



Full wwPDB EM Validation Report ⓘ

Nov 29, 2022 – 03:42 AM EST

PDB ID : 7UT4
EMDB ID : EMD-26754
Title : Gea2 closed/closed conformation (composite structure)
Authors : Muccini, A.; Fromme, J.C.
Deposited on : 2022-04-26
Resolution : 3.90 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) : 2.31.2

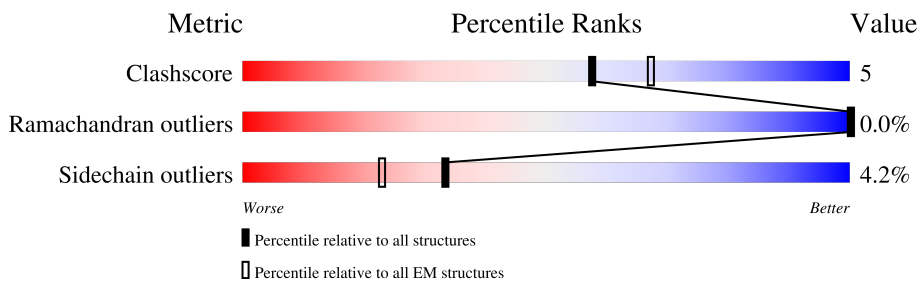
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.90 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	1459	
1	B	1459	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 39740 atoms, of which 20092 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.

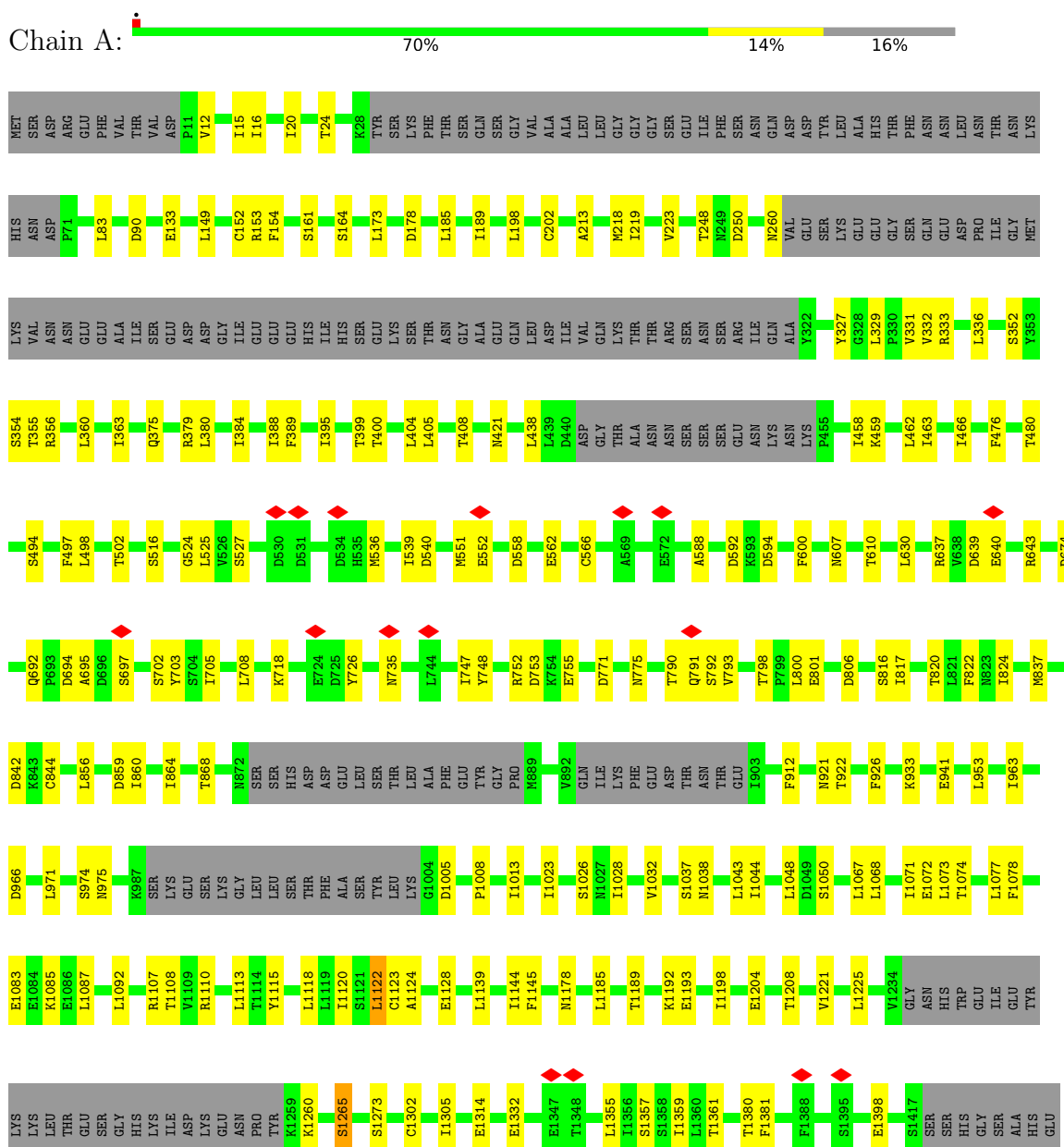
- Molecule 1 is a protein called GEA2 isoform 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	A	1224	Total	C	H	N	O	S	0	0
			19870	6306	10046	1603	1878	37		
1	B	1224	Total	C	H	N	O	S	0	0
			19870	6306	10046	1603	1878	37		

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: GEA2 isoform 1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	83665	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1.0	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	125.253	Depositor
Minimum map value	-109.998	Depositor
Average map value	0.002	Depositor
Map value standard deviation	1.500	Depositor
Recommended contour level	5.85	Depositor
Map size (\AA)	499.2, 499.2, 499.2	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.664, 1.664, 1.664	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	0.31	0/9981	0.61	1/13478 (0.0%)
1	B	0.31	0/9981	0.62	1/13478 (0.0%)
All	All	0.31	0/19962	0.61	2/26956 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	842	ASP	CB-CG-OD1	9.60	126.94	118.30
1	B	842	ASP	CB-CG-OD2	8.67	126.11	118.30

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	9824	10046	10046	96	0
1	B	9824	10046	10046	105	0
All	All	19648	20092	20092	200	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 5.

All (200) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:953:LEU:HD11	1:A:1023:ILE:HD13	1.64	0.79
1:A:1032:VAL:O	1:A:1038:ASN:ND2	2.24	0.71
1:A:498:LEU:O	1:A:502:THR:OG1	2.09	0.71
1:A:705:ILE:HD12	1:A:747:ILE:HD11	1.74	0.70
1:B:1032:VAL:O	1:B:1038:ASN:ND2	2.23	0.70
1:A:674:ASP:O	1:A:692:GLN:NE2	2.26	0.69
1:B:222:THR:OG1	1:B:335:TYR:OH	2.11	0.68
1:B:12:VAL:HG23	1:B:83:LEU:HD12	1.77	0.67
1:A:637:ARG:NH1	1:A:640:GLU:OE2	2.28	0.66
1:A:1110:ARG:NH1	1:A:1144:ILE:O	2.29	0.65
1:B:1286:GLU:N	1:B:1286:GLU:OE1	2.30	0.65
1:B:1229:SER:OG	1:B:1297:GLN:NE2	2.30	0.64
1:A:153:ARG:NH1	1:A:260:ASN:OD1	2.31	0.64
1:B:951:LEU:HD11	1:B:1070:ILE:HD11	1.79	0.64
1:A:352:SER:O	1:A:355:THR:OG1	2.15	0.64
1:B:246:SER:N	1:B:753:ASP:OD1	2.31	0.63
1:B:208:GLU:OE2	1:B:211:ARG:NH2	2.31	0.63
1:A:1067:LEU:O	1:A:1071:ILE:HD12	1.98	0.63
1:B:602:ASN:ND2	1:B:606:MET:SD	2.72	0.62
1:B:1110:ARG:NH1	1:B:1144:ILE:O	2.33	0.61
1:A:607:ASN:O	1:A:610:THR:OG1	2.19	0.60
1:A:856:LEU:O	1:A:860:ILE:HD12	2.01	0.60
1:A:20:ILE:O	1:A:24:THR:HG22	2.01	0.60
1:A:637:ARG:NE	1:A:639:ASP:OD2	2.35	0.60
1:B:185:LEU:HB3	1:B:189:ILE:HD11	1.83	0.59
1:B:1224:LEU:O	1:B:1228:ILE:HD12	2.02	0.59
1:A:185:LEU:HB3	1:A:189:ILE:HD11	1.84	0.59
1:A:12:VAL:O	1:A:16:ILE:HD12	2.02	0.58
1:A:438:LEU:HD21	1:A:463:ILE:HD11	1.85	0.58
1:A:198:LEU:HD22	1:A:218:MET:CE	2.33	0.58
1:A:798:THR:OG1	1:A:801:GLU:OE1	2.21	0.58
1:A:1083:GLU:OE2	1:A:1085:LYS:N	2.36	0.58
1:A:1204:GLU:O	1:A:1208:THR:HG23	2.03	0.58
1:B:480:THR:HG23	1:B:494:SER:OG	2.03	0.58
1:A:480:THR:HG23	1:A:494:SER:OG	2.04	0.57
1:A:1044:ILE:HD11	1:A:1077:LEU:HB3	1.87	0.57
1:A:551:MET:SD	1:A:552:GLU:N	2.78	0.57
1:A:933:LYS:O	1:A:1037:SER:OG	2.19	0.57
1:B:941:GLU:N	1:B:941:GLU:OE2	2.38	0.57
1:A:1192:LYS:NZ	1:A:1193:GLU:OE1	2.36	0.56
1:B:418:ILE:HG22	1:B:419:LEU:HD23	1.87	0.56
1:B:203:ASN:ND2	1:B:255:ASP:OD2	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1286:GLU:O	1:B:1290:ILE:HD12	2.05	0.56
1:A:771:ASP:O	1:A:775:ASN:ND2	2.38	0.56
1:A:1221:VAL:O	1:A:1225:LEU:HD22	2.06	0.56
1:B:943:TRP:O	1:B:947:VAL:HG12	2.05	0.55
1:B:943:TRP:CE3	1:B:946:ILE:HD11	2.40	0.55
1:B:486:CYS:SG	1:B:487:ASN:N	2.80	0.55
1:B:919:GLN:HE21	1:B:919:GLN:HA	1.72	0.55
1:A:941:GLU:N	1:A:941:GLU:OE1	2.39	0.55
1:B:694:ASP:N	1:B:697:SER:OG	2.40	0.55
1:B:1044:ILE:HD11	1:B:1077:LEU:HB3	1.89	0.55
1:A:375:GLN:NE2	1:A:790:THR:OG1	2.40	0.54
1:A:953:LEU:CD1	1:A:1023:ILE:HD13	2.36	0.54
1:B:476:PHE:O	1:B:480:THR:HG22	2.08	0.54
1:B:20:ILE:O	1:B:24:THR:HG22	2.08	0.53
1:B:1376:THR:OG1	1:B:1377:SER:N	2.41	0.53
1:A:219:ILE:O	1:A:223:VAL:HG23	2.09	0.53
1:A:1023:ILE:O	1:A:1026:SER:OG	2.25	0.53
1:A:1332:GLU:HG2	1:A:1380:THR:HG23	1.90	0.53
1:A:820:THR:O	1:A:824:ILE:HG22	2.08	0.53
1:B:856:LEU:O	1:B:860:ILE:HG22	2.08	0.53
1:B:166:LEU:HD12	1:B:200:LEU:HD21	1.90	0.53
1:A:133:GLU:N	1:A:133:GLU:OE1	2.41	0.53
1:A:588:ALA:N	1:A:594:ASP:OD2	2.41	0.53
1:B:645:LEU:HD23	1:B:646:LEU:HD12	1.91	0.53
1:B:920:LEU:HD12	1:B:920:LEU:O	2.10	0.52
1:B:170:VAL:CG1	1:B:217:THR:HG21	2.40	0.52
1:B:149:LEU:HD11	1:B:172:LEU:HD22	1.91	0.52
1:A:1357:SER:O	1:A:1361:THR:OG1	2.26	0.51
1:A:1072:GLU:OE2	1:A:1115:TYR:OH	2.24	0.51
1:B:1022:CYS:SG	1:B:1023:ILE:N	2.84	0.51
1:A:333:ARG:NH1	1:A:379:ARG:O	2.42	0.51
1:B:1011:GLU:N	1:B:1011:GLU:OE1	2.43	0.51
1:A:702:SER:HA	1:A:705:ILE:HG22	1.93	0.51
1:A:864:ILE:HG23	1:A:921:ASN:OD1	2.11	0.51
1:A:694:ASP:N	1:A:697:SER:OG	2.43	0.51
1:B:1360:LEU:HD12	1:B:1388:PHE:CZ	2.45	0.51
1:B:488:LEU:O	1:B:605:ARG:NH1	2.44	0.51
1:A:462:LEU:O	1:A:466:ILE:HG22	2.10	0.51
1:A:15:ILE:HG21	1:A:83:LEU:HD12	1.92	0.50
1:B:940:LYS:NZ	1:B:1042:ASP:OD1	2.44	0.50
1:A:1265:SER:OG	1:A:1305:ILE:HD13	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1325:THR:OG1	1:B:1328:MET:SD	2.68	0.50
1:A:792:SER:OG	1:A:793:VAL:N	2.45	0.50
1:B:502:THR:HG21	1:B:816:SER:HB3	1.94	0.50
1:B:1405:ILE:HD12	1:B:1408:LYS:HE3	1.94	0.50
1:A:476:PHE:O	1:A:480:THR:HG22	2.12	0.50
1:B:912:PHE:HZ	1:B:922:THR:HG21	1.76	0.50
1:B:716:GLN:O	1:B:716:GLN:NE2	2.45	0.49
1:B:198:LEU:HD21	1:B:342:LEU:HD21	1.94	0.49
1:A:404:LEU:O	1:A:408:THR:HG22	2.12	0.49
1:A:824:ILE:HG23	1:A:837:MET:CE	2.43	0.49
1:B:90:ASP:OD1	1:B:90:ASP:N	2.45	0.49
1:A:161:SER:O	1:A:164:SER:OG	2.30	0.49
1:A:395:ILE:HG23	1:A:399:THR:HG21	1.95	0.49
1:A:974:SER:OG	1:A:975:ASN:N	2.45	0.49
1:B:1009:THR:O	1:B:1013:ILE:HD12	2.13	0.48
1:A:438:LEU:O	1:A:459:LYS:NZ	2.46	0.48
1:B:926:PHE:CZ	1:B:1028:ILE:HD13	2.48	0.48
1:B:791:GLN:OE1	1:B:791:GLN:N	2.46	0.48
1:B:1064:GLU:HG3	1:B:1108:THR:HG21	1.95	0.48
1:A:748:TYR:OH	1:A:752:ARG:NH2	2.47	0.47
1:A:1123:CYS:SG	1:A:1124:ALA:N	2.87	0.47
1:B:919:GLN:HE21	1:B:919:GLN:CA	2.27	0.47
1:B:384:ILE:HA	1:B:388:ILE:HD12	1.97	0.47
1:A:643:ARG:NH1	1:A:755:GLU:O	2.47	0.47
1:B:89:LEU:HD22	1:B:92:ILE:HG21	1.96	0.47
1:A:336:LEU:HD12	1:A:380:LEU:HD23	1.96	0.47
1:A:1068:LEU:HD11	1:A:1108:THR:HG22	1.96	0.47
1:A:1092:LEU:HD11	1:A:1120:ILE:HD13	1.96	0.47
1:A:363:ILE:HD11	1:A:388:ILE:HD13	1.96	0.47
1:B:1123:CYS:SG	1:B:1124:ALA:N	2.88	0.47
1:A:458:ILE:O	1:A:462:LEU:HD23	2.16	0.46
1:B:1070:ILE:O	1:B:1074:THR:OG1	2.28	0.46
1:A:562:GLU:OE2	1:A:562:GLU:N	2.43	0.46
1:A:694:ASP:OD1	1:A:695:ALA:N	2.49	0.46
1:A:971:LEU:HD11	1:A:1118:LEU:HD22	1.96	0.46
1:B:912:PHE:CZ	1:B:922:THR:HG21	2.50	0.46
1:A:963:ILE:HD11	1:A:1073:LEU:CD2	2.45	0.46
1:A:1359:ILE:H	1:A:1359:ILE:HD12	1.80	0.46
1:B:820:THR:O	1:B:824:ILE:HG22	2.14	0.46
1:B:891:LEU:HD23	1:B:1061:ARG:HE	1.80	0.46
1:B:1149:PHE:O	1:B:1152:THR:HG22	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:822:PHE:CE2	1:B:841:LEU:HD21	2.51	0.46
1:B:571:ASN:ND2	1:B:607:ASN:OD1	2.49	0.46
1:B:15:ILE:HD11	1:B:130:ILE:HD12	1.97	0.46
1:B:487:ASN:OD1	1:B:488:LEU:N	2.49	0.46
1:B:1370:TYR:O	1:B:1374:GLY:N	2.48	0.46
1:B:943:TRP:HE3	1:B:946:ILE:HD11	1.79	0.45
1:B:1044:ILE:HG21	1:B:1087:LEU:HD11	1.97	0.45
1:B:159:GLN:N	1:B:159:GLN:OE1	2.50	0.45
1:B:891:LEU:HD21	1:B:957:LEU:CD1	2.46	0.45
1:B:1162:LEU:O	1:B:1162:LEU:HD23	2.16	0.45
1:A:525:LEU:HD22	1:A:817:ILE:HD12	1.99	0.45
1:B:136:GLN:OE1	1:B:137:ASN:N	2.49	0.45
1:A:213:ALA:HB2	1:B:167:LEU:HD21	1.99	0.45
1:B:1302:CYS:SG	1:B:1303:LYS:N	2.90	0.45
1:B:551:MET:SD	1:B:552:GLU:N	2.90	0.45
1:A:1048:LEU:HD13	1:A:1074:THR:HG21	1.99	0.44
1:B:919:GLN:HA	1:B:919:GLN:NE2	2.31	0.44
1:A:1044:ILE:HG21	1:A:1087:LEU:HD11	1.99	0.44
1:B:506:LEU:O	1:B:509:SER:OG	2.36	0.44
1:A:816:SER:O	1:A:820:THR:HG23	2.18	0.44
1:A:1078:PHE:CE2	1:A:1122:LEU:HD13	2.53	0.44
1:B:250:ASP:OD1	1:B:250:ASP:N	2.50	0.44
1:A:856:LEU:HD23	1:A:860:ILE:HD11	1.99	0.44
1:B:161:SER:O	1:B:164:SER:OG	2.35	0.43
1:B:1368:LEU:HD22	1:B:1411:ILE:CD1	2.48	0.43
1:B:540:ASP:OD2	1:B:542:GLU:N	2.49	0.43
1:B:1286:GLU:O	1:B:1289:ALA:HB3	2.18	0.43
1:A:1043:LEU:HD12	1:A:1043:LEU:O	2.18	0.43
1:A:1189:THR:HG22	1:A:1198:ILE:HG21	2.00	0.43
1:A:363:ILE:CD1	1:A:388:ILE:HD13	2.48	0.43
1:A:926:PHE:CZ	1:A:1028:ILE:HD13	2.54	0.43
1:A:868:THR:HG23	1:A:922:THR:HG22	2.01	0.43
1:B:538:ASP:OD1	1:B:539:ILE:N	2.51	0.43
1:B:892:VAL:HG23	1:B:983:ILE:O	2.19	0.43
1:B:694:ASP:OD2	1:B:695:ALA:N	2.51	0.43
1:B:1178:ASN:OD1	1:B:1181:PHE:N	2.48	0.43
1:B:1113:LEU:HD22	1:B:1145:PHE:CE2	2.53	0.42
1:A:1113:LEU:HD22	1:A:1145:PHE:CZ	2.54	0.42
1:B:1367:TYR:HH	1:B:1381:PHE:HD1	1.66	0.42
1:A:384:ILE:HA	1:A:388:ILE:HD12	2.01	0.42
1:B:489:ASP:OD1	1:B:490:ARG:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1206:ILE:HG21	1:B:1271:LEU:HD21	2.01	0.42
1:B:1263:GLU:HA	1:B:1266:LEU:HD12	2.02	0.42
1:A:149:LEU:HD13	1:A:173:LEU:HD12	2.02	0.42
1:B:194:LEU:HD23	1:B:335:TYR:CD1	2.55	0.42
1:B:352:SER:OG	1:B:353:TYR:N	2.53	0.42
1:B:438:LEU:HD21	1:B:463:ILE:HD11	2.02	0.42
1:B:963:ILE:HD12	1:B:964:PHE:CD2	2.55	0.42
1:A:329:LEU:O	1:A:332:VAL:HG22	2.20	0.41
1:A:912:PHE:CZ	1:A:922:THR:HG21	2.55	0.41
1:B:686:ASP:N	1:B:686:ASP:OD1	2.53	0.41
1:A:1185:LEU:O	1:A:1189:THR:HG23	2.20	0.41
1:B:395:ILE:O	1:B:399:THR:OG1	2.37	0.41
1:B:471:THR:OG1	1:B:472:ARG:N	2.53	0.41
1:B:891:LEU:HD21	1:B:957:LEU:HD11	2.01	0.41
1:B:1357:SER:O	1:B:1361:THR:OG1	2.36	0.41
1:A:539:ILE:HD11	1:A:800:LEU:HD22	2.03	0.41
1:B:1312:THR:O	1:B:1316:THR:OG1	2.35	0.41
1:A:708:LEU:HD23	1:A:708:LEU:O	2.20	0.41
1:A:1355:LEU:O	1:A:1359:ILE:HD12	2.19	0.41
1:A:356:ARG:O	1:A:360:LEU:HD23	2.21	0.41
1:A:524:GLY:O	1:A:527:SER:OG	2.32	0.41
1:A:824:ILE:HG23	1:A:837:MET:HE3	2.02	0.41
1:A:1008:PRO:HB2	1:A:1013:ILE:HD11	2.02	0.41
1:B:1039:ILE:H	1:B:1039:ILE:HD12	1.85	0.41
1:B:932:ASN:OD1	1:B:933:LYS:N	2.53	0.41
1:A:248:THR:O	1:A:248:THR:OG1	2.31	0.41
1:B:83:LEU:HD13	1:B:86:LEU:HG	2.03	0.41
1:B:162:ASP:OD2	1:B:162:ASP:N	2.53	0.41
1:B:367:LEU:HD12	1:B:384:ILE:HD11	2.03	0.41
1:B:645:LEU:HD13	1:B:665:PHE:HD1	1.86	0.41
1:B:1221:VAL:O	1:B:1225:LEU:HD23	2.21	0.41
1:A:327:TYR:CE1	1:A:331:VAL:HG11	2.56	0.41
1:A:336:LEU:CD1	1:A:380:LEU:HD23	2.51	0.40
1:A:399:THR:OG1	1:A:400:THR:N	2.54	0.40
1:B:385:SER:OG	1:B:426:GLN:NE2	2.55	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	1208/1459 (83%)	1150 (95%)	58 (5%)	0	100	100
1	B	1208/1459 (83%)	1143 (95%)	64 (5%)	1 (0%)	51	84
All	All	2416/2918 (83%)	2293 (95%)	122 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1233	ALA

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	1137/1346 (84%)	1093 (96%)	44 (4%)	32	59
1	B	1137/1346 (84%)	1085 (95%)	52 (5%)	27	55
All	All	2274/2692 (84%)	2178 (96%)	96 (4%)	33	57

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	ASP
1	A	152	CYS
1	A	154	PHE
1	A	178	ASP
1	A	202	CYS

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Mol	Chain	Res	Type
1	A	250	ASP
1	A	354	SER
1	A	389	PHE
1	A	405	LEU
1	A	421	ASN
1	A	497	PHE
1	A	516	SER
1	A	536	MET
1	A	540	ASP
1	A	558	ASP
1	A	566	CYS
1	A	592	ASP
1	A	600	PHE
1	A	630	LEU
1	A	703	TYR
1	A	718	LYS
1	A	726	TYR
1	A	735	ASN
1	A	753	ASP
1	A	791	GLN
1	A	806	ASP
1	A	822	PHE
1	A	844	CYS
1	A	859	ASP
1	A	966	ASP
1	A	1005	ASP
1	A	1050	SER
1	A	1107	ARG
1	A	1122	LEU
1	A	1128	GLU
1	A	1139	LEU
1	A	1178	ASN
1	A	1260	LYS
1	A	1265	SER
1	A	1273	SER
1	A	1302	CYS
1	A	1314	GLU
1	A	1381	PHE
1	A	1398	GLU
1	B	124	LYS
1	B	174	ARG
1	B	175	SER

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Mol	Chain	Res	Type
1	B	183	ASP
1	B	206	ARG
1	B	250	ASP
1	B	322	TYR
1	B	327	TYR
1	B	348	GLU
1	B	381	PHE
1	B	389	PHE
1	B	433	ARG
1	B	485	ASP
1	B	488	LEU
1	B	497	PHE
1	B	508	GLU
1	B	516	SER
1	B	527	SER
1	B	551	MET
1	B	592	ASP
1	B	605	ARG
1	B	630	LEU
1	B	639	ASP
1	B	686	ASP
1	B	702	SER
1	B	720	HIS
1	B	723	PHE
1	B	724	GLU
1	B	729	ASN
1	B	743	TYR
1	B	753	ASP
1	B	802	LEU
1	B	805	PHE
1	B	806	ASP
1	B	844	CYS
1	B	919	GLN
1	B	939	SER
1	B	940	LYS
1	B	986	ASN
1	B	1022	CYS
1	B	1027	ASN
1	B	1050	SER
1	B	1073	LEU
1	B	1107	ARG
1	B	1121	SER

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Mol	Chain	Res	Type
1	B	1139	LEU
1	B	1150	PHE
1	B	1314	GLU
1	B	1328	MET
1	B	1358	SER
1	B	1376	THR
1	B	1381	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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.

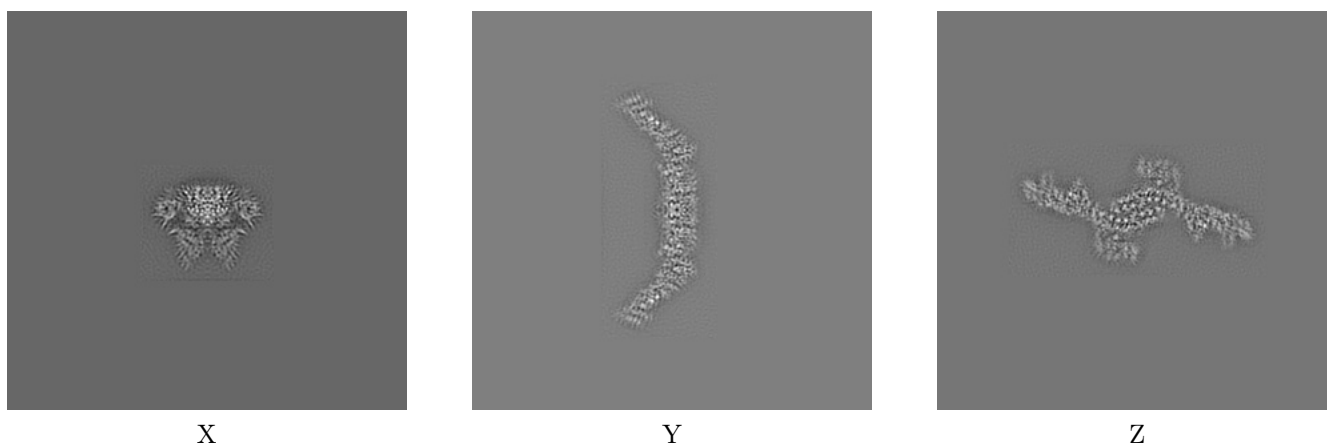
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26754. 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](#)

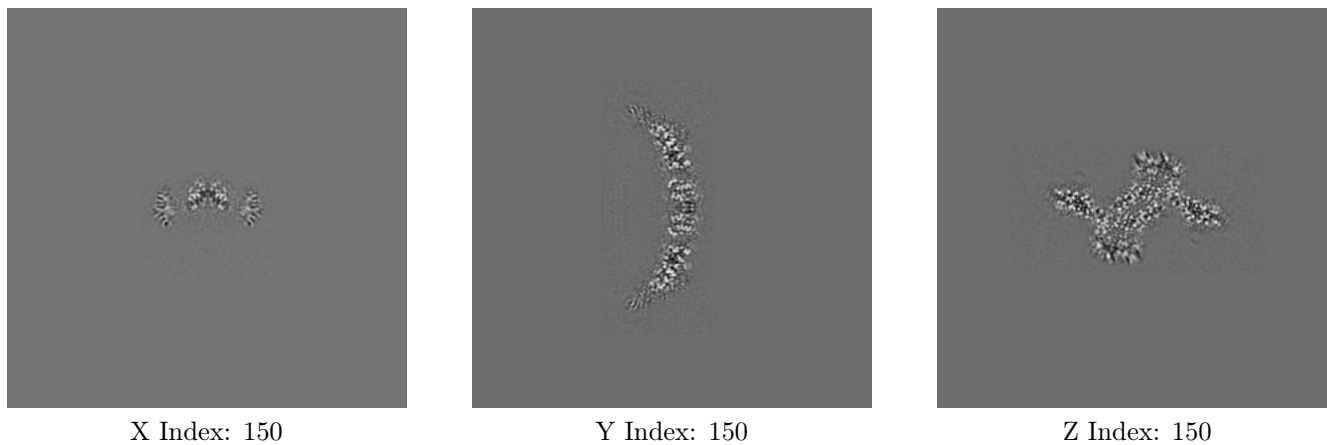
6.1.1 Primary map



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

6.2 Central slices [i](#)

6.2.1 Primary map



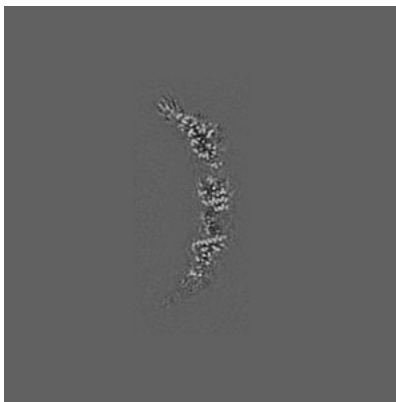
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

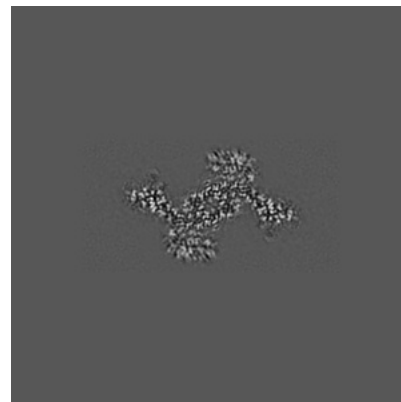
6.3.1 Primary map



X Index: 147



Y Index: 147

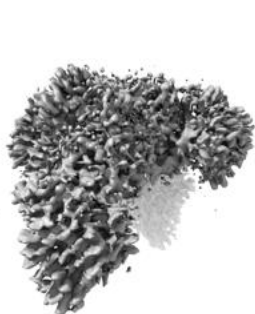


Z Index: 152

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

6.4 Orthogonal surface views [i](#)

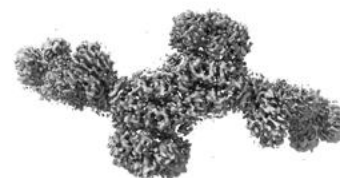
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 5.85. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

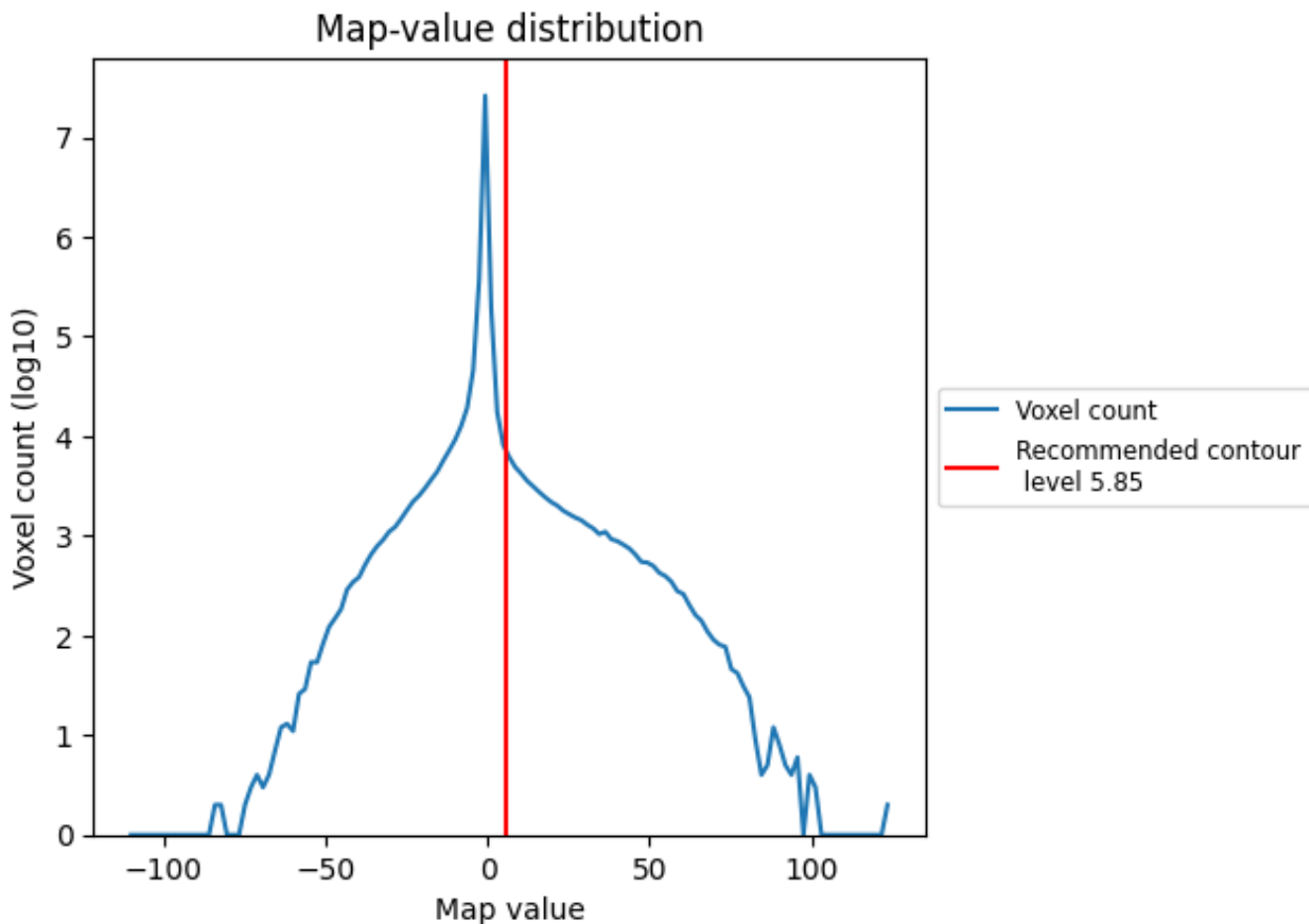
6.5 Mask visualisation

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

7 Map analysis [i](#)

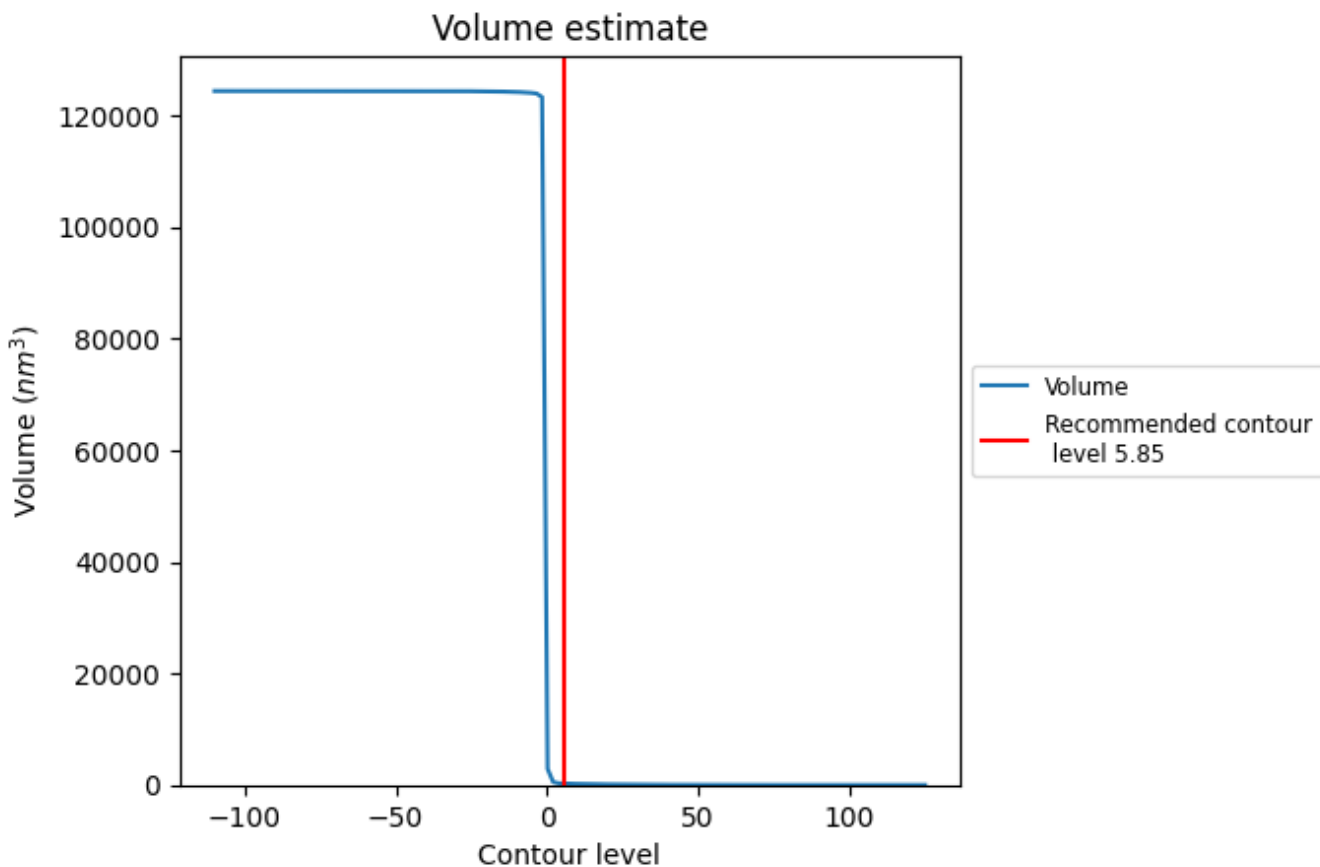
This section contains the results of statistical analysis of the map.

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.

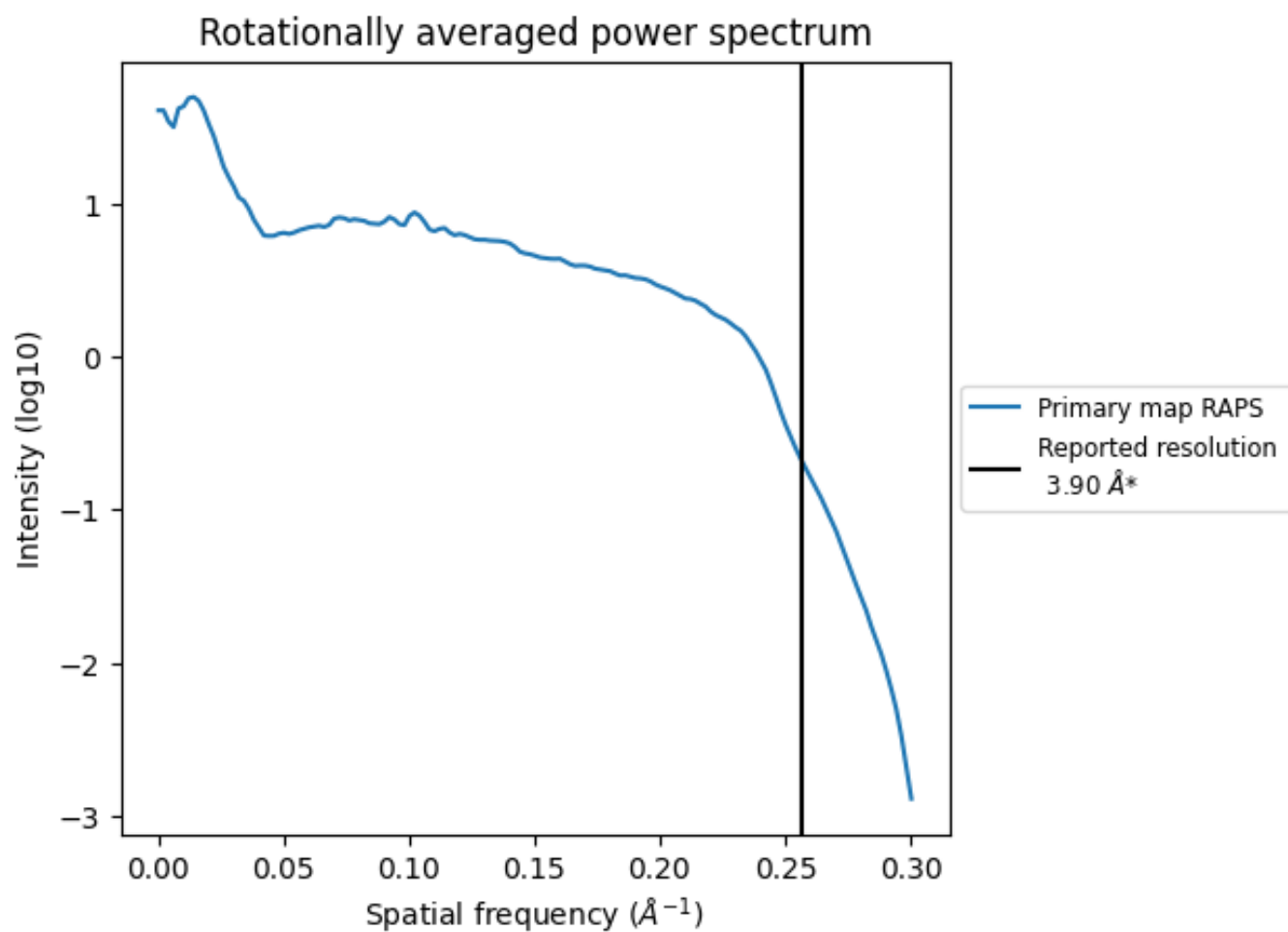
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 250 nm³; this corresponds to an approximate mass of 226 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 0.256 Å⁻¹

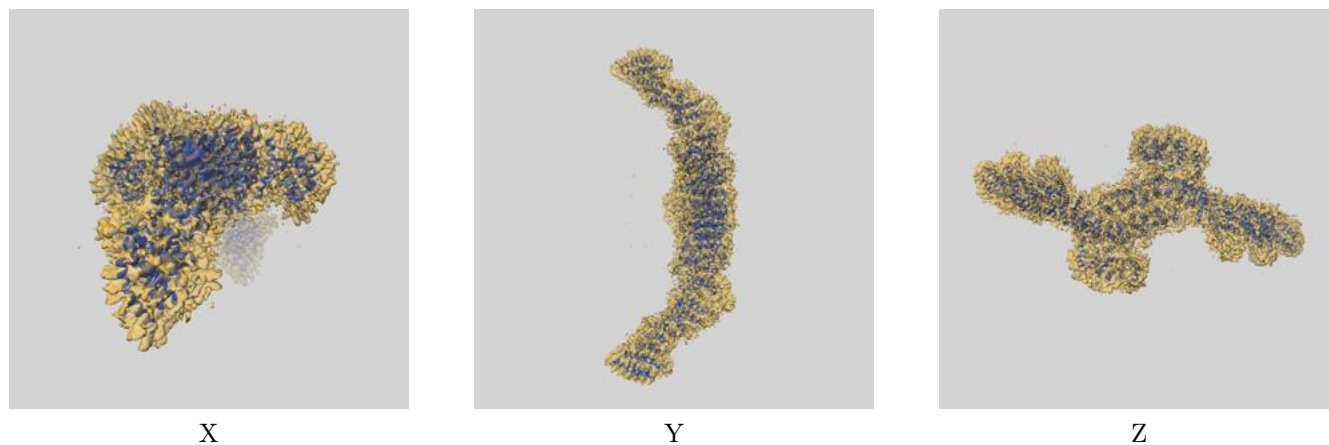
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

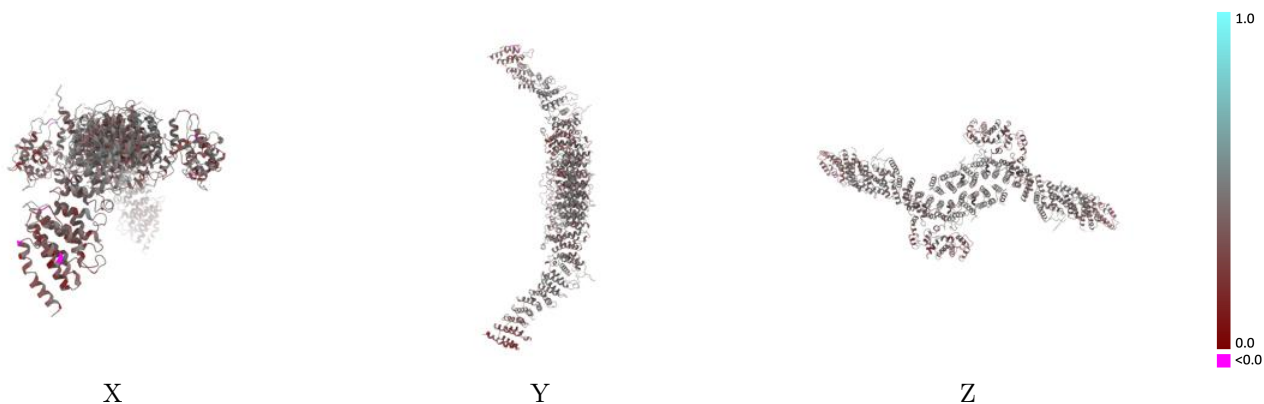
This section contains information regarding the fit between EMDB map EMD-26754 and PDB model 7UT4. Per-residue inclusion information can be found in section [3](#) on page [4](#).

9.1 Map-model overlay [i](#)



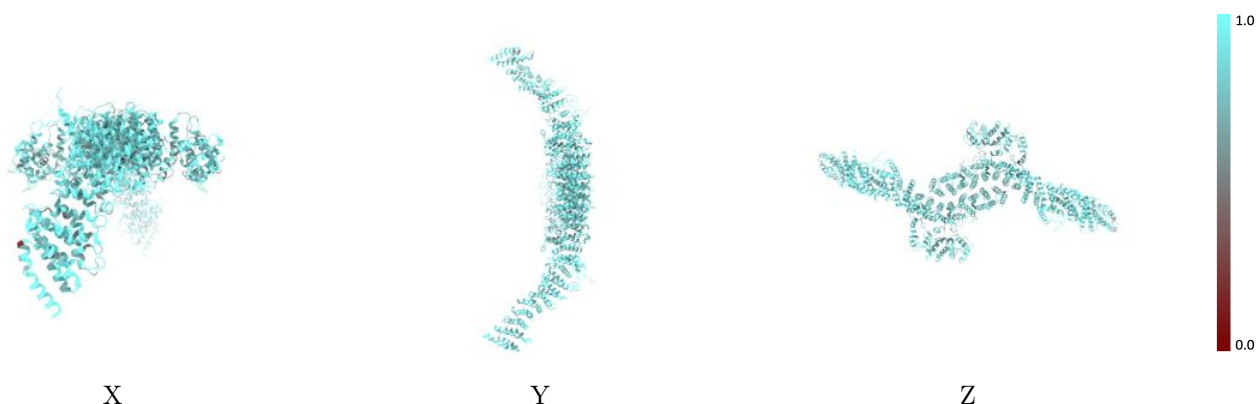
The images above show the 3D surface view of the map at the recommended contour level 5.85 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



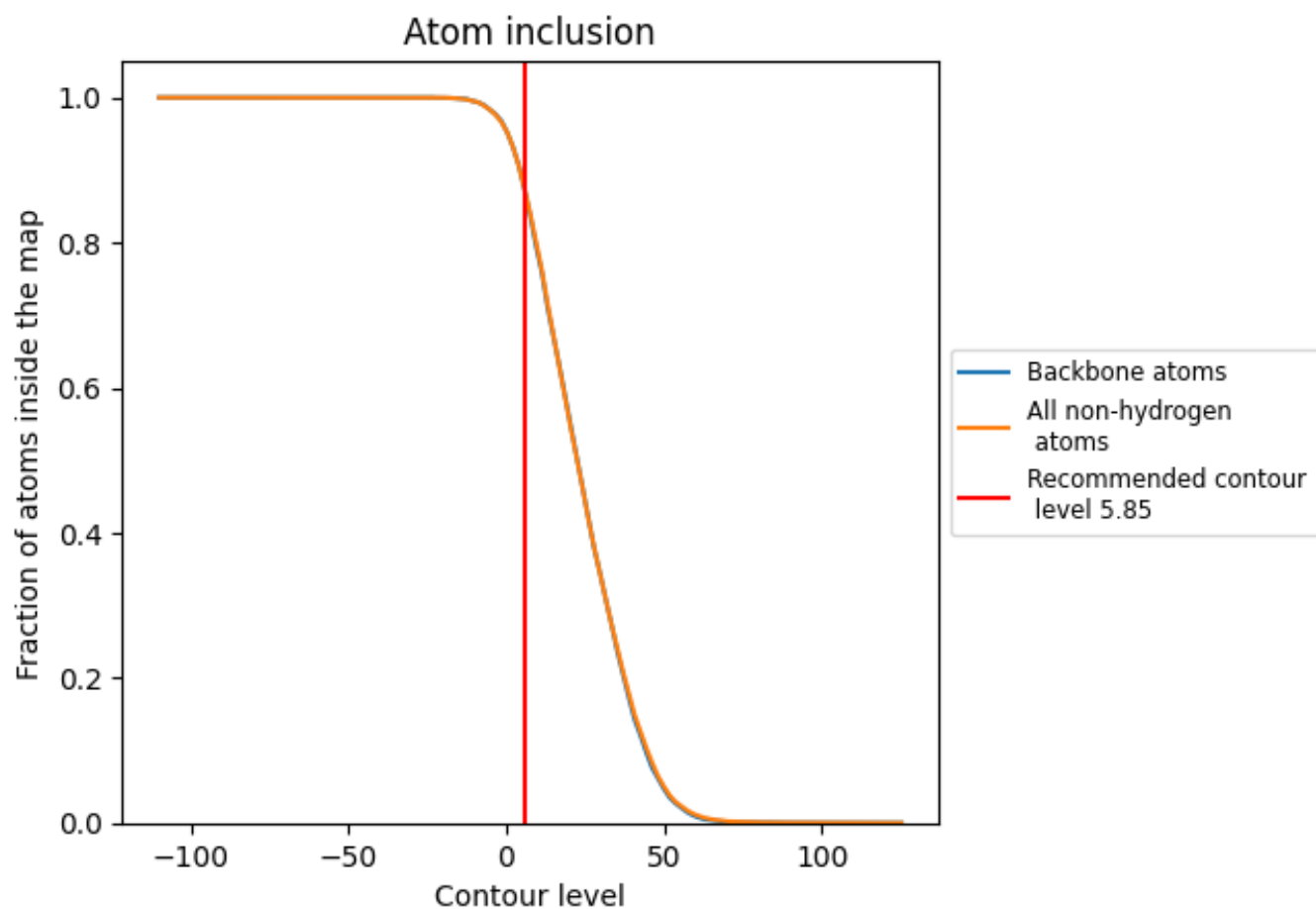
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (5.85).




9.4 Atom inclusion [i](#)



At the recommended contour level, 87% of all backbone atoms, 87% 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 (5.85) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8727	 0.4120
A	 0.8726	 0.4120
B	 0.8759	 0.4120

