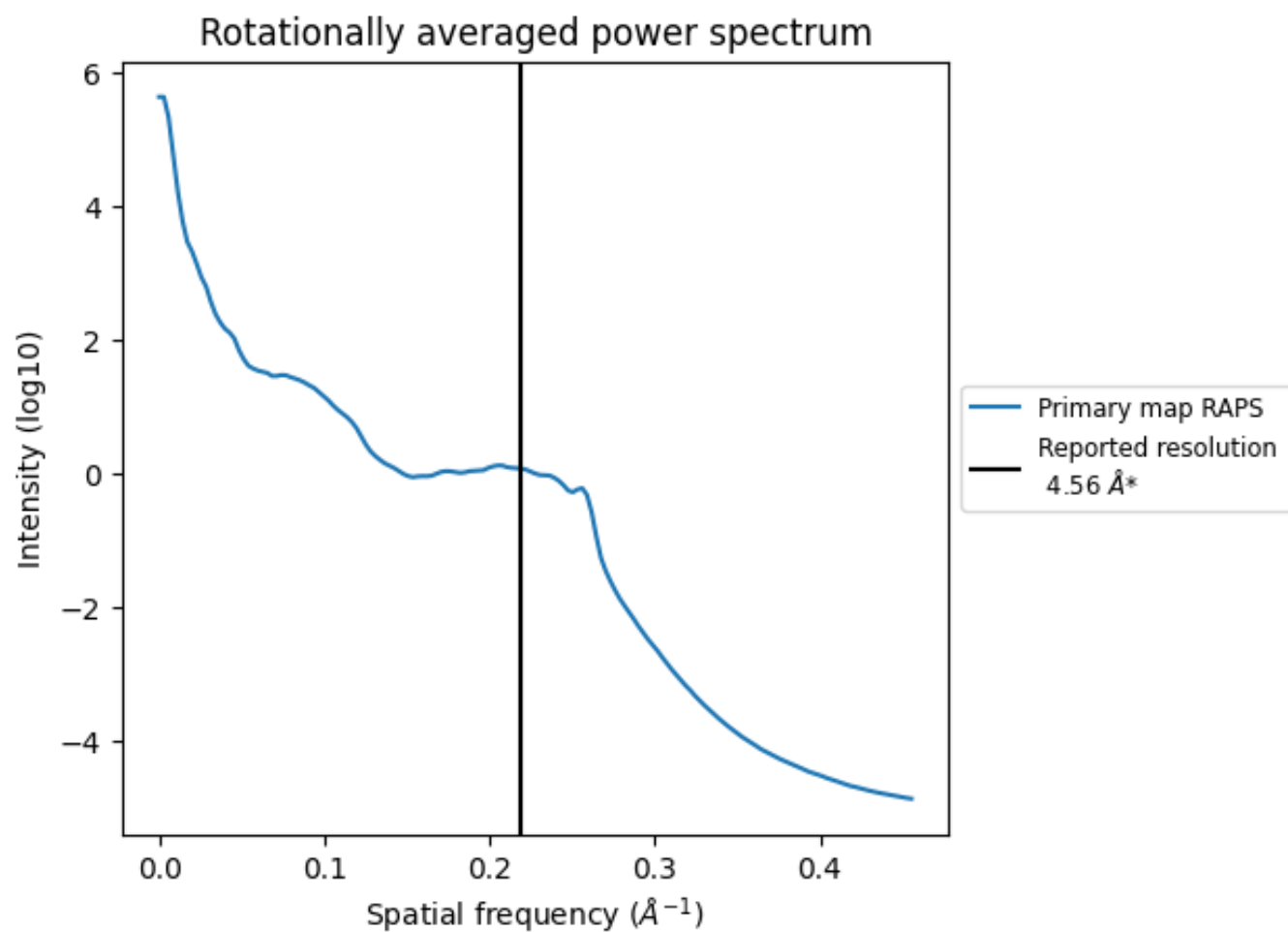
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 949 nm³; this corresponds to an approximate mass of 857 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 0.219 Å⁻¹

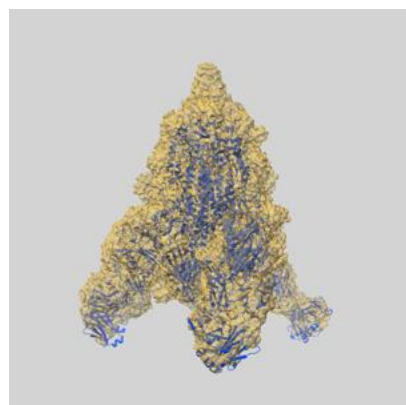
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

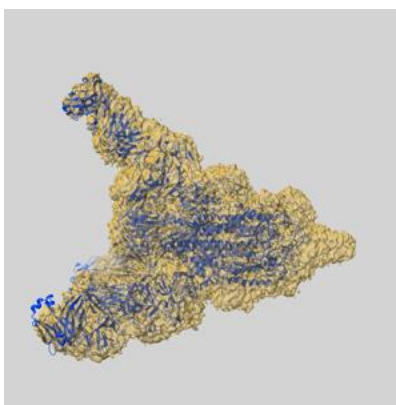
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-31744 and PDB model 7V6O. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

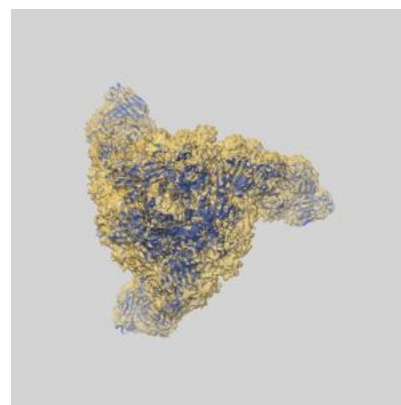
9.1 Map-model overlay [i](#)



X



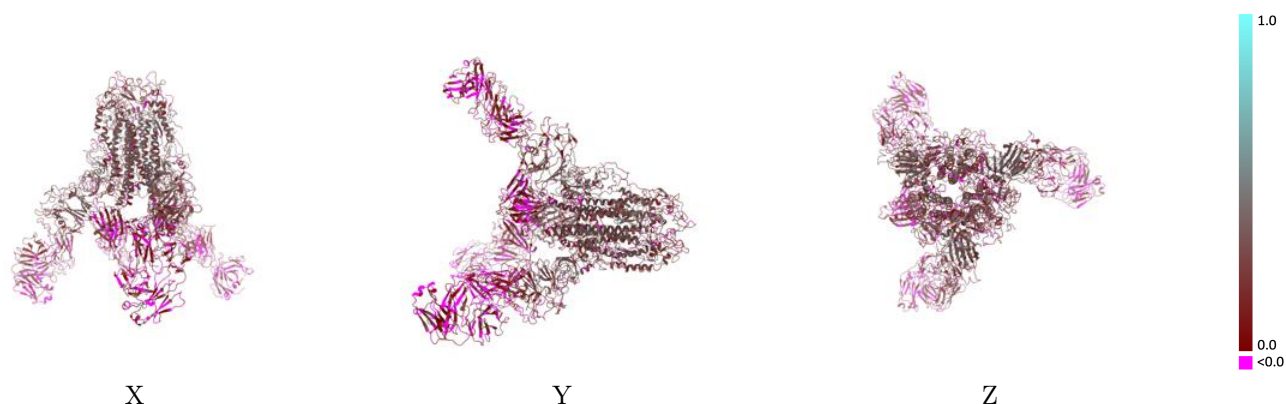
Y



Z

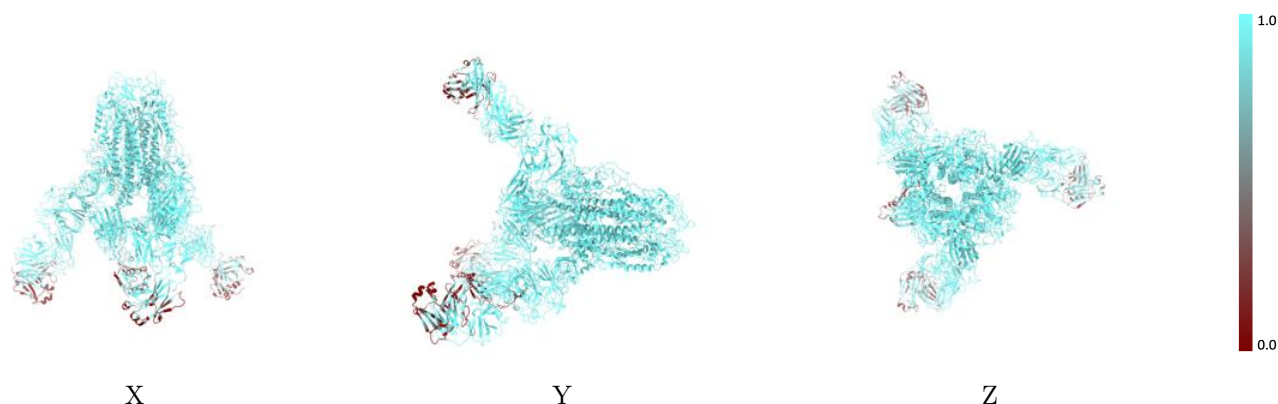
The images above show the 3D surface view of the map at the recommended contour level 0.004 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



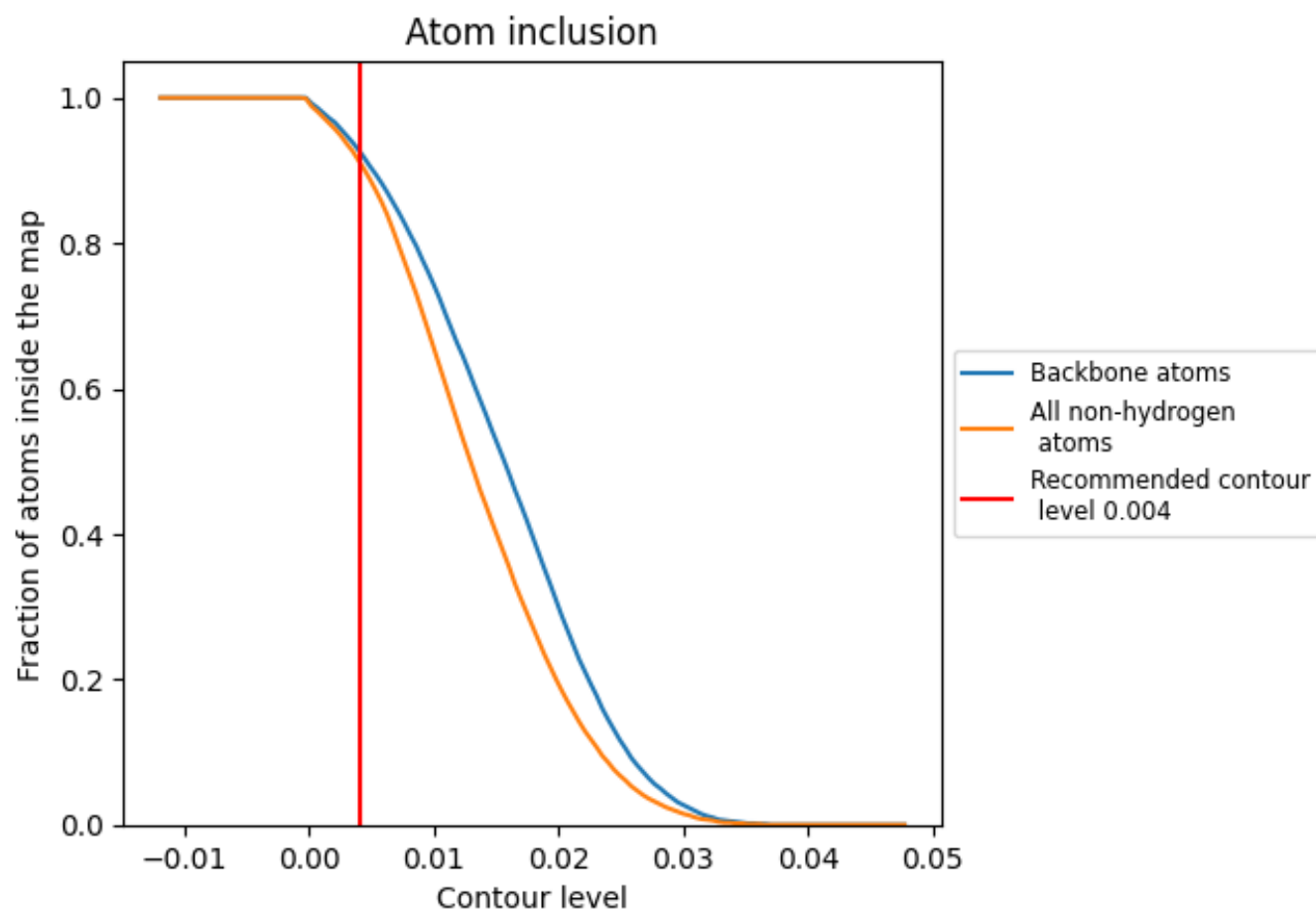
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.004).

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 91% 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 (0.004) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.9120	<div><div></div></div> 0.1850
A	<div><div></div></div> 0.9840	<div><div></div></div> 0.2230
B	<div><div></div></div> 0.9860	<div><div></div></div> 0.2300
C	<div><div></div></div> 0.9280	<div><div></div></div> 0.2210
D	<div><div></div></div> 0.7230	<div><div></div></div> 0.0790
E	<div><div></div></div> 0.7770	<div><div></div></div> 0.0460
F	<div><div></div></div> 0.7650	<div><div></div></div> 0.0970
G	<div><div></div></div> 0.8330	<div><div></div></div> 0.0880
H	<div><div></div></div> 0.7020	<div><div></div></div> 0.0770
I	<div><div></div></div> 0.8360	<div><div></div></div> 0.0980

1.0

0.0

<0.0