



Full wwPDB EM Validation Report ⓘ

Nov 19, 2022 – 02:34 PM EST

PDB ID : 7TPQ
EMDB ID : EMD-26061
Title : Cryo-em structure of human prothrombinase on a nanodisc at 5.3 Angstrom resolution
Authors : Di Cera, E.; Ruben, E.A.
Deposited on : 2022-01-25
Resolution : 5.30 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.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

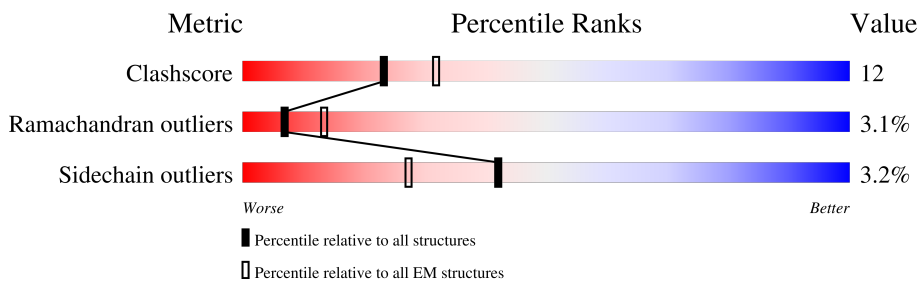
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 5.30 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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

Mol	Chain	Length	Quality of chain
1	A	139	29% (red), 73% (green), 22% (yellow), . (grey)
2	C	709	12% (red), 67% (green), 31% (yellow), . (grey)
3	D	651	7% (red), 67% (green), 30% (yellow), .. (grey)
4	B	254	20% (red), 59% (green), 35% (yellow), 6% (orange)

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 14020 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 Factor X light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	139	1089	657	184	231	17	0	0

- Molecule 2 is a protein called Coagulation factor Va.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	C	709	5715	3619	966	1100	30	0	0

- Molecule 3 is a protein called Coagulation factor Va.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	D	641	5217	3325	896	971	25	0	0

- Molecule 4 is a protein called Activated factor Xa heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	B	254	1999	1260	351	374	14	0	0

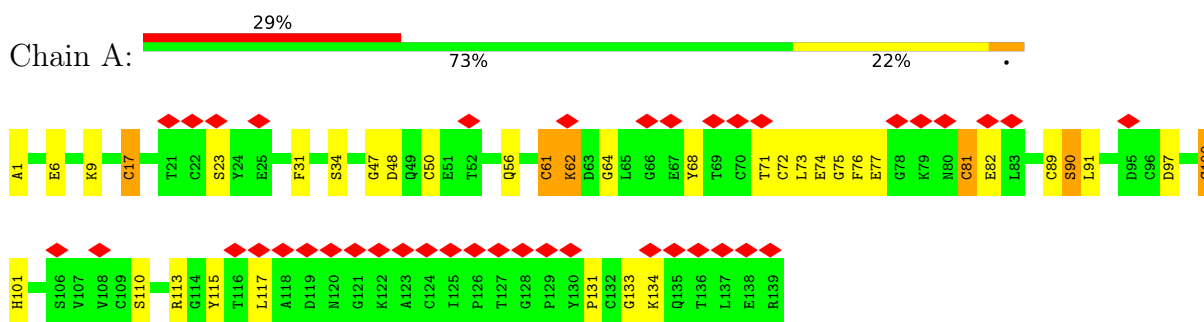
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	379	ALA	SER	engineered mutation	UNP P00742

3 Residue-property plots [i](#)

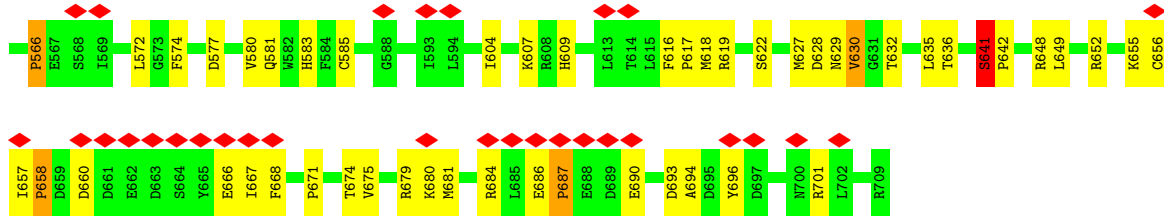
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Factor X light chain

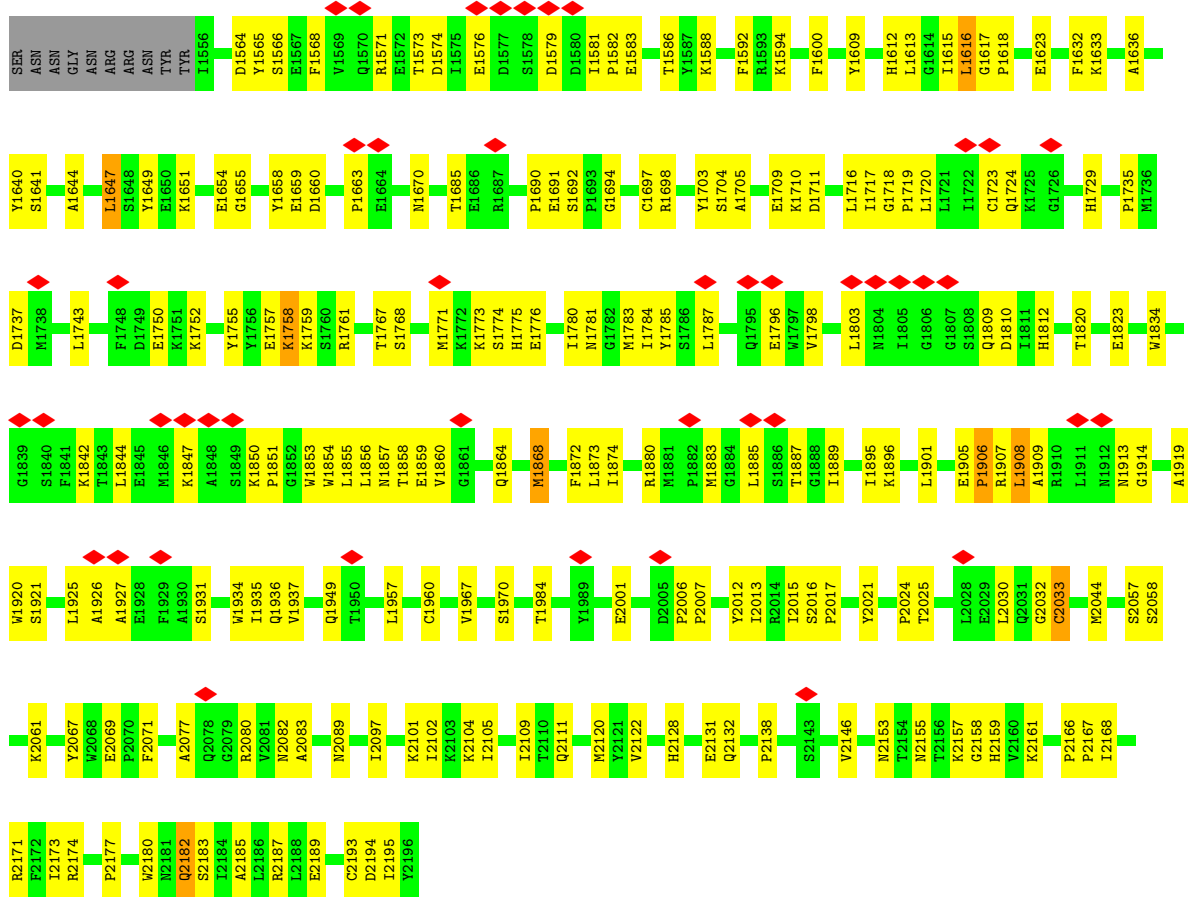


- Molecule 2: Coagulation factor Va

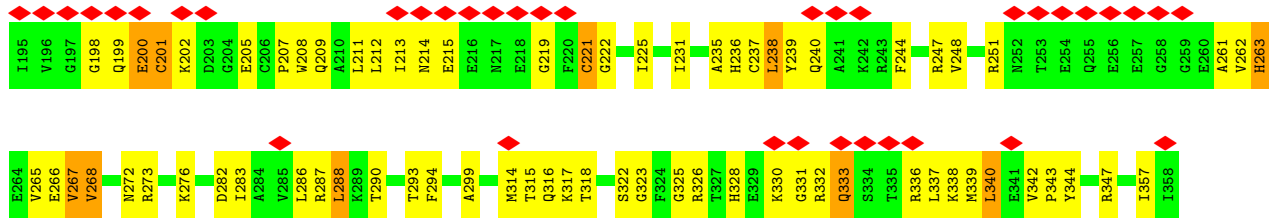


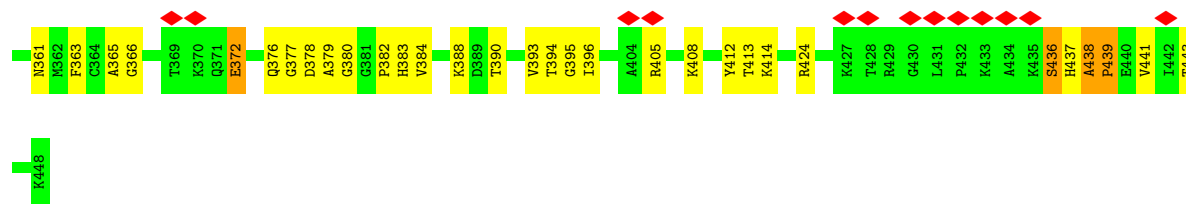


• Molecule 3: Coagulation factor Va



• Molecule 4: Activated factor Xa heavy chain





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	16835	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51.28	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.071	Depositor
Minimum map value	-0.504	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.030	Depositor
Recommended contour level	0.2	Depositor
Map size (\AA)	381.63998, 381.63998, 381.63998	wwPDB
Map dimensions	406, 406, 406	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.93999994, 0.93999994, 0.93999994	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	A	0.24	0/1107	0.43	0/1483
2	C	0.24	0/5861	0.43	0/7945
3	D	0.24	0/5358	0.43	0/7249
4	B	0.24	0/2041	0.45	0/2751
All	All	0.24	0/14367	0.43	0/19428

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1089	0	983	23	0
2	C	5715	0	5531	137	0
3	D	5217	0	5078	125	0
4	B	1999	0	1971	66	0
All	All	14020	0	13563	339	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 (339) 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)
4:B:437:HIS:HB3	4:B:438:ALA:HB3	1.60	0.82
4:B:380:GLY:HA2	4:B:396:ILE:HG23	1.62	0.81
2:C:313:ARG:HB3	2:C:701:ARG:HH22	1.53	0.74
2:C:325:PHE:HB2	2:C:371:TYR:HB2	1.70	0.73
2:C:37:VAL:HG12	2:C:157:GLY:HA3	1.71	0.71
4:B:318:THR:HA	4:B:343:PRO:HA	1.74	0.69
3:D:2128:HIS:HB2	3:D:2174:ARG:HB3	1.74	0.67
3:D:1612:HIS:HA	3:D:1781:ASN:HD21	1.60	0.66
2:C:420:HIS:HB3	2:C:475:ARG:HB3	1.78	0.66
4:B:436:SER:HA	4:B:439:PRO:HB3	1.76	0.66
2:C:581:GLN:NE2	2:C:622:SER:OG	2.28	0.66
3:D:2104:LYS:HB3	3:D:2193:CYS:HB3	1.77	0.65
2:C:628:ASP:OD2	4:B:347:ARG:NH2	2.30	0.65
3:D:1609:TYR:HB2	3:D:1984:THR:HG22	1.77	0.65
4:B:236:HIS:HE1	4:B:379:ALA:HB1	1.61	0.64
2:C:480:ASP:OD1	2:C:480:ASP:N	2.30	0.64
2:C:101:LEU:HD12	2:C:103:HIS:HE1	1.62	0.63
2:C:405:ILE:HG21	2:C:417:ILE:HG21	1.81	0.62
3:D:2193:CYS:SG	3:D:2194:ASP:N	2.72	0.62
4:B:283:ILE:HG12	4:B:396:ILE:HD12	1.81	0.62
3:D:1885:LEU:HD21	3:D:2030:LEU:H	1.65	0.62
3:D:1883:MET:HG3	3:D:2032:GLY:H	1.65	0.62
4:B:361:ASN:ND2	4:B:413:THR:OG1	2.33	0.62
3:D:1780:ILE:HG21	3:D:1868:MET:HB3	1.82	0.62
4:B:211:LEU:HD22	4:B:326:ARG:HH12	1.64	0.61
2:C:244:ASN:ND2	2:C:418:TYR:OH	2.32	0.61
2:C:472:CYS:HA	2:C:498:CYS:HA	1.81	0.61
2:C:316:ARG:NH1	2:C:500:SER:OG	2.34	0.61
2:C:23:ASN:HA	2:C:26:LEU:HD23	1.81	0.61
4:B:384:VAL:HA	4:B:393:VAL:HA	1.83	0.61
3:D:1564:ASP:HA	3:D:1586:THR:HA	1.82	0.61
4:B:396:ILE:H	4:B:413:THR:H	1.49	0.61
4:B:332:ARG:HH12	4:B:377:GLY:H	1.49	0.60
3:D:1970:SER:HB3	3:D:2012:TYR:HB2	1.82	0.60
2:C:604:ILE:HG13	2:C:609:HIS:HD1	1.67	0.60
4:B:221:CYS:SG	4:B:222:GLY:N	2.74	0.60
2:C:278:TRP:HB2	2:C:296:ILE:HB	1.84	0.59
3:D:1690:PRO:O	3:D:1724:GLN:NE2	2.35	0.59
3:D:1757:GLU:O	3:D:1759:LYS:N	2.34	0.59
4:B:199:GLN:NE2	4:B:339:MET:O	2.35	0.59
4:B:200:GLU:HA	4:B:338:LYS:HA	1.85	0.59
2:C:641:SER:H	2:C:642:PRO:HD2	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:2077:ALA:O	3:D:2082:ASN:ND2	2.35	0.59
2:C:675:VAL:HG21	4:B:273:ARG:HH21	1.68	0.59
4:B:328:HIS:HB3	4:B:332:ARG:HE	1.68	0.59
3:D:1616:LEU:HD21	3:D:1718:GLY:HA2	1.85	0.58
2:C:58:LEU:HB2	2:C:186:LEU:HD22	1.85	0.58
3:D:1949:GLN:HA	3:D:2001:GLU:HA	1.85	0.58
2:C:2:GLN:OE1	2:C:4:ARG:NH2	2.35	0.58
4:B:405:ARG:O	4:B:408:LYS:NZ	2.36	0.58
4:B:207:PRO:HG3	4:B:299:ALA:HB3	1.86	0.58
2:C:155:ASN:ND2	2:C:194:LYS:O	2.37	0.57
2:C:420:HIS:HD2	2:C:475:ARG:HD2	1.68	0.57
4:B:231:ILE:HB	4:B:286:LEU:HB2	1.84	0.57
4:B:342:VAL:HB	4:B:366:GLY:HA2	1.85	0.57
3:D:2111:GLN:HB3	3:D:2187:ARG:HD2	1.87	0.57
3:D:1880:ARG:NH1	3:D:2033:CYS:SG	2.78	0.56
1:A:1:ALA:N	1:A:6:GLU:OE1	2.36	0.56
2:C:185:LEU:HD11	2:C:227:TRP:HE3	1.70	0.56
3:D:1937:VAL:N	3:D:2013:ILE:O	2.38	0.56
2:C:277:LYS:HB3	2:C:435:PHE:HB2	1.87	0.56
2:C:316:ARG:NH2	2:C:693:ASP:OD2	2.39	0.56
2:C:326:ILE:HD13	2:C:405:ILE:HG23	1.87	0.56
1:A:6:GLU:HA	1:A:9:LYS:HE2	1.87	0.56
2:C:132:PRO:O	2:C:134:HIS:ND1	2.34	0.56
2:C:142:HIS:N	2:C:162:LEU:O	2.33	0.55
3:D:2120:MET:HG3	3:D:2183:SER:HB3	1.87	0.55
2:C:89:ILE:HB	3:D:1820:THR:HG23	1.88	0.55
3:D:1647:LEU:HG	3:D:1649:TYR:HD1	1.71	0.55
4:B:212:LEU:HD23	4:B:237:CYS:HB3	1.88	0.55
2:C:54:ILE:HD13	2:C:215:PRO:HG2	1.89	0.55
2:C:229:LEU:HD13	2:C:258:ILE:HG12	1.88	0.55
3:D:1565:TYR:HE1	3:D:1615:ILE:HD12	1.72	0.55
2:C:656:CYS:H	2:C:660:ASP:HB2	1.72	0.55
3:D:2128:HIS:HA	3:D:2138:PRO:HA	1.89	0.55
2:C:636:THR:HB	2:C:648:ARG:HG2	1.87	0.55
2:C:144:TYR:HD1	2:C:162:LEU:HB2	1.72	0.55
3:D:2083:ALA:HB2	3:D:2187:ARG:HG2	1.89	0.55
2:C:205:THR:HG22	2:C:210:VAL:HG22	1.88	0.55
2:C:572:LEU:HD13	2:C:649:LEU:HD21	1.89	0.55
3:D:1758:LYS:HG3	3:D:1759:LYS:HG3	1.89	0.55
2:C:485:ARG:NH2	2:C:528:SER:O	2.41	0.54
3:D:1690:PRO:HG3	3:D:1698:ARG:HB2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:325:GLY:N	4:B:378:ASP:OD1	2.39	0.54
3:D:1588:LYS:HE2	3:D:1755:TYR:HD1	1.72	0.54
3:D:1901:LEU:HA	3:D:1925:LEU:HB3	1.89	0.54
3:D:1750:GLU:HB3	3:D:1775:HIS:HB3	1.89	0.54
3:D:1594:LYS:HB2	3:D:1618:PRO:HG3	1.88	0.54
3:D:1641:SER:OG	3:D:1670:ASN:ND2	2.41	0.54
3:D:1773:LYS:HB3	3:D:1776:GLU:HB3	1.88	0.54
2:C:95:SER:O	2:C:124:TRP:NE1	2.39	0.54
3:D:1581:ILE:HD12	3:D:1582:PRO:HD2	1.89	0.54
2:C:153:ASP:HA	2:C:158:LEU:HD12	1.89	0.53
2:C:680:LYS:HD3	2:C:681:MET:HG3	1.89	0.53
2:C:316:ARG:NH1	2:C:500:SER:O	2.35	0.53
4:B:363:PHE:HE2	4:B:414:LYS:HB2	1.72	0.53
3:D:2111:GLN:HE21	3:D:2157:LYS:HD2	1.74	0.53
3:D:1649:TYR:HE2	3:D:1654:GLU:HG3	1.73	0.53
3:D:1655:GLY:O	3:D:1670:ASN:ND2	2.35	0.53
1:A:74:GLU:OE1	1:A:75:GLY:N	2.42	0.53
4:B:323:GLY:HA3	4:B:378:ASP:H	1.74	0.53
2:C:675:VAL:H	4:B:424:ARG:HH22	1.56	0.53
3:D:1857:ASN:ND2	3:D:1859:GLU:OE1	2.42	0.52
2:C:301:CYS:SG	2:C:303:LYS:NZ	2.82	0.52
2:C:428:GLU:HG2	2:C:456:LYS:H	1.73	0.52
3:D:1709:GLU:HB3	3:D:1752:LYS:HE3	1.90	0.52
2:C:616:PHE:HD2	2:C:619:ARG:HH11	1.57	0.52
2:C:313:ARG:O	2:C:701:ARG:NH2	2.43	0.52
3:D:1644:ALA:O	3:D:1649:TYR:OH	2.25	0.52
3:D:1697:CYS:HA	3:D:1723:CYS:HA	1.91	0.52
3:D:1592:PHE:HB2	3:D:1617:GLY:HA3	1.92	0.52
3:D:1967:VAL:HG13	3:D:2015:ILE:HG22	1.90	0.52
1:A:110:SER:HA	1:A:117:LEU:HD21	1.92	0.52
2:C:400:ARG:N	2:C:459:ILE:O	2.42	0.52
3:D:2061:LYS:HG3	3:D:2067:TYR:HB3	1.92	0.52
4:B:365:ALA:HB3	4:B:412:TYR:CD1	2.45	0.52
1:A:50:CYS:SG	1:A:61:CYS:N	2.84	0.51
1:A:72:CYS:SG	1:A:73:LEU:N	2.83	0.51
2:C:227:TRP:O	2:C:267:THR:OG1	2.24	0.51
2:C:351:HIS:O	2:C:363:TYR:OH	2.23	0.51
3:D:1780:ILE:HD13	3:D:1868:MET:HA	1.92	0.51
4:B:198:GLY:HA2	4:B:340:LEU:HB3	1.91	0.51
2:C:232:MET:HG3	2:C:262:SER:HA	1.91	0.51
1:A:89:CYS:O	1:A:91:LEU:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1571:ARG:HE	3:D:1574:ASP:HB3	1.76	0.51
3:D:1771:MET:HB2	3:D:1774:SER:HB3	1.92	0.51
4:B:328:HIS:HD2	4:B:330:LYS:HD3	1.76	0.51
1:A:133:GLY:HA3	4:B:390:THR:HA	1.93	0.51
2:C:684:ARG:NH2	4:B:272:ASN:O	2.42	0.51
2:C:484:MET:SD	2:C:484:MET:N	2.84	0.50
3:D:1581:ILE:HD11	3:D:1583:GLU:HG2	1.93	0.50
3:D:1851:PRO:HA	3:D:1874:ILE:HB	1.93	0.50
3:D:2132:GLN:HA	3:D:2171:ARG:HD2	1.93	0.50
2:C:18:ARG:NH1	2:C:211:ASN:OD1	2.41	0.50
3:D:1640:TYR:OH	3:D:1755:TYR:OH	2.20	0.50
2:C:334:ASP:HB3	2:C:347:TYR:HE2	1.75	0.50
3:D:1600:PHE:HE1	3:D:1633:LYS:HB3	1.76	0.50
2:C:139:CYS:HA	2:C:165:CYS:HA	1.94	0.50
1:A:76:PHE:HB2	1:A:81:CYS:HB2	1.93	0.50
2:C:574:PHE:HB2	2:C:627:MET:HB3	1.93	0.50
2:C:14:SER:HA	2:C:34:LYS:HA	1.93	0.49
2:C:36:ILE:HG22	2:C:158:LEU:HD21	1.94	0.49
2:C:58:LEU:HD11	2:C:161:PRO:HD3	1.94	0.49
3:D:1757:GLU:C	3:D:1759:LYS:H	2.15	0.49
2:C:100:TYR:CZ	3:D:1855:LEU:HB2	2.48	0.49
2:C:398:GLN:H	2:C:401:ASP:HB2	1.77	0.49
2:C:655:LYS:HA	2:C:660:ASP:HB2	1.94	0.49
4:B:202:LYS:HD2	4:B:205:GLU:HB3	1.94	0.49
4:B:383:HIS:HB3	4:B:395:GLY:HA3	1.94	0.49
2:C:629:ASN:OD1	2:C:630:VAL:N	2.44	0.49
3:D:1887:THR:HA	3:D:2159:HIS:HB2	1.94	0.49
1:A:89:CYS:SG	1:A:90:SER:N	2.86	0.49
2:C:693:ASP:OD1	2:C:694:ALA:N	2.45	0.49
3:D:1810:ASP:OD1	3:D:1812:HIS:NE2	2.46	0.49
4:B:225:ILE:HD11	4:B:248:VAL:HB	1.95	0.49
1:A:47:GLY:HA2	1:A:48:ASP:HA	1.52	0.49
3:D:2044:MET:HB2	3:D:2189:GLU:HG2	1.95	0.49
3:D:2120:MET:SD	3:D:2120:MET:N	2.86	0.49
2:C:96:GLU:HG3	2:C:98:ALA:H	1.78	0.48
2:C:346:LYS:NZ	2:C:555:SER:OG	2.46	0.48
2:C:40:GLU:HB2	2:C:60:PRO:HB3	1.96	0.48
2:C:476:PRO:HA	2:C:494:LEU:HA	1.96	0.48
3:D:1823:GLU:HA	3:D:1844:LEU:HD11	1.95	0.48
4:B:247:ARG:HB2	4:B:261:ALA:HB1	1.96	0.48
3:D:1710:LYS:HZ2	3:D:1752:LYS:HA	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1856:LEU:HB2	3:D:1872:PHE:HD1	1.78	0.48
2:C:632:THR:HB	2:C:652:ARG:HG2	1.95	0.48
4:B:214:ASN:OD1	4:B:215:GLU:N	2.44	0.48
2:C:63:TYR:HA	2:C:163:LEU:HB2	1.96	0.48
2:C:181:LYS:HB2	2:C:225:ILE:HA	1.96	0.48
2:C:617:PRO:O	2:C:619:ARG:N	2.43	0.48
3:D:1729:HIS:ND1	3:D:1737:ASP:OD1	2.47	0.48
3:D:1632:PHE:HZ	3:D:1640:TYR:HD2	1.62	0.48
2:C:315:GLN:HG3	2:C:316:ARG:H	1.78	0.47
4:B:372:GLU:O	4:B:376:GLN:NE2	2.47	0.47
2:C:635:LEU:HD12	2:C:649:LEU:HD23	1.95	0.47
3:D:1896:LYS:O	3:D:1936:GLN:N	2.47	0.47
3:D:2131:GLU:OE1	3:D:2131:GLU:N	2.46	0.47
4:B:239:TYR:HE2	4:B:276:LYS:HG3	1.79	0.47
4:B:322:SER:HB3	4:B:382:PRO:HG2	1.95	0.47
3:D:1925:LEU:HD22	3:D:2024:PRO:HD2	1.97	0.47
2:C:92:SER:HB3	3:D:1850:LYS:HE2	1.97	0.47
2:C:525:GLU:H	2:C:556:ASN:HB3	1.79	0.47
2:C:76:ASN:ND2	2:C:80:LYS:O	2.48	0.47
2:C:359:ILE:HD13	2:C:534:ASN:HD22	1.79	0.47
2:C:73:HIS:NE2	2:C:119:GLU:HB3	2.28	0.47
4:B:238:LEU:HD22	4:B:268:VAL:HG11	1.95	0.47
4:B:438:ALA:HB1	4:B:441:VAL:H	1.80	0.47
2:C:96:GLU:N	2:C:111:ASP:OD2	2.48	0.47
2:C:137:PRO:HD3	2:C:251:ASN:HA	1.97	0.47
3:D:1834:TRP:NE1	3:D:1842:LYS:HD3	2.30	0.47
4:B:201:CYS:HB2	4:B:337:LEU:HG	1.97	0.47
3:D:1651:LYS:NZ	3:D:1660:ASP:HB3	2.30	0.47
2:C:55:SER:OG	2:C:58:LEU:O	2.33	0.46
3:D:2122:VAL:O	3:D:2153:ASN:N	2.48	0.46
2:C:657:ILE:HB	2:C:658:PRO:HD3	1.97	0.46
3:D:1864:GLN:NE2	3:D:1868:MET:O	2.47	0.46
2:C:8:VAL:HA	2:C:40:GLU:HA	1.98	0.46
3:D:1960:CYS:SG	3:D:2025:THR:OG1	2.60	0.46
3:D:1780:ILE:N	3:D:1783:MET:O	2.49	0.46
2:C:519:VAL:HG23	2:C:561:ILE:HA	1.98	0.46
3:D:1703:TYR:HB2	3:D:1716:LEU:O	2.15	0.46
2:C:80:LYS:NZ	2:C:152:GLU:OE2	2.46	0.46
2:C:161:PRO:HB3	2:C:230:LEU:HD13	1.98	0.46
3:D:2097:ILE:HB	3:D:2173:ILE:HG23	1.98	0.46
3:D:2104:LYS:HG2	3:D:2167:PRO:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1568:PHE:HE2	3:D:1579:ASP:HB2	1.79	0.45
1:A:9:LYS:HB3	1:A:31:PHE:HD2	1.81	0.45
2:C:284:THR:OG1	2:C:287:HIS:O	2.28	0.45
2:C:398:GLN:NE2	2:C:505:ARG:O	2.50	0.45
2:C:220:CYS:HB3	2:C:299:LYS:HB2	1.98	0.45
2:C:690:GLU:OE2	4:B:357:ILE:N	2.49	0.45
3:D:2122:VAL:HG23	3:D:2185:ALA:H	1.82	0.45
4:B:317:LYS:O	4:B:344:TYR:N	2.38	0.45
1:A:77:GLU:O	1:A:82:GLU:N	2.50	0.45
2:C:101:LEU:HD23	2:C:101:LEU:H	1.81	0.45
3:D:1913:ASN:OD1	3:D:1914:GLY:N	2.49	0.45
4:B:263:HIS:CE1	4:B:288:LEU:HD13	2.52	0.45
3:D:2089:ASN:ND2	3:D:2182:GLN:O	2.49	0.45
4:B:326:ARG:HG2	4:B:333:GLN:HE22	1.81	0.45
1:A:115:TYR:HE1	1:A:134:LYS:HE2	1.82	0.45
2:C:445:ARG:HD3	2:C:445:ARG:H	1.82	0.45
2:C:519:VAL:HG12	2:C:585:CYS:HB3	1.99	0.44
3:D:1780:ILE:HD11	3:D:1785:TYR:H	1.80	0.44
3:D:1906:PRO:HA	3:D:1920:TRP:HA	1.99	0.44
3:D:2101:LYS:NZ	3:D:2102:ILE:O	2.51	0.44
3:D:2122:VAL:HG11	3:D:2177:PRO:HB3	1.99	0.44
4:B:209:GLN:HB3	4:B:251:ARG:HH21	1.81	0.44
2:C:35:LYS:NZ	2:C:155:ASN:O	2.40	0.44
2:C:222:HIS:HA	2:C:306:ARG:HH22	1.82	0.44
2:C:513:ASP:HB3	2:C:580:VAL:HG12	1.99	0.44
3:D:1640:TYR:HH	3:D:1755:TYR:HH	1.60	0.44
2:C:482:ASP:OD2	2:C:485:ARG:N	2.51	0.44
2:C:474:THR:OG1	2:C:622:SER:HB2	2.18	0.44
2:C:319:MET:HB2	2:C:400:ARG:HD3	1.99	0.44
3:D:1579:ASP:N	3:D:1579:ASP:OD1	2.51	0.44
4:B:363:PHE:CE2	4:B:414:LYS:HB2	2.53	0.44
2:C:38:TYR:HB2	2:C:59:GLY:HA3	2.00	0.43
2:C:111:ASP:HB2	2:C:122:TYR:HE1	1.84	0.43
3:D:2104:LYS:HB2	3:D:2195:ILE:HD11	1.99	0.43
2:C:191:ASP:OD1	2:C:193:SER:OG	2.32	0.43
2:C:547:ARG:HE	2:C:547:ARG:HA	1.83	0.43
3:D:1743:LEU:HD11	3:D:1787:LEU:HD22	2.01	0.43
3:D:1901:LEU:HD11	3:D:2017:PRO:HG3	2.00	0.43
3:D:1935:ILE:N	3:D:2015:ILE:O	2.47	0.43
3:D:2155:ASN:HD21	3:D:2158:GLY:HA3	1.83	0.43
2:C:90:ARG:HE	2:C:127:SER:HA	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:240:SER:H	2:C:284:THR:HG22	1.84	0.43
3:D:1568:PHE:O	3:D:1571:ARG:NH1	2.51	0.43
3:D:1905:GLU:HB2	3:D:1908:LEU:HB3	2.01	0.43
3:D:1960:CYS:HA	3:D:2021:TYR:HB3	2.01	0.43
4:B:265:VAL:HG23	4:B:287:ARG:O	2.18	0.43
2:C:92:SER:HA	3:D:1854:TRP:HH2	1.83	0.43
3:D:1568:PHE:HB3	3:D:1785:TYR:HB2	2.01	0.43
2:C:186:LEU:HD12	2:C:231:GLY:HA2	2.01	0.43
2:C:323:GLU:HB3	2:C:404:LYS:HE2	2.00	0.43
1:A:101:HIS:O	1:A:101:HIS:ND1	2.51	0.43
2:C:72:VAL:HG13	2:C:122:TYR:HB2	2.01	0.43
2:C:666:GLU:HG2	2:C:668:PHE:HD2	1.83	0.43
1:A:9:LYS:HD2	1:A:31:PHE:HA	2.00	0.43
2:C:629:ASN:H	2:C:655:LYS:HZ3	1.67	0.43
2:C:192:GLU:HB3	2:C:201:SER:HB3	1.99	0.42
2:C:342:ASN:HD21	2:C:566:PRO:HB2	1.84	0.42
4:B:437:HIS:H	4:B:438:ALA:C	2.22	0.42
1:A:62:LYS:NZ	1:A:64:GLY:HA3	2.34	0.42
2:C:141:THR:HG22	2:C:163:LEU:HG	2.02	0.42
2:C:310:LYS:HG2	2:C:460:LEU:HD13	2.00	0.42
2:C:686:GLU:HB3	2:C:687:PRO:HD3	2.01	0.42
4:B:248:VAL:HG22	4:B:263:HIS:HB2	2.00	0.42
1:A:48:ASP:N	1:A:48:ASP:OD1	2.53	0.42
2:C:409:ASN:ND2	2:C:413:ARG:O	2.50	0.42
4:B:248:VAL:H	4:B:263:HIS:HB2	1.85	0.42
4:B:314:MET:HG3	4:B:315:THR:HG23	2.01	0.42
2:C:607:LYS:HE2	2:C:607:LYS:HB2	1.91	0.42
3:D:1719:PRO:HB3	3:D:1803:LEU:HD22	2.02	0.42
4:B:328:HIS:CD2	4:B:330:LYS:HB2	2.54	0.42
3:D:1623:GLU:HG3	3:D:1735:PRO:HG3	2.01	0.42
3:D:1812:HIS:CD2	3:D:1860:VAL:HG11	2.54	0.42
3:D:1812:HIS:HB3	3:D:1858:THR:HB	2.02	0.42
3:D:2105:ILE:HB	3:D:2168:ILE:HD12	2.00	0.42
3:D:2109:ILE:HG22	3:D:2189:GLU:HB3	2.02	0.42
3:D:1907:ARG:O	3:D:1909:ALA:N	2.53	0.42
3:D:2057:SER:OG	3:D:2058:SER:N	2.52	0.42
3:D:2069:GLU:HB3	3:D:2071:PHE:CE1	2.55	0.42
4:B:213:ILE:HA	4:B:219:GLY:HA2	2.02	0.42
4:B:293:THR:OG1	4:B:294:PHE:N	2.52	0.42
2:C:64:ALA:N	2:C:163:LEU:O	2.52	0.42
3:D:1616:LEU:HD23	3:D:1617:GLY:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1934:TRP:HA	3:D:2016:SER:HB2	2.01	0.42
2:C:76:ASN:O	2:C:117:GLY:N	2.50	0.42
2:C:160:GLY:HA2	2:C:161:PRO:HD3	1.86	0.42
3:D:1566:SER:HA	3:D:1784:ILE:HD12	2.02	0.42
3:D:1796:GLU:HB3	3:D:1798:VAL:HG13	2.02	0.42
2:C:675:VAL:HG22	2:C:679:ARG:HG3	2.01	0.42
3:D:2122:VAL:HA	3:D:2180:TRP:HA	2.01	0.42
1:A:113:ARG:H	1:A:113:ARG:HG3	1.71	0.41
4:B:209:GLN:HE22	4:B:323:GLY:HA2	1.85	0.41
2:C:515:GLU:OE2	2:C:583:HIS:ND1	2.53	0.41
3:D:1889:ILE:HG13	3:D:2161:LYS:H	1.84	0.41
2:C:9:ALA:HB1	2:C:77:LYS:HE2	2.01	0.41
2:C:139:CYS:HB2	2:C:267:THR:HG21	2.01	0.41
2:C:170:LEU:HD13	2:C:176:GLN:HA	2.02	0.41
3:D:1647:LEU:HD23	3:D:1647:LEU:H	1.85	0.41
3:D:1895:ILE:HD11	3:D:1935:ILE:HG22	2.02	0.41
3:D:2166:PRO:HA	3:D:2167:PRO:HD3	1.91	0.41
4:B:336:ARG:NH1	4:B:337:LEU:HB3	2.35	0.41
4:B:221:CYS:HB2	4:B:379:ALA:HB3	2.01	0.41
4:B:288:LEU:HD23	4:B:288:LEU:H	1.85	0.41
1:A:90:SER:N	1:A:100:CYS:SG	2.93	0.41
2:C:233:SER:HB3	2:C:239:PHE:HE2	1.85	0.41
2:C:351:HIS:HB3	2:C:359:ILE:HD12	2.03	0.41
2:C:577:ASP:H	4:B:347:ARG:HH12	1.66	0.41
3:D:1704:SER:HB3	3:D:1711:ASP:HB3	2.03	0.41
3:D:1920:TRP:CH2	3:D:1935:ILE:HG12	2.56	0.41
4:B:240:GLN:HE22	4:B:276:LYS:HD3	1.85	0.41
2:C:144:TYR:CD1	2:C:162:LEU:HB2	2.55	0.41
4:B:235:ALA:N	4:B:282:ASP:OD1	2.37	0.41
3:D:1757:GLU:HB3	3:D:1761:ARG:HH12	1.86	0.41
2:C:283:LEU:HD22	2:C:483:ILE:HG21	2.03	0.41
2:C:425:SER:HB2	2:C:457:TRP:HE3	1.86	0.41
3:D:1704:SER:OG	3:D:1705:ALA:N	2.54	0.41
1:A:61:CYS:HB3	1:A:68:TYR:HB2	2.03	0.41
2:C:280:ILE:H	2:C:295:TYR:HA	1.86	0.41
3:D:2069:GLU:N	3:D:2069:GLU:OE1	2.53	0.41
4:B:267:VAL:HG12	4:B:268:VAL:H	1.86	0.41
1:A:56:GLN:H	1:A:56:GLN:HG3	1.74	0.41
2:C:417:ILE:HG23	2:C:430:GLU:HG3	2.03	0.41
4:B:326:ARG:NH1	4:B:378:ASP:OD2	2.53	0.41
2:C:157:GLY:HA2	2:C:190:PHE:HE1	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1600:PHE:CE1	3:D:1633:LYS:HB3	2.55	0.40
2:C:235:GLY:N	2:C:237:GLU:OE1	2.55	0.40
3:D:1906:PRO:HA	3:D:1921:SER:H	1.86	0.40
1:A:97:ASP:HA	4:B:388:LYS:HD3	2.03	0.40
3:D:1857:ASN:OD1	3:D:1857:ASN:N	2.54	0.40
2:C:526:ASN:HA	2:C:531:LEU:HD13	2.03	0.40
3:D:1658:TYR:HD2	3:D:1659:GLU:HG3	1.87	0.40
3:D:1853:TRP:CD1	3:D:1873:LEU:HD13	2.57	0.40
3:D:1925:LEU:O	3:D:1927:ALA:N	2.54	0.40
3:D:2006:PRO:HA	3:D:2007:PRO:HD3	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	137/139 (99%)	106 (77%)	26 (19%)	5 (4%)	3	27
2	C	707/709 (100%)	545 (77%)	139 (20%)	23 (3%)	4	29
3	D	639/651 (98%)	483 (76%)	139 (22%)	17 (3%)	5	33
4	B	252/254 (99%)	191 (76%)	52 (21%)	9 (4%)	3	27
All	All	1735/1753 (99%)	1325 (76%)	356 (20%)	54 (3%)	7	30

All (54) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	658	PRO
3	D	1663	PRO
3	D	1758	LYS
4	B	439	PRO
1	A	90	SER

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Mol	Chain	Res	Type
1	A	131	PRO
2	C	383	PRO
2	C	566	PRO
2	C	687	PRO
3	D	1926	ALA
4	B	208	TRP
2	C	51	GLN
2	C	641	SER
3	D	1636	ALA
3	D	1768	SER
3	D	1906	PRO
3	D	1957	LEU
4	B	333	GLN
1	A	17	CYS
1	A	81	CYS
2	C	265	SER
2	C	379	HIS
2	C	465	PRO
3	D	1692	SER
3	D	1931	SER
3	D	2080	ARG
3	D	2182	GLN
4	B	266	GLU
1	A	34	SER
2	C	148	GLU
2	C	393	PRO
2	C	499	LYS
2	C	618	MET
3	D	1767	THR
3	D	1847	LYS
3	D	1908	LEU
3	D	1919	ALA
4	B	263	HIS
4	B	438	ALA
2	C	263	ALA
2	C	339	ILE
2	C	432	ASN
4	B	436	SER
2	C	86	PRO
2	C	89	ILE
4	B	331	GLY
2	C	132	PRO

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Mol	Chain	Res	Type
2	C	514	ILE
3	D	2146	VAL
3	D	1694	GLY
4	B	262	VAL
2	C	116	PRO
2	C	667	ILE
2	C	671	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	122/122 (100%)	116 (95%)	6 (5%)	25	51
2	C	637/637 (100%)	619 (97%)	18 (3%)	43	65
3	D	567/576 (98%)	555 (98%)	12 (2%)	53	72
4	B	215/215 (100%)	201 (94%)	14 (6%)	17	43
All	All	1541/1550 (99%)	1491 (97%)	50 (3%)	42	61

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	CYS
1	A	23	SER
1	A	61	CYS
1	A	62	LYS
1	A	71	THR
1	A	100	CYS
2	C	31	THR
2	C	34	LYS
2	C	47	LYS
2	C	101	LEU
2	C	206	VAL
2	C	306	ARG
2	C	313	ARG
2	C	338	VAL

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Mol	Chain	Res	Type
2	C	358	GLN
2	C	379	HIS
2	C	445	ARG
2	C	480	ASP
2	C	506	ARG
2	C	551	LYS
2	C	630	VAL
2	C	641	SER
2	C	674	THR
2	C	696	TYR
3	D	1573	THR
3	D	1576	GLU
3	D	1613	LEU
3	D	1616	LEU
3	D	1647	LEU
3	D	1685	THR
3	D	1691	GLU
3	D	1717	ILE
3	D	1720	LEU
3	D	1809	GLN
3	D	1868	MET
3	D	2033	CYS
4	B	200	GLU
4	B	201	CYS
4	B	221	CYS
4	B	238	LEU
4	B	244	PHE
4	B	267	VAL
4	B	268	VAL
4	B	288	LEU
4	B	290	THR
4	B	316	GLN
4	B	340	LEU
4	B	372	GLU
4	B	394	THR
4	B	443	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
2	C	5	GLN

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Mol	Chain	Res	Type
2	C	244	ASN
2	C	300	ASN
2	C	315	GLN
2	C	398	GLN
2	C	420	HIS
2	C	534	ASN
2	C	581	GLN
2	C	597	HIS
3	D	1670	ASN
3	D	1724	GLN
3	D	1781	ASN
3	D	1804	ASN
3	D	1809	GLN
3	D	1815	HIS
3	D	1817	HIS
3	D	1824	ASN
3	D	1940	GLN
3	D	2031	GLN
3	D	2089	ASN
3	D	2128	HIS
4	B	236	HIS
4	B	240	GLN
4	B	316	GLN
4	B	328	HIS
4	B	361	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 monosaccharides 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.

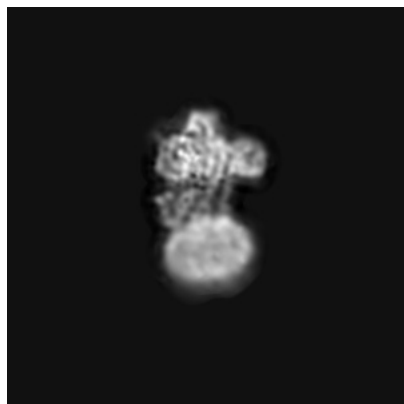
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26061. These allow visual inspection of the internal detail of the map and identification of artifacts.

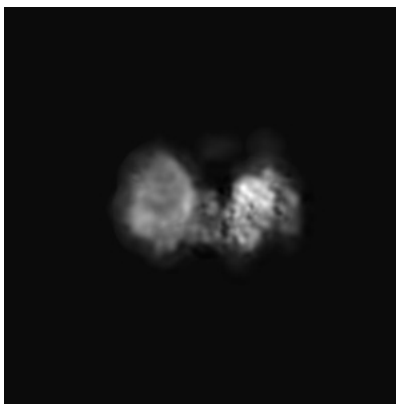
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

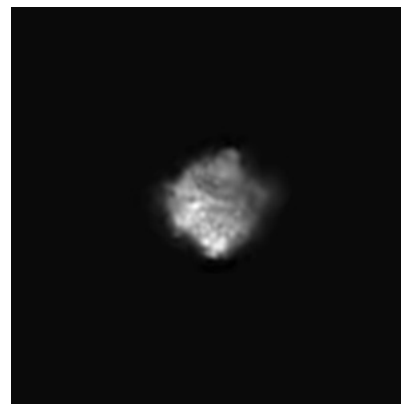
6.1.1 Primary map



X

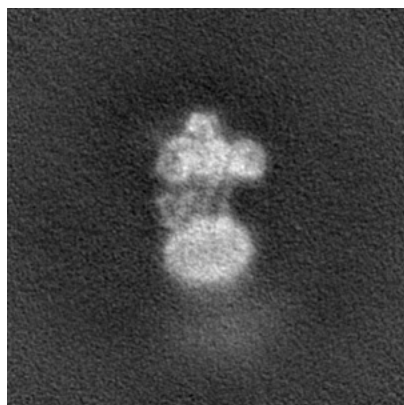


Y

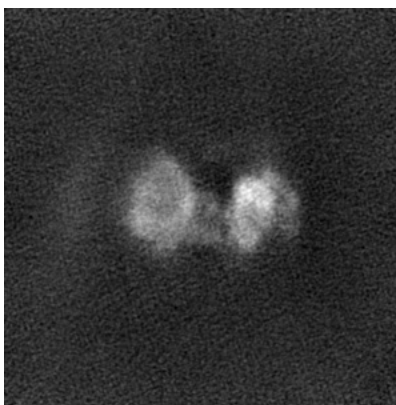


Z

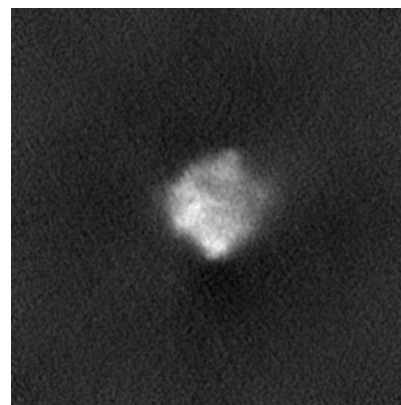
6.1.2 Raw map



X



Y



Z

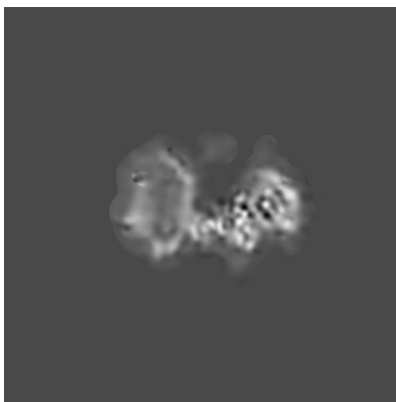
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

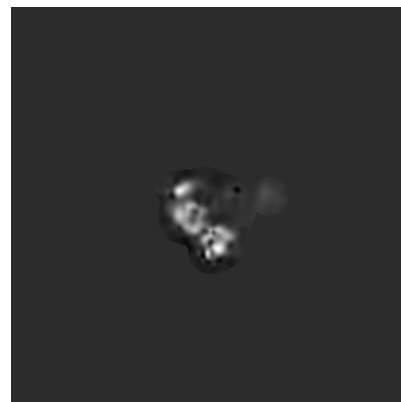
6.2.1 Primary map



X Index: 203

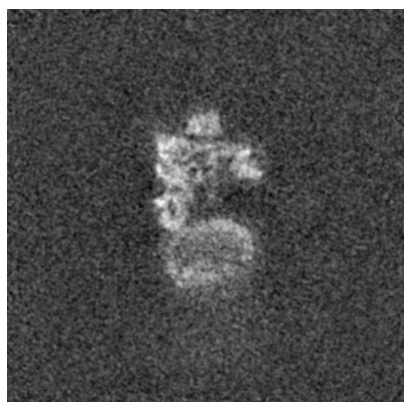


Y Index: 203



Z Index: 203

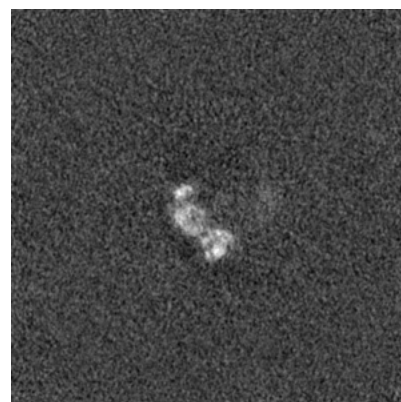
6.2.2 Raw map



X Index: 203



Y Index: 203



Z Index: 203

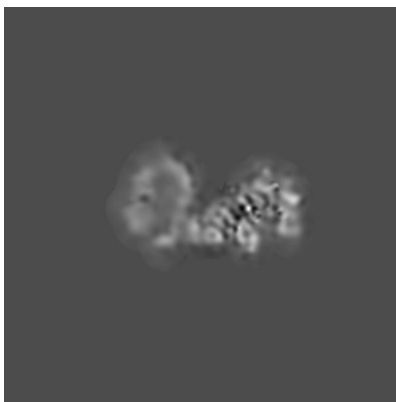
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 213

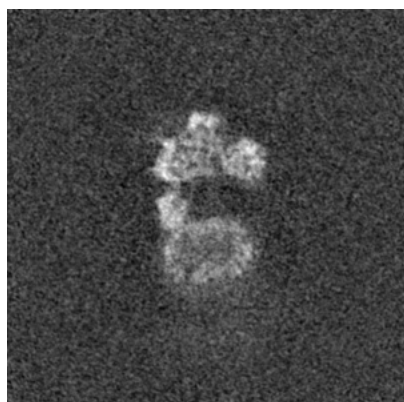


Y Index: 188

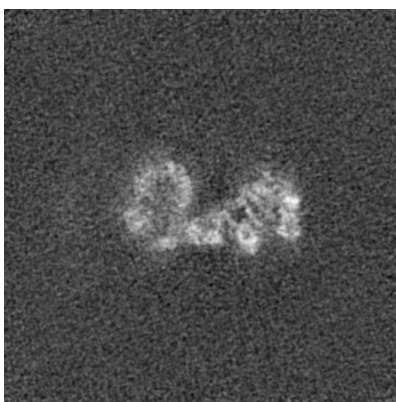


Z Index: 242

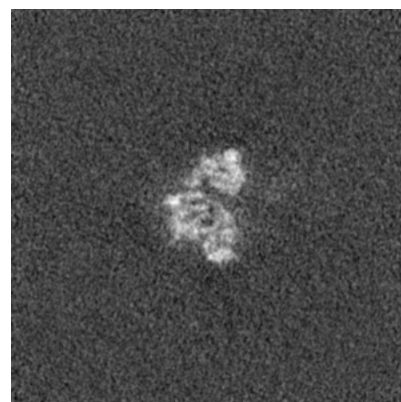
6.3.2 Raw map



X Index: 212



Y Index: 189

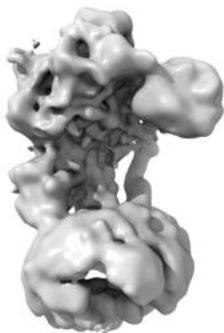


Z Index: 243

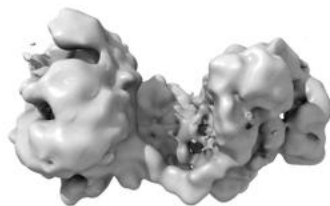
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

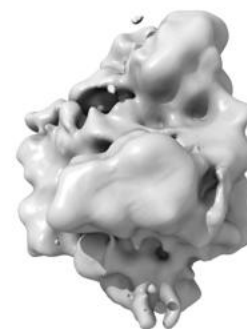
6.4.1 Primary map



X



Y



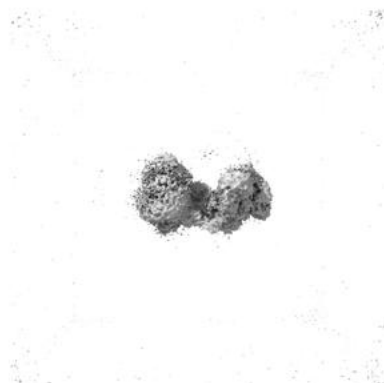
Z

The images above show the 3D surface view of the map at the recommended contour level 0.2. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

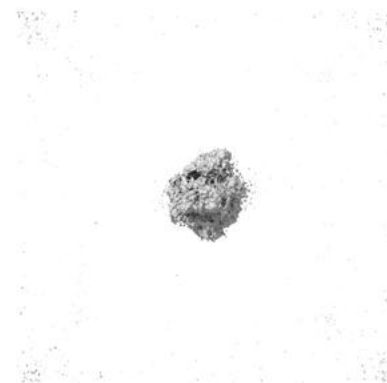
6.4.2 Raw map



X



Y



Z

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.

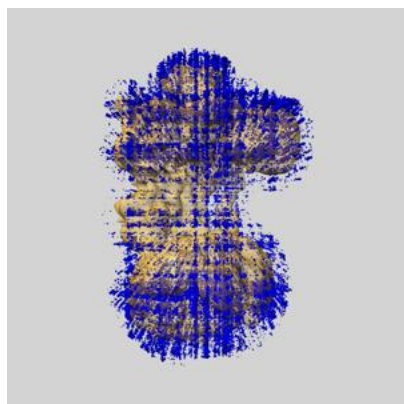
6.5 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

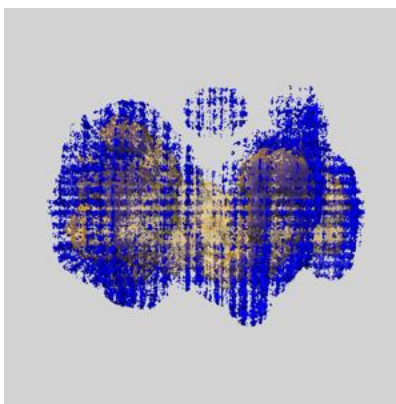
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

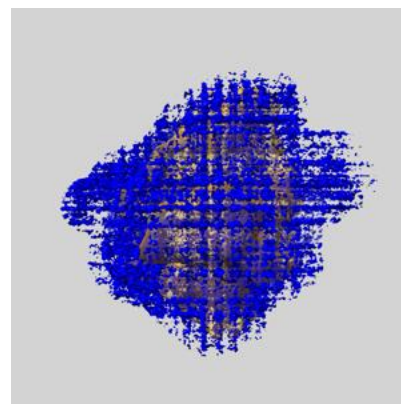
6.5.1 emd_26061_msk_1.map [i](#)



X



Y

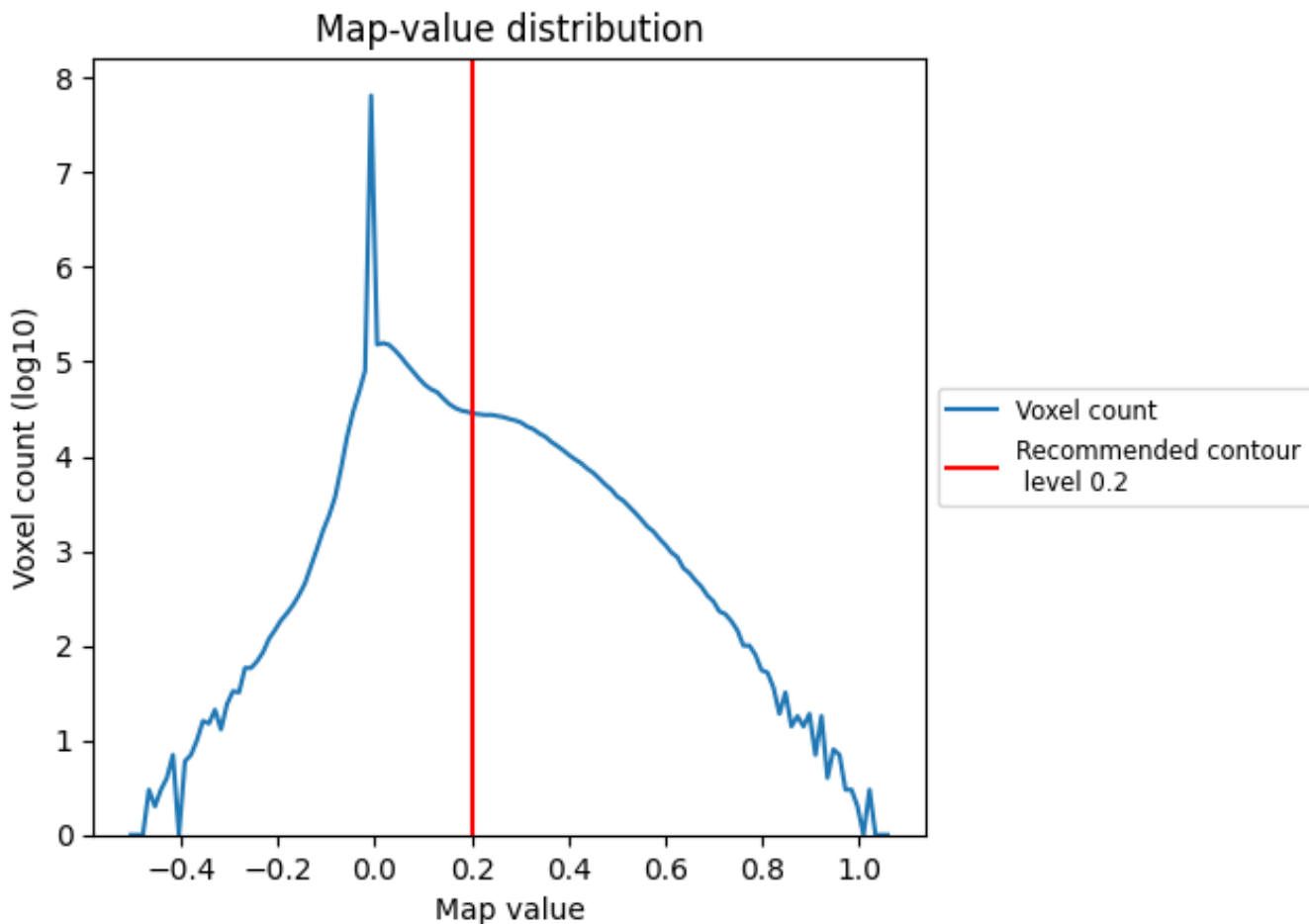


Z

7 Map analysis [i](#)

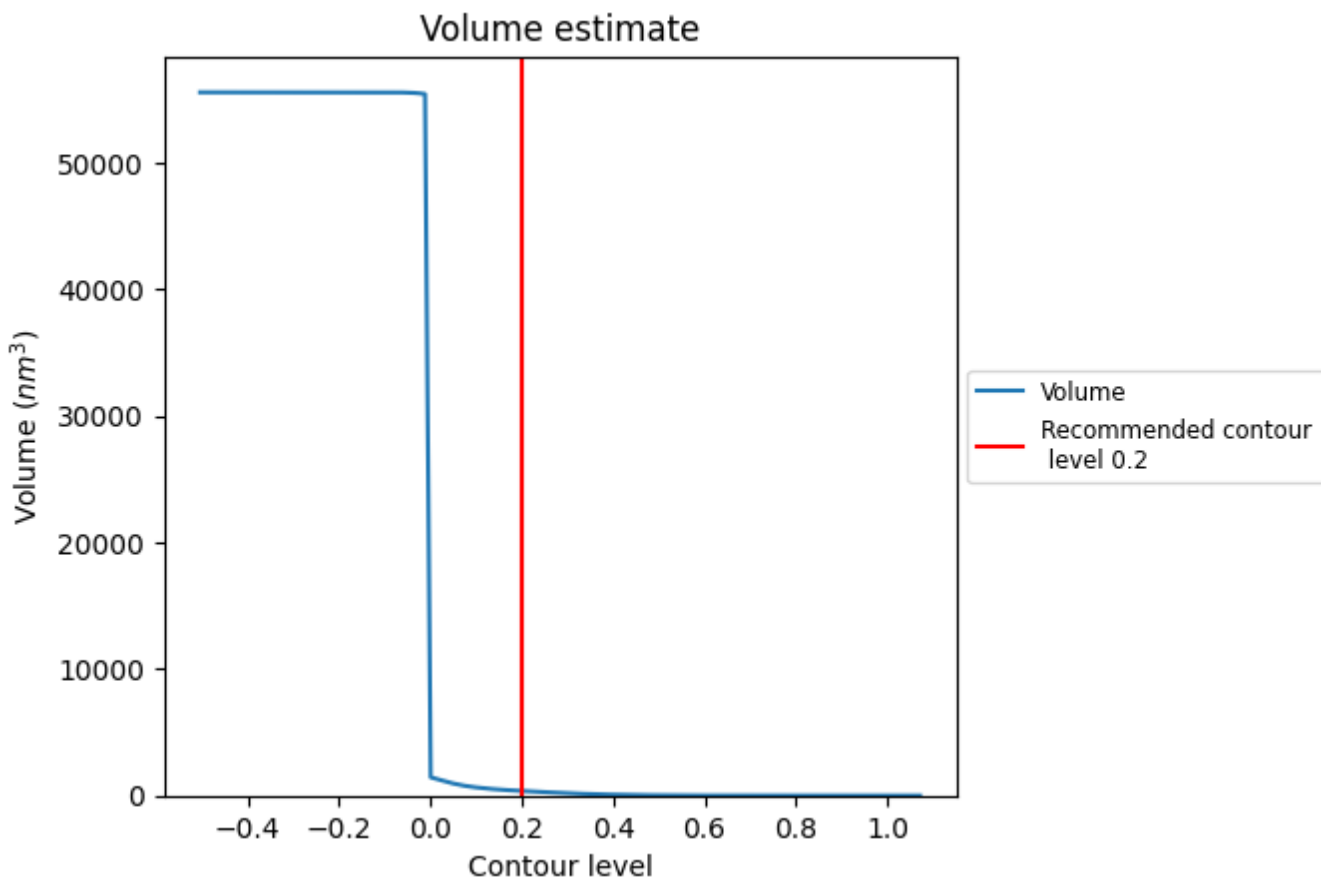
This section contains the results of statistical analysis of the map.

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.

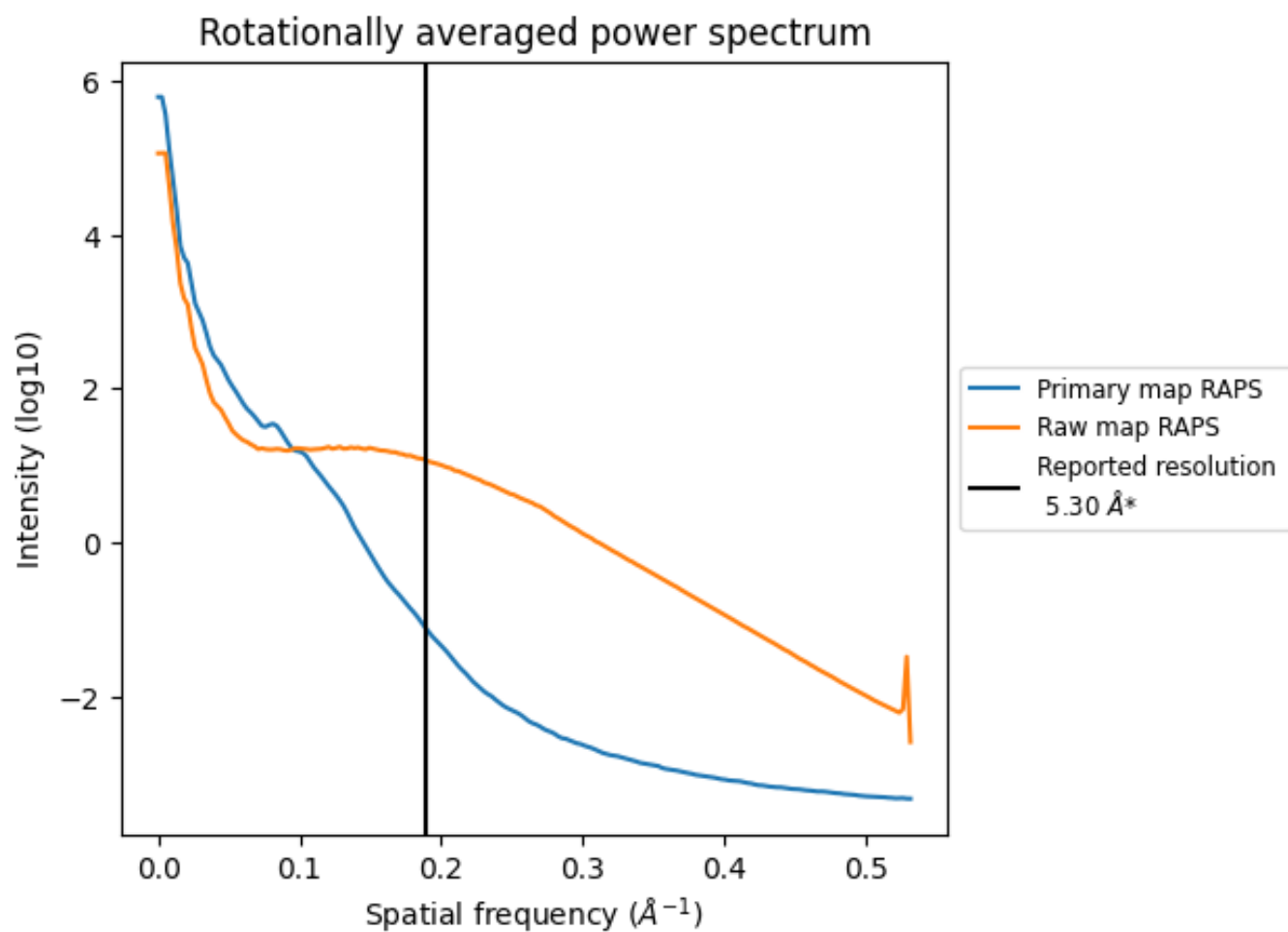
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 365 nm³; this corresponds to an approximate mass of 330 kDa.

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

7.3 Rotationally averaged power spectrum (i)

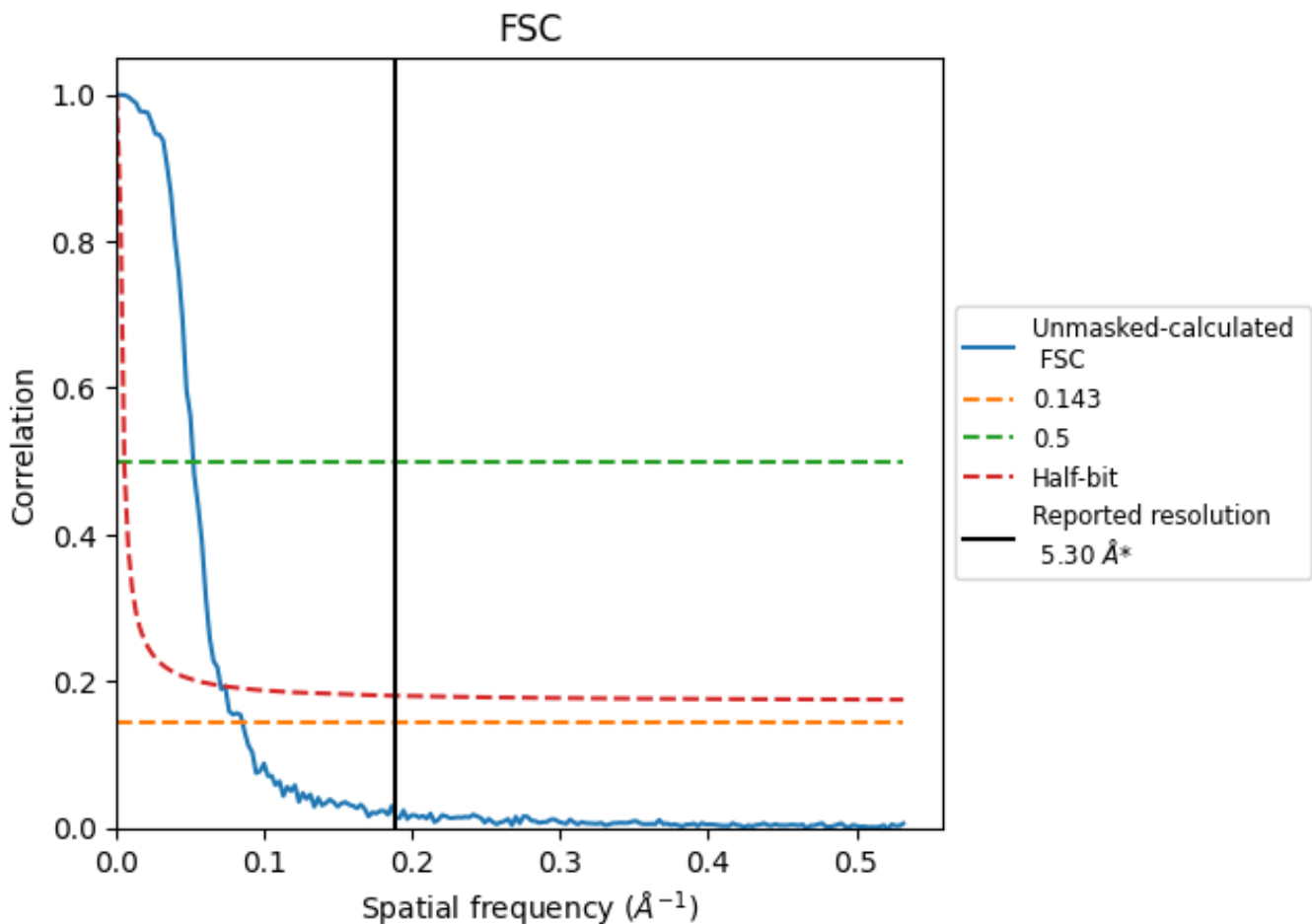


*Reported resolution corresponds to spatial frequency of 0.189 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.189 Å⁻¹

8.2 Resolution estimates [i](#)

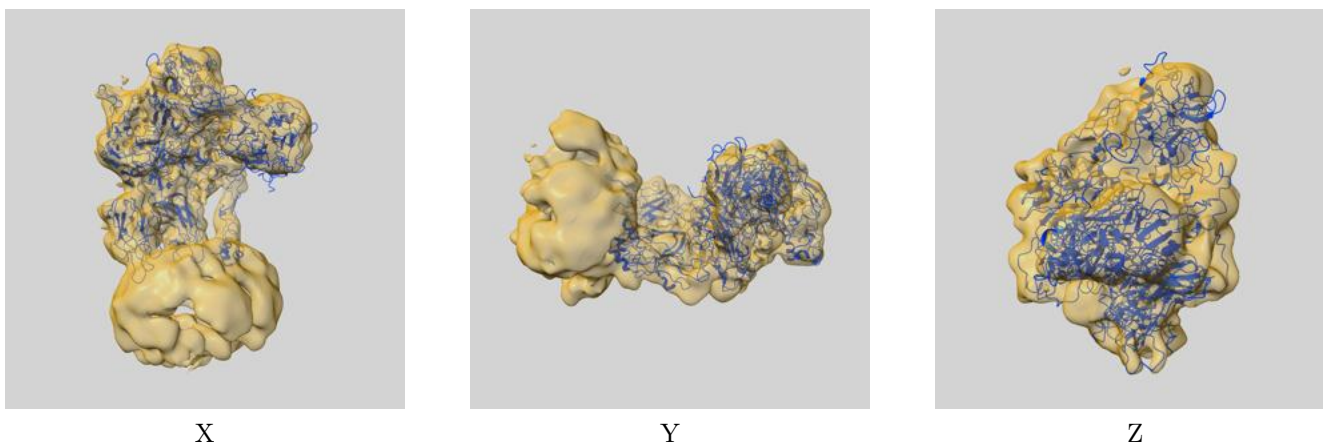
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	5.30	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	11.75	19.31	14.22

*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 11.75 differs from the reported value 5.3 by more than 10 %

9 Map-model fit [i](#)

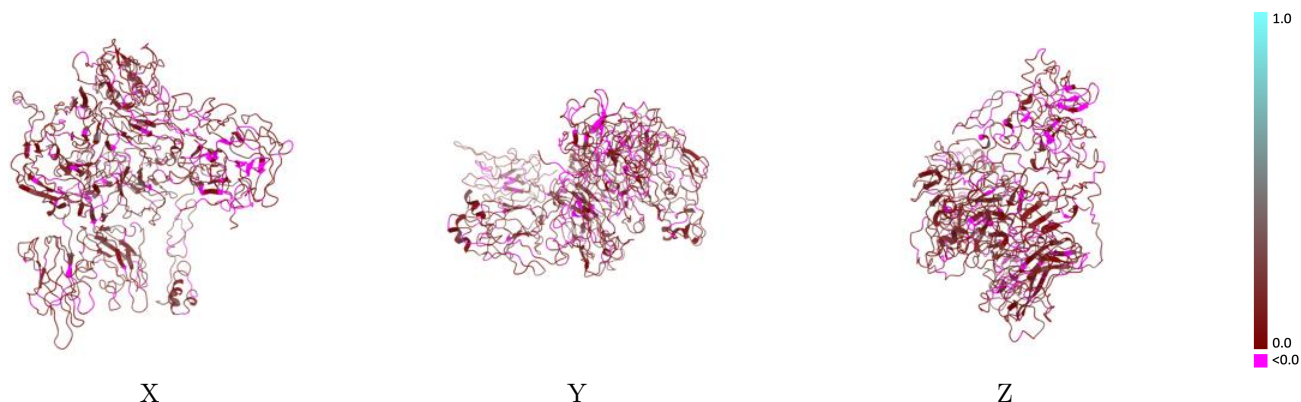
This section contains information regarding the fit between EMDB map EMD-26061 and PDB model 7TPQ. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)



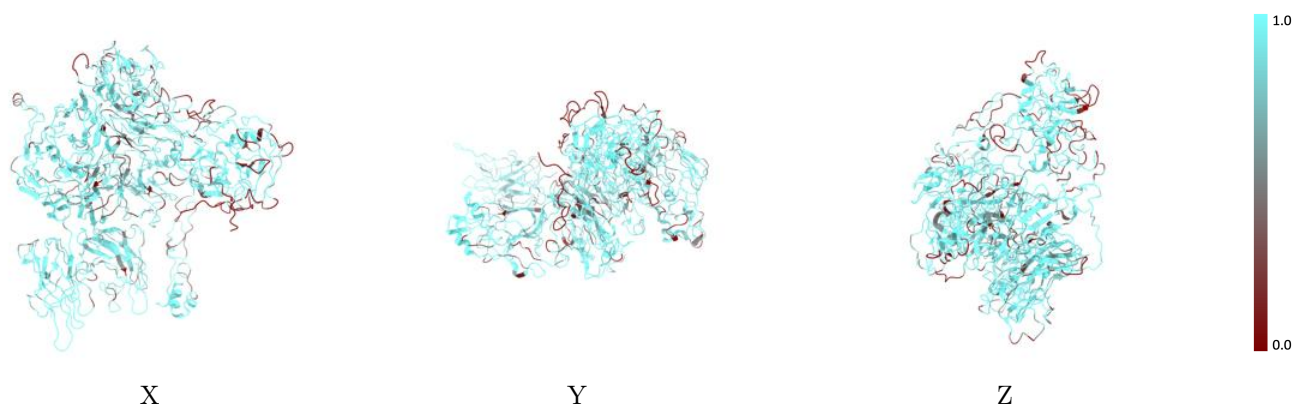
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



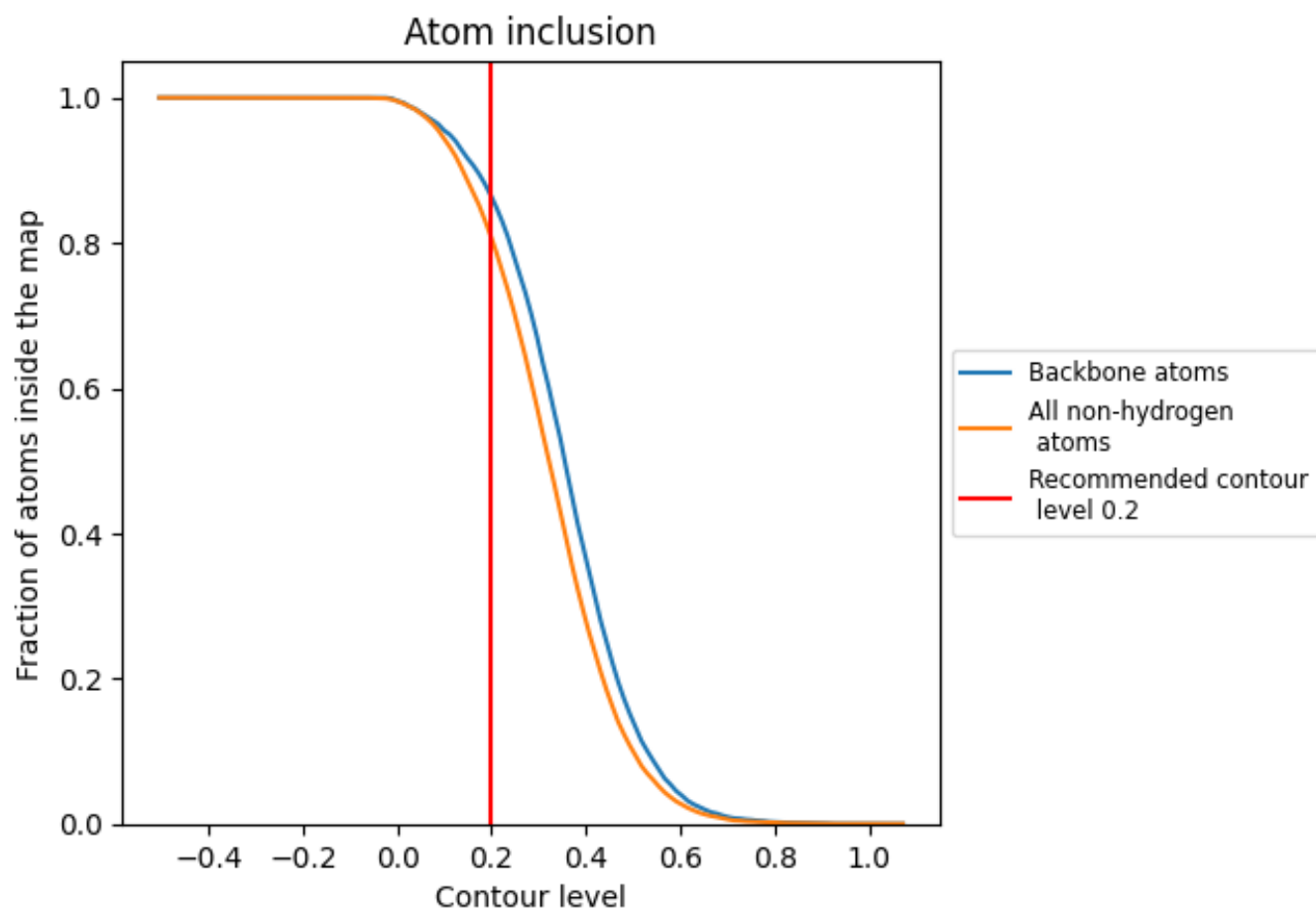
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).











9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 81% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8064	 0.1090
A	 0.6794	 0.0990
B	 0.7731	 0.0660
C	 0.8023	 0.1070
D	 0.8503	 0.1300

