Full wwPDB EM Validation Report ⓘ

Nov 29, 2022 – 03:47 AM EST

PDB ID	:	7UTH
EMDB ID	:	EMD-26770
Title	:	Gea2 open/open 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:

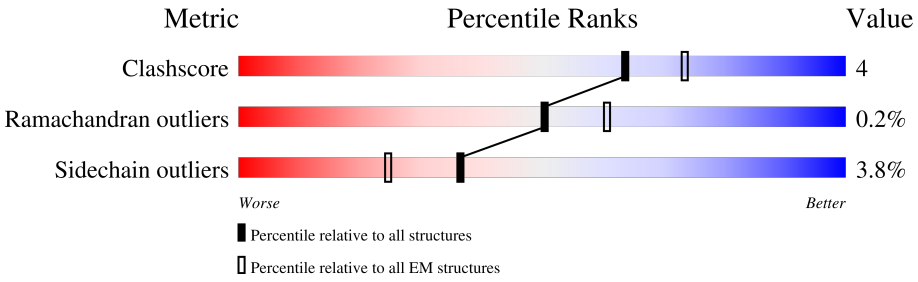
EMDB validation analysis	:	0.0.1.dev43
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

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 <=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	65% 12% 23%
1	B	1459	66% 11% 23%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 36590 atoms, of which 18572 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.

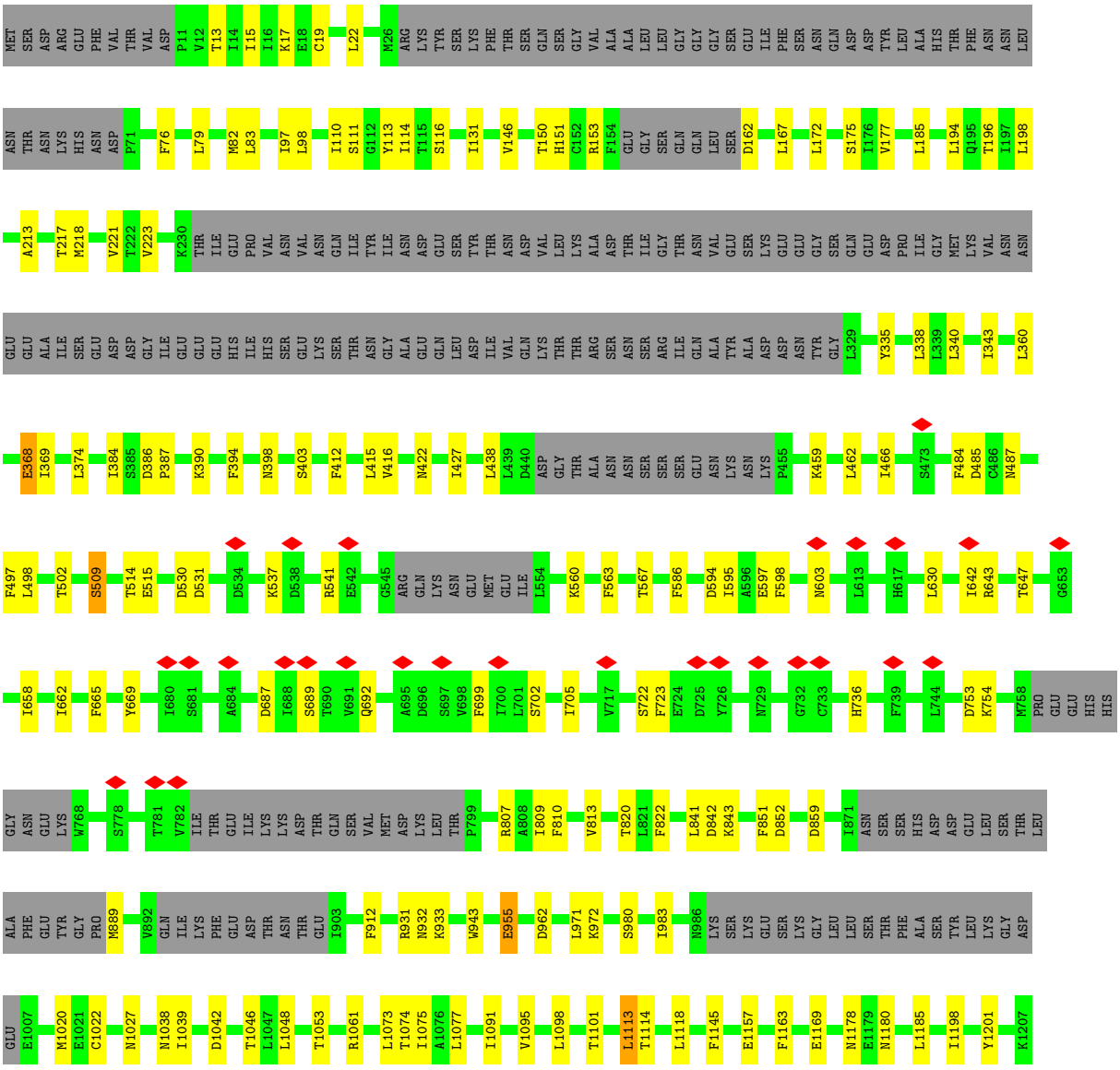
Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	1121	Total	C	H	N	O	S	0	0
			18295	5821	9286	1465	1688	35		
1	B	1121	Total	C	H	N	O	S	0	0
			18295	5821	9286	1465	1688	35		

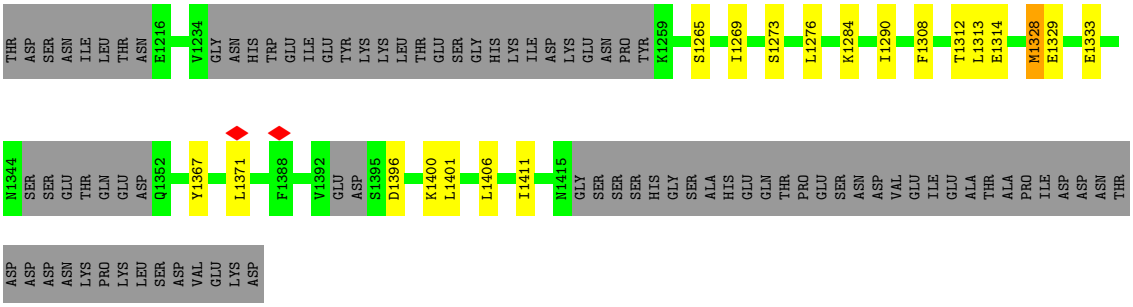
3 Residue-property plots (i)

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

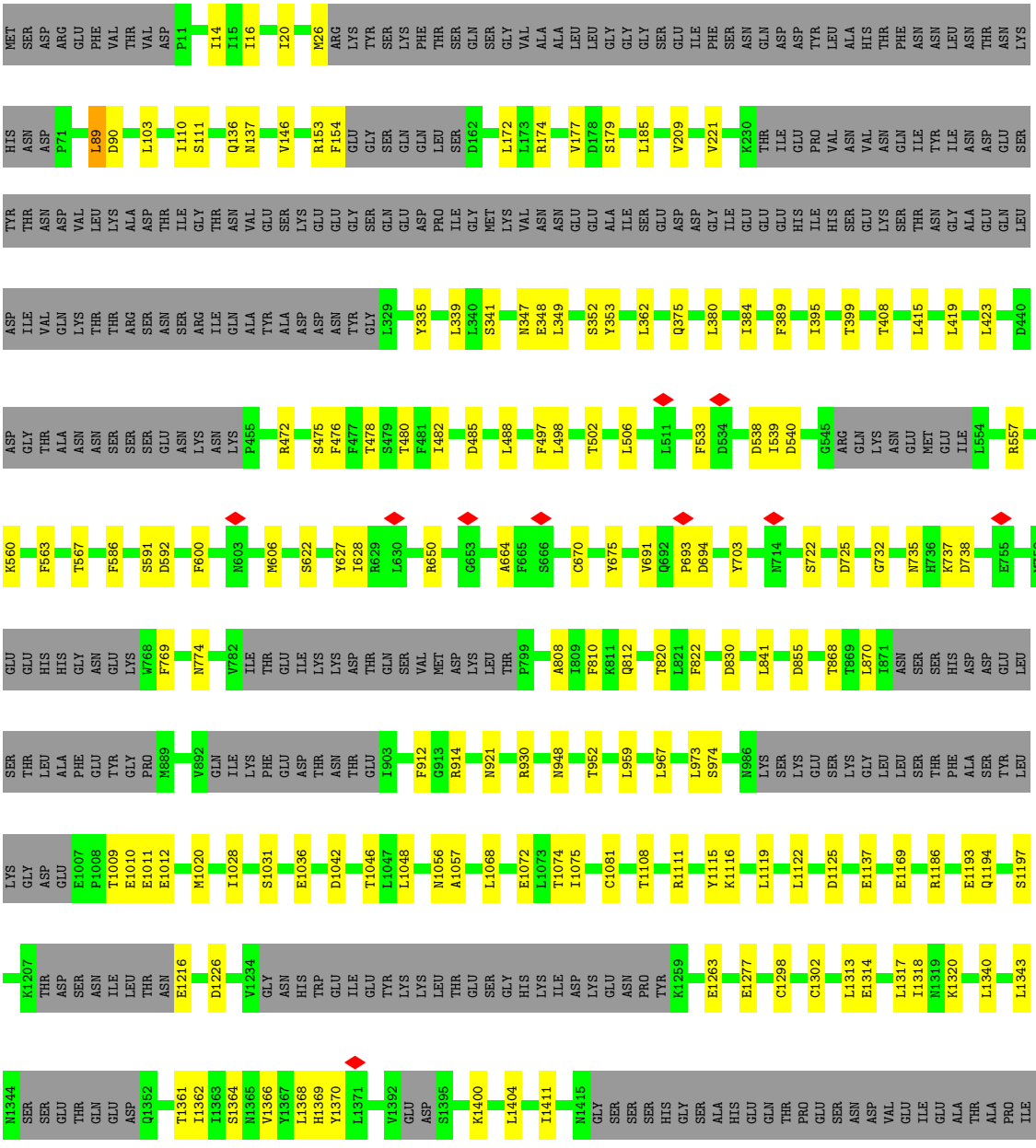
Chain A:





• Molecule 1: GEA2 isoform 1

Chain B:




ASP
ASP
ASN
THR
ASP
ASP
ASP
ASN
LYS
PRO
LYS
LEU
SER
ASP
VAL
GLU
LYS
ASP

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	74385	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 ($e^-/Å^2$)	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	115.814	Depositor
Minimum map value	-104.415	Depositor
Average map value	-0.069	Depositor
Map value standard deviation	1.599	Depositor
Recommended contour level	6.0	Depositor
Map size (Å)	499.2, 499.2, 499.2	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.664, 1.664, 1.664	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	0.32	0/9151	0.62	0/12344
1	B	0.32	0/9151	0.63	0/12344
All	All	0.32	0/18302	0.63	0/24688

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

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	9009	9286	9286	83	0
1	B	9009	9286	9286	76	0
All	All	18018	18572	18572	156	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 4.

All (156) 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:670:CYS:HG	1:B:675:TYR:HH	1.36	0.74
1:A:498:LEU:O	1:A:502:THR:OG1	2.09	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:488:LEU:O	1:B:560:LYS:NZ	2.24	0.69
1:A:151:HIS:O	1:A:153:ARG:NH1	2.26	0.69
1:B:557:ARG:NH2	1:B:812:GLN:OE1	2.26	0.68
1:B:592:ASP:OD2	1:B:622:SER:OG	2.11	0.68
1:B:694:ASP:OD2	1:B:735:ASN:ND2	2.26	0.68
1:A:1284:LYS:NZ	1:A:1333:GLU:OE1	2.27	0.68
1:A:530:ASP:OD1	1:A:843:LYS:NZ	2.27	0.68
1:B:1111:ARG:NH1	1:B:1115:TYR:OH	2.27	0.68
1:B:914:ARG:NH2	1:B:1011:GLU:O	2.26	0.67
1:B:1318:ILE:HD13	1:B:1366:VAL:HG23	1.75	0.67
1:B:868:THR:HG22	1:B:870:LEU:HD23	1.76	0.67
1:B:722:SER:OG	1:B:725:ASP:OD2	2.13	0.67
1:B:1072:GLU:OE1	1:B:1115:TYR:OH	2.14	0.65
1:A:394:PHE:O	1:A:398:ASN:ND2	2.29	0.65
1:B:339:LEU:HD11	1:B:362:LEU:HD23	1.77	0.65
1:B:341:SER:O	1:B:347:ASN:ND2	2.30	0.65
1:B:384:ILE:HD11	1:B:415:LEU:HD21	1.79	0.65
1:A:15:ILE:HD13	1:A:97:ILE:HD11	1.79	0.64
1:A:541:ARG:NH2	1:A:852:ASP:OD2	2.32	0.63
1:A:485:ASP:O	1:A:560:LYS:NZ	2.32	0.61
1:A:509:SER:OG	1:A:509:SER:O	2.17	0.61
1:B:1361:THR:OG1	1:B:1400:LYS:NZ	2.34	0.61
1:B:732:GLY:N	1:B:738:ASP:OD1	2.34	0.61
1:A:842:ASP:OD1	1:A:931:ARG:NH2	2.33	0.61
1:A:955:GLU:O	1:A:1061:ARG:NH2	2.33	0.60
1:A:412:PHE:O	1:A:416:VAL:HG23	2.01	0.60
1:B:1048:LEU:HD21	1:B:1074:THR:HG21	1.84	0.60
1:B:948:ASN:O	1:B:952:THR:HG22	2.02	0.59
1:A:889:MET:SD	1:A:889:MET:N	2.75	0.59
1:B:498:LEU:O	1:B:502:THR:HG22	2.02	0.59
1:A:1048:LEU:HD21	1:A:1074:THR:HG21	1.84	0.59
1:A:643:ARG:O	1:A:647:THR:HG23	2.03	0.58
1:A:1169:GLU:N	1:A:1169:GLU:OE1	2.37	0.58
1:A:1178:ASN:OD1	1:A:1180:ASN:N	2.36	0.58
1:B:1263:GLU:N	1:B:1263:GLU:OE2	2.37	0.57
1:B:1036:GLU:N	1:B:1036:GLU:OE1	2.38	0.57
1:A:162:ASP:N	1:A:162:ASP:OD1	2.37	0.57
1:A:1308:PHE:O	1:A:1312:THR:OG1	2.22	0.57
1:A:150:THR:O	1:A:151:HIS:ND1	2.37	0.57
1:A:1367:TYR:O	1:A:1371:LEU:HD12	2.05	0.57
1:A:150:THR:HB	1:A:196:THR:HG21	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:THR:OG1	1:A:515:GLU:N	2.38	0.56
1:A:972:LYS:NZ	1:A:1157:GLU:OE1	2.38	0.56
1:A:722:SER:OG	1:A:723:PHE:N	2.40	0.55
1:B:1075:ILE:HD11	1:B:1119:LEU:HD21	1.88	0.55
1:B:478:THR:O	1:B:482:ILE:HD12	2.06	0.55
1:B:1010:GLU:N	1:B:1010:GLU:OE1	2.40	0.54
1:A:809:ILE:O	1:A:813:VAL:HG12	2.07	0.54
1:A:689:SER:O	1:A:692:GLN:NE2	2.40	0.54
1:B:1216:GLU:N	1:B:1216:GLU:OE1	2.40	0.54
1:A:1073:LEU:HD12	1:A:1077:LEU:HD23	1.90	0.54
1:A:1091:ILE:O	1:A:1095:VAL:HG23	2.08	0.54
1:B:1277:GLU:OE1	1:B:1320:LYS:NZ	2.38	0.54
1:B:1009:THR:N	1:B:1012:GLU:OE2	2.41	0.54
1:A:177:VAL:HG11	1:A:221:VAL:HG23	1.89	0.54
1:B:475:SER:O	1:B:475:SER:OG	2.20	0.53
1:A:146:VAL:HG21	1:A:185:LEU:HD22	1.90	0.53
1:A:223:VAL:HG23	1:A:369:ILE:HG13	1.90	0.53
1:B:822:PHE:CE2	1:B:841:LEU:HD21	2.44	0.52
1:B:16:ILE:O	1:B:20:ILE:HD12	2.10	0.52
1:B:606:MET:SD	1:B:606:MET:N	2.83	0.52
1:A:110:ILE:HD11	1:A:114:ILE:CG1	2.39	0.52
1:A:1185:LEU:CD1	1:A:1198:ILE:HD11	2.40	0.52
1:B:136:GLN:OE1	1:B:137:ASN:N	2.43	0.52
1:B:973:LEU:HD12	1:B:974:SER:H	1.74	0.52
1:A:753:ASP:OD1	1:A:754:LYS:N	2.43	0.52
1:B:26:MET:SD	1:B:26:MET:N	2.82	0.52
1:A:416:VAL:HG22	1:A:427:ILE:HD11	1.92	0.51
1:B:1193:GLU:N	1:B:1193:GLU:OE1	2.43	0.51
1:A:1269:ILE:HG22	1:A:1312:THR:HG21	1.92	0.51
1:A:1371:LEU:HD13	1:A:1411:ILE:HD13	1.91	0.51
1:B:177:VAL:HG11	1:B:221:VAL:HG23	1.92	0.51
1:A:1113:LEU:HD12	1:A:1145:PHE:CZ	2.46	0.50
1:A:1098:LEU:O	1:A:1101:THR:HG23	2.12	0.50
1:A:1039:ILE:HD12	1:A:1077:LEU:HD13	1.93	0.49
1:B:563:PHE:O	1:B:567:THR:HG23	2.12	0.49
1:A:943:TRP:HB3	1:A:1046:THR:HG21	1.94	0.49
1:A:597:GLU:OE1	1:A:630:LEU:HD11	2.12	0.48
1:A:699:PHE:O	1:A:702:SER:OG	2.31	0.48
1:B:1313:LEU:HD12	1:B:1317:LEU:HD13	1.95	0.48
1:A:563:PHE:O	1:A:567:THR:HG23	2.14	0.48
1:A:736:HIS:O	1:A:736:HIS:ND1	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:VAL:HG22	1:A:427:ILE:CD1	2.43	0.48
1:B:476:PHE:O	1:B:480:THR:HG23	2.13	0.48
1:A:98:LEU:HD11	1:A:131:ILE:CD1	2.44	0.48
1:B:959:LEU:HD13	1:B:1028:ILE:CG2	2.43	0.47
1:B:347:ASN:O	1:B:349:LEU:N	2.47	0.47
1:B:1169:GLU:OE1	1:B:1169:GLU:N	2.47	0.47
1:B:538:ASP:OD1	1:B:539:ILE:N	2.47	0.47
1:B:670:CYS:SG	1:B:675:TYR:OH	2.55	0.47
1:B:973:LEU:HD12	1:B:974:SER:N	2.29	0.47
1:A:22:LEU:HD22	1:A:76:PHE:HE2	1.80	0.47
1:A:1042:ASP:O	1:A:1046:THR:HG22	2.15	0.47
1:A:980:SER:OG	1:A:1061:ARG:O	2.33	0.47
1:B:1318:ILE:HD11	1:B:1369:HIS:ND1	2.30	0.47
1:B:1194:GLN:O	1:B:1197:SER:OG	2.30	0.46
1:A:194:LEU:HD13	1:A:335:TYR:CD1	2.51	0.46
1:A:384:ILE:HD11	1:A:415:LEU:HD21	1.98	0.45
1:B:110:ILE:HD12	1:B:111:SER:H	1.80	0.45
1:B:1042:ASP:O	1:B:1046:THR:HG22	2.15	0.45
1:B:1364:SER:HB3	1:B:1404:LEU:HD21	1.99	0.45
1:A:368:GLU:HG3	1:B:14:ILE:HD11	1.98	0.45
1:B:691:VAL:HG12	1:B:693:PRO:HD3	1.98	0.45
1:A:79:LEU:HD12	1:A:83:LEU:HD23	1.99	0.45
1:B:395:ILE:CG2	1:B:408:THR:HG21	2.46	0.44
1:B:1340:LEU:HA	1:B:1343:LEU:HD13	2.00	0.44
1:B:1313:LEU:HD12	1:B:1317:LEU:CD1	2.48	0.44
1:B:1368:LEU:CD1	1:B:1411:ILE:HD11	2.48	0.44
1:A:167:LEU:HD22	1:B:209:VAL:CG2	2.48	0.43
1:A:822:PHE:CE1	1:A:841:LEU:HD11	2.53	0.43
1:B:415:LEU:HD22	1:B:423:LEU:CD2	2.48	0.43
1:B:506:LEU:HD11	1:B:820:THR:HG22	2.00	0.43
1:B:1116:LYS:NZ	1:B:1137:GLU:OE1	2.30	0.43
1:A:438:LEU:O	1:A:459:LYS:NZ	2.44	0.43
1:B:808:ALA:O	1:B:812:GLN:NE2	2.51	0.43
1:B:973:LEU:HD11	1:B:1111:ARG:CD	2.48	0.43
1:B:395:ILE:O	1:B:399:THR:HG22	2.19	0.43
1:A:1396:ASP:O	1:A:1400:LYS:N	2.48	0.43
1:A:110:ILE:HD12	1:A:111:SER:H	1.84	0.43
1:A:1276:LEU:CD2	1:A:1290:ILE:HG21	2.49	0.43
1:A:932:ASN:OD1	1:A:933:LYS:N	2.52	0.42
1:B:1056:ASN:OD1	1:B:1057:ALA:N	2.51	0.42
1:A:502:THR:HG22	1:A:820:THR:HG21	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:658:ILE:HG22	1:A:662:ILE:HD12	2.01	0.42
1:B:675:TYR:OH	1:B:693:PRO:O	2.37	0.42
1:A:22:LEU:HD23	1:A:22:LEU:O	2.20	0.42
1:A:79:LEU:HD12	1:A:79:LEU:O	2.19	0.42
1:B:1313:LEU:HG	1:B:1362:ILE:HD13	2.01	0.42
1:A:113:TYR:O	1:A:116:SER:OG	2.36	0.42
1:B:1314:GLU:O	1:B:1318:ILE:HG22	2.20	0.42
1:B:628:ILE:HD11	1:B:664:ALA:HB1	2.01	0.42
1:B:89:LEU:HD12	1:B:90:ASP:H	1.84	0.42
1:A:374:LEU:HD23	1:A:374:LEU:O	2.19	0.42
1:A:594:ASP:OD1	1:A:595:ILE:N	2.53	0.42
1:A:213:ALA:O	1:A:217:THR:OG1	2.34	0.42
1:A:1328:MET:SD	1:A:1329:GLU:N	2.93	0.42
1:A:343:ILE:HD13	1:A:360:LEU:HD22	2.02	0.41
1:A:462:LEU:O	1:A:466:ILE:HG22	2.20	0.41
1:A:1075:ILE:HG21	1:A:1118:LEU:HD23	2.01	0.41
1:A:13:THR:O	1:B:472:ARG:NH2	2.54	0.41
1:B:485:ASP:OD1	1:B:557:ARG:NH1	2.52	0.41
1:B:352:SER:OG	1:B:353:TYR:N	2.54	0.41
1:A:1406:LEU:HD23	1:A:1406:LEU:O	2.21	0.41
1:B:335:TYR:OH	1:B:362:LEU:HD21	2.21	0.41
1:A:198:LEU:HD12	1:A:338:LEU:HD21	2.02	0.41
1:A:340:LEU:HD11	1:A:387:PRO:HG2	2.02	0.41
1:B:103:LEU:HD23	1:B:103:LEU:O	2.21	0.41
1:A:971:LEU:HD12	1:A:1114:THR:CG2	2.50	0.41
1:A:1329:GLU:N	1:A:1329:GLU:OE1	2.54	0.40
1:A:1053:THR:HG23	1:A:1098:LEU:HD23	2.03	0.40
1:A:642:ILE:HD13	1:A:705:ILE:HG21	2.04	0.40
1:B:146:VAL:HG21	1:B:185:LEU:HD22	2.02	0.40
1:A:15:ILE:HG21	1:A:97:ILE:HD11	2.03	0.40
1:B:1068:LEU:HD11	1:B:1108:THR:HG22	2.02	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	1091/1459 (75%)	1025 (94%)	64 (6%)	2 (0%)	47 79
1	B	1091/1459 (75%)	1037 (95%)	52 (5%)	2 (0%)	47 79
All	All	2182/2918 (75%)	2062 (94%)	116 (5%)	4 (0%)	50 79

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	830	ASP
1	A	983	ILE
1	A	1201	TYR
1	B	348	GLU

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	1042/1346 (77%)	1000 (96%)	42 (4%)	31 58
1	B	1042/1346 (77%)	1004 (96%)	38 (4%)	35 61
All	All	2084/2692 (77%)	2004 (96%)	80 (4%)	36 59

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	19	CYS
1	A	82	MET
1	A	172	LEU
1	A	175	SER
1	A	218	MET
1	A	368	GLU
1	A	386	ASP
1	A	390	LYS

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Mol	Chain	Res	Type
1	A	403	SER
1	A	422	ASN
1	A	484	PHE
1	A	487	ASN
1	A	497	PHE
1	A	509	SER
1	A	531	ASP
1	A	537	LYS
1	A	586	PHE
1	A	598	PHE
1	A	603	ASN
1	A	665	PHE
1	A	669	TYR
1	A	687	ASP
1	A	807	ARG
1	A	810	PHE
1	A	851	PHE
1	A	859	ASP
1	A	912	PHE
1	A	955	GLU
1	A	962	ASP
1	A	1020	MET
1	A	1022	CYS
1	A	1027	ASN
1	A	1038	ASN
1	A	1113	LEU
1	A	1163	PHE
1	A	1265	SER
1	A	1273	SER
1	A	1313	LEU
1	A	1314	GLU
1	A	1328	MET
1	A	1401	LEU
1	B	89	LEU
1	B	153	ARG
1	B	154	PHE
1	B	172	LEU
1	B	174	ARG
1	B	179	SER
1	B	375	GLN
1	B	380	LEU
1	B	389	PHE

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Mol	Chain	Res	Type
1	B	419	LEU
1	B	497	PHE
1	B	533	PHE
1	B	540	ASP
1	B	586	PHE
1	B	591	SER
1	B	600	PHE
1	B	627	TYR
1	B	650	ARG
1	B	703	TYR
1	B	737	LYS
1	B	769	PHE
1	B	774	ASN
1	B	810	PHE
1	B	855	ASP
1	B	912	PHE
1	B	921	ASN
1	B	930	ARG
1	B	967	LEU
1	B	1020	MET
1	B	1031	SER
1	B	1081	CYS
1	B	1122	LEU
1	B	1125	ASP
1	B	1186	ARG
1	B	1226	ASP
1	B	1298	CYS
1	B	1302	CYS
1	B	1370	TYR

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

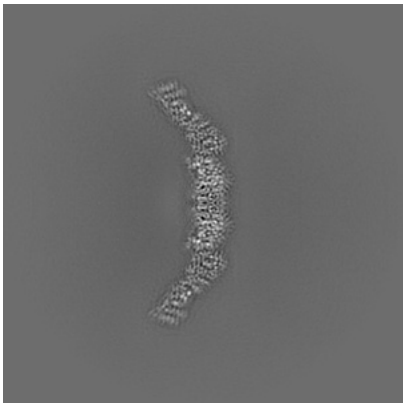
This section contains visualisations of the EMDB entry EMD-26770. These allow visual inspection of the internal detail of the map and identification of artifacts.

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.

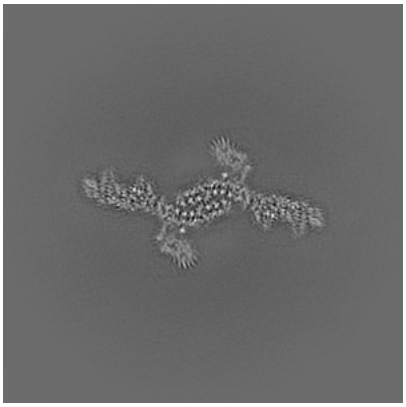
6.1 Orthogonal projections

6.1.1 Primary map

X



Y

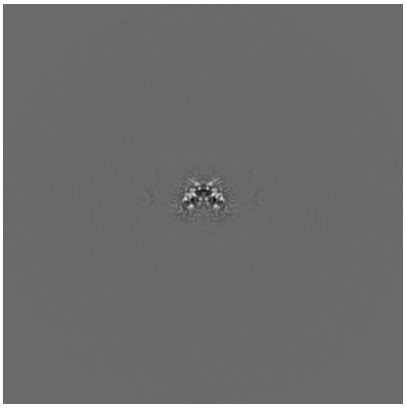


Z

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

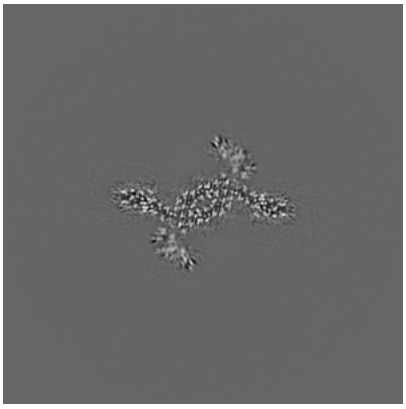
6.2 Central slices

6.2.1 Primary map



X Index: 150

Y Index: 150

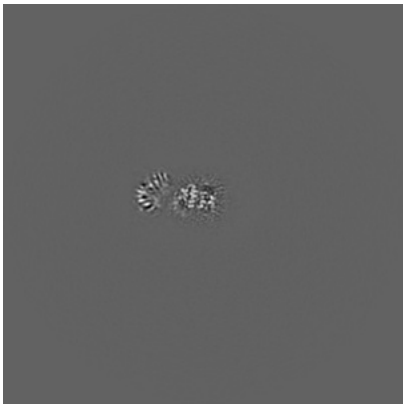


Z Index: 150

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

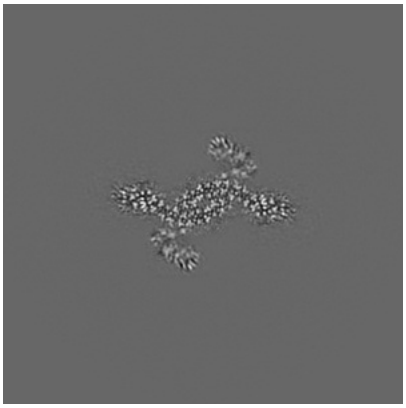
6.3 Largest variance slices (i)

6.3.1 Primary map



X Index: 136

Y Index: 147



Z Index: 151

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

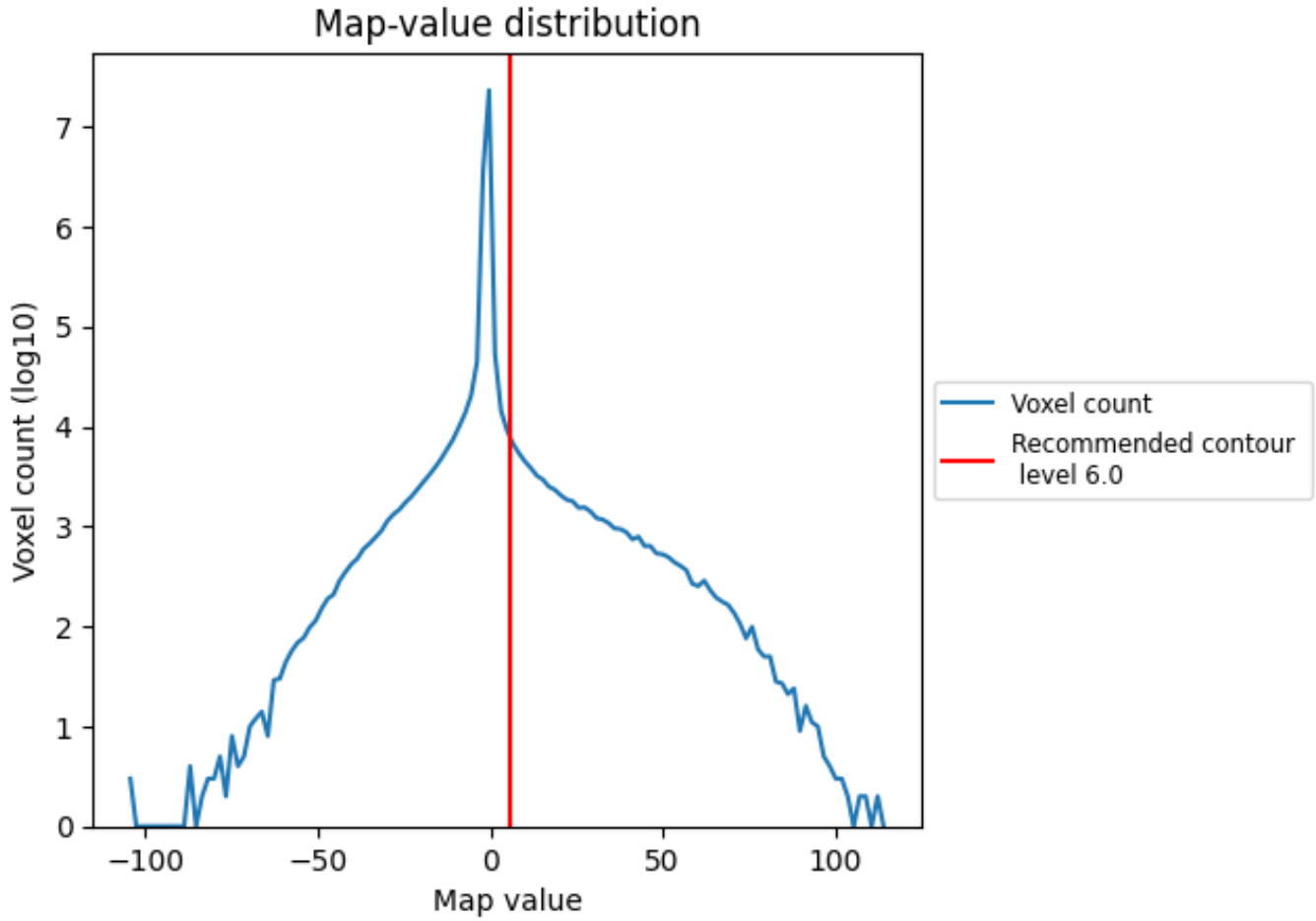
6.5 Mask visualisation (i)

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

7 Map analysis

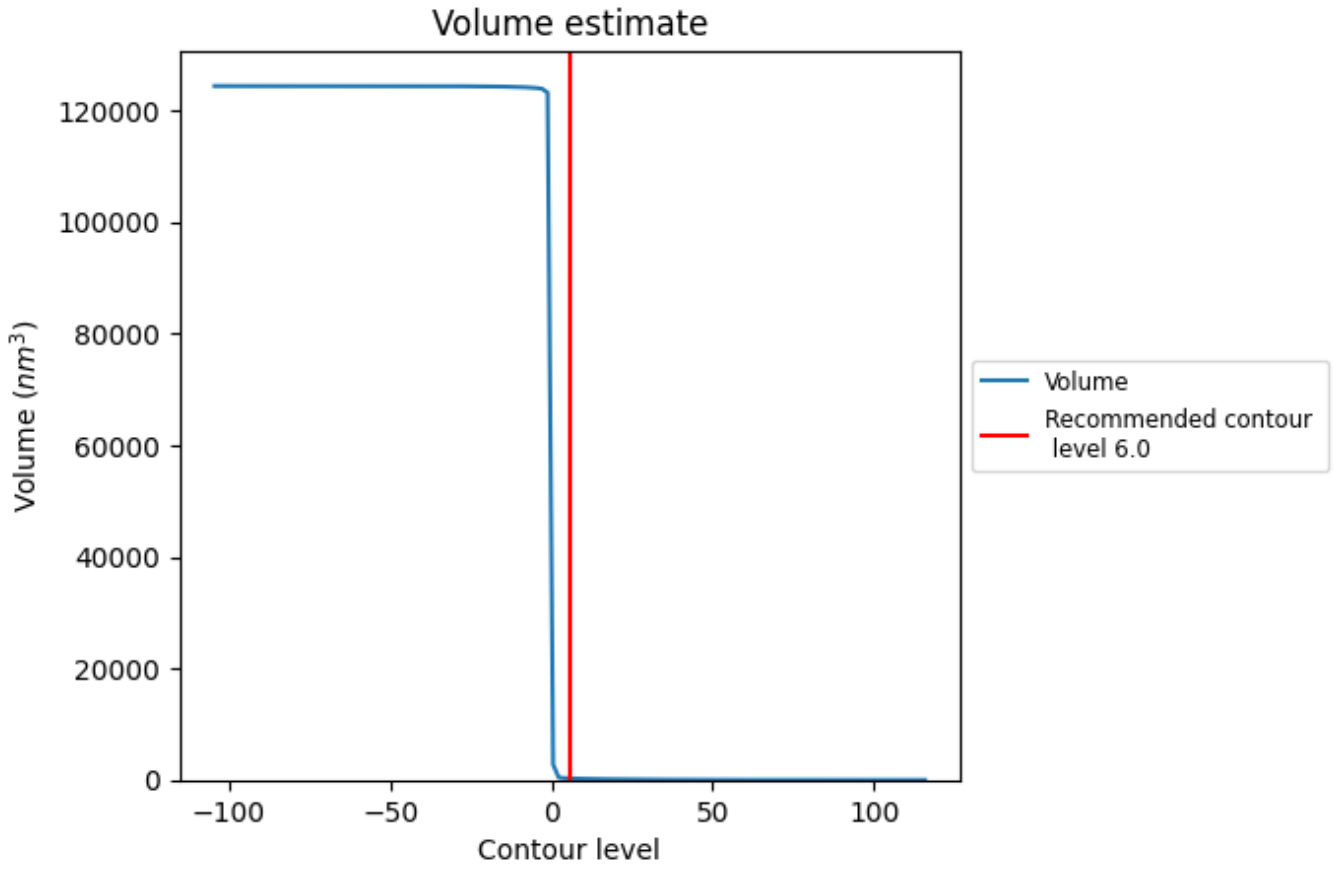
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

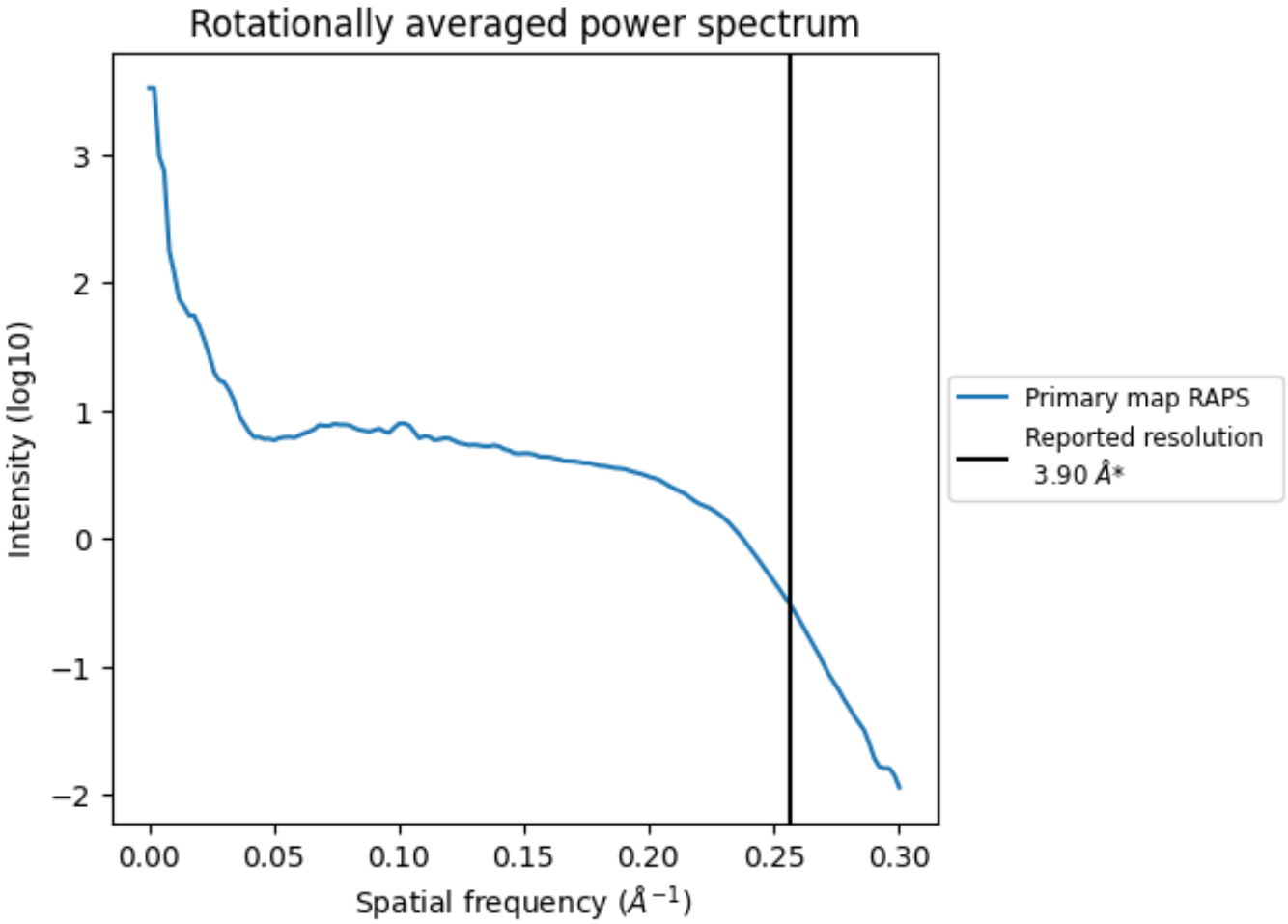
7.2 Volume estimate



The volume at the recommended contour level is 270 nm^3; this corresponds to an approximate mass of 244 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 $Å^{-1}$

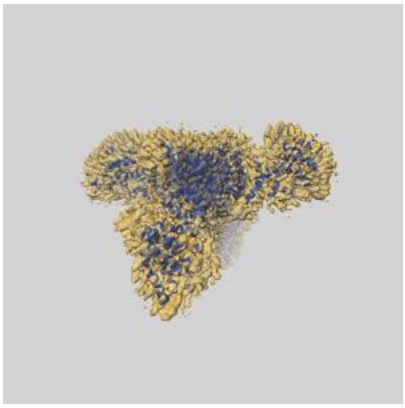
8 Fourier-Shell correlation (i)

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

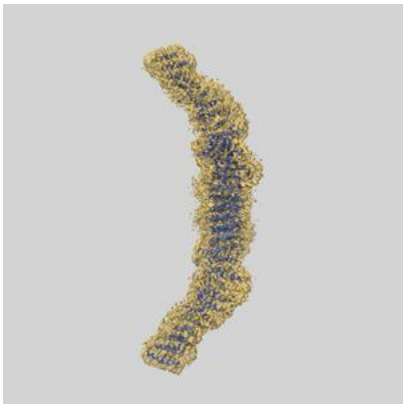
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-26770 and PDB model 7UTH. Per-residue inclusion information can be found in section 3 on page 4.

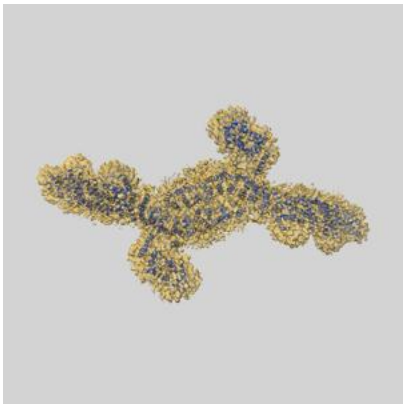
9.1 Map-model overlay



X



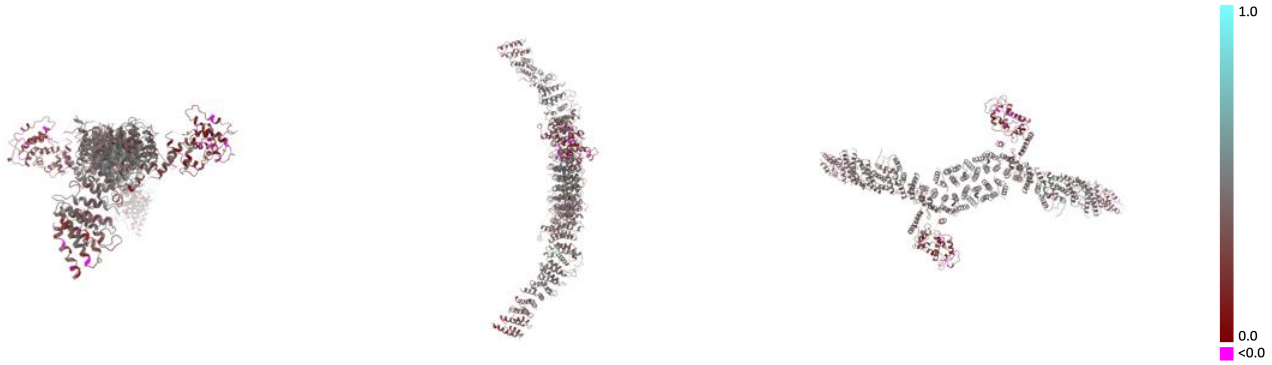
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 6.0 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.

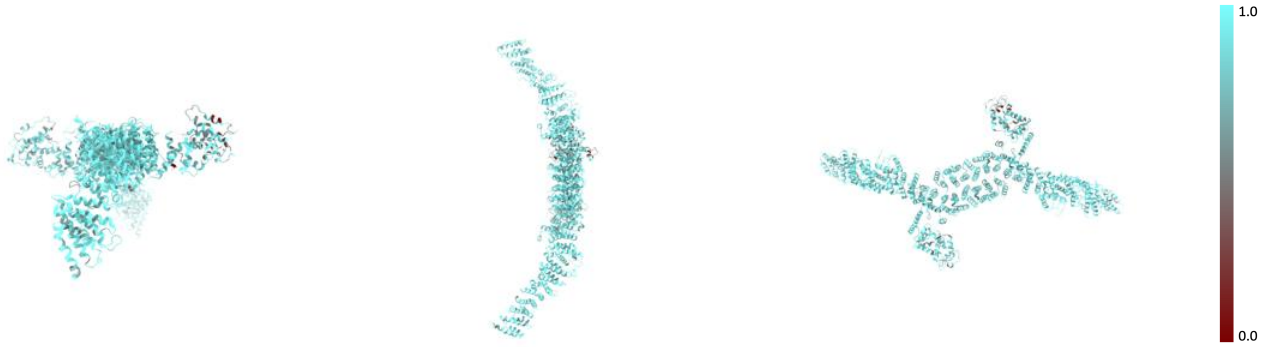
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

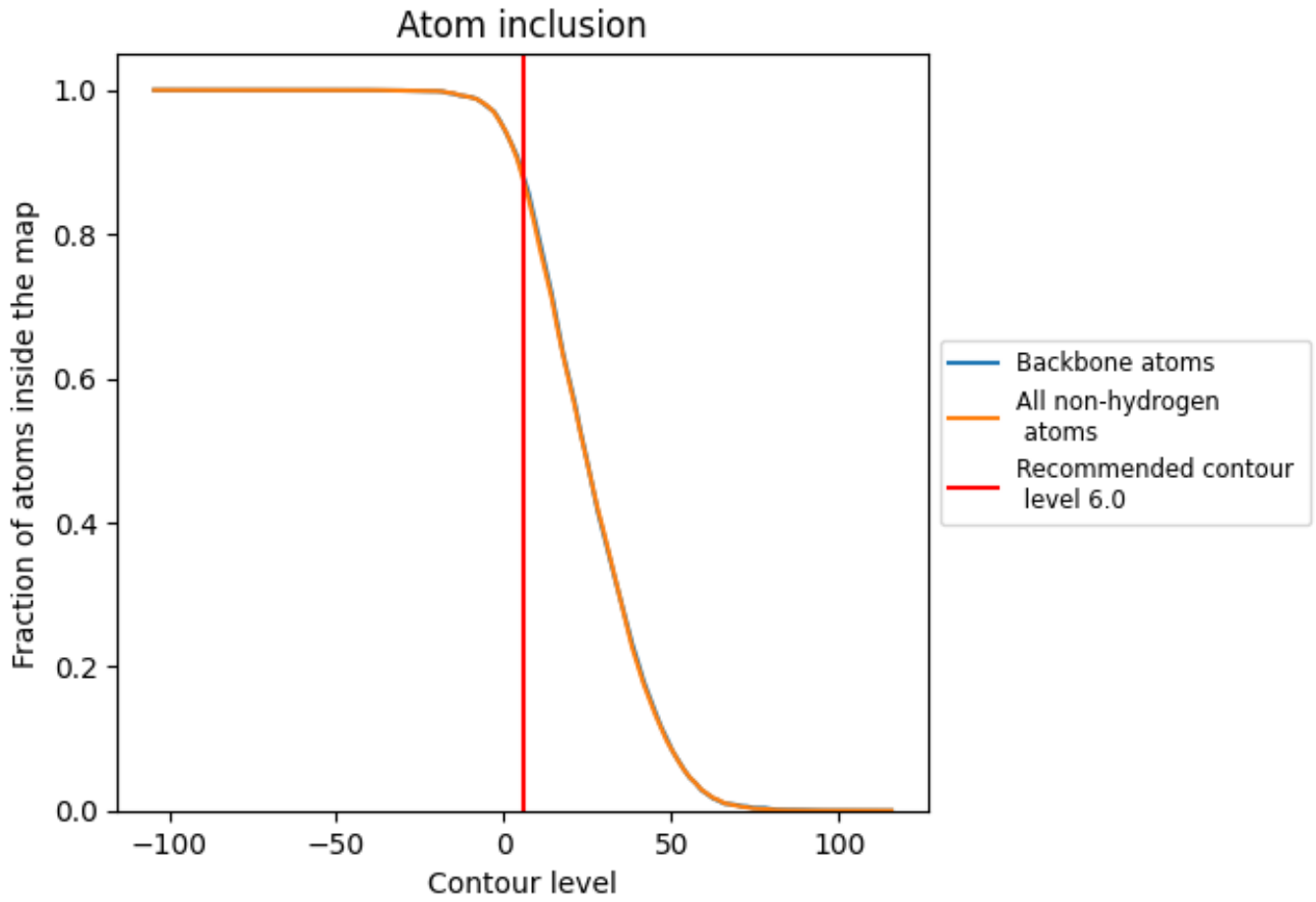
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (6.0).

9.4 Atom inclusion (i)



At the recommended contour level, 88% of all backbone atoms, 87% 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 (6.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8729	0.3800
A	0.8660	0.3790
B	0.8796	0.3810