



## Full wwPDB EM Validation Report ⓘ

Nov 10, 2024 – 01:46 AM EST

PDB ID : 6UM1  
EMDB ID : EMD-20815  
Title : Structure of M-6-P/IGFII Receptor at pH 4.5  
Authors : Wang, R.; Qi, X.; Li, X.  
Deposited on : 2019-10-08  
Resolution : 3.46 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](mailto: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>.

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

EMDB validation analysis	:	0.0.1.dev113
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4.02b-467
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

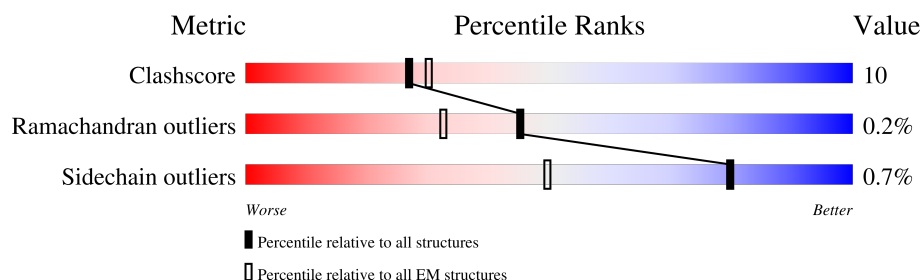
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.46 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	2499	
2	B	2	
3	C	3	

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	2508	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 17134 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 Cation-independent mannose-6-phosphate receptor.

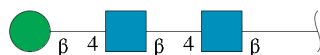
Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	2208	Total	C	N	O	S	0	0
			17025	10631	2916	3342	136		

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
2	B	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
3	C	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

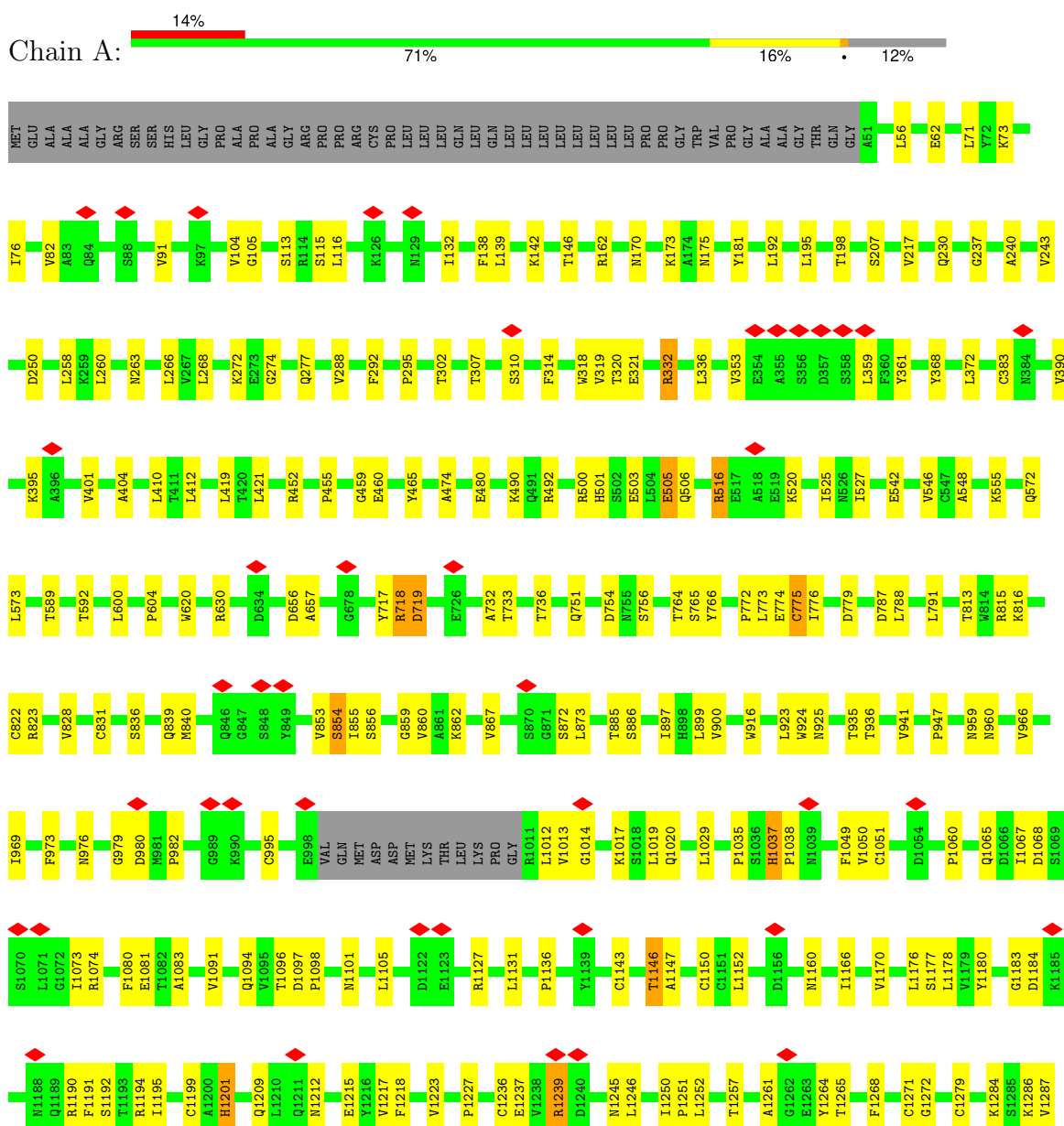


Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	

### 3 Residue-property plots

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

- Molecule 1: Cation-independent mannose-6-phosphate receptor





- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B:  100%

MAG1  
MAG2

- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  33% 67%

MAG1  
MAG2  
BMA3

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	128789	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	80	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	0.137	Depositor
Minimum map value	-0.096	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0145	Depositor
Map size ( $\text{\AA}$ )	280.0, 280.0, 280.0	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.0, 1.0, 1.0	Depositor



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, 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	0.69	0/17415	0.87	7/23648 (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	20

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	656	ASP	CB-CA-C	-8.75	92.90	110.40
1	A	1425	GLY	N-CA-C	-8.25	92.48	113.10
1	A	1037	HIS	C-N-CD	-6.15	107.07	120.60
1	A	1914	PHE	CB-CA-C	6.07	122.53	110.40
1	A	1423	GLN	N-CA-C	-5.92	95.00	111.00
1	A	1239	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	A	791	LEU	CB-CA-C	-5.34	100.05	110.20

There are no chirality outliers.

All (20) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1035	PRO	Peptide
1	A	1037	HIS	Mainchain
1	A	1051	CYS	Peptide

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Mol	Chain	Res	Type	Group
1	A	1180	TYR	Peptide
1	A	1199	CYS	Peptide
1	A	1201	HIS	Peptide
1	A	1272	GLY	Peptide
1	A	1304	ALA	Peptide
1	A	1576	PHE	Peptide
1	A	1656	GLU	Peptide
1	A	1766	ASN	Peptide
1	A	1914	PHE	Peptide
1	A	2186	LEU	Peptide
1	A	310	SER	Peptide
1	A	505	GLU	Peptide
1	A	813	THR	Peptide
1	A	823	ARG	Peptide
1	A	854	SER	Peptide
1	A	855	ILE	Peptide
1	A	995	CYS	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	17025	0	16238	323	0
2	B	28	0	25	0	0
3	C	39	0	34	0	0
4	A	42	0	37	8	0
All	All	17134	0	16334	323	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 10.

All (323) 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:1239:ARG:NH2	1:A:1424:ALA:HB2	1.59	1.17
1:A:718:ARG:HB3	1:A:732:ALA:HB1	1.28	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1239:ARG:HH22	1:A:1424:ALA:CB	1.71	1.02
1:A:1663:VAL:HG13	1:A:1670:ILE:HG23	1.44	1.00
1:A:718:ARG:HB3	1:A:732:ALA:CB	1.93	0.98
1:A:1663:VAL:CG1	1:A:1670:ILE:HG23	1.95	0.96
1:A:897:ILE:HG12	1:A:924:TRP:CD1	2.02	0.95
1:A:1414:ILE:HD11	1:A:1494:PRO:CG	1.97	0.95
1:A:764:THR:HG22	1:A:766:TYR:H	1.31	0.93
1:A:1257:THR:HG21	1:A:1364:TRP:HZ2	1.33	0.93
1:A:146:THR:HG21	1:A:181:TYR:HB3	1.53	0.90
1:A:295:PRO:HG2	1:A:501:HIS:CE1	2.07	0.90
1:A:1287:VAL:HG13	1:A:1307:PHE:HB3	1.56	0.87
1:A:1323:THR:HB	4:A:2508:NAG:H62	1.57	0.86
1:A:1321:ASN:OD1	4:A:2508:NAG:C1	2.25	0.84
1:A:600:LEU:HB3	1:A:772:PRO:HB3	1.59	0.84
1:A:764:THR:HG22	1:A:766:TYR:N	1.92	0.83
1:A:1414:ILE:HD11	1:A:1494:PRO:HG3	1.61	0.83
1:A:274:GLY:O	1:A:277:GLN:NE2	2.14	0.81
1:A:1414:ILE:HG22	1:A:1435:VAL:HG23	1.63	0.80
1:A:192:LEU:CD2	1:A:217:VAL:HG11	2.13	0.79
1:A:2053:SER:OG	1:A:2067:GLY:O	2.01	0.78
1:A:718:ARG:CB	1:A:732:ALA:CB	2.61	0.78
1:A:192:LEU:HD22	1:A:217:VAL:HG11	1.65	0.77
1:A:1414:ILE:HD11	1:A:1494:PRO:CB	2.17	0.74
1:A:192:LEU:HD23	1:A:192:LEU:O	1.87	0.72
1:A:1321:ASN:HD21	4:A:2508:NAG:C1	2.02	0.72
1:A:91:VAL:HG13	1:A:104:VAL:HB	1.71	0.71
1:A:897:ILE:CD1	1:A:924:TRP:NE1	2.53	0.71
1:A:754:ASP:CG	1:A:756:SER:OG	2.29	0.71
1:A:1257:THR:HG21	1:A:1364:TRP:CZ2	2.22	0.71
1:A:516:ARG:HB2	1:A:520:LYS:HB3	1.73	0.70
1:A:2272:THR:HB	1:A:2276:GLN:O	1.93	0.67
1:A:1083:ALA:HB1	1:A:1146:THR:HB	1.77	0.67
1:A:1094:GLN:HE22	1:A:1315:ASN:HB2	1.58	0.67
1:A:717:TYR:O	1:A:733:THR:N	2.27	0.67
1:A:1257:THR:CG2	1:A:1364:TRP:HZ2	2.06	0.66
1:A:1195:ILE:HG22	1:A:1218:PHE:HB2	1.77	0.66
1:A:828:VAL:HG23	1:A:831:CYS:SG	2.35	0.66
1:A:1427:ASP:OD1	1:A:1427:ASP:N	2.28	0.66
1:A:2017:LEU:HD22	1:A:2037:LEU:HB2	1.77	0.66
1:A:525:ILE:HD12	1:A:527:ILE:CD1	2.25	0.65
1:A:1287:VAL:CG1	1:A:1307:PHE:HB3	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1663:VAL:HG13	1:A:1670:ILE:CG2	2.24	0.64
1:A:2053:SER:CB	1:A:2067:GLY:O	2.46	0.64
1:A:321:GLU:HG3	1:A:501:HIS:CE1	2.33	0.64
1:A:718:ARG:CB	1:A:732:ALA:HB2	2.27	0.64
1:A:1321:ASN:ND2	4:A:2508:NAG:C1	2.60	0.64
1:A:2017:LEU:HD12	1:A:2128:TRP:NE1	2.13	0.63
1:A:899:LEU:HA	1:A:924:TRP:HB2	1.80	0.63
1:A:1495:MET:SD	1:A:1888:GLN:NE2	2.72	0.62
1:A:500:ARG:NH2	1:A:506:GLN:O	2.31	0.62
1:A:1742:LEU:HD12	1:A:1748:GLU:O	1.98	0.62
1:A:899:LEU:HG	1:A:924:TRP:HB2	1.79	0.62
1:A:195:LEU:HD21	1:A:320:THR:HG21	1.82	0.61
1:A:773:LEU:HD11	1:A:822:CYS:HB3	1.83	0.61
1:A:718:ARG:HG2	1:A:718:ARG:HH11	1.64	0.61
1:A:1696:SER:OG	1:A:1697:PRO:HD3	2.01	0.61
1:A:288:VAL:HG22	1:A:314:PHE:HB2	1.82	0.60
1:A:1323:THR:CB	4:A:2508:NAG:H62	2.28	0.60
1:A:1350:VAL:CG2	1:A:1363:GLU:HB3	2.32	0.60
1:A:1250:ILE:N	1:A:1251:PRO:HD2	2.16	0.60
1:A:815:ARG:HD2	1:A:839:GLN:HE21	1.66	0.59
1:A:773:LEU:CD1	1:A:822:CYS:HB3	2.33	0.59
1:A:1291:CYS:SG	1:A:1302:LYS:O	2.61	0.59
1:A:897:ILE:HG12	1:A:924:TRP:NE1	2.16	0.59
1:A:146:THR:HG21	1:A:181:TYR:CB	2.31	0.59
1:A:899:LEU:HG	1:A:924:TRP:CD1	2.38	0.59
1:A:959:ASN:HD21	1:A:980:ASP:HB2	1.68	0.59
1:A:1624:ARG:NH1	1:A:1627:VAL:O	2.35	0.59
1:A:966:VAL:HB	1:A:973:PHE:HB2	1.85	0.58
1:A:897:ILE:CG1	1:A:924:TRP:CD1	2.82	0.58
1:A:2090:TYR:HB3	1:A:2092:GLY:O	2.02	0.58
1:A:718:ARG:HA	1:A:732:ALA:HA	1.85	0.58
1:A:525:ILE:HD12	1:A:527:ILE:HD11	1.83	0.57
1:A:773:LEU:HD11	1:A:822:CYS:CB	2.33	0.57
1:A:116:LEU:HD22	1:A:459:GLY:O	2.04	0.57
1:A:525:ILE:HD13	1:A:604:PRO:HG3	1.85	0.57
1:A:1416:VAL:HG12	1:A:1509:TRP:CZ3	2.40	0.57
1:A:295:PRO:HG2	1:A:501:HIS:ND1	2.19	0.57
1:A:718:ARG:HB2	1:A:732:ALA:HB2	1.86	0.57
1:A:1183:GLY:HA3	1:A:1191:PHE:HB2	1.86	0.57
1:A:1663:VAL:HG12	1:A:1670:ILE:O	2.04	0.57
1:A:91:VAL:HG12	1:A:105:GLY:HA3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:717:TYR:O	1:A:733:THR:O	2.23	0.57
1:A:321:GLU:HG3	1:A:501:HIS:HE1	1.68	0.57
1:A:1494:PRO:HB3	1:A:1509:TRP:CE3	2.40	0.57
1:A:899:LEU:HG	1:A:924:TRP:CG	2.41	0.56
1:A:1287:VAL:HG13	1:A:1287:VAL:O	2.04	0.56
1:A:525:ILE:CD1	1:A:527:ILE:HD11	2.35	0.56
1:A:1759:CYS:HB3	1:A:1761:ALA:O	2.05	0.56
1:A:195:LEU:CD2	1:A:320:THR:HG21	2.35	0.56
1:A:1771:ILE:HG22	1:A:1796:PHE:HB2	1.86	0.56
1:A:899:LEU:HG	1:A:924:TRP:CB	2.35	0.55
1:A:1060:PRO:HA	1:A:1080:PHE:HB3	1.88	0.55
1:A:1597:LEU:HB2	1:A:1621:PHE:HB2	1.87	0.55
1:A:260:LEU:HD23	1:A:260:LEU:O	2.06	0.55
1:A:1321:ASN:CG	4:A:2508:NAG:C1	2.73	0.55
1:A:1318:LEU:HB2	1:A:1340:PHE:HB2	1.89	0.55
1:A:764:THR:HG22	1:A:765:SER:N	2.23	0.54
1:A:1321:ASN:HD21	4:A:2508:NAG:C2	2.19	0.54
1:A:899:LEU:CD1	1:A:924:TRP:CD1	2.91	0.54
1:A:1178:LEU:HB3	1:A:1195:ILE:HG13	1.87	0.54
1:A:2017:LEU:HD12	1:A:2128:TRP:CE2	2.43	0.54
1:A:516:ARG:HB2	1:A:520:LYS:CB	2.38	0.54
1:A:361:TYR:OH	1:A:452:ARG:NH1	2.41	0.54
1:A:1414:ILE:CG2	1:A:1435:VAL:HG23	2.37	0.54
1:A:2268:PHE:HE1	1:A:2277:TYR:HB3	1.73	0.54
1:A:1535:LEU:HD23	1:A:1708:ASN:HD22	1.73	0.53
1:A:1355:THR:HG23	1:A:1356:SER:N	2.24	0.53
1:A:2065:VAL:O	1:A:2089:HIS:HB2	2.09	0.53
1:A:2173:VAL:HG21	1:A:2266:PRO:HG2	1.90	0.53
1:A:935:THR:HG21	1:A:947:PRO:HG3	1.91	0.53
1:A:1019:LEU:HB3	1:A:1029:LEU:HG	1.91	0.53
1:A:885:THR:HG23	1:A:886:SER:O	2.09	0.52
1:A:2053:SER:HB3	1:A:2067:GLY:O	2.09	0.52
1:A:897:ILE:HD13	1:A:924:TRP:NE1	2.25	0.52
1:A:1236:CYS:O	1:A:1311:LEU:HD21	2.09	0.52
1:A:1338:ILE:HG21	1:A:1340:PHE:CZ	2.44	0.52
1:A:816:LYS:HZ2	1:A:1096:THR:HG21	1.74	0.52
1:A:872:SER:HB3	1:A:900:VAL:HG12	1.91	0.52
1:A:1029:LEU:HD13	1:A:1049:PHE:HE1	1.74	0.52
1:A:1875:PHE:HA	1:A:1958:ARG:HB3	1.92	0.51
1:A:525:ILE:HD12	1:A:527:ILE:HD13	1.91	0.51
1:A:969:ILE:HD11	1:A:1065:GLN:HE21	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1393:LEU:HD11	1:A:1511:THR:HG21	1.92	0.51
1:A:2017:LEU:CD2	1:A:2037:LEU:HB2	2.40	0.51
1:A:1143:CYS:SG	1:A:1160:ASN:ND2	2.83	0.51
1:A:1695:THR:O	1:A:1698:ASP:HB2	2.11	0.51
1:A:1883:LEU:HB3	1:A:1894:LEU:HA	1.93	0.51
1:A:1955:THR:HG22	1:A:1955:THR:O	2.11	0.50
1:A:525:ILE:CD1	1:A:604:PRO:HG3	2.41	0.50
1:A:840:MET:HB3	1:A:853:VAL:HB	1.92	0.50
1:A:460:GLU:HG3	1:A:465:TYR:CE1	2.45	0.50
1:A:1288:ILE:HG23	1:A:1304:ALA:HB2	1.94	0.50
1:A:1416:VAL:CG1	1:A:1509:TRP:CH2	2.95	0.50
1:A:774:GLU:O	1:A:776:ILE:N	2.44	0.50
1:A:542:GLU:OE1	1:A:542:GLU:N	2.39	0.50
1:A:754:ASP:CG	1:A:756:SER:HG	2.15	0.50
1:A:1239:ARG:HH22	1:A:1424:ALA:HB2	0.74	0.50
1:A:1387:THR:HG23	1:A:1517:VAL:HG22	1.93	0.50
1:A:82:VAL:HG13	1:A:82:VAL:O	2.12	0.49
1:A:302:THR:HA	1:A:319:VAL:O	2.11	0.49
1:A:897:ILE:CG1	1:A:924:TRP:NE1	2.76	0.49
1:A:1973:GLY:O	1:A:1974:ARG:NH1	2.45	0.49
1:A:1098:PRO:HD2	1:A:1170:VAL:HG21	1.94	0.49
1:A:1201:HIS:CD2	1:A:1223:VAL:HG21	2.47	0.49
1:A:1201:HIS:CD2	1:A:1396:TYR:CD2	3.01	0.49
1:A:91:VAL:HG13	1:A:104:VAL:CB	2.41	0.49
1:A:1314:GLU:O	1:A:1317:VAL:HG12	2.13	0.49
1:A:1416:VAL:HG13	1:A:1509:TRP:CH2	2.48	0.49
1:A:2216:VAL:HG23	1:A:2275:CYS:O	2.13	0.49
1:A:113:SER:HB2	1:A:116:LEU:HB3	1.95	0.49
1:A:1423:GLN:NE2	1:A:1429:CYS:H	2.11	0.49
1:A:460:GLU:HG3	1:A:465:TYR:HE1	1.77	0.48
1:A:1271:CYS:N	1:A:1322:TYR:OH	2.44	0.48
1:A:751:GLN:OE1	1:A:751:GLN:N	2.46	0.48
1:A:923:LEU:HD22	1:A:1352:LEU:HD11	1.96	0.48
1:A:1252:LEU:HD22	1:A:1364:TRP:HZ3	1.78	0.48
1:A:1416:VAL:CG1	1:A:1509:TRP:CZ3	2.96	0.48
1:A:1657:GLN:HG3	1:A:1706:PRO:HD3	1.96	0.48
1:A:775:CYS:O	1:A:776:ILE:HG12	2.14	0.48
1:A:2083:VAL:HG23	1:A:2101:ILE:HB	1.95	0.48
1:A:1387:THR:HG23	1:A:1517:VAL:CG2	2.44	0.48
1:A:1425:GLY:C	1:A:1426:SER:O	2.46	0.47
1:A:2110:GLY:HA3	1:A:2129:ASP:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:839:GLN:HB3	1:A:856:SER:HB3	1.95	0.47
1:A:1732:ASP:OD1	1:A:1734:GLY:O	2.33	0.47
1:A:266:LEU:HB2	1:A:292:PHE:HB2	1.96	0.47
1:A:1904:PRO:HG2	1:A:1953:HIS:HA	1.95	0.47
1:A:2101:ILE:HG23	1:A:2126:PHE:HB3	1.96	0.47
1:A:836:SER:N	1:A:859:GLY:O	2.44	0.47
1:A:1201:HIS:NE2	1:A:1396:TYR:HD2	2.12	0.47
1:A:1670:ILE:HD11	1:A:1803:VAL:HA	1.97	0.47
1:A:1131:LEU:HD12	1:A:1150:CYS:HB3	1.96	0.47
1:A:1184:ASP:N	1:A:1190:ARG:O	2.48	0.47
1:A:1830:SER:HB2	1:A:1849:THR:HG23	1.96	0.47
1:A:368:TYR:HA	1:A:395:LYS:HG2	1.97	0.47
1:A:779:ASP:HA	1:A:867:VAL:HG11	1.96	0.47
1:A:1271:CYS:H	1:A:1322:TYR:HH	1.61	0.47
1:A:1013:VAL:HG22	1:A:1014:GLY:O	2.15	0.47
1:A:1201:HIS:CD2	1:A:1223:VAL:CG2	2.98	0.47
1:A:318:TRP:NE1	1:A:320:THR:HG22	2.30	0.46
1:A:490:LYS:O	1:A:492:ARG:NH2	2.46	0.46
1:A:1194:ARG:HB2	1:A:1217:VAL:HG23	1.97	0.46
1:A:1672:LEU:HD12	1:A:1675:LEU:HD12	1.96	0.46
1:A:62:GLU:HG2	1:A:73:LYS:HG2	1.97	0.46
1:A:336:LEU:HD23	1:A:412:LEU:HG	1.98	0.46
1:A:1029:LEU:HD13	1:A:1049:PHE:CE1	2.50	0.46
1:A:1509:TRP:CD1	1:A:1511:THR:HB	2.50	0.46
1:A:1551:TYR:HB3	1:A:1635:LEU:HD23	1.97	0.46
1:A:1569:ALA:O	1:A:1572:VAL:HG12	2.15	0.46
1:A:170:ASN:HB2	1:A:173:LYS:HB2	1.97	0.46
1:A:2266:PRO:HB3	1:A:2281:TRP:CG	2.51	0.46
1:A:372:LEU:HB2	1:A:390:VAL:HG12	1.97	0.46
1:A:1127:ARG:HD2	1:A:1152:LEU:HD21	1.97	0.46
1:A:1067:ILE:HG22	1:A:1074:ARG:HB2	1.97	0.46
1:A:230:GLN:HB2	1:A:243:VAL:HG11	1.96	0.46
1:A:1177:SER:OG	1:A:1194:ARG:NH1	2.48	0.46
1:A:332:ARG:HA	1:A:332:ARG:HD3	1.51	0.46
1:A:1019:LEU:HD22	1:A:1029:LEU:HD11	1.98	0.46
1:A:1091:VAL:HG13	1:A:1166:ILE:HG23	1.98	0.46
1:A:1406:THR:HG23	1:A:1806:ASP:HB2	1.98	0.46
1:A:1689:GLU:HB2	1:A:2225:TYR:HD2	1.82	0.46
1:A:873:LEU:HD12	1:A:873:LEU:O	2.16	0.45
1:A:115:SER:O	1:A:138:PHE:HB2	2.16	0.45
1:A:383:CYS:HA	1:A:401:VAL:HG11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1355:THR:HG21	1:A:1361:LEU:CD2	2.45	0.45
1:A:1624:ARG:HE	1:A:1650:HIS:HB3	1.80	0.45
1:A:258:LEU:HB3	1:A:268:LEU:HB2	1.98	0.45
1:A:1600:VAL:HA	1:A:1618:VAL:HG12	1.98	0.45
1:A:104:VAL:HG13	1:A:132:ILE:HG22	1.98	0.45
1:A:410:LEU:HD23	1:A:421:LEU:HD11	1.98	0.45
1:A:546:VAL:HG21	1:A:589:THR:HG21	1.97	0.45
1:A:941:VAL:O	1:A:1019:LEU:HD21	2.16	0.45
1:A:1735:ARG:N	1:A:1754:GLU:O	2.46	0.45
1:A:1068:ASP:O	1:A:1073:ILE:O	2.34	0.45
1:A:1268:PHE:HA	1:A:1290:SER:HA	1.99	0.45
1:A:372:LEU:HD11	1:A:455:PRO:HG3	1.98	0.44
1:A:548:ALA:O	1:A:555:LYS:N	2.51	0.44
1:A:2268:PHE:HA	1:A:2279:PHE:CB	2.47	0.44
1:A:1450:ARG:HB2	1:A:1465:VAL:HG22	1.99	0.44
1:A:175:ASN:HD21	1:A:237:GLY:HA2	1.82	0.44
1:A:1402:ALA:HB3	1:A:1412:TYR:HB2	2.00	0.44
1:A:548:ALA:HB3	1:A:555:LYS:HB2	1.99	0.44
1:A:1105:LEU:HD11	1:A:1176:LEU:HD13	1.99	0.44
1:A:1265:THR:OG1	1:A:1295:ARG:NH2	2.50	0.44
1:A:307:THR:HB	1:A:736:THR:HG21	1.99	0.44
1:A:1212:ASN:O	1:A:1215:GLU:HG2	2.18	0.44
1:A:1245:ASN:HD22	1:A:1372:PRO:HD2	1.81	0.44
1:A:764:THR:CG2	1:A:766:TYR:H	2.17	0.44
1:A:1261:ALA:HB3	1:A:1264:TYR:HB2	1.99	0.44
1:A:1696:SER:OG	1:A:2009:HIS:N	2.31	0.44
1:A:1786:LEU:HA	1:A:1796:PHE:HA	1.99	0.44
1:A:1866:LEU:HB3	1:A:1873:ALA:HB3	1.99	0.44
1:A:1773:PHE:HB3	1:A:1803:VAL:HG21	2.00	0.43
1:A:2035:ILE:HA	1:A:2054:VAL:HG22	2.00	0.43
1:A:527:ILE:HD11	1:A:620:TRP:CZ3	2.53	0.43
1:A:1321:ASN:HD21	4:A:2508:NAG:H2	1.81	0.43
1:A:2115:THR:HB	1:A:2125:HIS:C	2.39	0.43
1:A:816:LYS:HE3	1:A:1096:THR:HG21	2.00	0.43
1:A:2120:ASP:OD1	1:A:2121:SER:N	2.52	0.43
1:A:1252:LEU:HD21	1:A:1368:TYR:HE2	1.83	0.43
1:A:1423:GLN:NE2	1:A:1429:CYS:SG	2.87	0.43
1:A:2259:PRO:HA	1:A:2284:SER:HA	2.01	0.43
1:A:916:TRP:CG	1:A:1356:SER:HB3	2.54	0.43
1:A:1292:GLN:HE21	1:A:1302:LYS:HB2	1.83	0.43
1:A:718:ARG:HH11	1:A:718:ARG:CG	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:899:LEU:HD11	1:A:924:TRP:CD1	2.54	0.43
1:A:1414:ILE:CD1	1:A:1494:PRO:HG3	2.42	0.43
1:A:2081:VAL:HG12	1:A:2103:LEU:HB3	2.00	0.43
1:A:2254:PHE:HE2	1:A:2278:LEU:HD21	1.84	0.43
1:A:207:SER:OG	1:A:630:ARG:NH1	2.51	0.43
1:A:1192:SER:O	1:A:1215:GLU:HA	2.19	0.43
1:A:527:ILE:HD11	1:A:620:TRP:HZ3	1.83	0.43
1:A:776:ILE:HD13	1:A:787:ASP:HA	2.01	0.43
1:A:1314:GLU:HB2	1:A:1317:VAL:HG13	2.01	0.43
1:A:1426:SER:O	1:A:1427:ASP:C	2.57	0.42
1:A:56:LEU:HB3	1:A:76:ILE:HD13	2.00	0.42
1:A:897:ILE:HD11	1:A:924:TRP:CE2	2.54	0.42
1:A:1536:PHE:HB3	1:A:1597:LEU:HD11	2.02	0.42
1:A:1981:VAL:HG22	1:A:1982:ARG:HD2	2.01	0.42
1:A:62:GLU:HB3	1:A:71:LEU:HD11	2.01	0.42
1:A:419:LEU:HD23	1:A:474:ALA:HB1	2.01	0.42
1:A:1209:GLN:HB3	1:A:1217:VAL:HG13	1.99	0.42
1:A:718:ARG:CG	1:A:718:ARG:NH1	2.81	0.42
1:A:897:ILE:CD1	1:A:924:TRP:CE2	3.02	0.42
1:A:1201:HIS:CD2	1:A:1396:TYR:HD2	2.37	0.42
1:A:1411:HIS:O	1:A:1438:LEU:N	2.44	0.42
1:A:1877:ARG:HB3	1:A:1880:SER:HB3	2.01	0.42
1:A:353:VAL:HG21	1:A:359:LEU:O	2.20	0.42
1:A:1517:VAL:HG23	1:A:1517:VAL:O	2.19	0.42
1:A:2241:CYS:HB2	1:A:2275:CYS:HB2	1.89	0.42
1:A:788:LEU:HD11	1:A:873:LEU:CD1	2.50	0.42
1:A:250:ASP:O	1:A:272:LYS:HD3	2.20	0.42
1:A:390:VAL:HG23	1:A:404:ALA:HB3	2.01	0.41
1:A:840:MET:HA	1:A:854:SER:O	2.21	0.41
1:A:979:GLY:HA3	1:A:1017:LYS:HE2	2.01	0.41
1:A:1050:VAL:HG23	1:A:1081:GLU:HG2	2.01	0.41
1:A:1788:ARG:NH2	1:A:1789:THR:O	2.53	0.41
1:A:1284:LYS:HB2	1:A:1286:LYS:HG2	2.02	0.41
1:A:1338:ILE:HG21	1:A:1340:PHE:CE1	2.55	0.41
1:A:1338:ILE:CG2	1:A:1340:PHE:CZ	3.02	0.41
1:A:936:THR:HG22	1:A:1020:GLN:HB3	2.02	0.41
1:A:1012:LEU:HD22	1:A:1013:VAL:HG12	2.01	0.41
1:A:2057:LYS:HA	1:A:2063:VAL:HG13	2.02	0.41
1:A:2214:ARG:HD2	1:A:2214:ARG:HA	1.85	0.41
1:A:503:GLU:HG3	1:A:505:GLU:H	1.86	0.41
1:A:2159:LEU:HG	1:A:2163:TYR:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ALA:HB1	1:A:288:VAL:HG21	2.02	0.41
1:A:2075:ASP:HB3	1:A:2082:ILE:HD12	2.02	0.41
1:A:925:ASN:ND2	1:A:1363:GLU:OE2	2.39	0.41
1:A:1583:VAL:HG21	1:A:1645:LEU:HD12	2.03	0.41
1:A:572:GLN:HG2	1:A:592:THR:HG22	2.01	0.41
1:A:718:ARG:HG2	1:A:718:ARG:NH1	2.33	0.41
1:A:860:VAL:HG13	1:A:862:LYS:HE2	2.03	0.41
1:A:872:SER:CB	1:A:900:VAL:HG12	2.51	0.41
1:A:1136:PRO:HB3	1:A:1147:ALA:HA	2.01	0.41
1:A:1435:VAL:HG21	1:A:1479:THR:HG21	2.02	0.41
1:A:1623:CYS:SG	1:A:1624:ARG:N	2.93	0.41
1:A:1201:HIS:CD2	1:A:1396:TYR:CE2	3.09	0.40
1:A:1686:ASP:HA	1:A:1786:LEU:HB3	2.04	0.40
1:A:1743:ASN:HB2	1:A:1750:TYR:HE2	1.86	0.40
1:A:1067:ILE:HG13	1:A:1067:ILE:O	2.21	0.40
1:A:2185:GLN:NE2	1:A:2188:SER:OG	2.54	0.40
1:A:139:LEU:HB2	1:A:162:ARG:HB3	2.03	0.40
1:A:941:VAL:O	1:A:1019:LEU:CD2	2.70	0.40
1:A:1414:ILE:HD11	1:A:1494:PRO:HB2	2.00	0.40
1:A:1757:THR:HA	1:A:1758:PRO:HD3	1.97	0.40
1:A:1787:LEU:HD13	1:A:2278:LEU:HD21	2.04	0.40
1:A:2103:LEU:HD13	1:A:2128:TRP:HD1	1.87	0.40
1:A:527:ILE:O	1:A:573:LEU:HD21	2.21	0.40
1:A:717:TYR:HD1	1:A:717:TYR:HA	1.65	0.40
1:A:900:VAL:HG11	1:A:1350:VAL:HG21	2.03	0.40
1:A:923:LEU:HD13	1:A:1352:LEU:HD11	2.03	0.40
1:A:1101:ASN:HB3	1:A:1227:PRO:HB3	2.03	0.40
1:A:1237:GLU:HB3	1:A:1246:LEU:HD12	2.03	0.40
1:A:2000:LYS:HD2	1:A:2039:GLN:HA	2.02	0.40
1:A:2010:ARG:HD2	1:A:2134:CYS:HA	2.03	0.40
1:A:142:LYS:HB3	1:A:198:THR:HG21	2.04	0.40
1:A:960:ASN:O	1:A:976:ASN:ND2	2.55	0.40
1:A:980:ASP:O	1:A:982:PRO:HD3	2.21	0.40
1:A:1845:VAL:HG12	1:A:1864:VAL:HG13	2.03	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	2200/2499 (88%)	2044 (93%)	151 (7%)	5 (0%)	44 76

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	ASN
1	A	657	ALA
1	A	719	ASP
1	A	1038	PRO
1	A	775	CYS

### 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	1904/2142 (89%)	1891 (99%)	13 (1%)	81 89

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	332	ARG
1	A	480	GLU
1	A	516	ARG
1	A	718	ARG
1	A	719	ASP
1	A	1097	ASP

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Mol	Chain	Res	Type
1	A	1146	THR
1	A	1279	CYS
1	A	1423	GLN
1	A	1426	SER
1	A	1427	ASP
1	A	1902	CYS
1	A	2083	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	127	GLN
1	A	311	ASN
1	A	348	GLN
1	A	501	HIS
1	A	688	GLN
1	A	801	ASN
1	A	839	GLN
1	A	857	ASN
1	A	909	HIS
1	A	959	ASN
1	A	1039	ASN
1	A	1094	GLN
1	A	1201	HIS
1	A	1292	GLN
1	A	1329	HIS
1	A	1399	ASN
1	A	1567	ASN
1	A	1708	ASN
1	A	1887	HIS
1	A	2064	GLN
1	A	2089	HIS
1	A	2145	ASN
1	A	2153	ASN
1	A	2227	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

5 monosaccharides 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$
2	NAG	B	1	1,2	14,14,15	0.37	0	17,19,21	0.70	0
2	NAG	B	2	2	14,14,15	0.36	0	17,19,21	0.74	0
3	NAG	C	1	1,3	14,14,15	0.29	0	17,19,21	1.03	2 (11%)
3	NAG	C	2	3	14,14,15	0.31	0	17,19,21	1.05	2 (11%)
3	BMA	C	3	3	11,11,12	0.39	0	15,15,17	0.65	0

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
2	NAG	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
3	NAG	C	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1
3	BMA	C	3	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	NAG	C2-N2-C7	2.37	126.08	122.90
3	C	2	NAG	C8-C7-N2	2.31	119.95	116.12
3	C	1	NAG	C8-C7-N2	2.28	119.90	116.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	NAG	C2-N2-C7	2.04	125.63	122.90

There are no chirality outliers.

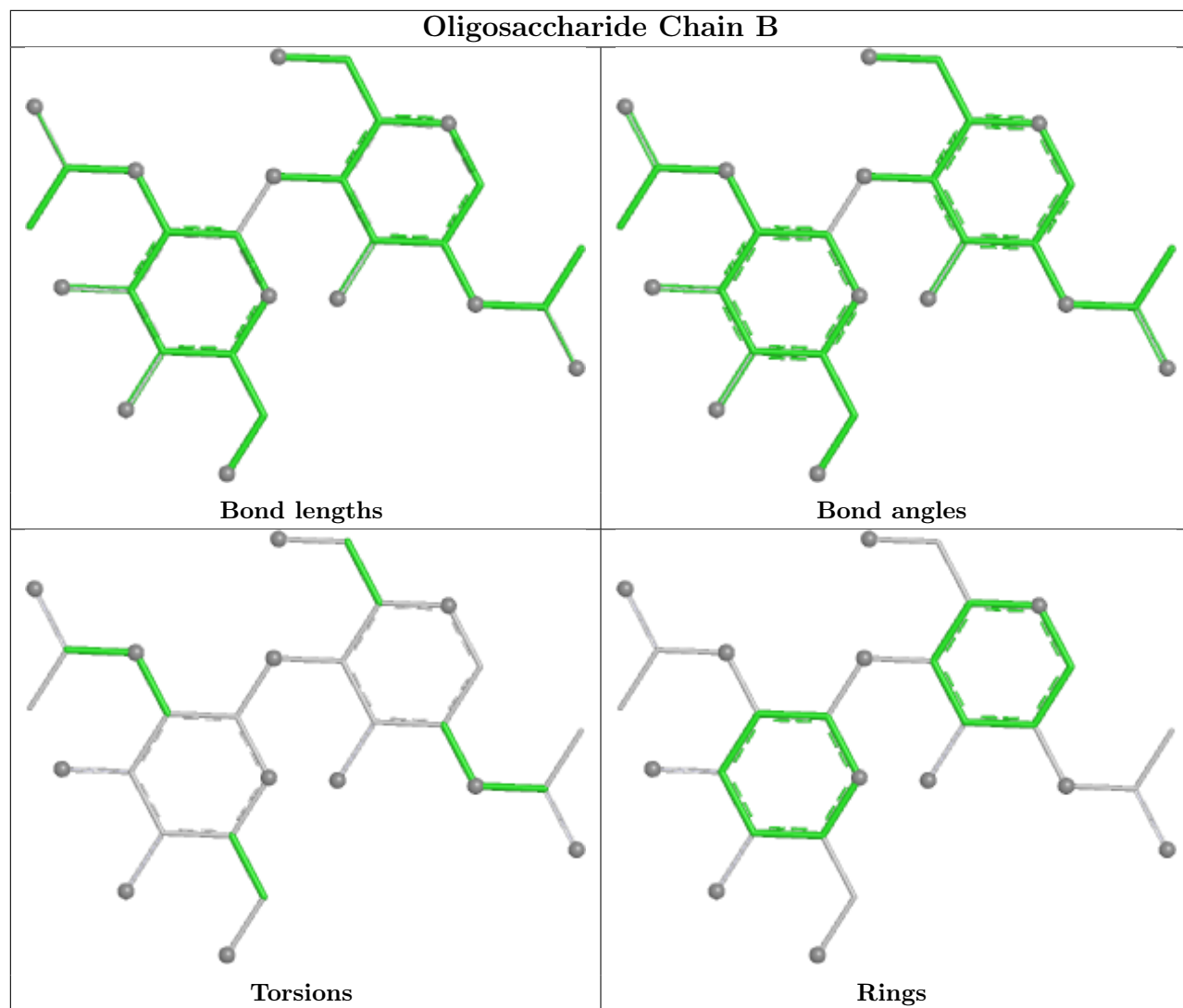
All (6) torsion outliers are listed below:

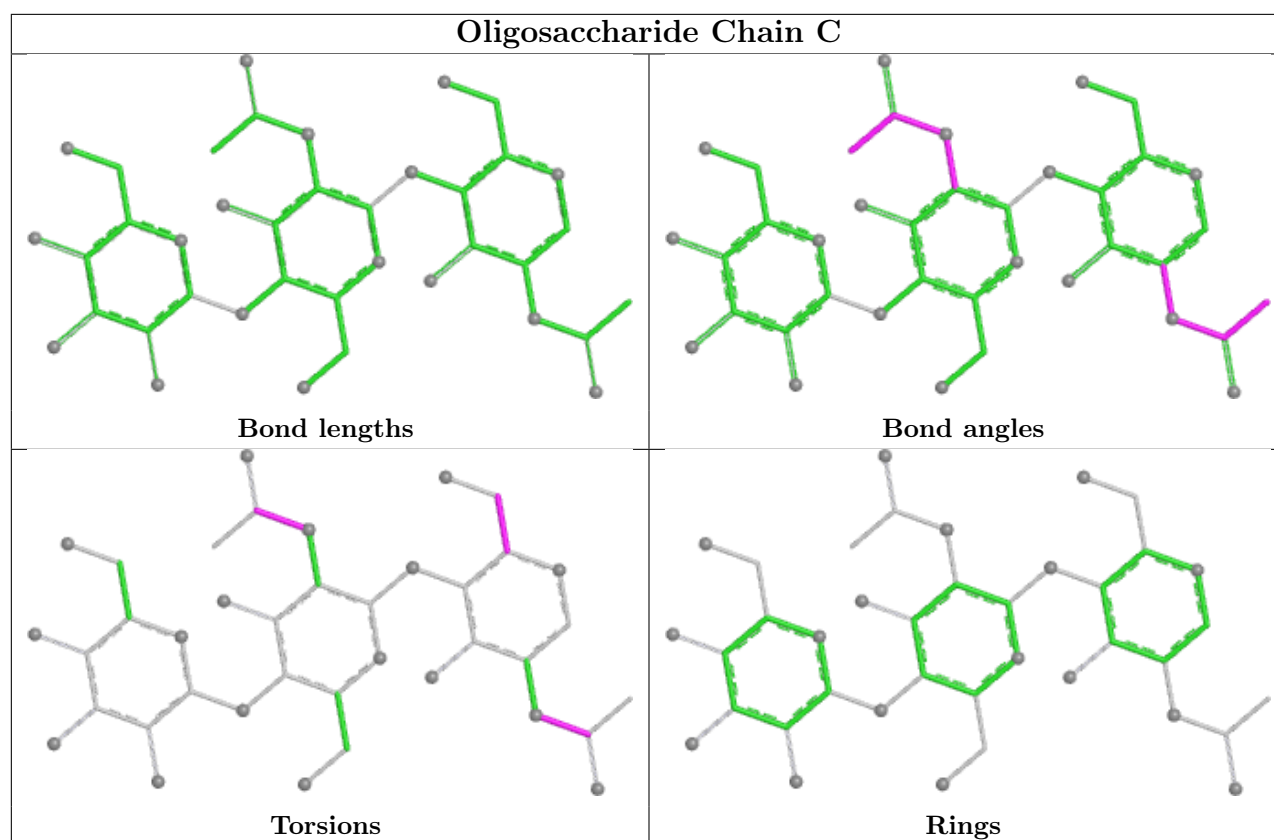
Mol	Chain	Res	Type	Atoms
3	C	1	NAG	O5-C5-C6-O6
3	C	1	NAG	C4-C5-C6-O6
3	C	1	NAG	C8-C7-N2-C2
3	C	1	NAG	O7-C7-N2-C2
3	C	2	NAG	C8-C7-N2-C2
3	C	2	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





## 5.6 Ligand geometry [i](#)

3 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	$\# Z  > 2$
4	NAG	A	2508	-	14,14,15	0.40	0	17,19,21	1.17	1 (5%)
4	NAG	A	2501	1	14,14,15	0.59	0	17,19,21	0.94	1 (5%)
4	NAG	A	2507	1	14,14,15	0.44	0	17,19,21	1.31	3 (17%)

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
4	NAG	A	2508	-	-	3/6/23/26	0/1/1/1
4	NAG	A	2501	1	-	0/6/23/26	0/1/1/1
4	NAG	A	2507	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2507	NAG	C1-O5-C5	3.07	116.30	112.19
4	A	2507	NAG	C2-N2-C7	2.50	126.25	122.90
4	A	2508	NAG	C4-C3-C2	-2.44	107.44	111.02
4	A	2507	NAG	O5-C1-C2	-2.43	107.53	111.29
4	A	2501	NAG	O5-C1-C2	2.05	114.46	111.29

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2508	NAG	C1-C2-N2-C7
4	A	2507	NAG	O5-C5-C6-O6
4	A	2507	NAG	C4-C5-C6-O6
4	A	2508	NAG	O5-C5-C6-O6
4	A	2508	NAG	C4-C5-C6-O6
4	A	2507	NAG	C1-C2-N2-C7
4	A	2507	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2508	NAG	8	0

## 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.

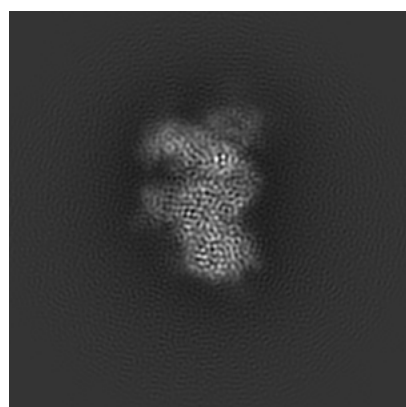
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20815. These allow visual inspection of the internal detail of the map and identification of artifacts.

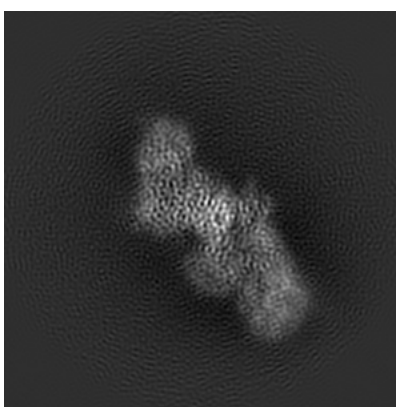
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

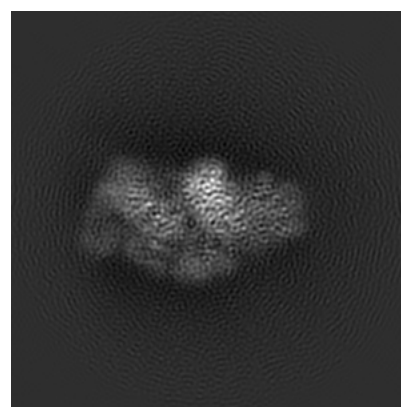
#### 6.1.1 Primary map



X



Y

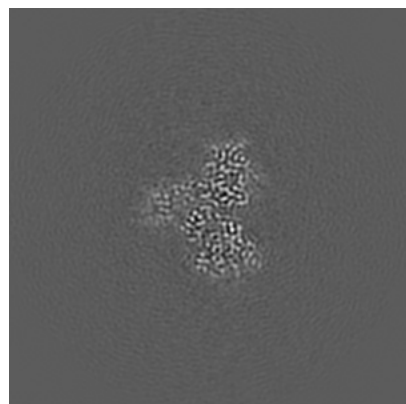


Z

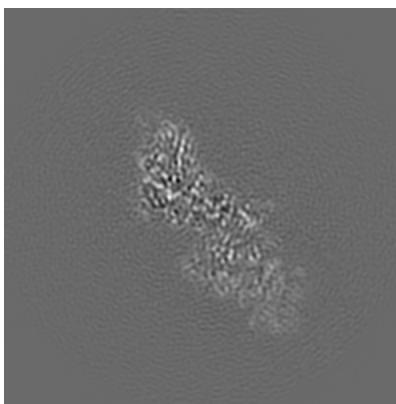
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

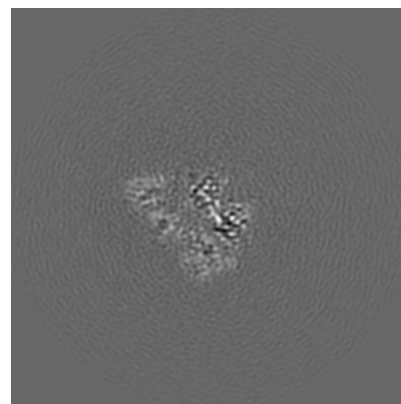
#### 6.2.1 Primary map



X Index: 140



Y Index: 140

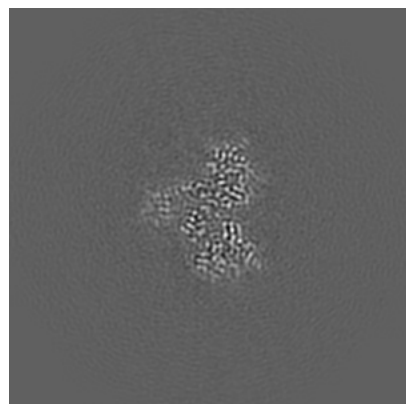


Z Index: 140

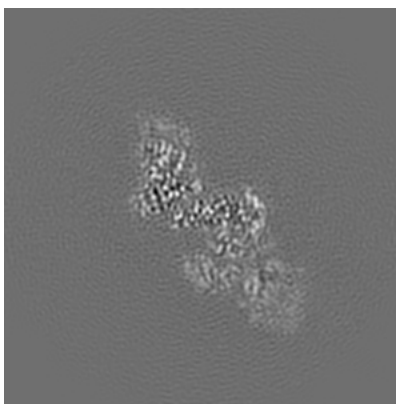
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

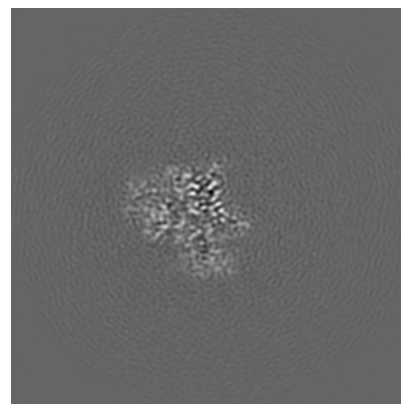
### 6.3.1 Primary map



X Index: 141



Y Index: 145

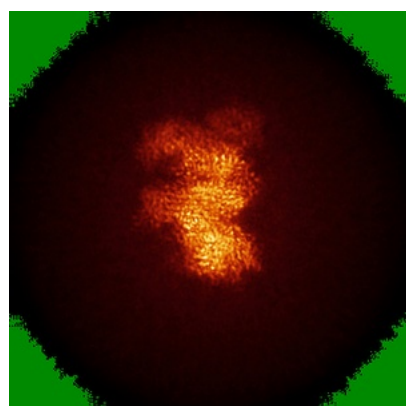


Z Index: 149

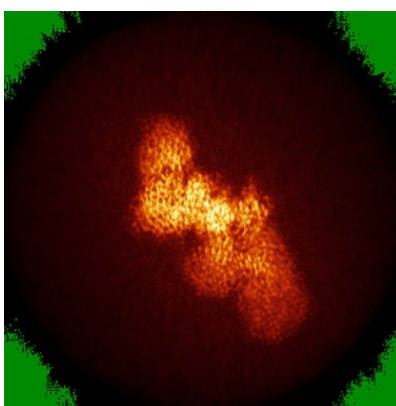
The images above show the largest variance slices of the map in three orthogonal directions.

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

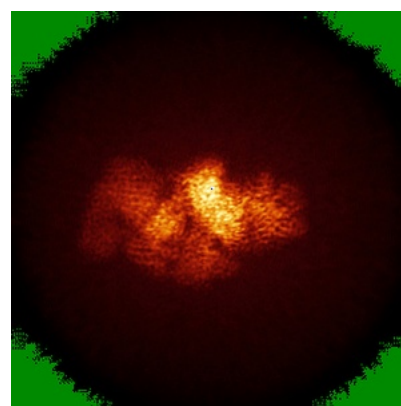
### 6.4.1 Primary map



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 0.0145. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

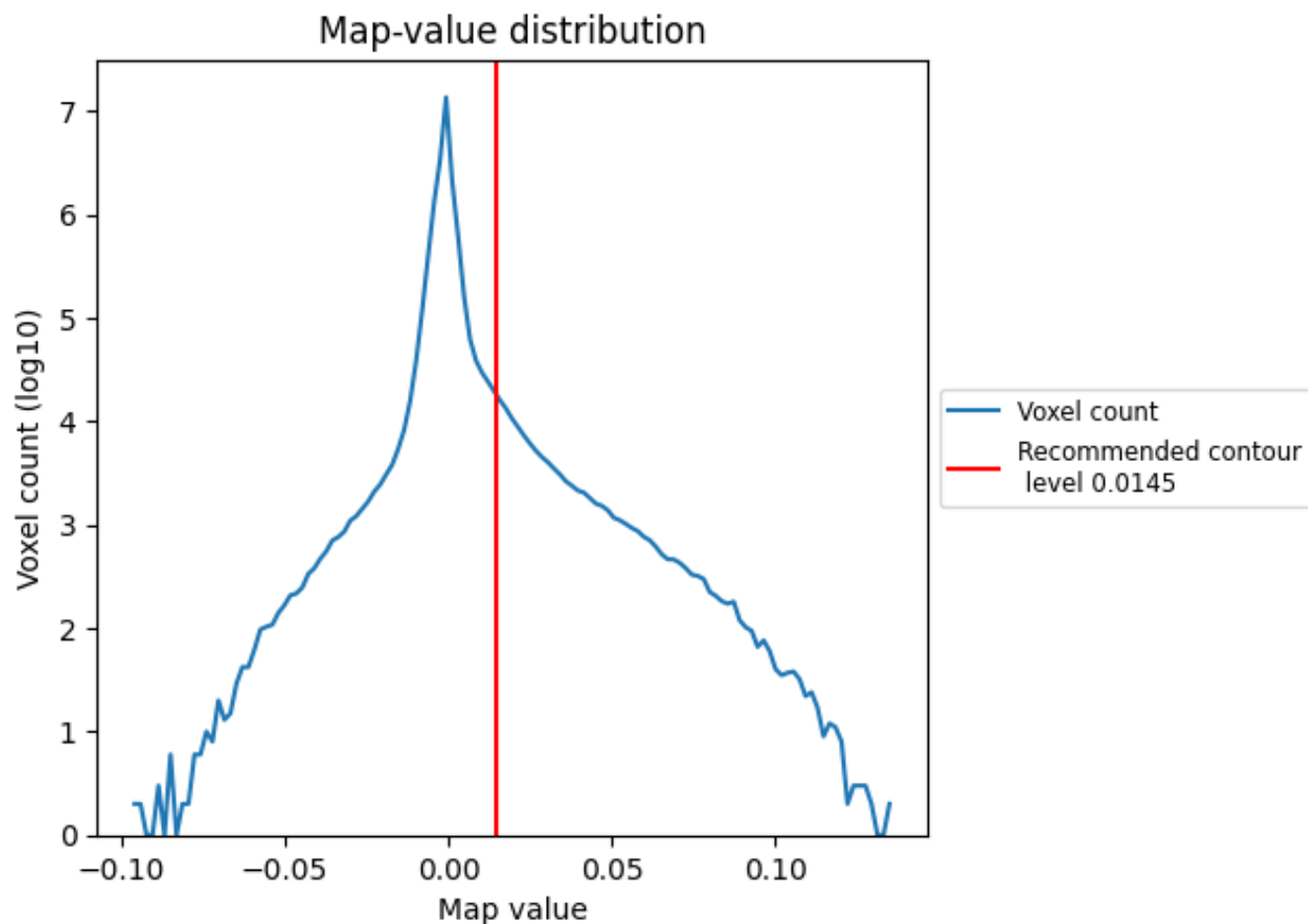
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

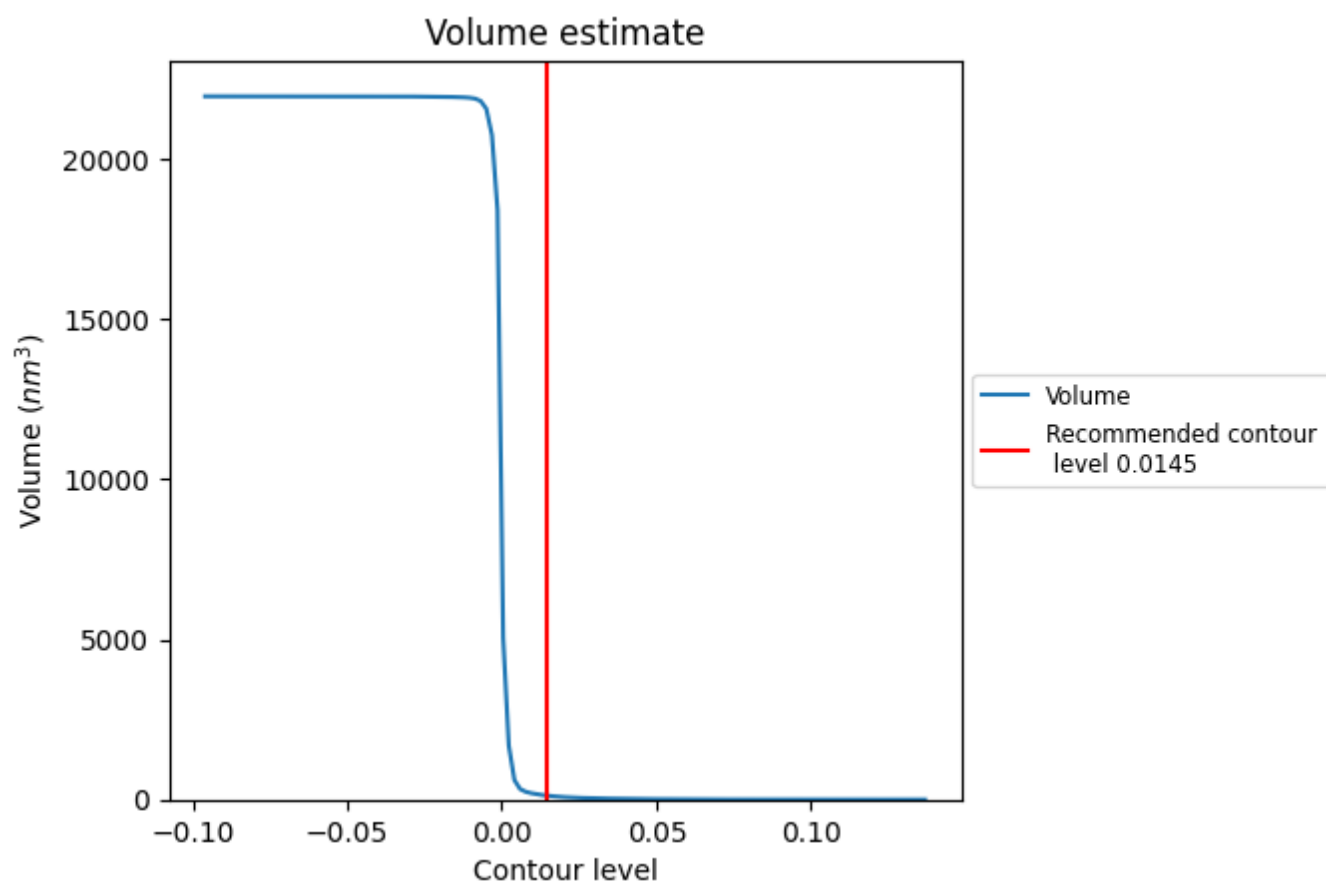
This section contains the results of statistical analysis of the map.

### 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.

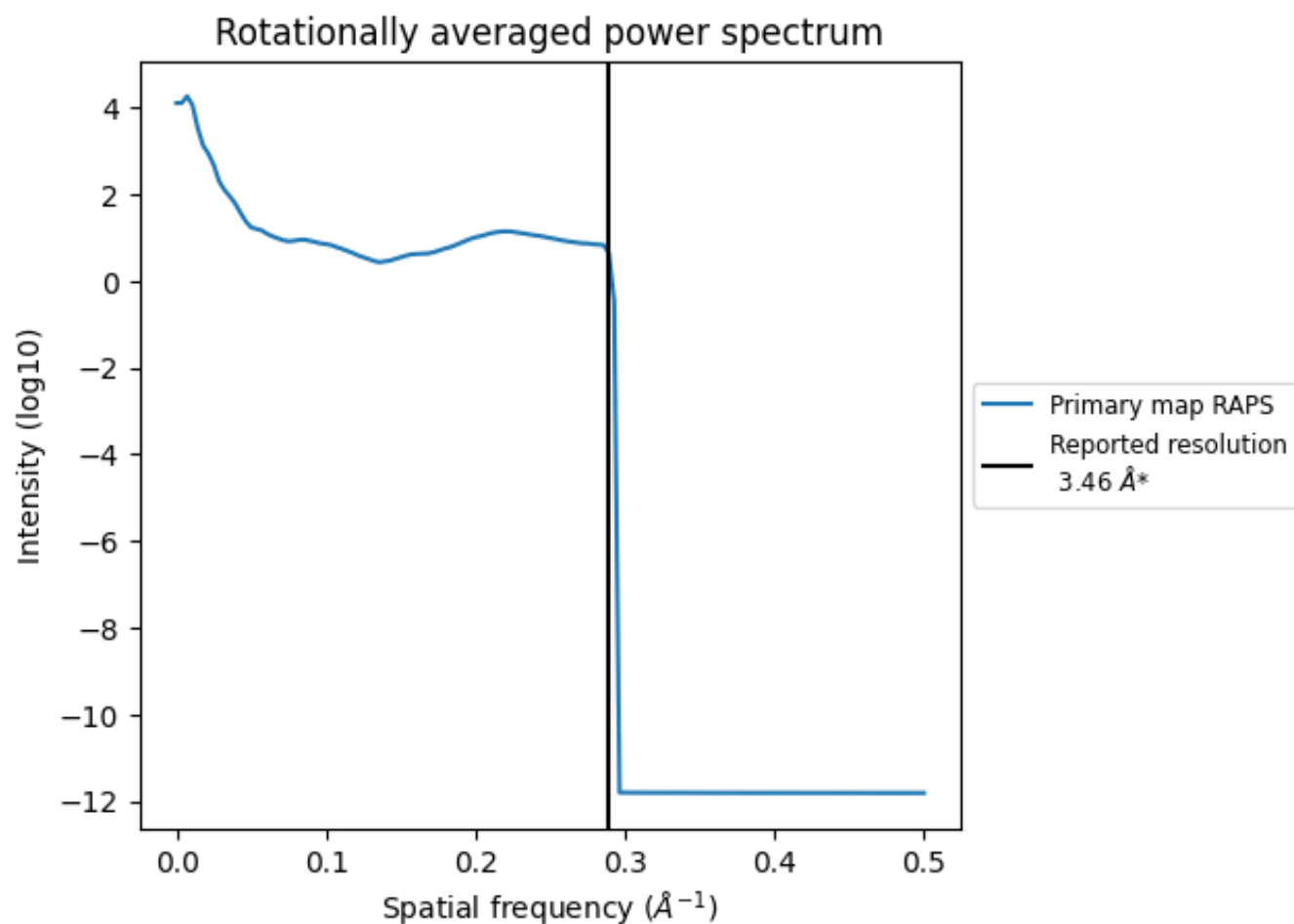
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 124 nm<sup>3</sup>; this corresponds to an approximate mass of 112 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 0.289  $\text{\AA}^{-1}$

## 8 Fourier-Shell correlation ⓘ

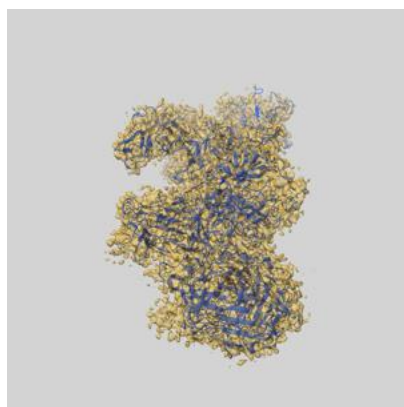
This section was not generated. No FSC curve or half-maps provided.



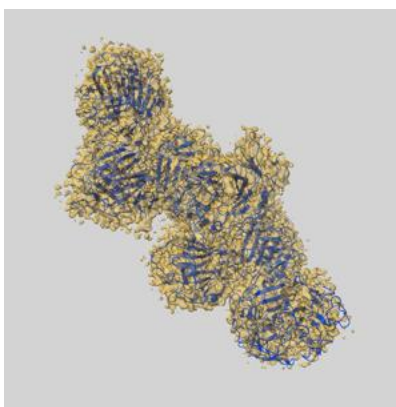
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-20815 and PDB model 6UM1. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

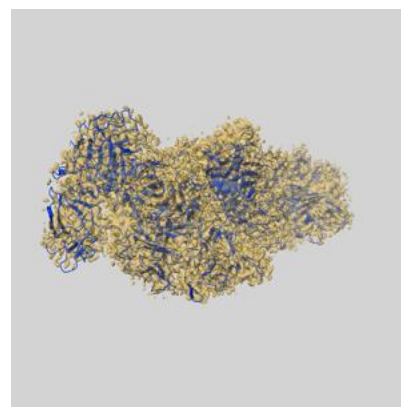
### 9.1 Map-model overlay [i](#)



X



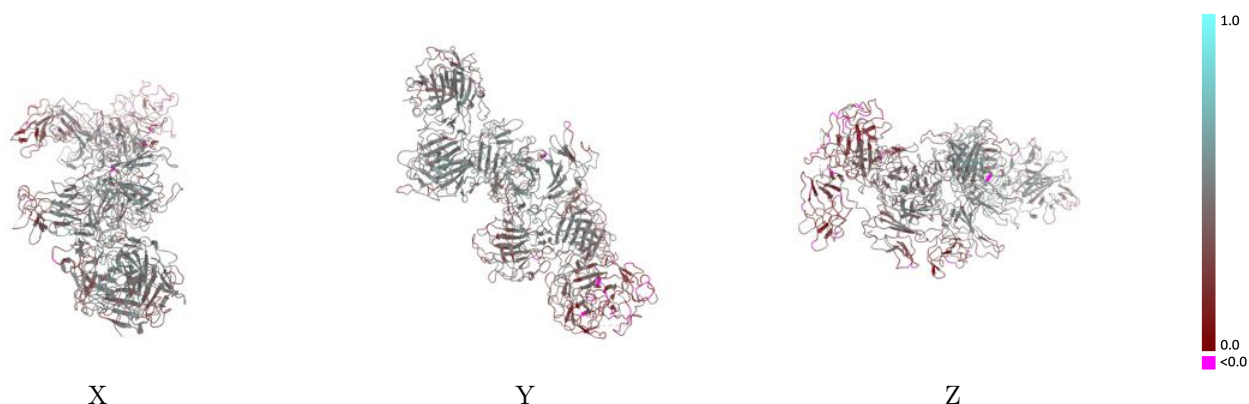
Y



Z

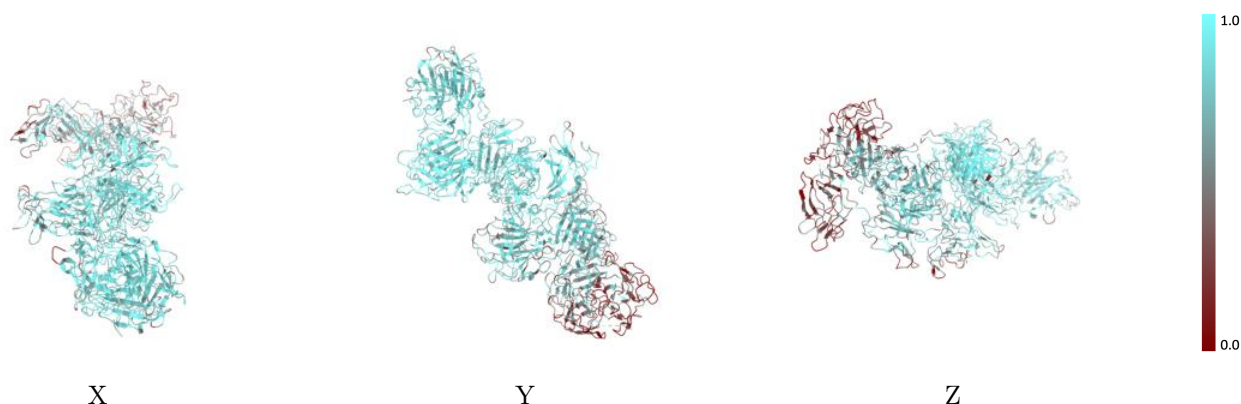
The images above show the 3D surface view of the map at the recommended contour level 0.0145 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



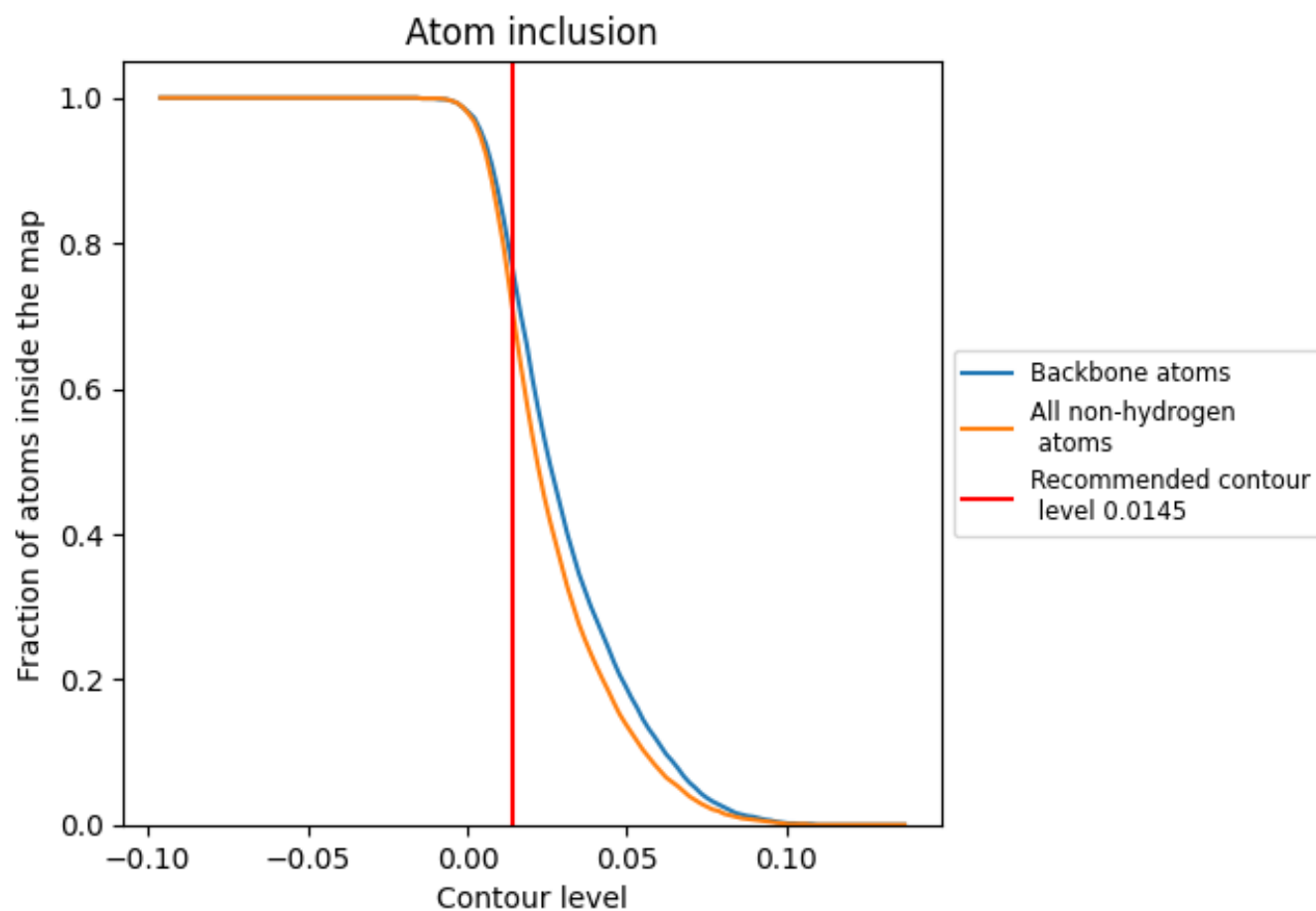
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0145).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 70% 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 (0.0145) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6990	<div></div> 0.4010
A	<div></div> 0.7000	<div></div> 0.4010
B	<div></div> 0.5710	<div></div> 0.3220
C	<div></div> 0.5900	<div></div> 0.3430

