



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 03:56 AM UTC

PDB ID : 6FN4 / pdb_00006fn4
EMDB ID : EMD-4282
Title : Apo form of UIC2 Fab complex of human-mouse chimeric ABCB1 (ABCB1HM)
Authors : Alam, A.; Locher, K.P.
Deposited on : 2018-02-02
Resolution : 4.14 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.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

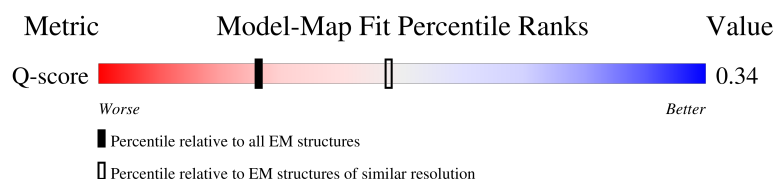
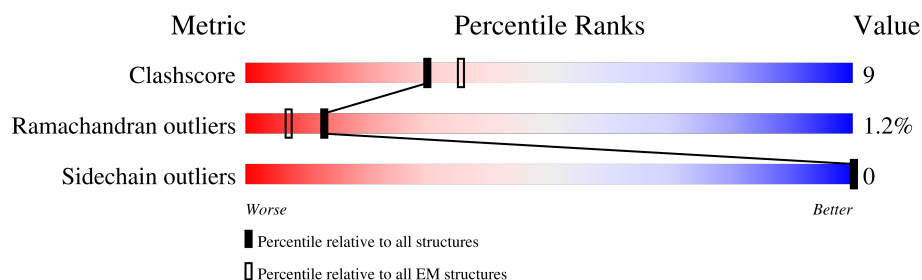
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.14 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	5577 (3.64 - 4.64)

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	1279	
2	B	220	
3	C	225	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	A	1301	-	-	X	-
4	NAG	A	1302	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12702 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 Apo form of Human-mouse chimeric ABCB1 (ABCB1HM) in complex with Antigen binding fragment of UIC2..

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1181	9182	5913	1552	1684	33	0	0

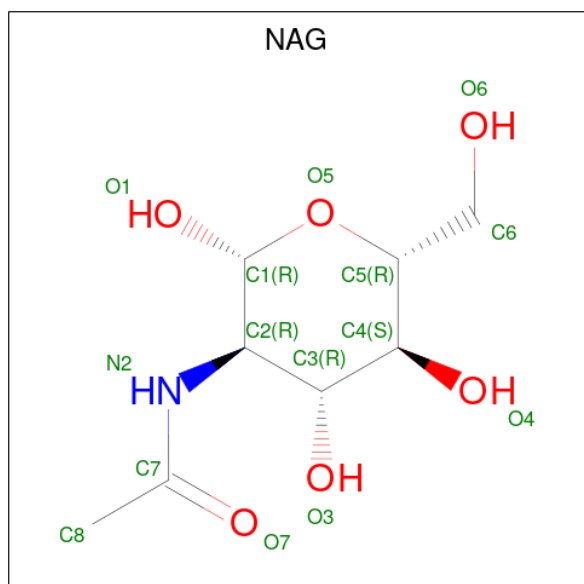
- Molecule 2 is a protein called UIC2 Antigen Binding Fragment Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	220	1710	1069	292	342	7	0	0

- Molecule 3 is a protein called UIC2 Antigen binding fragment Heavy chain.

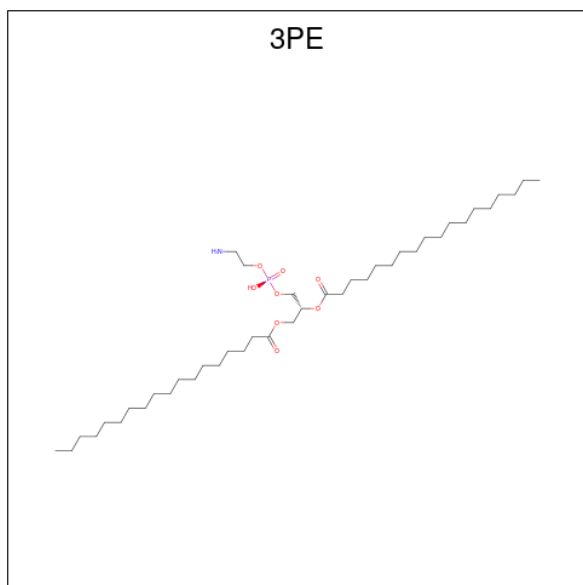
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	225	1717	1098	274	337	8	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

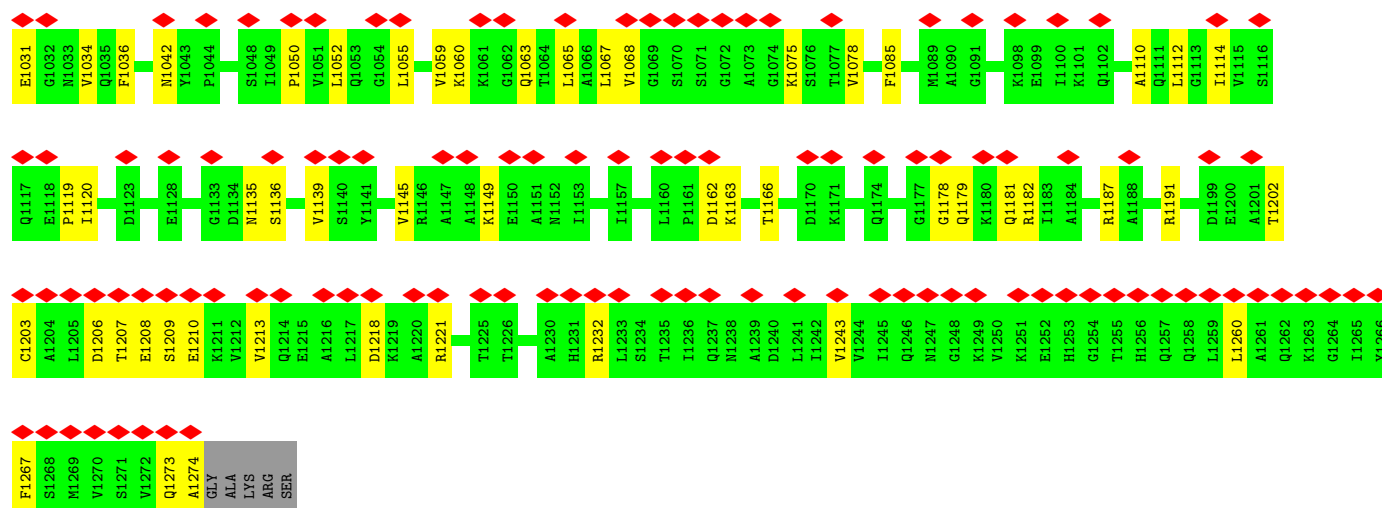


Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	

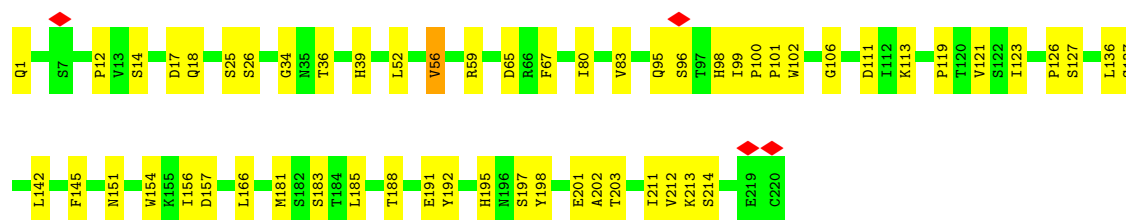
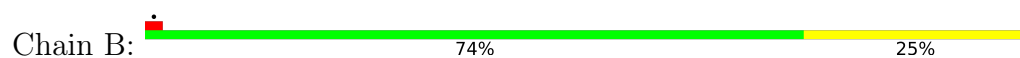
- Molecule 5 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (CCD ID: 3PE) (formula: $C_{41}H_{82}NO_8P$).



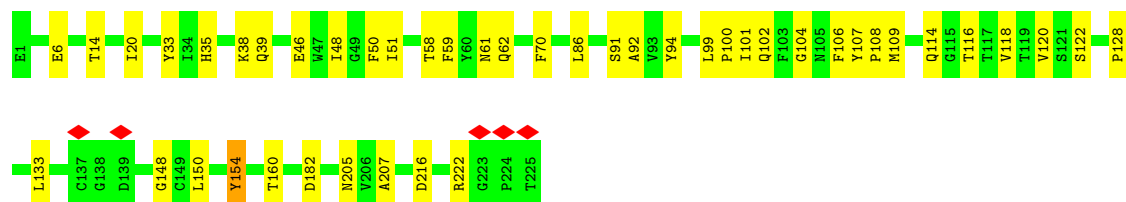
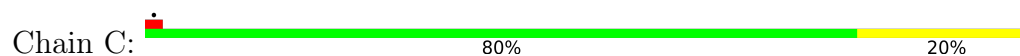
Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total	C	N	O	P	0
			51	41	1	8	1	



• Molecule 2: UIC2 Antigen Binding Fragment Light chain



• Molecule 3: UIC2 Antigen binding fragment Heavy chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	517053	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	0.9	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.690	Depositor
Minimum map value	-0.464	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.015	Depositor
Recommended contour level	0.07	Depositor
Map size (Å)	277.4, 277.4, 277.4	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.387, 1.387, 1.387	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3PE, NAG

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	0.18	0/9352	0.45	4/12645 (0.0%)
2	B	0.16	0/1751	0.43	0/2376
3	C	0.16	0/1767	0.39	0/2412
All	All	0.17	0/12870	0.44	4/17433 (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	A	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ASP	CA-C-N	6.38	133.19	121.70
1	A	375	ASP	C-N-CA	6.38	133.19	121.70
1	A	237	SER	CA-C-N	5.46	131.52	121.70
1	A	237	SER	C-N-CA	5.46	131.52	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	488	ARG	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9182	0	9355	171	0
2	B	1710	0	1652	38	0
3	C	1717	0	1656	31	0
4	A	42	0	39	18	0
5	A	51	0	82	2	0
All	All	12702	0	12784	232	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 9.

All (232) 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:94:ARG:HD2	4:A:1301:NAG:C8	1.74	1.16
1:A:90:ASN:ND2	4:A:1302:NAG:C1	2.10	1.15
1:A:94:ARG:CD	4:A:1301:NAG:H82	1.87	1.04
1:A:94:ARG:HD2	4:A:1301:NAG:H82	0.96	0.94
1:A:587:ARG:HA	1:A:628:GLN:HE22	1.36	0.90
1:A:93:ASN:HD22	4:A:1301:NAG:C1	1.85	0.87
1:A:375:ASP:HA	1:A:376:SER:HB3	1.61	0.83
1:A:93:ASN:ND2	4:A:1301:NAG:C1	2.44	0.80
1:A:33:VAL:H	1:A:34:SER:HA	1.48	0.78
1:A:742:ASP:OD1	1:A:747:LYS:NZ	2.17	0.78
3:C:133:LEU:HD11	3:C:150:LEU:HG	1.68	0.76
1:A:546:ARG:NH1	1:A:909:LEU:O	2.19	0.75
1:A:159:ILE:HG21	1:A:442:LEU:HB3	1.69	0.75
1:A:491:ARG:HA	1:A:911:ARG:NH2	2.02	0.75
1:A:271:LYS:HE3	1:A:1136:SER:H	1.52	0.74
1:A:790:MET:HG2	1:A:1011:ILE:HD11	1.71	0.72
1:A:376:SER:H	1:A:462:VAL:HB	1.52	0.72
1:A:1218:ASP:OD1	1:A:1221:ARG:NH2	2.23	0.71
1:A:1060:LYS:H	1:A:1063:GLN:HE21	1.36	0.71
1:A:259:ALA:HB2	1:A:1120:ILE:HG21	1.73	0.71
1:A:432:LYS:HB3	1:A:584:ILE:HD12	1.72	0.70
1:A:797:ARG:HG2	1:A:1015:PRO:HG2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:LYS:HD3	1:A:785:LYS:HG3	1.75	0.68
2:B:119:PRO:HB3	2:B:145:PHE:HB3	1.76	0.67
1:A:1179:GLN:HG2	1:A:1182:ARG:NH2	2.12	0.65
1:A:425:VAL:HG11	1:A:588:LEU:HD11	1.79	0.64
1:A:555:GLU:N	1:A:584:ILE:O	2.29	0.64
1:A:610:ASN:OD1	1:A:611:HIS:N	2.31	0.64
1:A:557:THR:HG23	1:A:560:LEU:HD12	1.80	0.63
1:A:1008:ILE:HA	1:A:1011:ILE:HG22	1.81	0.63
1:A:481:THR:HG22	1:A:482:THR:H	1.63	0.62
1:A:736:GLY:HA3	1:A:971:GLU:HG2	1.81	0.62
2:B:126:PRO:HG2	2:B:136:LEU:HD12	1.82	0.62
1:A:97:ILE:HG12	4:A:1303:NAG:O7	1.99	0.62
3:C:205:ASN:ND2	3:C:216:ASP:OD1	2.30	0.61
1:A:539:ALA:HA	1:A:542:ARG:HE	1.66	0.61
1:A:702:LEU:HD11	1:A:1008:ILE:HD11	1.83	0.61
1:A:1210:GLU:OE2	1:A:1232:ARG:NH1	2.34	0.61
1:A:838:ASN:OD1	1:A:839:ILE:N	2.34	0.60
1:A:97:ILE:HG12	4:A:1303:NAG:H2	1.81	0.60
1:A:629:THR:HB	1:A:1274:ALA:HA	1.83	0.60
1:A:1202:THR:OG1	1:A:1203:CYS:SG	2.60	0.60
1:A:716:VAL:HG11	1:A:771:LEU:HD11	1.83	0.59
1:A:436:VAL:HG13	1:A:552:LEU:HD23	1.85	0.59
3:C:128:PRO:HB3	3:C:154:TYR:HB3	1.84	0.59
2:B:14:SER:OG	2:B:17:ASP:OD2	2.21	0.59
1:A:439:MET:HB3	1:A:552:LEU:HD21	1.84	0.58
1:A:424:LEU:HD12	1:A:584:ILE:HD11	1.85	0.58
1:A:1112:LEU:O	1:A:1191:ARG:NH1	2.37	0.58
2:B:18:GLN:HA	2:B:80:ILE:O	2.03	0.58
2:B:201:GLU:HG2	2:B:212:VAL:HG22	1.87	0.57
1:A:237:SER:HB2	1:A:239:THR:HG23	1.86	0.57
1:A:83:ASN:ND2	3:C:101:ILE:O	2.38	0.57
1:A:588:LEU:HD23	1:A:621:TYR:CE1	2.40	0.57
3:C:38:LYS:HB2	3:C:48:ILE:HD11	1.86	0.57
1:A:507:ASN:O	1:A:537:ARG:NH2	2.38	0.56
1:A:1110:ALA:HA	1:A:1191:ARG:NH2	2.20	0.56
1:A:90:ASN:HD21	4:A:1302:NAG:C1	2.15	0.56
1:A:33:VAL:HG22	1:A:358:ARG:HG3	1.88	0.56
1:A:395:ASN:N	1:A:413:ASN:OD1	2.33	0.56
1:A:827:ALA:HB2	1:A:1000:ALA:HB1	1.87	0.56
1:A:484:ALA:O	1:A:488:ARG:HD2	2.07	0.55
2:B:39:HIS:ND1	3:C:108:PRO:HB3	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:ASN:HD22	4:A:1302:NAG:C1	2.13	0.54
1:A:424:LEU:HB2	1:A:584:ILE:HD13	1.89	0.54
3:C:99:LEU:HD12	3:C:100:PRO:HD2	1.89	0.54
1:A:702:LEU:O	1:A:786:ARG:NH2	2.41	0.54
2:B:102:TRP:HE1	3:C:109:MET:HE3	1.73	0.54
1:A:816:ARG:O	1:A:820:ASP:HB2	2.07	0.54
2:B:1:GLN:HA	2:B:101:PRO:HD3	1.88	0.54
2:B:59:ARG:NH2	2:B:65:ASP:OD1	2.36	0.54
1:A:991:SER:HB3	5:A:1304:3PE:H31	1.89	0.53
1:A:33:VAL:HG23	1:A:34:SER:C	2.33	0.53
2:B:151:ASN:HB2	2:B:203:THR:HB	1.90	0.53
1:A:460:ILE:HD11	1:A:465:LEU:HD13	1.91	0.53
2:B:188:THR:OG1	2:B:191:GLU:OE1	2.27	0.53
1:A:90:ASN:CG	4:A:1302:NAG:C1	2.81	0.53
1:A:425:VAL:O	1:A:599:GLY:HA2	2.09	0.52
1:A:396:ILE:HB	1:A:412:LEU:HB3	1.91	0.52
1:A:1034:VAL:HB	1:A:1059:VAL:HB	1.91	0.52
1:A:38:MET:HE1	1:A:358:ARG:HA	1.92	0.52
3:C:160:THR:OG1	3:C:207:ALA:HB3	2.10	0.52
1:A:317:THR:HG23	1:A:738:PHE:HZ	1.73	0.52
1:A:738:PHE:O	3:C:102:GLN:NE2	2.43	0.52
2:B:17:ASP:H	2:B:83:VAL:HG12	1.75	0.52
1:A:94:ARG:NH2	4:A:1303:NAG:O6	2.40	0.52
1:A:376:SER:N	1:A:462:VAL:HB	2.23	0.51
1:A:73:GLU:OE2	1:A:209:ARG:NH2	2.41	0.51
1:A:143:ILE:HA	1:A:146:ILE:HG22	1.92	0.51
1:A:467:GLU:HA	1:A:546:ARG:HH21	1.74	0.51
2:B:99:ILE:HG21	3:C:59:PHE:HE2	1.75	0.51
1:A:1052:LEU:HD22	1:A:1055:LEU:HD22	1.92	0.51
1:A:440:GLN:HB2	1:A:442:LEU:HG	1.93	0.51
1:A:93:ASN:ND2	4:A:1301:NAG:O5	2.37	0.51
1:A:266:PHE:CE1	1:A:1114:ILE:HG22	2.46	0.50
1:A:491:ARG:HA	1:A:911:ARG:HH21	1.75	0.50
1:A:628:GLN:HB3	1:A:1273:GLN:HB2	1.93	0.50
1:A:425:VAL:HG11	1:A:588:LEU:HD21	1.92	0.50
2:B:136:LEU:HD11	2:B:192:TYR:CG	2.46	0.50
2:B:123:ILE:HD11	2:B:213:LYS:HB3	1.93	0.50
3:C:14:THR:OG1	3:C:122:SER:OG	2.21	0.50
1:A:257:LEU:HA	1:A:260:ILE:HG23	1.94	0.50
2:B:157:ASP:HA	2:B:197:SER:HB3	1.94	0.49
1:A:987:VAL:O	1:A:991:SER:CB	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:ASN:ND2	4:A:1302:NAG:O5	2.45	0.49
3:C:86:LEU:HD21	3:C:120:VAL:HG22	1.95	0.49
1:A:295:ASN:ND2	1:A:770:PHE:O	2.45	0.49
1:A:150:PHE:CD1	1:A:364:VAL:HG13	2.48	0.49
1:A:1016:GLU:HB3	1:A:1017:ILE:HD12	1.94	0.48
1:A:231:TRP:CD1	1:A:298:MET:HG3	2.48	0.48
1:A:1065:LEU:HD11	1:A:1243:VAL:HG23	1.96	0.48
3:C:38:LYS:HG3	3:C:94:TYR:CE1	2.48	0.48
1:A:492:GLU:HG3	1:A:911:ARG:NH1	2.28	0.48
1:A:555:GLU:HG2	1:A:585:ALA:HA	1.95	0.48
1:A:1036:PHE:CE2	1:A:1078:VAL:HG13	2.48	0.48
1:A:1178:GLY:O	1:A:1182:ARG:HG3	2.14	0.48
2:B:12:PRO:HA	2:B:111:ASP:HB3	1.95	0.48
1:A:147:ARG:HD2	1:A:923:LEU:HD11	1.95	0.48
1:A:370:ASN:OD1	1:A:371:LYS:N	2.46	0.47
1:A:205:ILE:HA	1:A:208:THR:HG22	1.97	0.47
1:A:1006:HIS:O	1:A:1010:ILE:HG13	2.13	0.47
2:B:121:VAL:HG11	2:B:213:LYS:HG3	1.96	0.47
1:A:422:VAL:HG23	1:A:596:VAL:HG13	1.95	0.47
1:A:425:VAL:HG12	1:A:426:GLY:N	2.29	0.47
1:A:391:LEU:HB2	1:A:416:VAL:HB	1.96	0.47
2:B:157:ASP:OD2	2:B:195:HIS:ND1	2.47	0.47
1:A:523:VAL:HG22	1:A:524:GLY:H	1.79	0.47
1:A:809:THR:H	1:A:812:ALA:HB3	1.78	0.47
1:A:213:LEU:HD22	1:A:333:VAL:HG13	1.96	0.47
1:A:1145:VAL:HG22	1:A:1149:LYS:HE3	1.97	0.47
1:A:999:LYS:HA	1:A:1002:VAL:HG22	1.95	0.47
1:A:1206:ASP:C	1:A:1208:GLU:H	2.22	0.47
1:A:131:GLN:HG3	1:A:132:VAL:N	2.29	0.46
1:A:479:PHE:O	1:A:481:THR:N	2.43	0.46
1:A:1119:PRO:HG3	1:A:1181:GLN:HA	1.97	0.46
1:A:849:SER:OG	1:A:857:THR:OG1	2.26	0.46
2:B:39:HIS:CG	3:C:108:PRO:HB3	2.51	0.46
1:A:272:GLU:OE1	1:A:275:ARG:NH2	2.48	0.46
1:A:795:MET:HE1	1:A:813:LEU:HB3	1.96	0.46
1:A:171:ASN:OD1	1:A:897:THR:HG22	2.16	0.46
2:B:156:ILE:HG23	2:B:198:TYR:HE1	1.80	0.46
1:A:427:ASN:O	1:A:432:LYS:NZ	2.48	0.46
2:B:99:ILE:HG21	3:C:59:PHE:CE2	2.50	0.46
1:A:1162:ASP:O	1:A:1166:THR:OG1	2.32	0.46
1:A:432:LYS:HB3	1:A:584:ILE:HG23	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:PHE:CE2	1:A:435:THR:HG23	2.51	0.46
1:A:600:PHE:CE1	1:A:605:ILE:HG12	2.51	0.46
1:A:92:THR:HG21	4:A:1302:NAG:H81	1.97	0.46
1:A:469:ILE:HG12	1:A:550:ILE:HB	1.97	0.46
1:A:1042:ASN:OD1	1:A:1050:PRO:HA	2.16	0.46
3:C:104:GLY:C	3:C:106:PHE:H	2.24	0.46
1:A:231:TRP:O	1:A:235:LEU:HG	2.16	0.45
1:A:724:GLN:HB3	1:A:725:PRO:HD3	1.99	0.45
2:B:202:ALA:HB3	2:B:211:ILE:HB	1.99	0.45
2:B:127:SER:HA	3:C:222:ARG:HH22	1.82	0.45
3:C:6:GLU:H	3:C:114:GLN:NE2	2.14	0.45
1:A:136:LEU:HD23	1:A:934:ALA:HA	1.98	0.45
1:A:437:GLN:O	1:A:442:LEU:HB2	2.16	0.45
1:A:251:ALA:O	1:A:254:GLU:HG3	2.17	0.45
1:A:1260:LEU:HD12	1:A:1267:PHE:HD1	1.81	0.44
2:B:52:LEU:HD11	2:B:67:PHE:CE2	2.52	0.44
1:A:110:MET:HG2	1:A:960:ALA:HB2	1.99	0.44
2:B:99:ILE:HA	2:B:100:PRO:HA	1.62	0.44
1:A:391:LEU:HD21	1:A:550:ILE:HD13	1.99	0.44
2:B:59:ARG:NH1	2:B:65:ASP:HA	2.33	0.44
3:C:20:ILE:HD11	3:C:116:THR:HB	1.99	0.44
1:A:240:ASP:HA	1:A:243:LEU:HD13	2.00	0.44
1:A:597:ILE:HG21	1:A:621:TYR:CE1	2.53	0.44
1:A:375:ASP:HA	1:A:376:SER:CB	2.38	0.44
1:A:755:SER:O	1:A:758:PHE:HB2	2.17	0.44
1:A:94:ARG:HD2	4:A:1301:NAG:C7	2.44	0.44
1:A:479:PHE:HB2	1:A:486:ASN:OD1	2.17	0.44
3:C:51:ILE:HG23	3:C:70:PHE:HD2	1.83	0.43
3:C:107:TYR:CG	3:C:108:PRO:HD2	2.53	0.43
1:A:97:ILE:HG21	4:A:1303:NAG:H4	2.00	0.43
1:A:492:GLU:OE2	1:A:914:LYS:HB2	2.19	0.43
1:A:866:ILE:HD13	1:A:947:MET:HG3	1.99	0.43
1:A:927:TYR:O	1:A:931:MET:HG2	2.19	0.43
2:B:154:TRP:CE2	2:B:185:LEU:HD13	2.53	0.43
2:B:166:LEU:O	2:B:183:SER:HA	2.18	0.43
3:C:6:GLU:H	3:C:114:GLN:HE22	1.66	0.43
1:A:721:GLY:HA2	1:A:837:GLN:HG3	2.00	0.43
1:A:866:ILE:HD12	1:A:867:ILE:N	2.33	0.43
1:A:178:VAL:HA	1:A:181:ILE:HG22	2.00	0.43
1:A:232:ALA:HA	1:A:235:LEU:HD12	2.01	0.43
1:A:349:PRO:O	1:A:352:GLU:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:126:PRO:HG3	2:B:137:SER:O	2.18	0.43
1:A:558:CYS:HA	1:A:587:ARG:NH2	2.34	0.43
1:A:1067:LEU:HD23	1:A:1243:VAL:HB	2.00	0.43
2:B:36:THR:HG23	2:B:56:VAL:HG21	2.01	0.43
1:A:1145:VAL:O	1:A:1149:LYS:HG3	2.18	0.42
1:A:487:ILE:O	1:A:490:GLY:N	2.38	0.42
1:A:984:ALA:HA	1:A:987:VAL:HG22	2.01	0.42
5:A:1304:3PE:H252	5:A:1304:3PE:H282	1.70	0.42
2:B:198:TYR:O	2:B:214:SER:HA	2.19	0.42
3:C:39:GLN:O	3:C:92:ALA:HB1	2.20	0.42
1:A:96:ASP:OD1	1:A:97:ILE:N	2.52	0.42
1:A:237:SER:HA	1:A:238:PHE:HB3	2.01	0.42
1:A:925:ILE:HB	1:A:926:PRO:HD3	2.00	0.42
1:A:1209:SER:O	1:A:1213:VAL:HG23	2.19	0.42
2:B:52:LEU:HD12	2:B:52:LEU:HA	1.84	0.42
1:A:93:ASN:CG	1:A:94:ARG:H	2.27	0.41
2:B:25:SER:OG	2:B:26:SER:N	2.53	0.41
1:A:88:MET:HE1	3:C:59:PHE:CG	2.54	0.41
3:C:33:TYR:CE2	3:C:101:ILE:HG13	2.55	0.41
3:C:133:LEU:HB2	3:C:148:GLY:C	2.45	0.41
1:A:800:VAL:HG21	1:A:1085:PHE:CG	2.55	0.41
3:C:61:ASN:OD1	3:C:62:GLN:N	2.50	0.41
3:C:91:SER:HA	3:C:118:VAL:O	2.20	0.41
1:A:1068:VAL:HA	1:A:1075:LYS:HZ3	1.85	0.41
1:A:92:THR:OG1	4:A:1302:NAG:H82	2.20	0.41
1:A:588:LEU:HD12	1:A:588:LEU:N	2.35	0.41
1:A:987:VAL:O	1:A:991:SER:HB2	2.20	0.41
2:B:12:PRO:HB2	2:B:113:LYS:HG2	2.02	0.41
2:B:95:GLN:CD	2:B:96:SER:H	2.28	0.41
1:A:266:PHE:HE2	1:A:1187:ARG:NE	2.19	0.41
1:A:892:SER:O	1:A:895:ILE:HG22	2.20	0.41
1:A:271:LYS:HD2	1:A:1135:ASN:OD1	2.21	0.41
1:A:315:TYR:HA	1:A:318:THR:HG22	2.03	0.41
1:A:788:ARG:HA	1:A:791:VAL:HG12	2.01	0.41
3:C:35:HIS:ND1	3:C:50:PHE:HB3	2.36	0.41
1:A:743:ASP:OD1	1:A:744:PRO:HD2	2.20	0.41
1:A:911:ARG:HA	1:A:911:ARG:HD3	1.54	0.41
1:A:158:GLU:HG3	1:A:160:GLY:H	1.86	0.41
1:A:987:VAL:O	1:A:991:SER:OG	2.29	0.40
2:B:100:PRO:HA	2:B:101:PRO:HA	1.88	0.40
1:A:729:VAL:HG23	1:A:975:LEU:HD11	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:LEU:HA	1:A:993:PHE:HE1	1.85	0.40
1:A:271:LYS:NZ	1:A:1135:ASN:HA	2.36	0.40
1:A:64:LEU:O	1:A:67:MET:HG2	2.20	0.40
2:B:142:LEU:HB2	2:B:181:MET:HG2	2.04	0.40
3:C:51:ILE:HG22	3:C:58:THR:HG22	2.04	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	1177/1279 (92%)	1088 (92%)	77 (6%)	12 (1%)	12	46
2	B	218/220 (99%)	191 (88%)	23 (11%)	4 (2%)	6	34
3	C	223/225 (99%)	198 (89%)	22 (10%)	3 (1%)	9	41
All	All	1618/1724 (94%)	1477 (91%)	122 (8%)	19 (1%)	13	42

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1163	LYS
3	C	182	ASP
1	A	42	ALA
1	A	589	SER
1	A	1031	GLU
3	C	154	TYR
1	A	166	ASP
3	C	46	GLU
1	A	480	ALA
1	A	587	ARG
1	A	1139	VAL
1	A	1207	THR

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Mol	Chain	Res	Type
2	B	98	HIS
1	A	163	ASP
1	A	1020	TYR
1	A	431	GLY
2	B	56	VAL
2	B	34	GLY
2	B	106	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	972/1056 (92%)	972 (100%)	0	100	100
2	B	199/199 (100%)	199 (100%)	0	100	100
3	C	195/195 (100%)	195 (100%)	0	100	100
All	All	1366/1450 (94%)	1366 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	ASN
1	A	329	GLN
1	A	356	ASN
1	A	586	HIS
1	A	611	HIS
1	A	628	GLN
1	A	724	GLN
1	A	924	GLN
1	A	1063	GLN
1	A	1106	GLN
1	A	1231	HIS
1	A	1256	HIS
2	B	43	GLN

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Mol	Chain	Res	Type
2	B	98	HIS
2	B	196	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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	3PE	A	1304	-	50,50,50	0.93	4 (8%)	53,55,55	1.17	3 (5%)
4	NAG	A	1302	-	14,14,15	2.24	3 (21%)	17,19,21	1.27	2 (11%)
4	NAG	A	1303	-	14,14,15	2.44	4 (28%)	17,19,21	1.49	3 (17%)
4	NAG	A	1301	-	14,14,15	2.23	3 (21%)	17,19,21	1.27	2 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	3PE	A	1304	-	-	19/54/54/54	-
4	NAG	A	1302	-	-	2/6/23/26	0/1/1/1
4	NAG	A	1303	-	-	2/6/23/26	0/1/1/1
4	NAG	A	1301	-	-	2/6/23/26	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1302	NAG	O5-C1	5.29	1.52	1.43
4	A	1301	NAG	O5-C1	5.28	1.52	1.43
4	A	1303	NAG	O5-C1	5.17	1.52	1.43
4	A	1303	NAG	C7-N2	4.70	1.49	1.34
4	A	1301	NAG	C7-N2	4.06	1.47	1.34
4	A	1302	NAG	C7-N2	4.04	1.47	1.34
4	A	1303	NAG	O7-C7	-3.27	1.15	1.23
4	A	1303	NAG	C2-N2	3.19	1.51	1.46
5	A	1304	3PE	O21-C2	-2.74	1.40	1.46
4	A	1302	NAG	O7-C7	-2.51	1.17	1.23
4	A	1301	NAG	O7-C7	-2.50	1.17	1.23
5	A	1304	3PE	O31-C31	2.44	1.40	1.33
5	A	1304	3PE	O31-C3	-2.11	1.40	1.45
5	A	1304	3PE	O21-C21	2.09	1.40	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1304	3PE	O21-C21-C22	4.05	120.25	111.48
4	A	1303	NAG	C1-O5-C5	3.20	116.47	112.19
4	A	1303	NAG	C8-C7-N2	3.13	121.31	116.12
4	A	1303	NAG	O7-C7-C8	-3.04	116.64	122.05
4	A	1301	NAG	C1-C2-N2	2.81	114.86	110.43
4	A	1302	NAG	C1-C2-N2	2.81	114.86	110.43
5	A	1304	3PE	O31-C31-C32	2.63	119.87	111.83
4	A	1301	NAG	C6-C5-C4	-2.57	106.71	113.02
4	A	1302	NAG	C6-C5-C4	-2.57	106.72	113.02
5	A	1304	3PE	O12-P-O14	-2.40	101.27	112.44

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1304	3PE	C22-C21-O21-C2
5	A	1304	3PE	O22-C21-O21-C2
4	A	1301	NAG	O5-C5-C6-O6
4	A	1302	NAG	O5-C5-C6-O6
4	A	1303	NAG	O5-C5-C6-O6
4	A	1303	NAG	C4-C5-C6-O6
4	A	1301	NAG	C4-C5-C6-O6
4	A	1302	NAG	C4-C5-C6-O6
5	A	1304	3PE	C23-C24-C25-C26
5	A	1304	3PE	C36-C37-C38-C39
5	A	1304	3PE	C27-C28-C29-C2A
5	A	1304	3PE	C24-C25-C26-C27
5	A	1304	3PE	C28-C29-C2A-C2B
5	A	1304	3PE	C2B-C2C-C2D-C2E
5	A	1304	3PE	C22-C23-C24-C25
5	A	1304	3PE	O11-C1-C2-O21
5	A	1304	3PE	C29-C2A-C2B-C2C
5	A	1304	3PE	O11-C1-C2-C3
5	A	1304	3PE	C39-C3A-C3B-C3C
5	A	1304	3PE	C1-O11-P-O14
5	A	1304	3PE	C33-C34-C35-C36
5	A	1304	3PE	C35-C36-C37-C38
5	A	1304	3PE	O31-C31-C32-C33
5	A	1304	3PE	C2F-C2G-C2H-C2I
5	A	1304	3PE	C3B-C3C-C3D-C3E

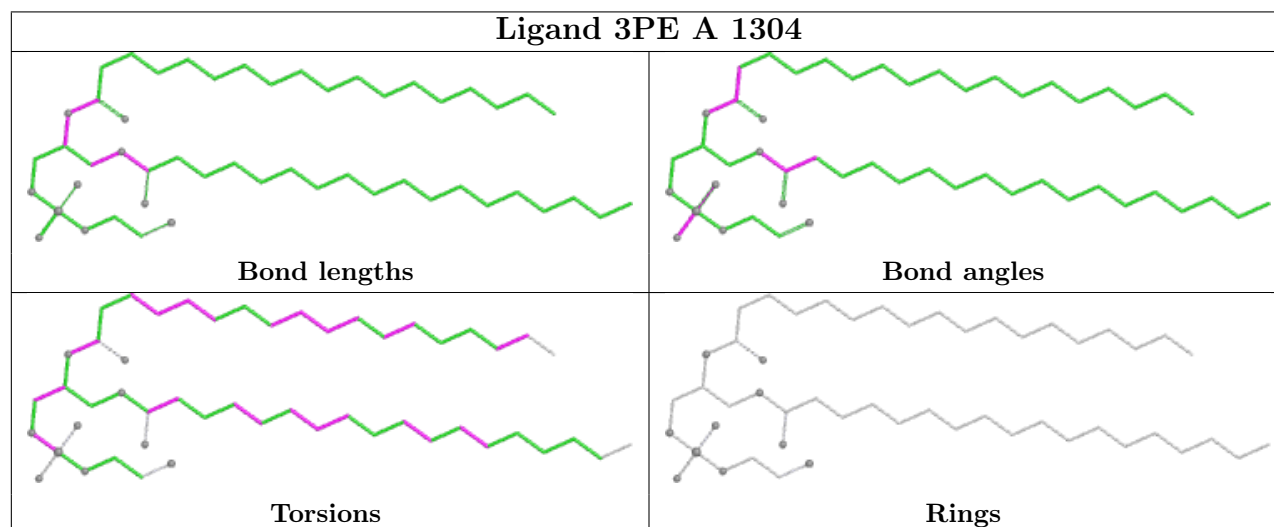
There are no ring outliers.

4 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1304	3PE	2	0
4	A	1302	NAG	7	0
4	A	1303	NAG	4	0
4	A	1301	NAG	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.

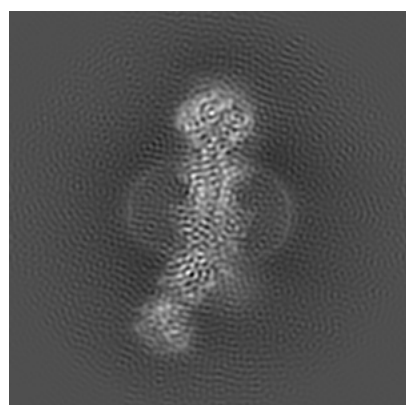
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4282. These allow visual inspection of the internal detail of the map and identification of artifacts.

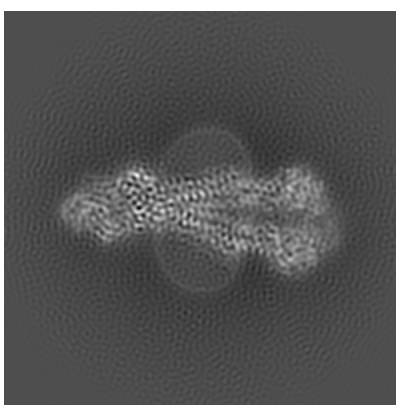
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

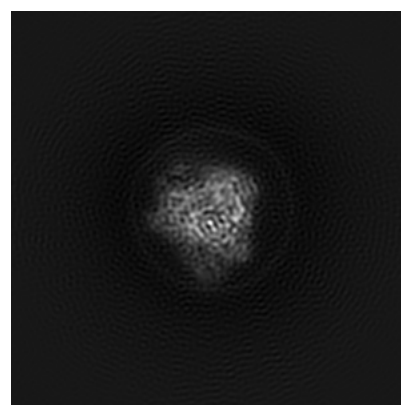
6.1.1 Primary map



X



Y

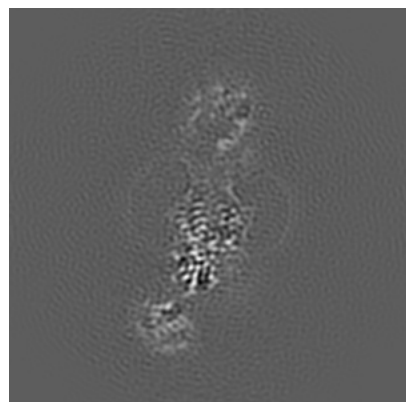


Z

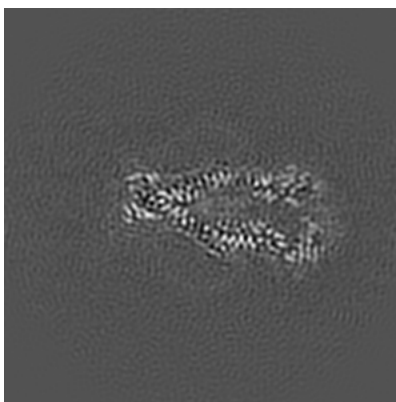
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

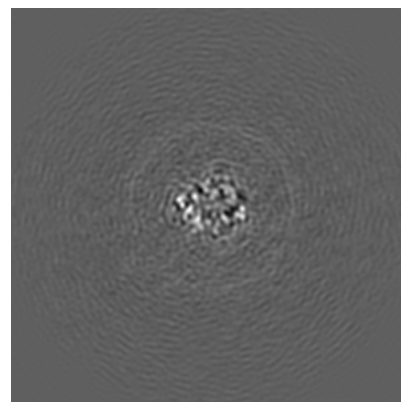
6.2.1 Primary map



X Index: 100



Y Index: 100

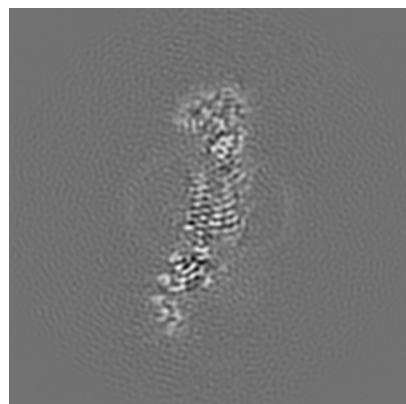


Z Index: 100

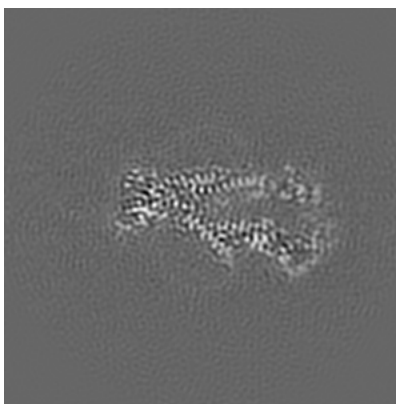
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

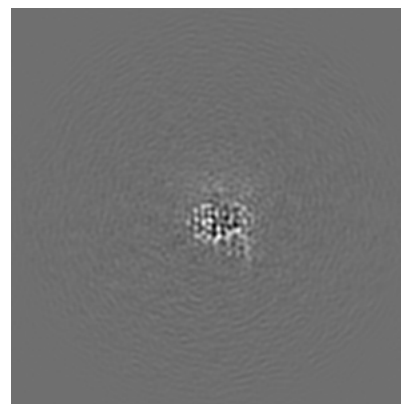
6.3.1 Primary map



X Index: 110



Y Index: 97

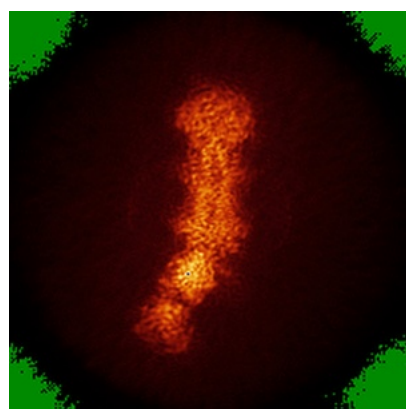


Z Index: 67

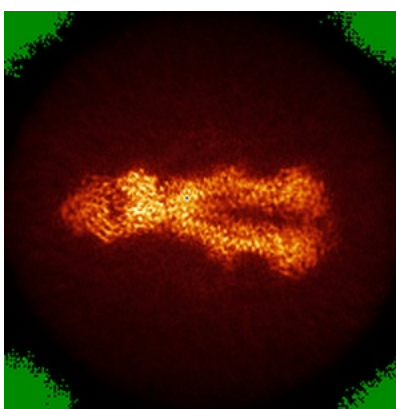
The images above show the largest variance slices of the map in three orthogonal directions.

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

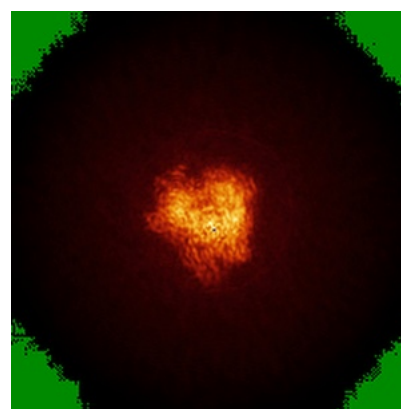
6.4.1 Primary map



X



Y

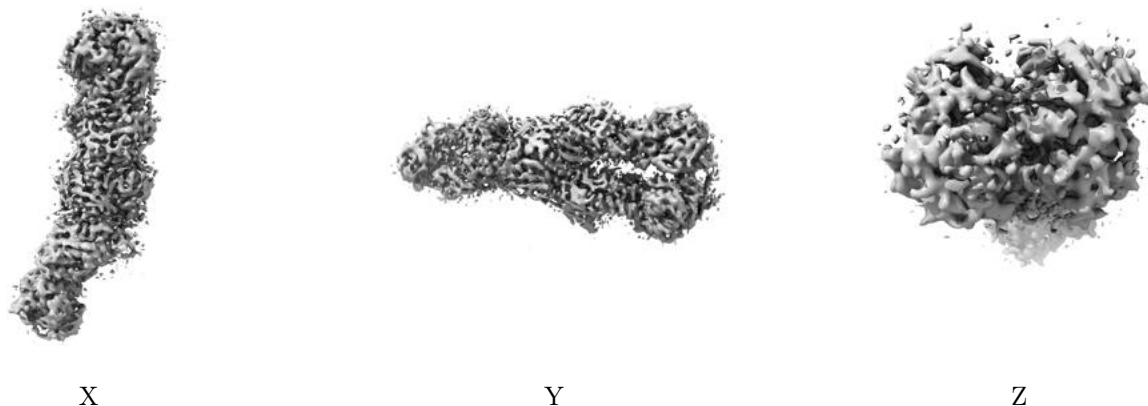


Z

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

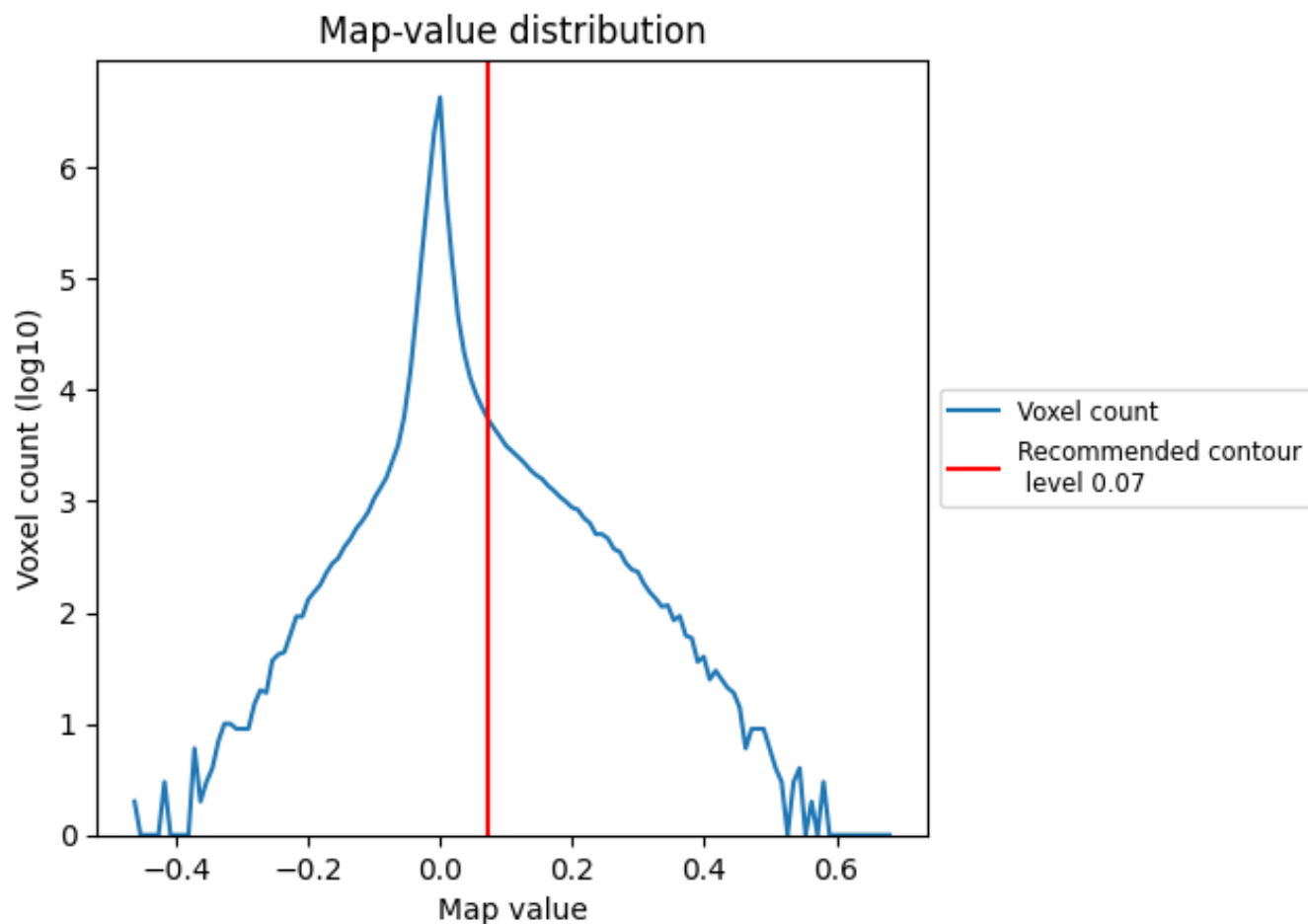
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

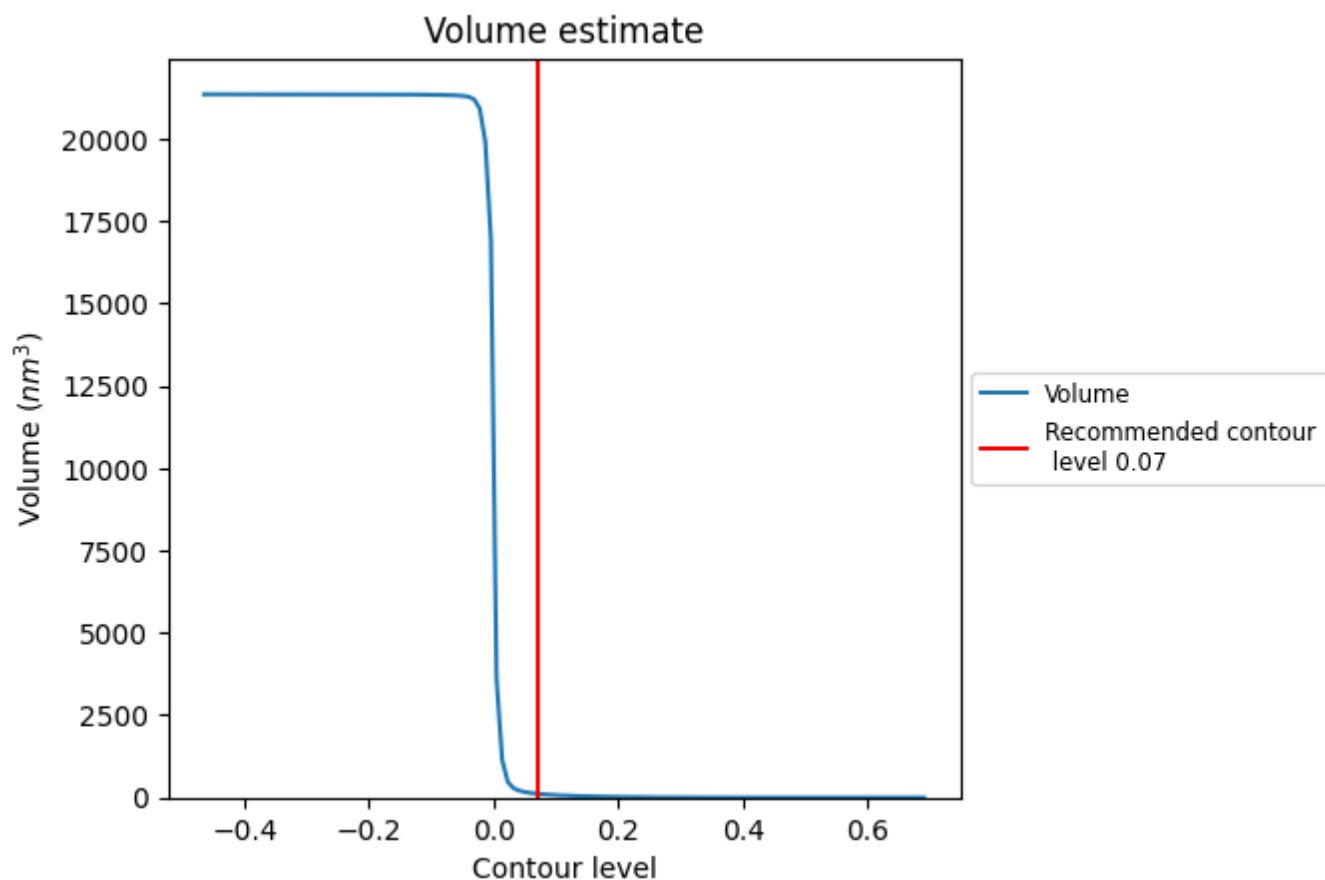
This section contains the results of statistical analysis of the map.

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.

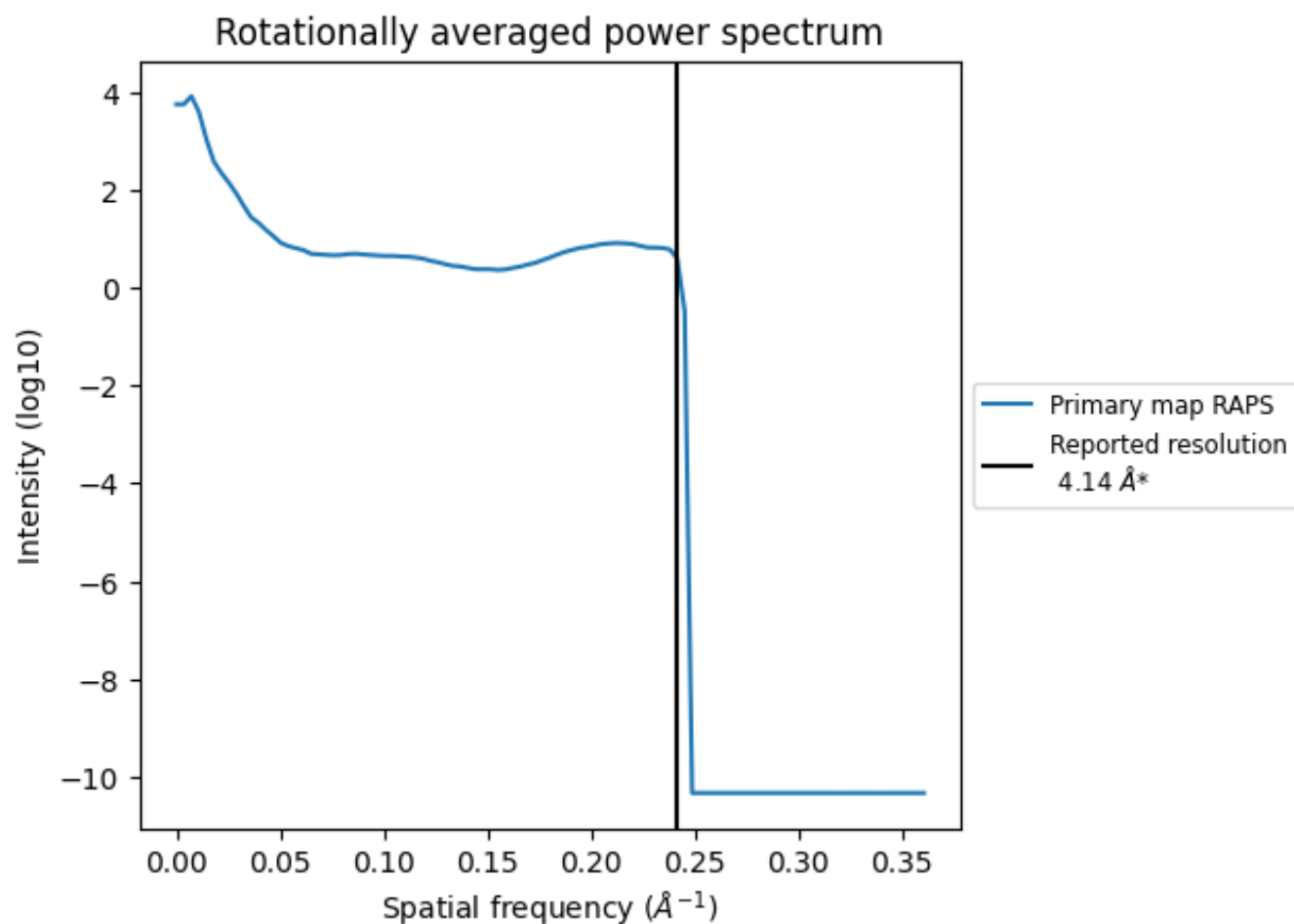
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 115 nm³; this corresponds to an approximate mass of 104 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 0.242 $\text{\AA}^{-1}$

8 Fourier-Shell correlation ⓘ

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

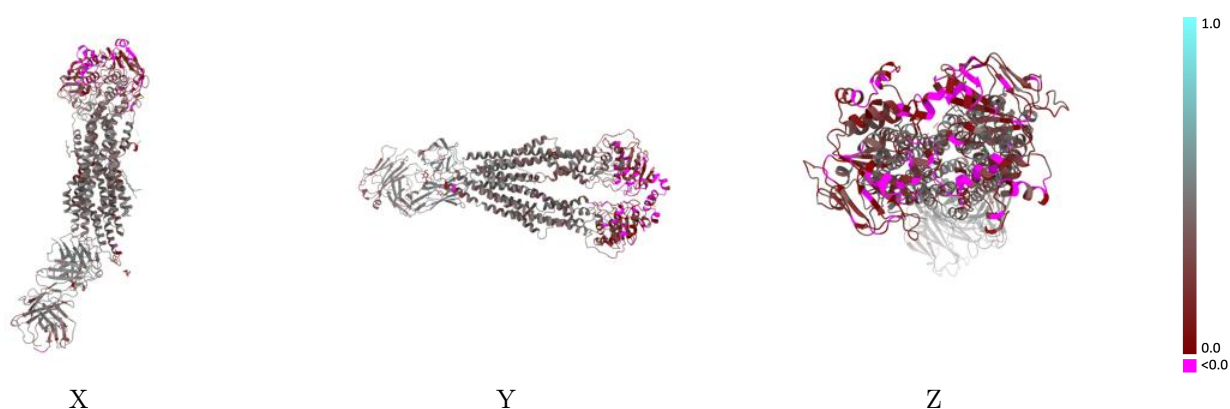
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-4282 and PDB model 6FN4. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

This section was not generated.

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

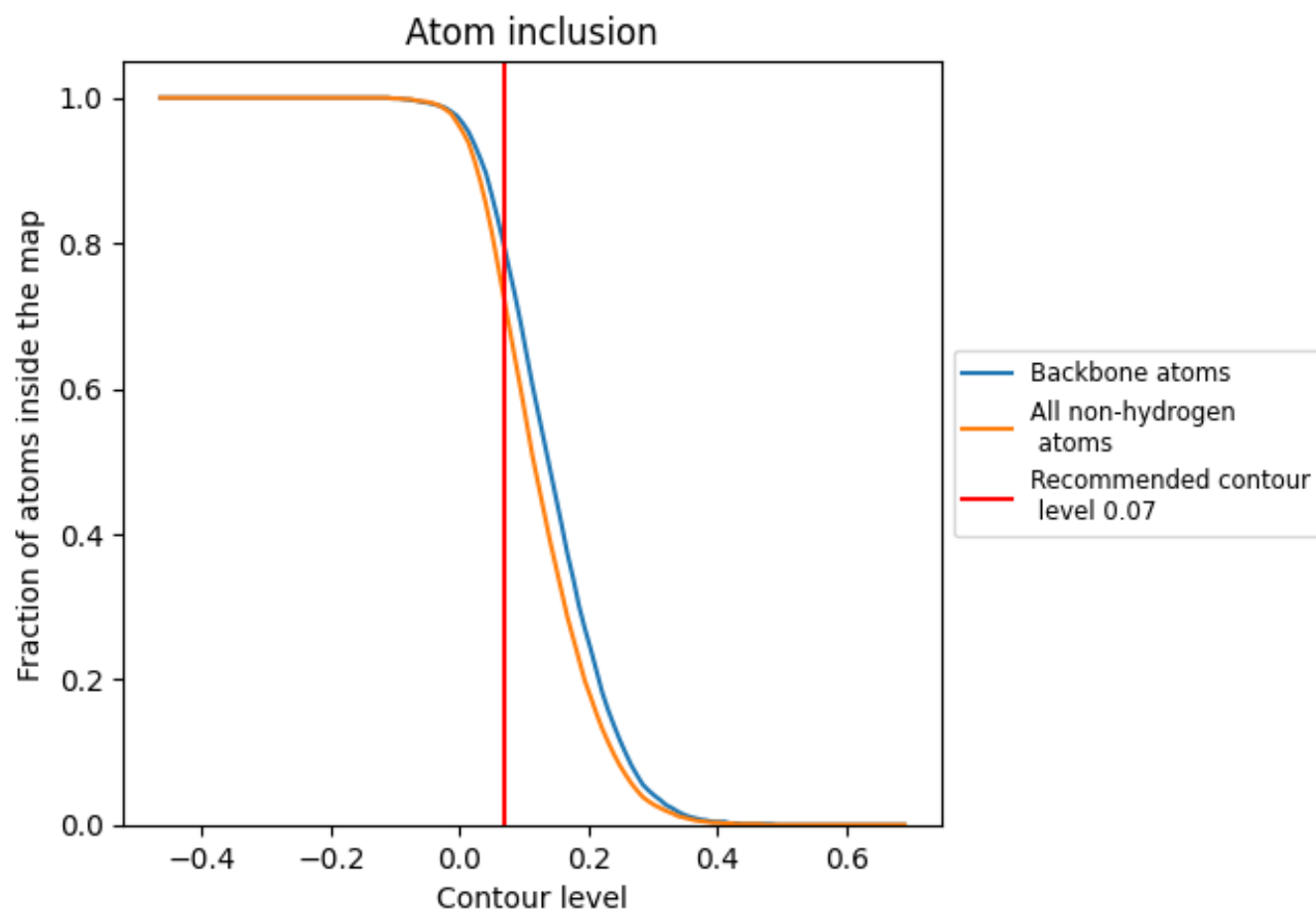


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 72% 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 (0.07) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7170	0.3400
A	0.6720	0.3060
B	0.8340	0.4180
C	0.8420	0.4460

