



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 12:42 AM UTC

PDB ID : 7AIN / pdb_00007ain
EMDB ID : EMD-11799
Title : Structure of Human Potassium Chloride Transporter KCC3 S45D/T940D/T997D in NaCl (Reference Map)
Authors : Chi, G.; Man, H.; Ebenhoch, R.; Reggiano, G.; Pike, A.C.W.; Wang, D.; McKinley, G.; Mukhopadhyay, S.M.M.; MacLean, E.M.; Chalk, R.; Moreau, C.; Snee, M.; Bohstedt, T.; Singh, N.K.; Abrusci, P.; Arrowsmith, C.H.; Bountra, C.; Edwards, A.M.; Marsden, B.D.; Burgess-Brown, N.A.; DiMaio, F.; Duerr, K.L.; Structural Genomics Consortium (SGC)
Deposited on : 2020-09-28
Resolution : 3.20 Å(reported)
Based on initial model : 6Y5R

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
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)

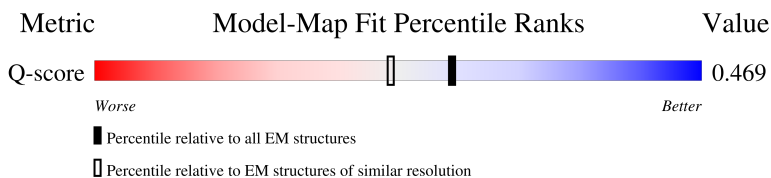
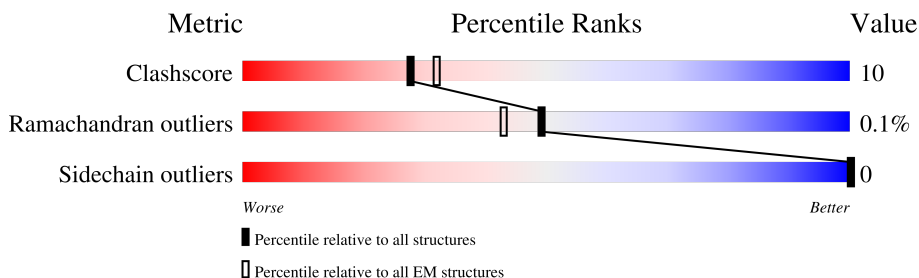
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 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	15020 (2.70 - 3.70)

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	1099	 64% 18% 18%
1	B	1099	 64% 18% 18%

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
2	C	3	33% 67%
2	G	3	67% 33%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13904 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 Isoform 2 of Solute carrier family 12 member 6.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	B	904	Total	C	N	O	P	S	0	0
			6899	4480	1147	1223	2	47		
1	A	904	Total	C	N	O	P	S	0	0
			6899	4480	1147	1223	2	47		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	45	ASP	SER	engineered mutation	UNP Q9UHW9
B	940	ASP	THR	engineered mutation	UNP Q9UHW9
B	997	ASP	THR	engineered mutation	UNP Q9UHW9
A	45	ASP	SER	engineered mutation	UNP Q9UHW9
A	940	ASP	THR	engineered mutation	UNP Q9UHW9
A	997	ASP	THR	engineered mutation	UNP Q9UHW9

- Molecule 2 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.



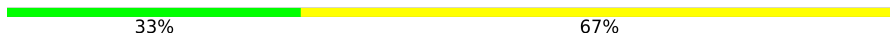
Mol	Chain	Residues	Atoms				AltConf	Trace
2	C	3	Total	C	N	O	0	0
			39	22	2	15		
2	G	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

-
- | Frequency | Percentage |
|-----------|------------|
| Daily | 64% |
| Weekly | 18% |
| Monthly | 18% |





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	408039	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.789	Depositor
Minimum map value	-1.207	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.044	Depositor
Recommended contour level	0.17	Depositor
Map size ($\text{\AA}$)	260.40002, 260.40002, 260.40002	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.6510001, 0.6510001, 0.6510001	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, SEP, 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	0.43	0/7033	0.52	0/9561
1	B	0.44	0/7033	0.53	0/9561
All	All	0.44	0/14066	0.52	0/19122

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6899	0	6910	138	0
1	B	6899	0	6910	138	0
2	C	39	0	34	0	0
2	G	39	0	34	0	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
All	All	13904	0	13914	270	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 (270) 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:873:MET:HG2	1:B:909:LEU:HD11	1.57	0.87
1:A:873:MET:HG2	1:A:909:LEU:HD11	1.57	0.85
1:A:862:VAL:HG21	1:A:876:PRO:HG3	1.76	0.67
1:B:600:LEU:HD22	1:B:640:CYS:HB2	1.77	0.66
1:A:600:LEU:HD22	1:A:640:CYS:HB2	1.77	0.66
1:A:813:TRP:HB2	1:A:822:TRP:HB3	1.77	0.66
1:B:813:TRP:HB2	1:B:822:TRP:HB3	1.77	0.66
1:B:862:VAL:HG21	1:B:876:PRO:HG3	1.76	0.65
1:A:826:ILE:HD12	1:A:1067:ASP:HB3	1.79	0.64
1:B:826:ILE:HD12	1:B:1067:ASP:HB3	1.79	0.64
1:A:675:TRP:NE1	1:A:685:SER:OG	2.31	0.64
1:B:155:LEU:HD21	1:B:239:ILE:HG12	1.81	0.63
1:B:422:ASN:OD1	1:B:423:HIS:N	2.31	0.63
1:A:163:THR:HG23	1:A:437:LEU:HD11	1.82	0.62
1:A:240:PHE:HA	1:A:244:ILE:HD12	1.81	0.62
1:A:422:ASN:OD1	1:A:423:HIS:N	2.31	0.62
1:B:675:TRP:NE1	1:B:685:SER:OG	2.31	0.62
1:B:865:ILE:O	1:B:867:HIS:N	2.32	0.62
1:A:865:ILE:O	1:A:867:HIS:N	2.32	0.62
1:B:163:THR:HG23	1:B:437:LEU:HD11	1.82	0.62
1:B:240:PHE:HA	1:B:244:ILE:HD12	1.81	0.62
1:A:155:LEU:HD21	1:A:239:ILE:HG12	1.81	0.62
1:B:840:LEU:HD21	1:B:871:MET:HE3	1.82	0.61
1:B:877:PHE:CE2	1:B:916:LEU:HD13	2.36	0.61
1:A:877:PHE:CE2	1:A:916:LEU:HD13	2.36	0.60
1:B:226:THR:HG22	1:B:575:THR:HG23	1.82	0.60
1:B:1040:ASN:ND2	1:B:1076:VAL:O	2.29	0.60
1:B:721:LYS:O	1:B:722:HIS:ND1	2.35	0.60
1:A:917:ARG:HG3	1:A:917:ARG:HH11	1.67	0.60
1:A:840:LEU:HD21	1:A:871:MET:HE3	1.82	0.60
1:A:226:THR:HG22	1:A:575:THR:HG23	1.82	0.59
1:B:917:ARG:HG3	1:B:917:ARG:HH11	1.66	0.59
1:A:1040:ASN:ND2	1:A:1076:VAL:O	2.29	0.59
1:B:698:PRO:O	1:A:680:ARG:NH1	2.36	0.59
1:A:721:LYS:O	1:A:722:HIS:ND1	2.35	0.59
1:B:550:ASP:HB3	1:B:552:ILE:HG13	1.85	0.59
1:B:804:THR:HG22	1:B:838:ALA:HB3	1.85	0.59
1:A:550:ASP:HB3	1:A:552:ILE:HG13	1.85	0.59
1:A:1063:ASN:HB3	1:A:1065:GLU:HG2	1.85	0.58
1:B:341:LYS:HA	1:B:354:LEU:HD11	1.85	0.58
1:B:601:MET:HE1	1:B:656:MET:HG2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:SER:HB2	1:B:466:PRO:HB3	1.85	0.58
1:B:877:PHE:CD2	1:B:916:LEU:HD13	2.38	0.58
1:A:934:ALA:HB2	1:A:1032:ARG:HD2	1.85	0.58
1:A:601:MET:HE1	1:A:656:MET:HG2	1.85	0.58
1:A:877:PHE:CD2	1:A:916:LEU:HD13	2.38	0.58
1:A:804:THR:HG22	1:A:838:ALA:HB3	1.85	0.58
1:A:755:GLU:OE1	1:A:755:GLU:N	2.37	0.58
1:B:132:PRO:O	1:B:134:MET:HG3	2.05	0.57
1:A:132:PRO:O	1:A:134:MET:HG3	2.04	0.57
1:A:188:SER:HB2	1:A:466:PRO:HB3	1.85	0.57
1:A:341:LYS:HA	1:A:354:LEU:HD11	1.85	0.57
1:B:755:GLU:N	1:B:755:GLU:OE1	2.37	0.56
1:B:1063:ASN:HB3	1:B:1065:GLU:HG2	1.85	0.56
1:B:364:PHE:HA	1:B:367:ALA:HB2	1.87	0.56
1:B:674:GLU:OE1	1:A:736:LYS:HB2	2.04	0.56
1:B:934:ALA:HB2	1:B:1032:ARG:HD2	1.85	0.56
1:B:443:PRO:HB3	1:B:595:LEU:HD22	1.88	0.55
1:A:364:PHE:HA	1:A:367:ALA:HB2	1.87	0.55
1:B:861:ASP:OD1	1:B:893:PHE:HE1	1.90	0.55
1:A:861:ASP:OD1	1:A:893:PHE:HE1	1.90	0.55
1:A:825:PHE:HD2	1:A:1071:MET:HE1	1.72	0.55
1:B:825:PHE:HD2	1:B:1071:MET:HE1	1.72	0.54
1:B:645:PHE:CD1	1:A:653:ILE:HD11	2.42	0.54
1:A:376:ASN:OD1	1:A:415:SER:N	2.33	0.54
1:A:443:PRO:HB3	1:A:595:LEU:HD22	1.88	0.54
1:B:519:TRP:HA	1:B:522:VAL:HG12	1.90	0.54
1:A:231:MET:HE1	1:A:443:PRO:HG3	1.90	0.54
1:A:162:GLY:HA3	1:A:496:LEU:HB3	1.89	0.53
1:B:160:VAL:HG22	1:B:437:LEU:HD22	1.90	0.53
1:B:376:ASN:OD1	1:B:415:SER:N	2.33	0.53
1:A:640:CYS:HA	1:A:643:LEU:HB2	1.91	0.53
1:B:231:MET:HE1	1:B:443:PRO:HG3	1.90	0.53
1:A:813:TRP:H	1:A:813:TRP:CD1	2.26	0.53
1:B:640:CYS:HA	1:B:643:LEU:HB2	1.91	0.53
1:A:519:TRP:HA	1:A:522:VAL:HG12	1.90	0.53
1:B:568:PRO:HB2	1:B:571:ALA:HB3	1.91	0.53
1:B:813:TRP:CD1	1:B:813:TRP:H	2.26	0.53
1:A:568:PRO:HB2	1:A:571:ALA:HB3	1.91	0.53
1:A:160:VAL:HG22	1:A:437:LEU:HD22	1.90	0.52
1:A:199:GLY:HA3	1:A:203:PHE:CD2	2.45	0.52
1:B:162:GLY:HA3	1:B:496:LEU:HB3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:790:HIS:HE2	1:A:781:ALA:HA	1.74	0.52
1:A:939:ARG:HH11	1:A:943:MET:HG3	1.76	0.51
1:B:206:SER:OG	1:B:207:ARG:NH1	2.43	0.51
1:A:206:SER:OG	1:A:207:ARG:NH1	2.43	0.51
1:B:789:SER:O	1:B:793:GLN:NE2	2.30	0.51
1:A:928:HIS:HB3	1:A:930:SEP:O3P	2.11	0.51
1:B:825:PHE:CD2	1:B:1071:MET:HE1	2.47	0.50
1:B:917:ARG:HG3	1:B:917:ARG:NH1	2.27	0.50
1:B:1092:SER:O	1:B:1092:SER:OG	2.29	0.50
1:B:939:ARG:HH11	1:B:943:MET:HG3	1.75	0.50
1:A:159:TRP:O	1:A:163:THR:HG22	2.11	0.50
1:A:789:SER:O	1:A:793:GLN:NE2	2.30	0.50
1:B:199:GLY:HA3	1:B:203:PHE:CD2	2.45	0.50
1:A:825:PHE:CD2	1:A:1071:MET:HE1	2.47	0.50
1:B:862:VAL:HG22	1:B:1054:LEU:HD22	1.94	0.50
1:B:928:HIS:HB3	1:B:930:SEP:O3P	2.11	0.49
1:A:862:VAL:HG22	1:A:1054:LEU:HD22	1.94	0.49
1:B:847:PHE:O	1:B:1083:ARG:HD3	2.12	0.49
1:B:763:ILE:HD12	1:B:778:LEU:HD21	1.94	0.49
1:A:763:ILE:HD12	1:A:778:LEU:HD21	1.94	0.49
1:B:258:LYS:O	1:B:261:ALA:N	2.46	0.48
1:B:674:GLU:HG3	1:B:675:TRP:CD1	2.49	0.48
1:A:934:ALA:HB2	1:A:1032:ARG:CD	2.43	0.48
1:B:159:TRP:O	1:B:163:THR:HG22	2.11	0.48
1:B:934:ALA:HB2	1:B:1032:ARG:CD	2.43	0.48
1:A:847:PHE:O	1:A:1083:ARG:HD3	2.12	0.48
1:B:237:ILE:HD12	1:B:270:TYR:HB3	1.95	0.48
1:B:1074:LEU:O	1:B:1078:THR:HG22	2.14	0.48
1:A:231:MET:HG3	1:A:588:LEU:HD22	1.95	0.48
1:A:249:ALA:HB1	1:A:266:ASN:ND2	2.29	0.48
1:A:917:ARG:HG3	1:A:917:ARG:NH1	2.27	0.48
1:A:674:GLU:HG3	1:A:675:TRP:CD1	2.49	0.48
1:B:231:MET:HG3	1:B:588:LEU:HD22	1.95	0.48
1:A:845:ILE:HA	1:A:848:PHE:CD1	2.49	0.48
1:B:845:ILE:HA	1:B:848:PHE:CD1	2.49	0.48
1:B:249:ALA:HB1	1:B:266:ASN:ND2	2.29	0.48
1:A:237:ILE:HD12	1:A:270:TYR:HB3	1.95	0.48
1:A:495:VAL:HG13	1:A:504:VAL:HG21	1.96	0.47
1:B:495:VAL:HG13	1:B:504:VAL:HG21	1.96	0.47
1:A:375:HIS:ND1	1:A:417:VAL:HG11	2.29	0.47
1:B:147:ASN:HD22	1:B:450:ALA:HB1	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:LEU:HD11	1:A:239:ILE:HG13	1.95	0.47
1:A:912:PHE:CE1	1:A:916:LEU:HD11	2.50	0.47
1:B:155:LEU:HD11	1:B:239:ILE:HG13	1.95	0.47
1:A:205:ILE:HG23	1:A:209:LEU:HD12	1.97	0.47
1:A:511:GLY:C	1:A:518:PRO:HB3	2.40	0.47
1:B:205:ILE:HG23	1:B:209:LEU:HD12	1.97	0.47
1:B:375:HIS:ND1	1:B:417:VAL:HG11	2.29	0.47
1:B:864:TRP:HB2	1:B:873:MET:HE1	1.96	0.47
1:B:399:ASN:OD1	1:B:497:ARG:NH1	2.48	0.47
1:A:147:ASN:HD22	1:A:450:ALA:HB1	1.80	0.47
1:A:241:LEU:HD21	1:A:266:ASN:HB3	1.97	0.47
1:A:864:TRP:HB2	1:A:873:MET:HE1	1.95	0.47
1:B:219:LEU:HD12	1:B:662:ILE:HG21	1.97	0.46
1:A:1074:LEU:O	1:A:1078:THR:HG22	2.14	0.46
1:A:219:LEU:HD12	1:A:662:ILE:HG21	1.97	0.46
1:A:399:ASN:OD1	1:A:497:ARG:NH1	2.48	0.46
1:B:841:VAL:HB	1:B:1086:LEU:HB2	1.98	0.46
1:B:912:PHE:CE1	1:B:916:LEU:HD11	2.50	0.46
1:A:541:PRO:HB2	1:A:568:PRO:CG	2.46	0.46
1:A:841:VAL:HB	1:A:1086:LEU:HB2	1.98	0.46
1:B:158:THR:CG2	1:B:510:VAL:HB	2.45	0.46
1:B:614:THR:HG21	1:B:630:TRP:HB3	1.98	0.46
1:B:241:LEU:HD21	1:B:266:ASN:HB3	1.97	0.46
1:B:541:PRO:HB2	1:B:568:PRO:CG	2.46	0.46
1:B:519:TRP:O	1:B:523:ILE:HG12	2.16	0.46
1:B:845:ILE:HA	1:B:848:PHE:HD1	1.81	0.46
1:A:158:THR:CG2	1:A:510:VAL:HB	2.45	0.46
1:A:845:ILE:HA	1:A:848:PHE:HD1	1.81	0.46
1:B:407:ILE:HG22	1:B:494:VAL:HG23	1.98	0.46
1:B:511:GLY:C	1:B:518:PRO:HB3	2.40	0.46
1:A:177:CYS:HA	1:A:180:THR:HG22	1.98	0.46
1:A:766:LEU:O	1:A:770:GLU:HG2	2.17	0.45
1:B:177:CYS:HA	1:B:180:THR:HG22	1.98	0.45
1:A:519:TRP:O	1:A:523:ILE:HG12	2.16	0.45
1:B:725:LEU:HG	1:B:845:ILE:HD13	1.98	0.45
1:A:725:LEU:HG	1:A:845:ILE:HD13	1.99	0.45
1:A:830:ARG:NH1	1:A:1061:PRO:O	2.43	0.45
1:A:600:LEU:HB3	1:A:644:MET:HG2	1.98	0.45
1:A:320:HIS:O	1:A:322:PRO:HD3	2.17	0.45
1:A:407:ILE:HG22	1:A:494:VAL:HG23	1.98	0.45
1:B:733:LYS:HE3	1:B:736:LYS:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:ALA:HB2	1:A:473:ILE:HD12	1.98	0.45
1:A:268:ARG:NE	1:A:584:LEU:O	2.50	0.45
1:B:184:ALA:HB2	1:B:473:ILE:CD1	2.47	0.44
1:B:733:LYS:HA	1:B:733:LYS:HD2	1.78	0.44
1:B:766:LEU:O	1:B:770:GLU:HG2	2.17	0.44
1:A:733:LYS:HD2	1:A:733:LYS:HA	1.78	0.44
1:B:184:ALA:HB2	1:B:473:ILE:HD12	1.98	0.44
1:B:933:SER:OG	1:B:1032:ARG:NH1	2.50	0.44
1:A:692:LEU:HD23	1:A:692:LEU:HA	1.80	0.44
1:A:807:MET:HB3	1:A:807:MET:HE2	1.64	0.44
1:A:933:SER:OG	1:A:1032:ARG:NH1	2.50	0.44
1:B:138:MET:HE3	1:B:138:MET:HB2	1.77	0.44
1:B:144:CYS:O	1:B:148:ILE:HG13	2.17	0.44
1:B:320:HIS:O	1:B:322:PRO:HD3	2.17	0.44
1:B:553:ILE:HG23	1:B:665:TYR:HE2	1.82	0.44
1:B:110:GLN:O	1:B:200:GLY:N	2.50	0.44
1:B:538:THR:O	1:B:542:ARG:HG3	2.18	0.44
1:A:733:LYS:HE3	1:A:736:LYS:O	2.17	0.44
1:B:499:LYS:NZ	1:B:589:ASP:OD2	2.49	0.44
1:A:144:CYS:O	1:A:148:ILE:HG13	2.17	0.44
1:A:330:THR:HG23	1:A:426:VAL:HB	2.00	0.44
1:A:614:THR:HG21	1:A:630:TRP:HB3	1.98	0.44
1:B:131:THR:HB	1:B:132:PRO:HD3	2.00	0.44
1:B:192:THR:O	1:B:624:ARG:NH2	2.34	0.44
1:B:404:GLY:O	1:B:427:LEU:HD22	2.18	0.44
1:B:600:LEU:CD2	1:B:640:CYS:HB2	2.46	0.44
1:B:644:MET:HE2	1:B:644:MET:HB3	1.75	0.44
1:A:131:THR:HB	1:A:132:PRO:HD3	2.00	0.44
1:A:600:LEU:CD2	1:A:640:CYS:HB2	2.46	0.44
1:A:184:ALA:HB2	1:A:473:ILE:CD1	2.47	0.44
1:A:330:THR:HG22	1:A:407:ILE:O	2.17	0.44
1:A:615:LEU:HD21	1:A:630:TRP:CE2	2.53	0.44
1:A:437:LEU:HD23	1:A:437:LEU:HA	1.77	0.43
1:B:600:LEU:HB3	1:B:644:MET:HG2	1.98	0.43
1:B:615:LEU:HD21	1:B:630:TRP:CE2	2.53	0.43
1:B:330:THR:HG22	1:B:407:ILE:O	2.17	0.43
1:A:791:LEU:HD22	1:A:795:CYS:SG	2.59	0.43
1:B:268:ARG:NE	1:B:584:LEU:O	2.50	0.43
1:B:330:THR:HG23	1:B:426:VAL:HB	2.00	0.43
1:A:110:GLN:O	1:A:200:GLY:N	2.50	0.43
1:B:807:MET:HB3	1:B:807:MET:HE2	1.64	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:873:MET:O	1:B:876:PRO:HD2	2.19	0.43
1:B:673:LYS:HA	1:B:673:LYS:HD2	1.73	0.43
1:A:771:LYS:HE2	1:A:771:LYS:HB2	1.82	0.43
1:A:553:ILE:HG23	1:A:665:TYR:HE2	1.82	0.43
1:A:873:MET:O	1:A:876:PRO:HD2	2.19	0.43
1:B:279:VAL:HG21	1:B:576:ALA:HB2	2.01	0.42
1:B:771:LYS:HE2	1:B:771:LYS:HB2	1.82	0.42
1:B:930:SEP:O	1:B:1035:THR:HG21	2.19	0.42
1:A:279:VAL:HG21	1:A:576:ALA:HB2	2.01	0.42
1:B:595:LEU:HD23	1:B:599:PHE:HD2	1.84	0.42
1:B:1028:SER:O	1:B:1032:ARG:HG2	2.19	0.42
1:A:930:SEP:O	1:A:1035:THR:HG21	2.19	0.42
1:A:538:THR:O	1:A:542:ARG:HG3	2.18	0.42
1:B:353:LYS:O	1:B:356:GLY:N	2.52	0.42
1:A:404:GLY:O	1:A:427:LEU:HD22	2.18	0.42
1:A:600:LEU:HB3	1:A:644:MET:CG	2.50	0.42
1:A:673:LYS:HD2	1:A:673:LYS:HA	1.73	0.42
1:B:136:THR:HA	1:B:464:SER:HB2	2.02	0.42
1:A:1028:SER:O	1:A:1032:ARG:HG2	2.19	0.42
1:B:791:LEU:HD22	1:B:795:CYS:SG	2.59	0.42
1:A:353:LYS:O	1:A:356:GLY:N	2.52	0.42
1:A:401:LEU:HB2	1:A:428:VAL:HG21	2.02	0.42
1:A:138:MET:HE3	1:A:138:MET:HB2	1.77	0.42
1:B:645:PHE:CG	1:A:653:ILE:HD11	2.55	0.41
1:A:472:ALA:O	1:A:476:THR:HG23	2.21	0.41
1:A:595:LEU:HD23	1:A:599:PHE:HD2	1.84	0.41
1:B:401:LEU:HB2	1:B:428:VAL:HG21	2.02	0.41
1:B:472:ALA:O	1:B:476:THR:HG23	2.21	0.41
1:B:861:ASP:OD1	1:B:893:PHE:CE1	2.73	0.41
1:A:409:LYS:HE2	1:A:492:GLU:OE2	2.20	0.41
1:B:247:ARG:HA	1:B:247:ARG:HD2	1.92	0.41
1:B:600:LEU:HB3	1:B:644:MET:CG	2.50	0.41
1:A:820:ARG:HD2	1:A:820:ARG:HA	1.84	0.41
1:B:249:ALA:HB1	1:B:266:ASN:CG	2.46	0.41
1:B:392:ILE:HD12	1:B:489:ALA:HB2	2.02	0.41
1:A:249:ALA:HB1	1:A:266:ASN:CG	2.46	0.41
1:B:692:LEU:HD23	1:B:692:LEU:HA	1.80	0.41
1:A:192:THR:O	1:A:624:ARG:NH2	2.34	0.41
1:A:219:LEU:HA	1:A:219:LEU:HD23	1.86	0.41
1:A:632:LEU:HD23	1:A:632:LEU:HA	1.83	0.41
1:A:861:ASP:OD1	1:A:893:PHE:CE1	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1040:ASN:O	1:B:1044:VAL:HG22	2.21	0.41
1:B:1071:MET:HE2	1:B:1071:MET:HB2	1.95	0.41
1:B:315:SER:OG	1:B:316:PHE:N	2.54	0.41
1:B:409:LYS:HE2	1:B:492:GLU:OE2	2.20	0.41
1:B:643:LEU:HA	1:B:643:LEU:HD23	1.82	0.41
1:B:653:ILE:HD11	1:A:645:PHE:CD1	2.55	0.41
1:B:830:ARG:NH2	1:B:1067:ASP:OD1	2.54	0.41
1:A:471:LEU:HD23	1:A:471:LEU:HA	1.90	0.41
1:A:791:LEU:HD23	1:A:791:LEU:HA	1.85	0.41
1:B:632:LEU:HD23	1:B:632:LEU:HA	1.83	0.41
1:A:615:LEU:O	1:A:617:ARG:HD2	2.21	0.40
1:A:314:SER:O	1:A:314:SER:OG	2.39	0.40
1:A:323:VAL:HG23	1:A:503:ALA:HB1	2.02	0.40
1:A:691:LEU:HA	1:A:691:LEU:HD23	1.88	0.40
1:A:806:VAL:HG22	1:A:840:LEU:HD12	2.03	0.40
1:B:806:VAL:HG22	1:B:840:LEU:HD12	2.03	0.40
1:A:153:LEU:HB2	1:A:480:TYR:CE1	2.56	0.40
1:A:392:ILE:O	1:A:396:LEU:HG	2.22	0.40
1:B:615:LEU:O	1:B:617:ARG:HD2	2.21	0.40
1:B:892:ILE:HB	1:B:922:VAL:HG22	2.04	0.40
1:A:136:THR:HA	1:A:464:SER:HB2	2.02	0.40
1:A:392:ILE:HD12	1:A:489:ALA:HB2	2.02	0.40
1:B:153:LEU:HB2	1:B:480:TYR:CE1	2.56	0.40
1:A:677:ASP:HB3	1:A:680:ARG:HB3	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	894/1099 (81%)	868 (97%)	25 (3%)	1 (0%)	48 79

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	894/1099 (81%)	868 (97%)	25 (3%)	1 (0%)	48 79
All	All	1788/2198 (81%)	1736 (97%)	50 (3%)	2 (0%)	49 79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	866	VAL
1	A	866	VAL

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	722/929 (78%)	722 (100%)	0	100 100
1	B	722/929 (78%)	722 (100%)	0	100 100
All	All	1444/1858 (78%)	1444 (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 (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	115	HIS
1	B	133	GLN
1	B	147	ASN
1	B	347	ASN
1	B	423	HIS
1	B	453	ASN
1	B	620	ASN
1	B	629	HIS
1	B	752	ASN
1	B	777	GLN
1	B	882	HIS
1	A	115	HIS

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Mol	Chain	Res	Type
1	A	147	ASN
1	A	347	ASN
1	A	453	ASN
1	A	752	ASN
1	A	882	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	TPO	A	727	1	8,10,11	1.77	1 (12%)	10,14,16	1.79	3 (30%)
1	TPO	B	727	1	8,10,11	1.76	1 (12%)	10,14,16	1.79	3 (30%)
1	SEP	A	930	1	8,9,10	1.57	1 (12%)	7,12,14	1.10	1 (14%)
1	SEP	B	930	1	8,9,10	1.58	1 (12%)	7,12,14	1.10	1 (14%)

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
1	TPO	A	727	1	-	4/9/11/13	-
1	TPO	B	727	1	-	4/9/11/13	-
1	SEP	A	930	1	-	1/6/8/10	-
1	SEP	B	930	1	-	1/6/8/10	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	727	TPO	P-O1P	3.67	1.61	1.50
1	B	727	TPO	P-O1P	3.66	1.61	1.50
1	B	930	SEP	P-O1P	3.47	1.61	1.50
1	A	930	SEP	P-O1P	3.46	1.61	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	727	TPO	CG2-CB-CA	-3.39	106.67	113.26
1	A	727	TPO	CG2-CB-CA	-3.37	106.69	113.26
1	A	727	TPO	P-OG1-CB	-3.29	114.38	123.33
1	B	727	TPO	P-OG1-CB	-3.28	114.42	123.33
1	A	930	SEP	OG-CB-CA	2.16	110.25	108.14
1	B	930	SEP	OG-CB-CA	2.14	110.23	108.14
1	A	727	TPO	O2P-P-OG1	2.12	114.09	105.85
1	B	727	TPO	O2P-P-OG1	2.11	114.06	105.85

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	727	TPO	CG2-CB-OG1-P
1	A	727	TPO	CG2-CB-OG1-P
1	B	930	SEP	CA-CB-OG-P
1	A	930	SEP	CA-CB-OG-P
1	B	727	TPO	CB-OG1-P-O1P
1	A	727	TPO	CB-OG1-P-O1P
1	B	727	TPO	C-CA-CB-CG2
1	B	727	TPO	O-C-CA-CB
1	A	727	TPO	C-CA-CB-CG2
1	A	727	TPO	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	930	SEP	2	0
1	B	930	SEP	2	0

5.5 Carbohydrates

6 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	C	1	1,2	14,14,15	0.28	0	17,19,21	0.77	1 (5%)
2	NAG	C	2	2	14,14,15	0.34	0	17,19,21	0.87	1 (5%)
2	BMA	C	3	2	11,11,12	0.50	0	15,15,17	0.70	0
2	NAG	G	1	1,2	14,14,15	0.29	0	17,19,21	0.77	0
2	NAG	G	2	2	14,14,15	0.35	0	17,19,21	0.87	1 (5%)
2	BMA	G	3	2	11,11,12	0.51	0	15,15,17	0.69	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	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	2/2/19/22	0/1/1/1
2	NAG	G	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	BMA	G	3	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	NAG	C1-O5-C5	2.80	115.93	112.19
2	C	2	NAG	C1-O5-C5	2.79	115.92	112.19
2	C	1	NAG	C4-C3-C2	-2.00	108.08	111.02

There are no chirality outliers.

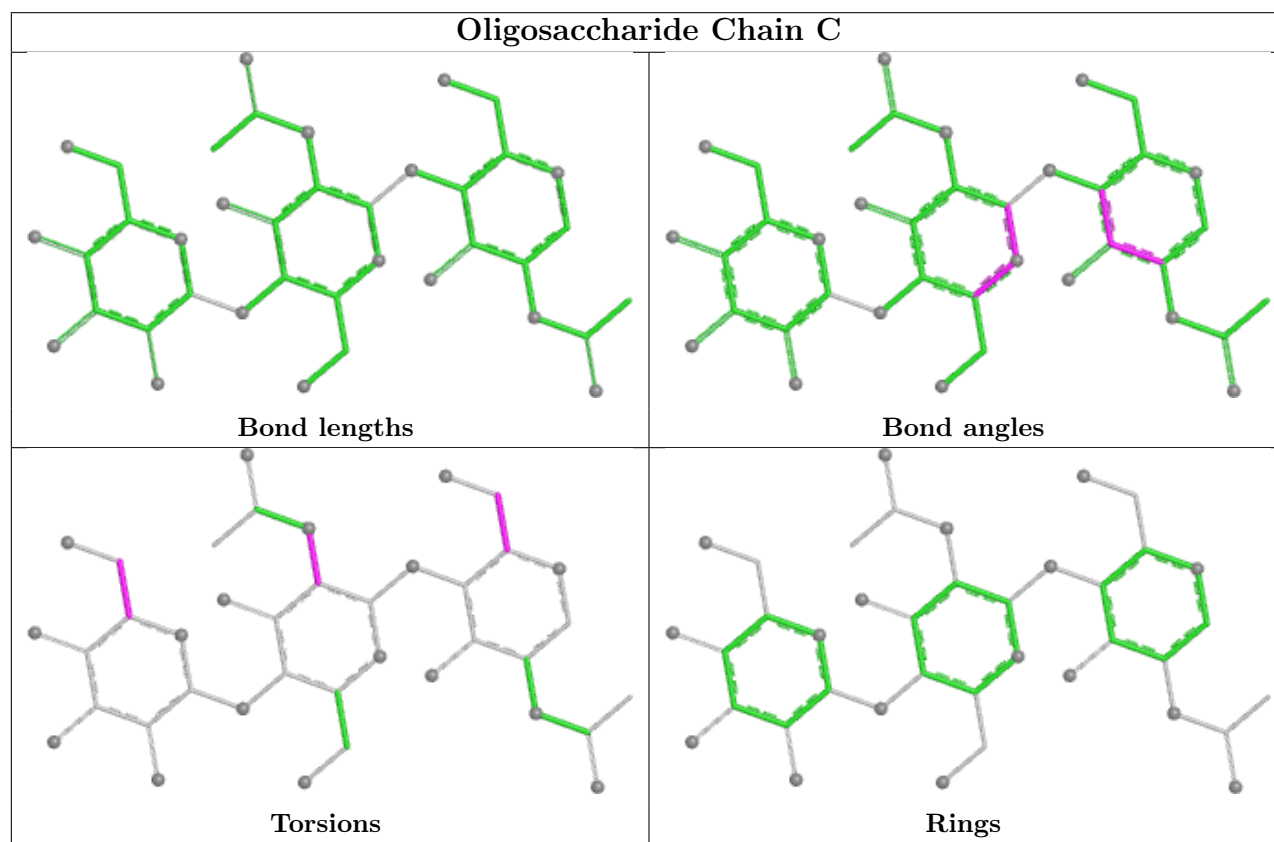
All (12) torsion outliers are listed below:

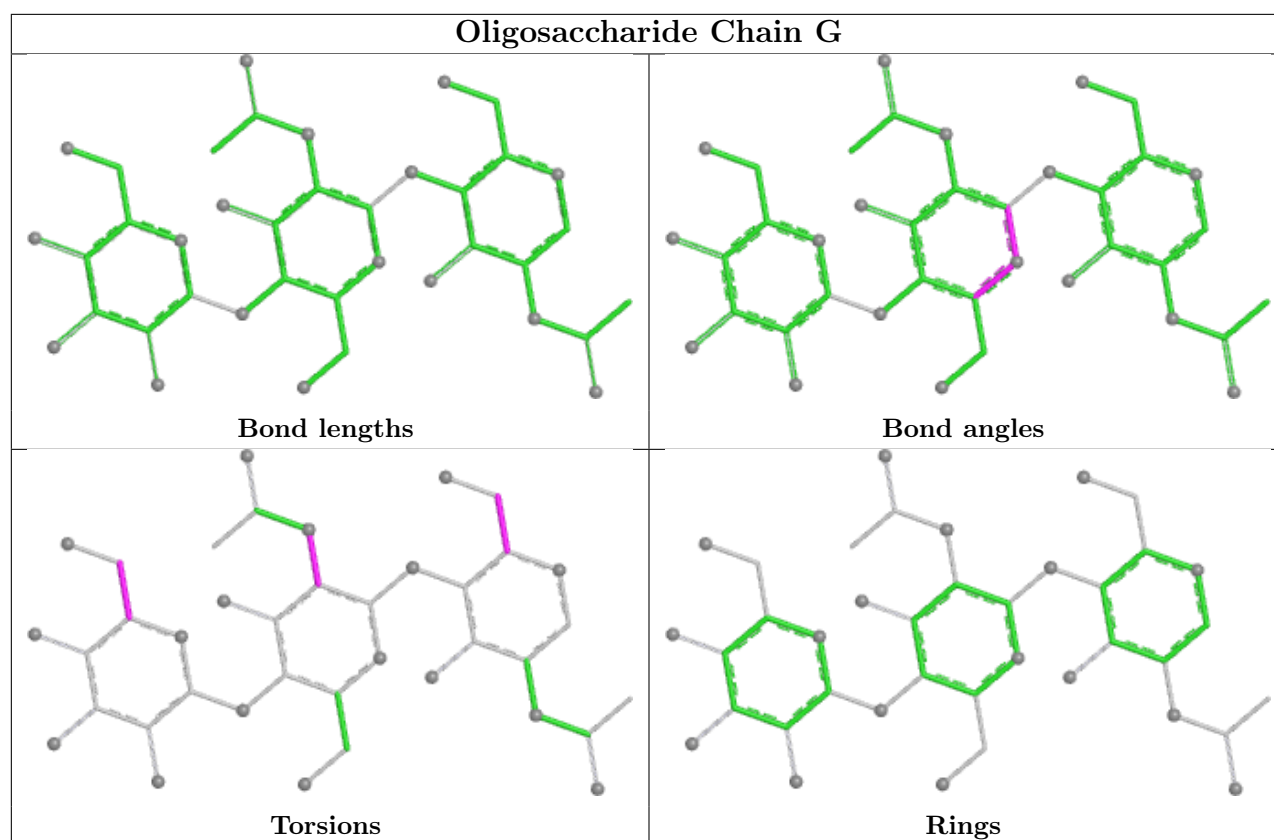
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C4-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	C	3	BMA	O5-C5-C6-O6
2	G	3	BMA	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	C	3	BMA	C4-C5-C6-O6
2	G	3	BMA	C4-C5-C6-O6
2	C	2	NAG	C1-C2-N2-C7
2	G	2	NAG	C1-C2-N2-C7
2	C	2	NAG	C3-C2-N2-C7
2	G	2	NAG	C3-C2-N2-C7

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](#)

2 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	$\# Z > 2$
3	NAG	B	1101	1	14,14,15	0.15	0	17,19,21	0.59	0
3	NAG	A	1101	1	14,14,15	0.14	0	17,19,21	0.59	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
3	NAG	B	1101	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1101	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1101	NAG	O5-C5-C6-O6
3	A	1101	NAG	O5-C5-C6-O6
3	B	1101	NAG	C4-C5-C6-O6
3	A	1101	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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.

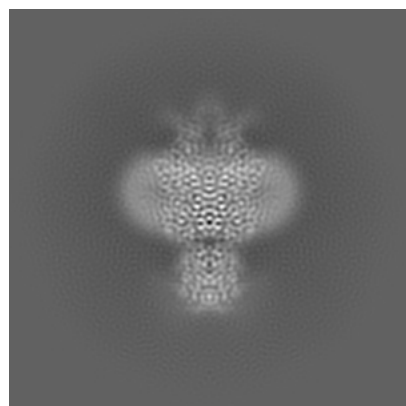
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-11799. These allow visual inspection of the internal detail of the map and identification of artifacts.

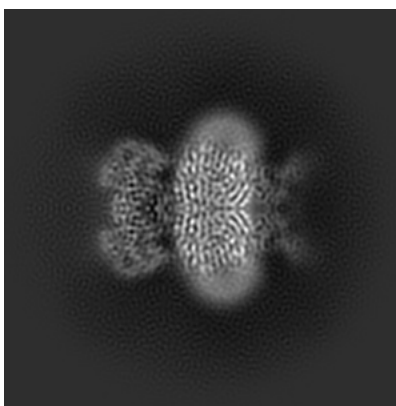
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

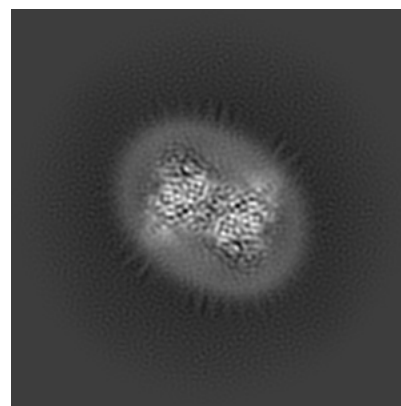
6.1.1 Primary map



X

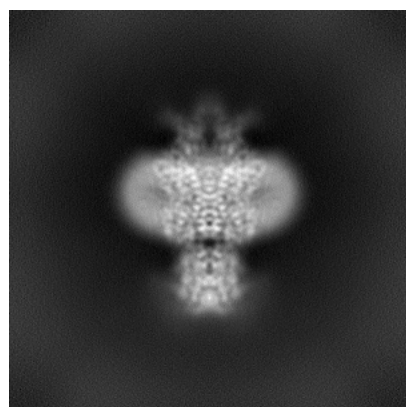


Y

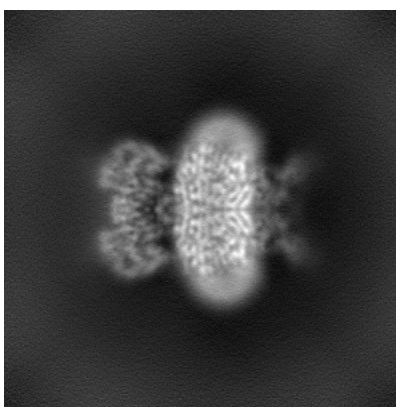


Z

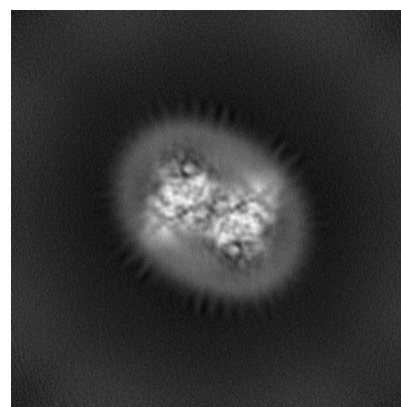
6.1.2 Raw map



X



Y

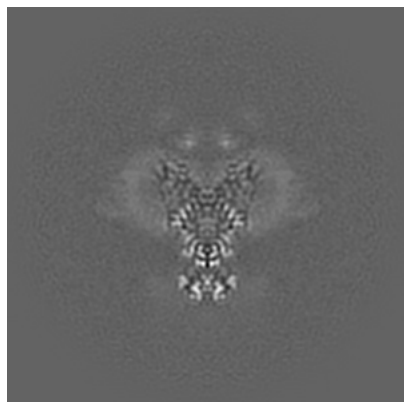


Z

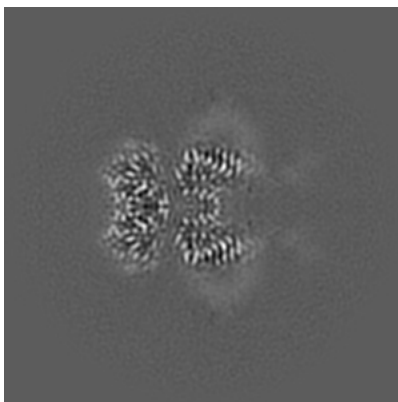
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

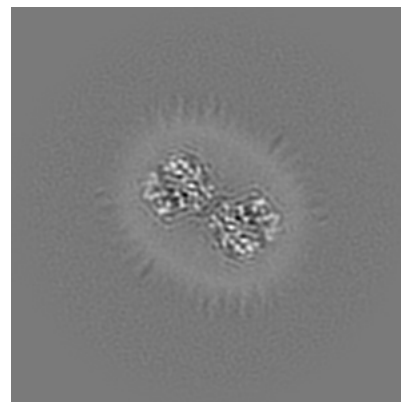
6.2.1 Primary map



X Index: 200

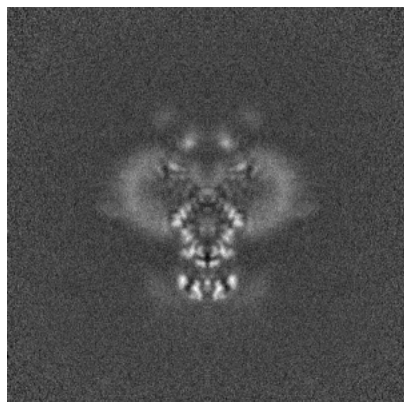


Y Index: 200

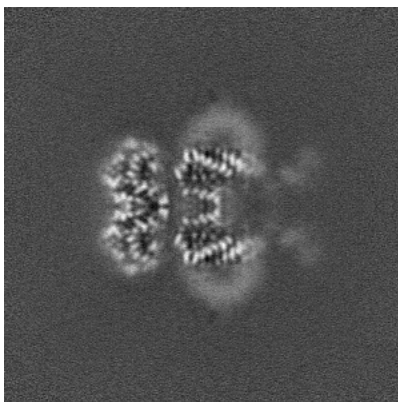


Z Index: 200

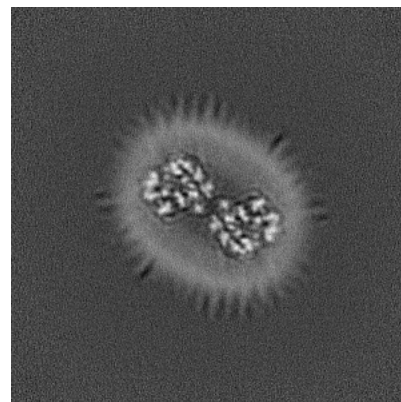
6.2.2 Raw map



X Index: 200



Y Index: 200

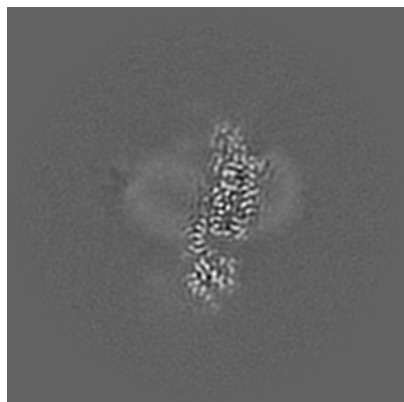


Z Index: 200

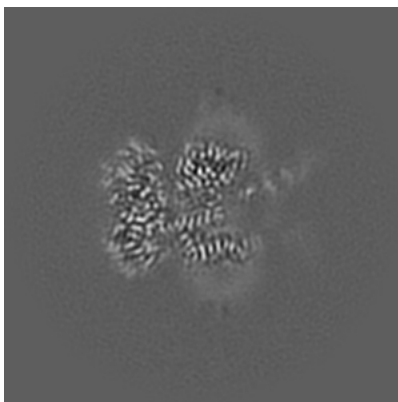
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

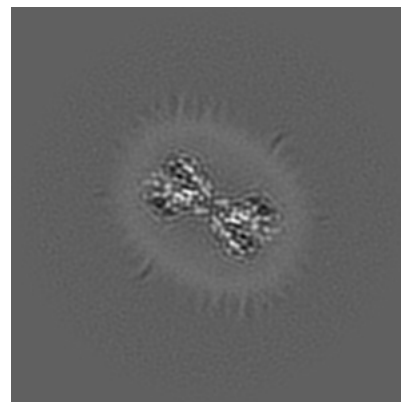
6.3.1 Primary map



X Index: 180

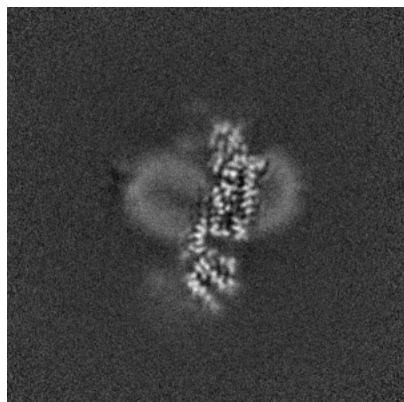


Y Index: 195

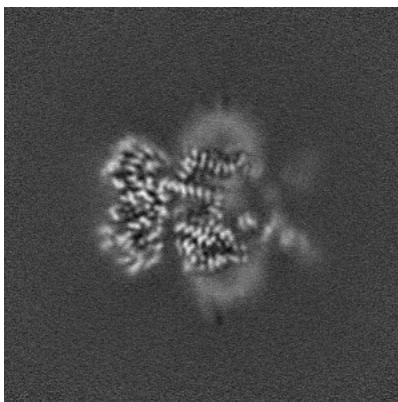


Z Index: 202

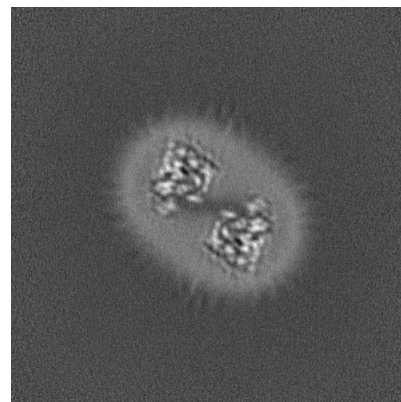
6.3.2 Raw map



X Index: 180



Y Index: 206

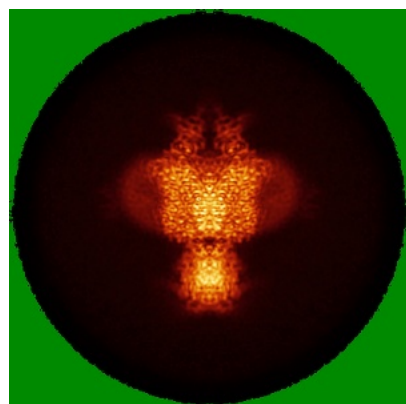


Z Index: 238

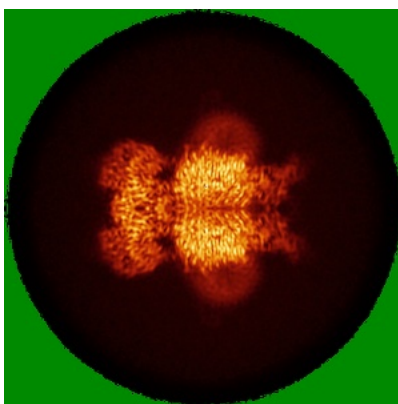
The images above show the largest variance slices of the map in three orthogonal directions.

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

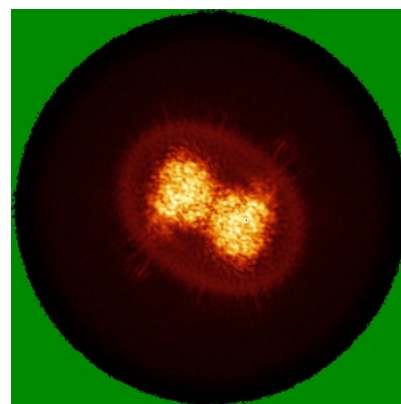
6.4.1 Primary map



X

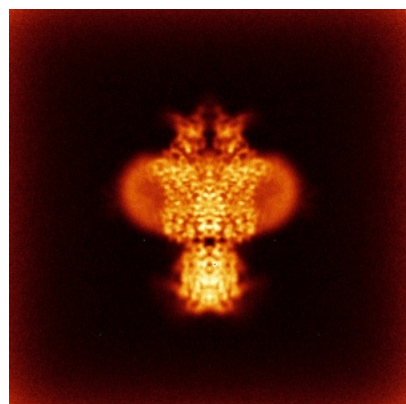


Y

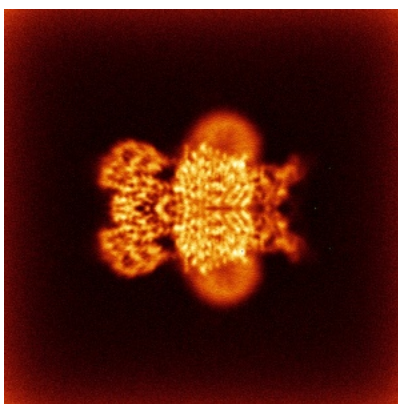


Z

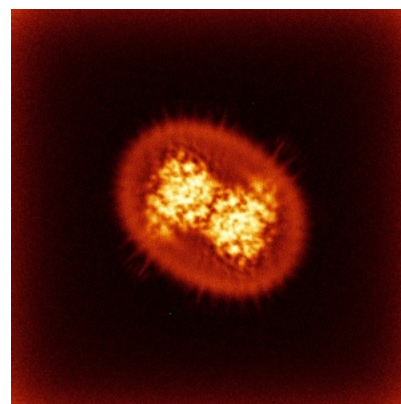
6.4.2 Raw map



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](#)

This section was not generated.

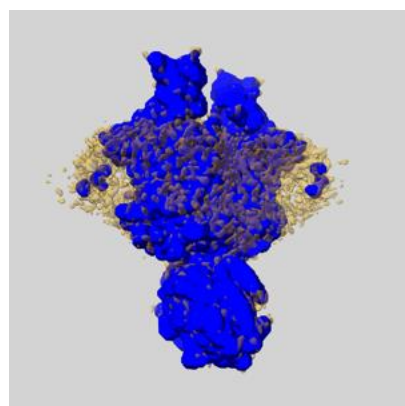
6.6 Mask visualisation [i](#)

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

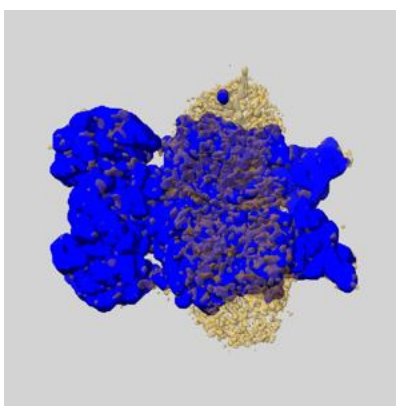
A mask typically either:

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

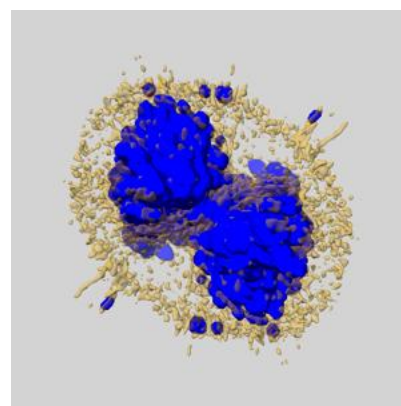
6.6.1 emd_11799_msk_1.map [i](#)



X



Y

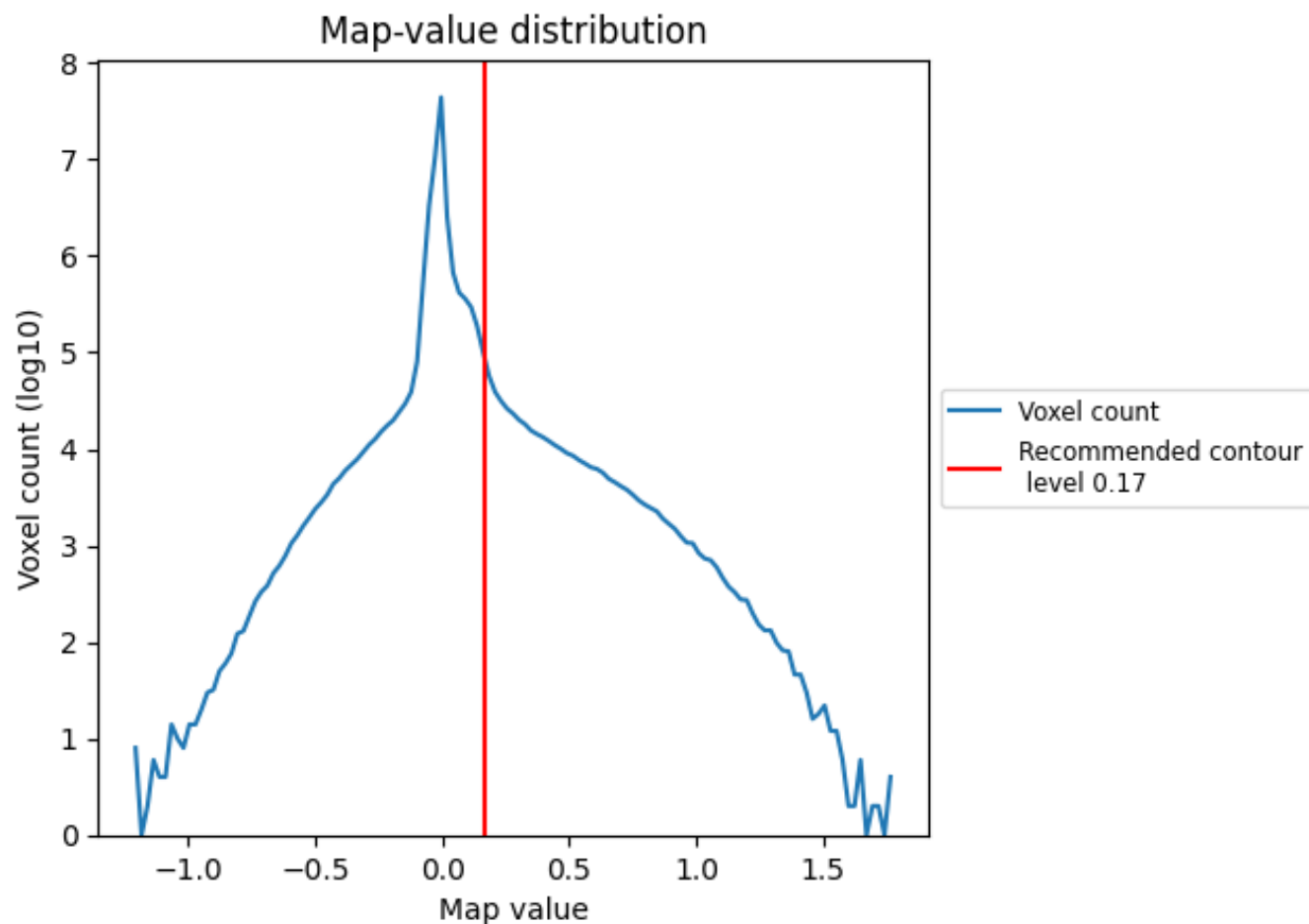


Z

7 Map analysis [i](#)

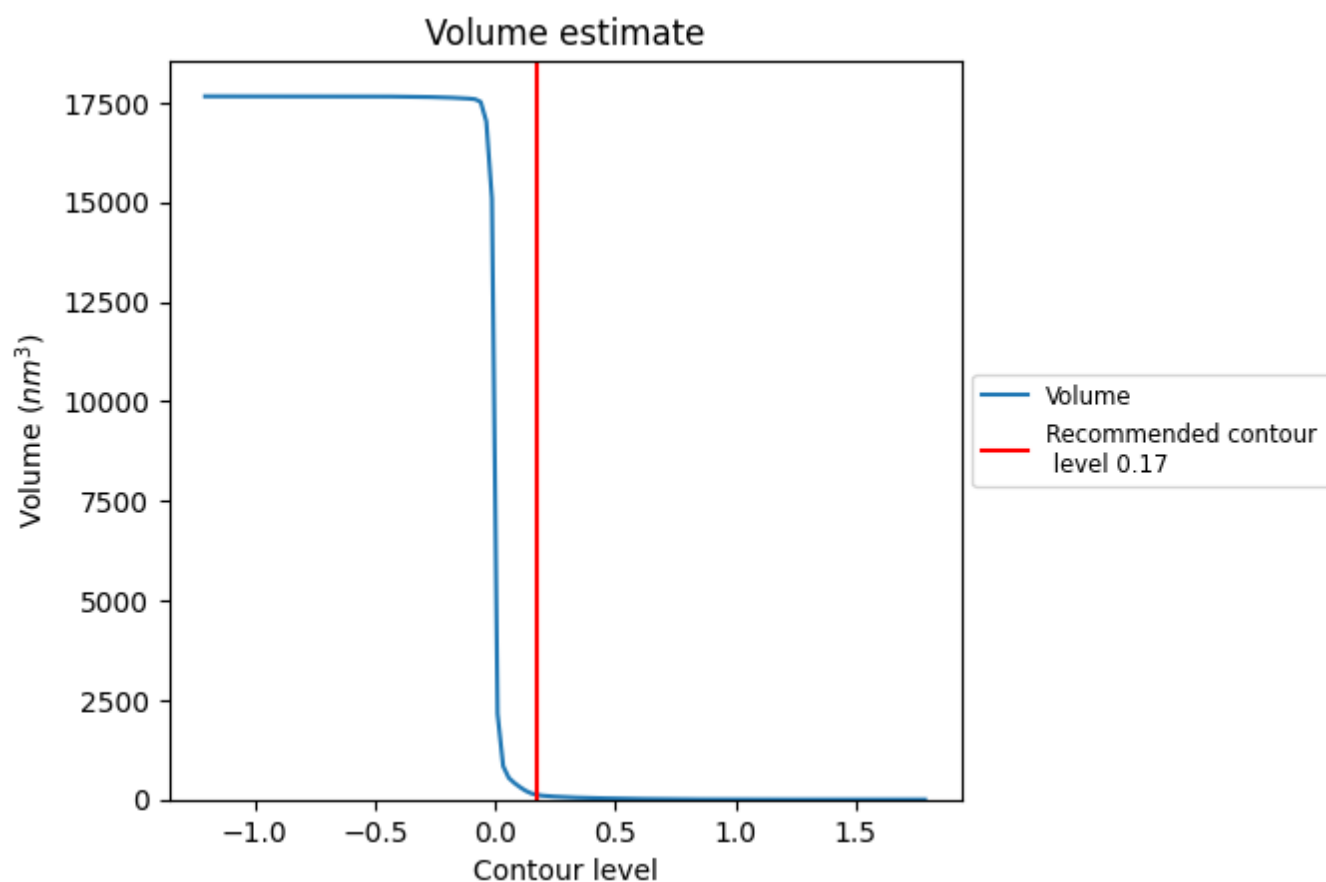
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

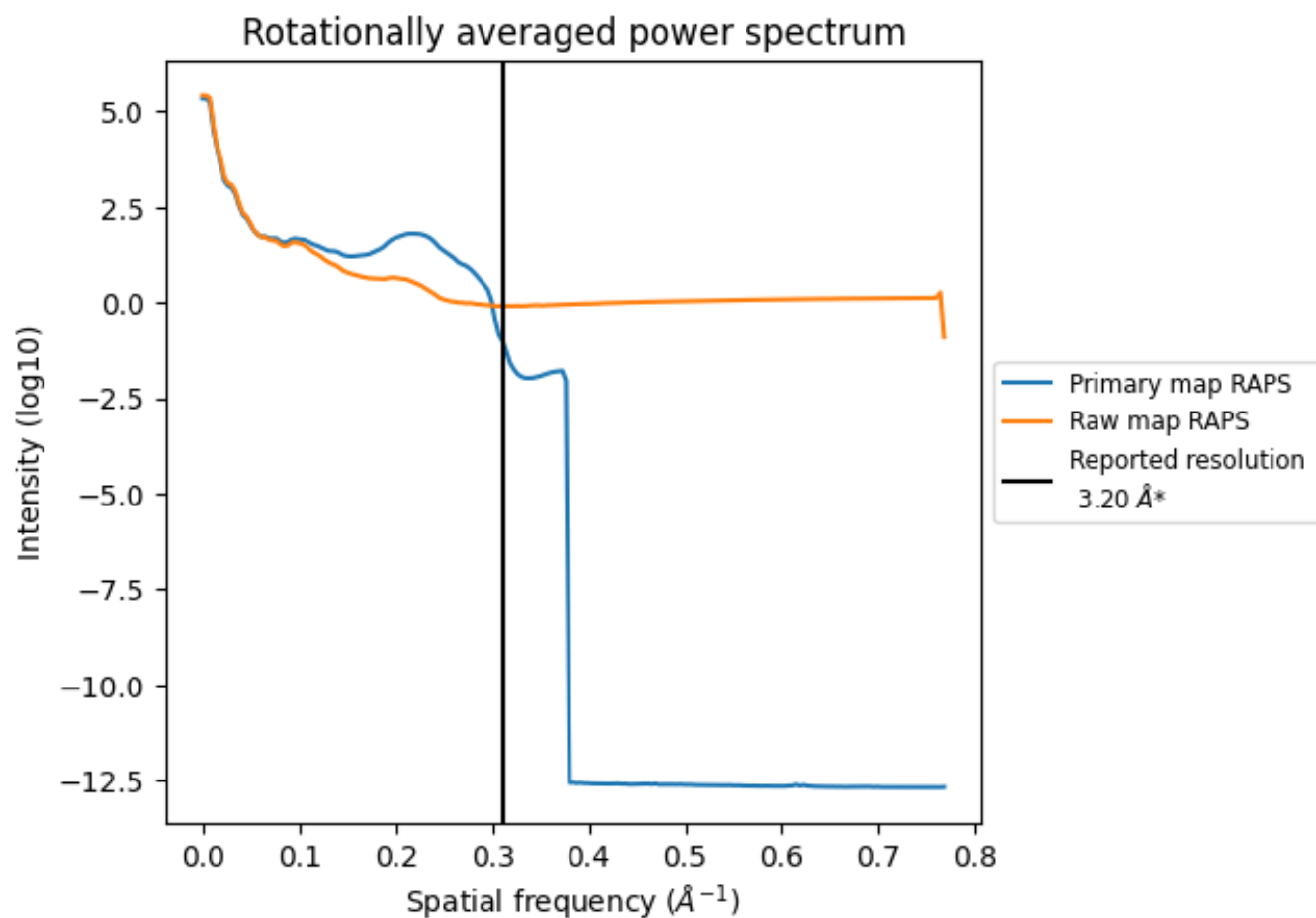
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 124 nm³; 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 ⓘ

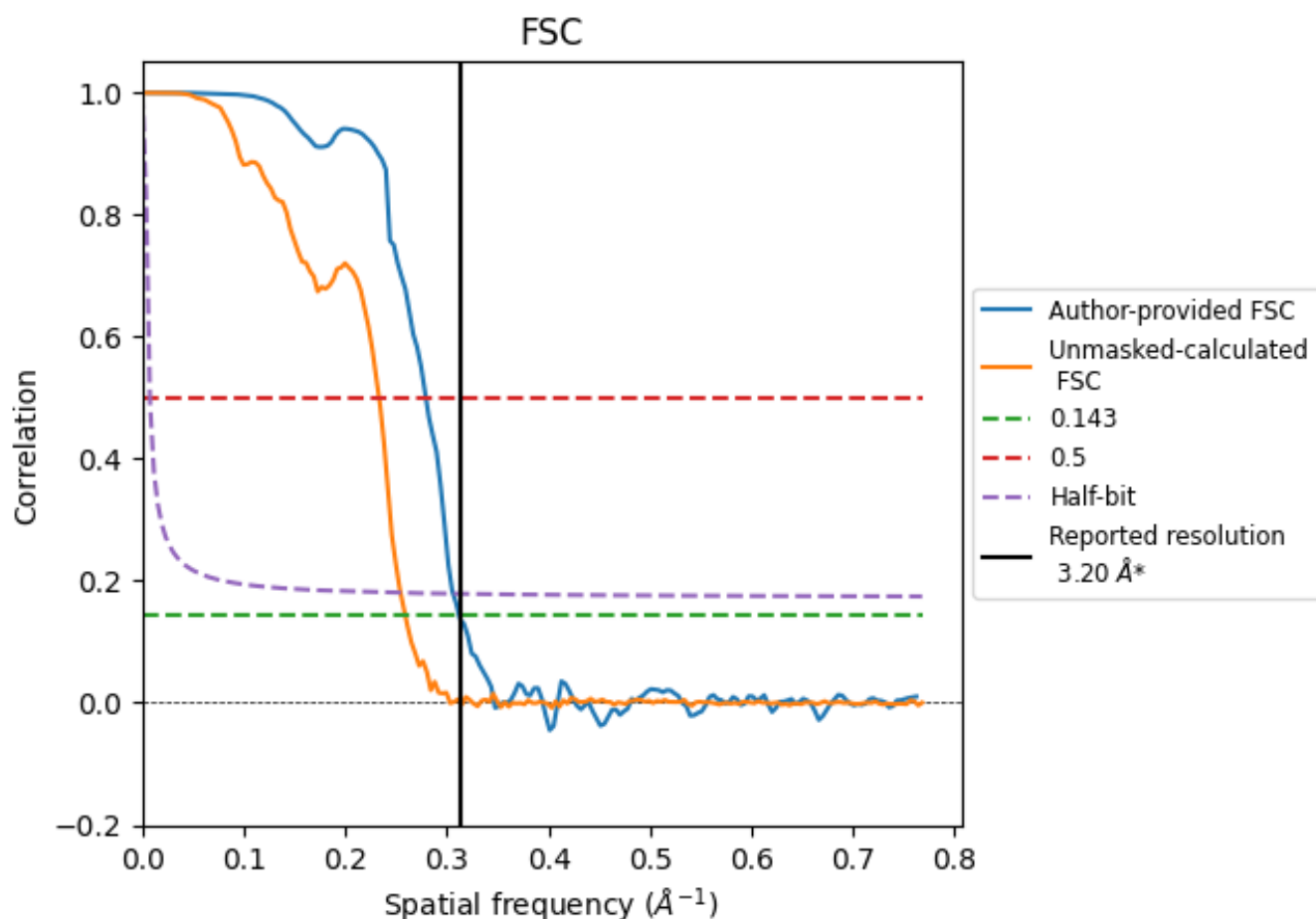


*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.312 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.20	3.58	3.27
Unmasked-calculated*	3.86	4.29	3.93

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.86 differs from the reported value 3.2 by more than 10 %

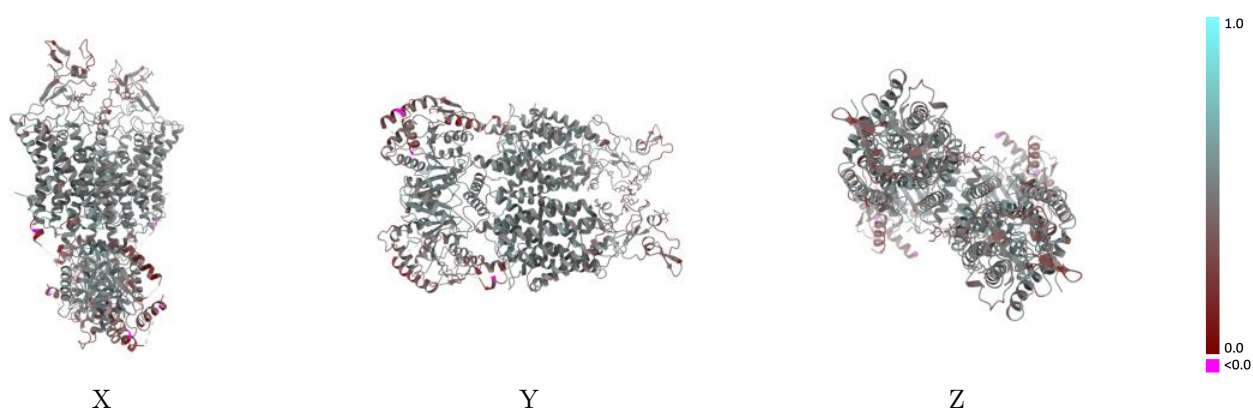
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-11799 and PDB model 7AIN. 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](#)

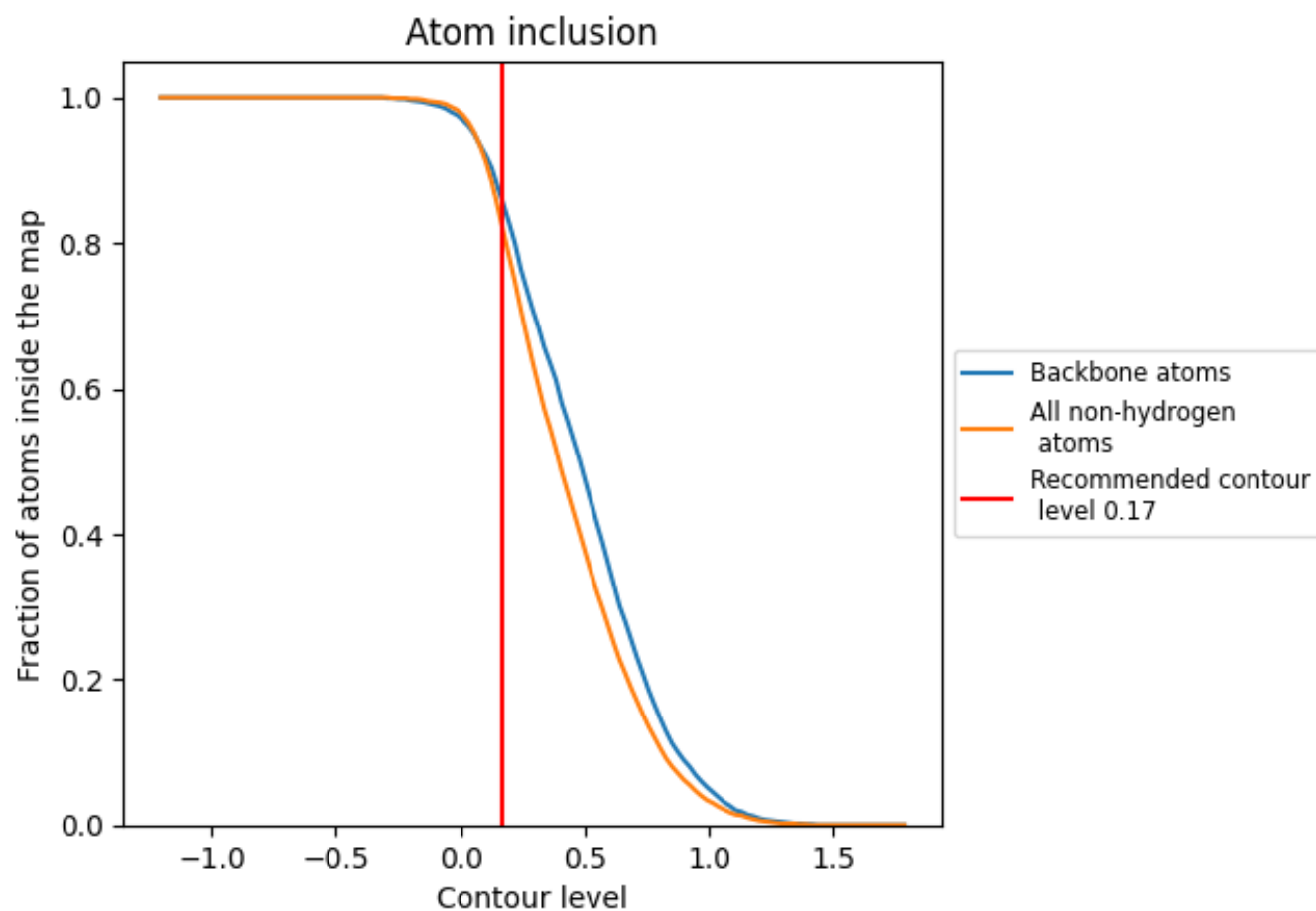


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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, 86% of all backbone atoms, 82% 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 (0.17) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8220	0.4690
A	0.8240	0.4700
B	0.8230	0.4690
C	0.5900	0.3520
G	0.5640	0.3630

