



Full wwPDB EM Validation Report ⓘ

Apr 23, 2026 – 06:27 PM EDT

PDB ID : 9P6X / pdb_00009p6x
EMDB ID : EMD-71328
Title : Structure of EBOV glycoprotein in complex with Nanosota-EB1 and EB2
Authors : Bu, F.; Ye, G.; Li, F.
Deposited on : 2025-06-20
Resolution : 2.92 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

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<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:

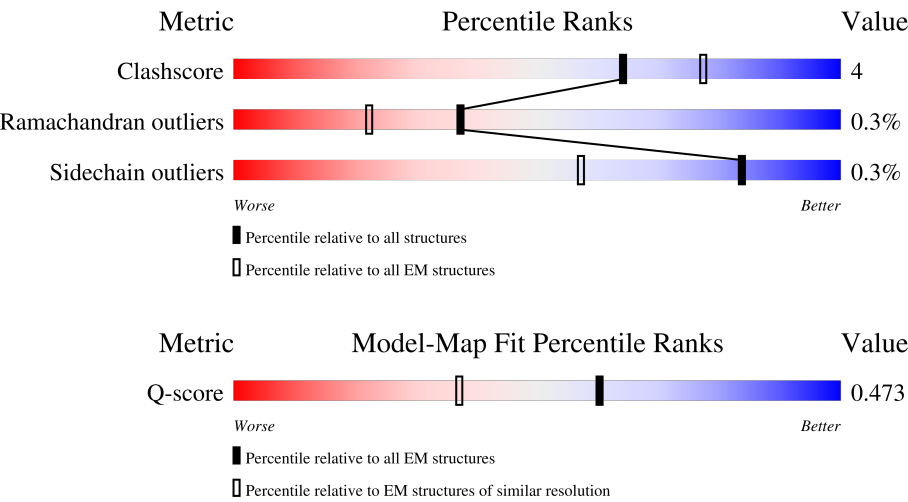
EMDB validation analysis	:	0.0.1.dev132
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.92 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	13007 (2.42 - 3.42)

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	509	61%7%31%
1	B	509	60%9%31%
1	C	509	56%10%34%
2	D	143	17% 76%8%15%

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Mol	Chain	Length	Quality of chain
2	E	143	
2	F	143	
3	G	143	
3	H	143	
3	I	143	
4	J	6	
4	K	6	
5	L	5	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13948 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 Envelope glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	349	Total	C	N	O	S	0	0
			2734	1744	469	510	11		
1	B	351	Total	C	N	O	S	0	0
			2748	1752	471	514	11		
1	C	338	Total	C	N	O	S	0	0
			2651	1697	453	490	11		

There are 177 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	MET	-	initiating methionine	UNP Q05320
A	10	ASP	-	expression tag	UNP Q05320
A	11	ALA	-	expression tag	UNP Q05320
A	12	MET	-	expression tag	UNP Q05320
A	13	LYS	-	expression tag	UNP Q05320
A	14	ARG	-	expression tag	UNP Q05320
A	15	GLY	-	expression tag	UNP Q05320
A	16	LEU	-	expression tag	UNP Q05320
A	17	CYS	-	expression tag	UNP Q05320
A	18	CYS	-	expression tag	UNP Q05320
A	19	VAL	-	expression tag	UNP Q05320
A	20	LEU	-	expression tag	UNP Q05320
A	21	LEU	-	expression tag	UNP Q05320
A	22	LEU	-	expression tag	UNP Q05320
A	23	CYS	-	expression tag	UNP Q05320
A	24	GLY	-	expression tag	UNP Q05320
A	25	ALA	-	expression tag	UNP Q05320
A	26	VAL	-	expression tag	UNP Q05320
A	27	PHE	-	expression tag	UNP Q05320
A	28	VAL	-	expression tag	UNP Q05320
A	29	SER	-	expression tag	UNP Q05320
A	30	ALA	-	expression tag	UNP Q05320
A	31	SER	-	expression tag	UNP Q05320
A	633	GLY	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
A	634	SER	-	expression tag	UNP Q05320
A	635	GLY	-	expression tag	UNP Q05320
A	636	TYR	-	expression tag	UNP Q05320
A	637	ILE	-	expression tag	UNP Q05320
A	638	PRO	-	expression tag	UNP Q05320
A	639	GLU	-	expression tag	UNP Q05320
A	640	ALA	-	expression tag	UNP Q05320
A	641	PRO	-	expression tag	UNP Q05320
A	642	ARG	-	expression tag	UNP Q05320
A	643	ASP	-	expression tag	UNP Q05320
A	644	GLY	-	expression tag	UNP Q05320
A	645	GLN	-	expression tag	UNP Q05320
A	646	ALA	-	expression tag	UNP Q05320
A	647	TYR	-	expression tag	UNP Q05320
A	648	VAL	-	expression tag	UNP Q05320
A	649	ARG	-	expression tag	UNP Q05320
A	650	LYS	-	expression tag	UNP Q05320
A	651	ASP	-	expression tag	UNP Q05320
A	652	GLY	-	expression tag	UNP Q05320
A	653	GLU	-	expression tag	UNP Q05320
A	654	TRP	-	expression tag	UNP Q05320
A	655	VAL	-	expression tag	UNP Q05320
A	656	LEU	-	expression tag	UNP Q05320
A	657	LEU	-	expression tag	UNP Q05320
A	658	SER	-	expression tag	UNP Q05320
A	659	THR	-	expression tag	UNP Q05320
A	660	PHE	-	expression tag	UNP Q05320
A	661	LEU	-	expression tag	UNP Q05320
A	662	GLY	-	expression tag	UNP Q05320
A	663	HIS	-	expression tag	UNP Q05320
A	664	HIS	-	expression tag	UNP Q05320
A	665	HIS	-	expression tag	UNP Q05320
A	666	HIS	-	expression tag	UNP Q05320
A	667	HIS	-	expression tag	UNP Q05320
A	668	HIS	-	expression tag	UNP Q05320
B	9	MET	-	initiating methionine	UNP Q05320
B	10	ASP	-	expression tag	UNP Q05320
B	11	ALA	-	expression tag	UNP Q05320
B	12	MET	-	expression tag	UNP Q05320
B	13	LYS	-	expression tag	UNP Q05320
B	14	ARG	-	expression tag	UNP Q05320
B	15	GLY	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
B	16	LEU	-	expression tag	UNP Q05320
B	17	CYS	-	expression tag	UNP Q05320
B	18	CYS	-	expression tag	UNP Q05320
B	19	VAL	-	expression tag	UNP Q05320
B	20	LEU	-	expression tag	UNP Q05320
B	21	LEU	-	expression tag	UNP Q05320
B	22	LEU	-	expression tag	UNP Q05320
B	23	CYS	-	expression tag	UNP Q05320
B	24	GLY	-	expression tag	UNP Q05320
B	25	ALA	-	expression tag	UNP Q05320
B	26	VAL	-	expression tag	UNP Q05320
B	27	PHE	-	expression tag	UNP Q05320
B	28	VAL	-	expression tag	UNP Q05320
B	29	SER	-	expression tag	UNP Q05320
B	30	ALA	-	expression tag	UNP Q05320
B	31	SER	-	expression tag	UNP Q05320
B	633	GLY	-	expression tag	UNP Q05320
B	634	SER	-	expression tag	UNP Q05320
B	635	GLY	-	expression tag	UNP Q05320
B	636	TYR	-	expression tag	UNP Q05320
B	637	ILE	-	expression tag	UNP Q05320
B	638	PRO	-	expression tag	UNP Q05320
B	639	GLU	-	expression tag	UNP Q05320
B	640	ALA	-	expression tag	UNP Q05320
B	641	PRO	-	expression tag	UNP Q05320
B	642	ARG	-	expression tag	UNP Q05320
B	643	ASP	-	expression tag	UNP Q05320
B	644	GLY	-	expression tag	UNP Q05320
B	645	GLN	-	expression tag	UNP Q05320
B	646	ALA	-	expression tag	UNP Q05320
B	647	TYR	-	expression tag	UNP Q05320
B	648	VAL	-	expression tag	UNP Q05320
B	649	ARG	-	expression tag	UNP Q05320
B	650	LYS	-	expression tag	UNP Q05320
B	651	ASP	-	expression tag	UNP Q05320
B	652	GLY	-	expression tag	UNP Q05320
B	653	GLU	-	expression tag	UNP Q05320
B	654	TRP	-	expression tag	UNP Q05320
B	655	VAL	-	expression tag	UNP Q05320
B	656	LEU	-	expression tag	UNP Q05320
B	657	LEU	-	expression tag	UNP Q05320
B	658	SER	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
B	659	THR	-	expression tag	UNP Q05320
B	660	PHE	-	expression tag	UNP Q05320
B	661	LEU	-	expression tag	UNP Q05320
B	662	GLY	-	expression tag	UNP Q05320
B	663	HIS	-	expression tag	UNP Q05320
B	664	HIS	-	expression tag	UNP Q05320
B	665	HIS	-	expression tag	UNP Q05320
B	666	HIS	-	expression tag	UNP Q05320
B	667	HIS	-	expression tag	UNP Q05320
B	668	HIS	-	expression tag	UNP Q05320
C	9	MET	-	initiating methionine	UNP Q05320
C	10	ASP	-	expression tag	UNP Q05320
C	11	ALA	-	expression tag	UNP Q05320
C	12	MET	-	expression tag	UNP Q05320
C	13	LYS	-	expression tag	UNP Q05320
C	14	ARG	-	expression tag	UNP Q05320
C	15	GLY	-	expression tag	UNP Q05320
C	16	LEU	-	expression tag	UNP Q05320
C	17	CYS	-	expression tag	UNP Q05320
C	18	CYS	-	expression tag	UNP Q05320
C	19	VAL	-	expression tag	UNP Q05320
C	20	LEU	-	expression tag	UNP Q05320
C	21	LEU	-	expression tag	UNP Q05320
C	22	LEU	-	expression tag	UNP Q05320
C	23	CYS	-	expression tag	UNP Q05320
C	24	GLY	-	expression tag	UNP Q05320
C	25	ALA	-	expression tag	UNP Q05320
C	26	VAL	-	expression tag	UNP Q05320
C	27	PHE	-	expression tag	UNP Q05320
C	28	VAL	-	expression tag	UNP Q05320
C	29	SER	-	expression tag	UNP Q05320
C	30	ALA	-	expression tag	UNP Q05320
C	31	SER	-	expression tag	UNP Q05320
C	633	GLY	-	expression tag	UNP Q05320
C	634	SER	-	expression tag	UNP Q05320
C	635	GLY	-	expression tag	UNP Q05320
C	636	TYR	-	expression tag	UNP Q05320
C	637	ILE	-	expression tag	UNP Q05320
C	638	PRO	-	expression tag	UNP Q05320
C	639	GLU	-	expression tag	UNP Q05320
C	640	ALA	-	expression tag	UNP Q05320
C	641	PRO	-	expression tag	UNP Q05320

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Chain	Residue	Modelled	Actual	Comment	Reference
C	642	ARG	-	expression tag	UNP Q05320
C	643	ASP	-	expression tag	UNP Q05320
C	644	GLY	-	expression tag	UNP Q05320
C	645	GLN	-	expression tag	UNP Q05320
C	646	ALA	-	expression tag	UNP Q05320
C	647	TYR	-	expression tag	UNP Q05320
C	648	VAL	-	expression tag	UNP Q05320
C	649	ARG	-	expression tag	UNP Q05320
C	650	LYS	-	expression tag	UNP Q05320
C	651	ASP	-	expression tag	UNP Q05320
C	652	GLY	-	expression tag	UNP Q05320
C	653	GLU	-	expression tag	UNP Q05320
C	654	TRP	-	expression tag	UNP Q05320
C	655	VAL	-	expression tag	UNP Q05320
C	656	LEU	-	expression tag	UNP Q05320
C	657	LEU	-	expression tag	UNP Q05320
C	658	SER	-	expression tag	UNP Q05320
C	659	THR	-	expression tag	UNP Q05320
C	660	PHE	-	expression tag	UNP Q05320
C	661	LEU	-	expression tag	UNP Q05320
C	662	GLY	-	expression tag	UNP Q05320
C	663	HIS	-	expression tag	UNP Q05320
C	664	HIS	-	expression tag	UNP Q05320
C	665	HIS	-	expression tag	UNP Q05320
C	666	HIS	-	expression tag	UNP Q05320
C	667	HIS	-	expression tag	UNP Q05320
C	668	HIS	-	expression tag	UNP Q05320

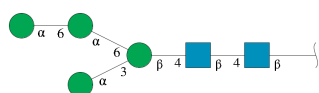
- Molecule 2 is a protein called Nanosota-EB1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	122	Total	C	N	O	S	0	0
			927	576	164	182	5		
2	E	122	Total	C	N	O	S	0	0
			927	576	164	182	5		
2	F	122	Total	C	N	O	S	0	0
			927	576	164	182	5		

- Molecule 3 is a protein called Nanosota-EB2.

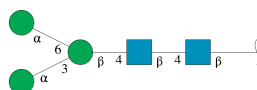
Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	122	Total	C	N	O	S	0	0
			929	575	166	184	4		
3	H	122	Total	C	N	O	S	0	0
			929	575	166	184	4		
3	I	122	Total	C	N	O	S	0	0
			929	575	166	184	4		

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



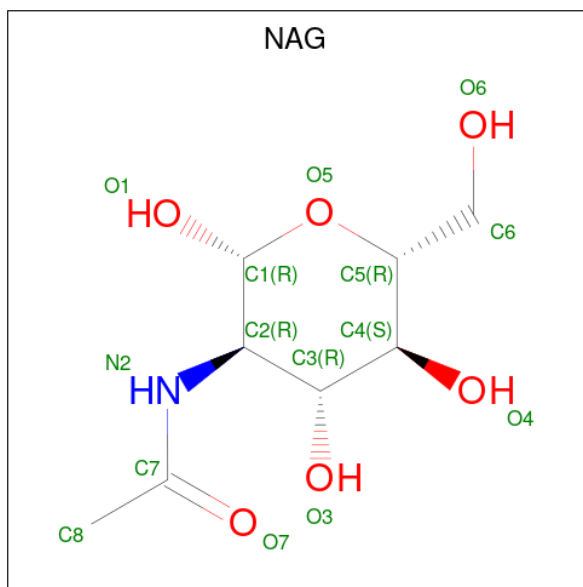
Mol	Chain	Residues	Atoms				AltConf	Trace
4	J	6	Total	C	N	O	0	0
			72	40	2	30		
4	K	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



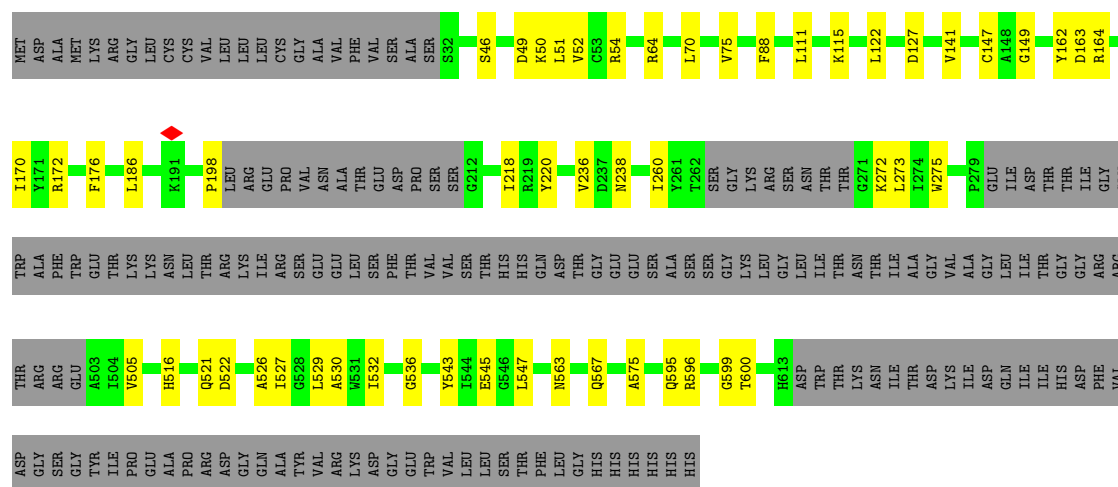
Mol	Chain	Residues	Atoms				AltConf	Trace
5	L	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



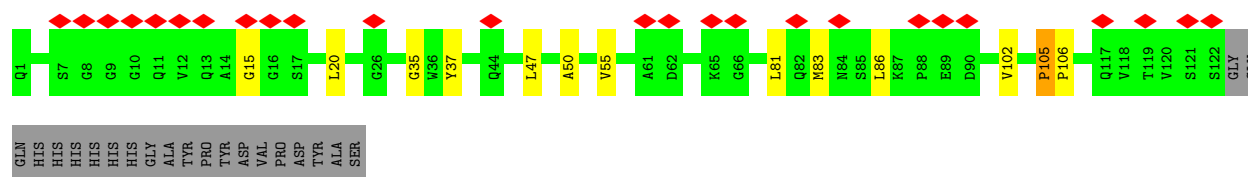
Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	

Chain C:  56% 10% 34%



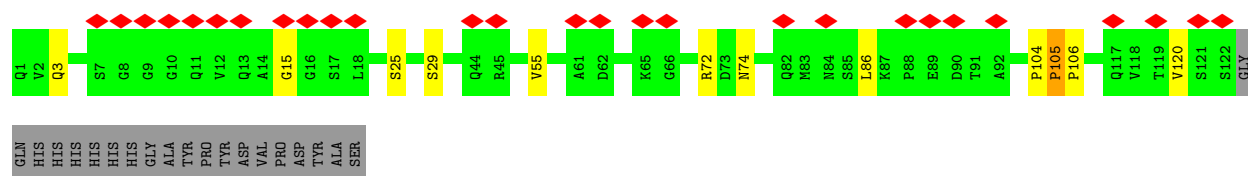
• Molecule 2: Nanosota-EB1

Chain D:  17% 76% 8% 15%



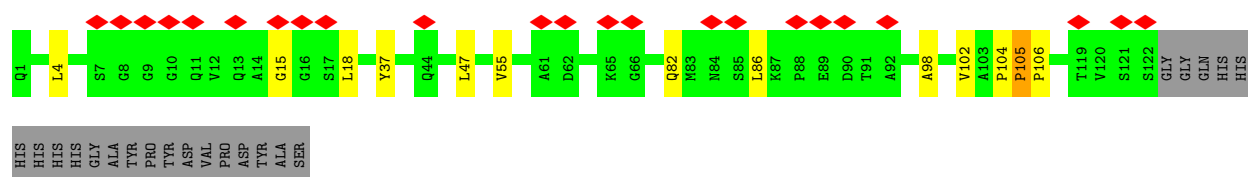
• Molecule 2: Nanosota-EB1

Chain E:  19% 77% 8% 15%



• Molecule 2: Nanosota-EB1

Chain F:  16% 76% 8% 15%

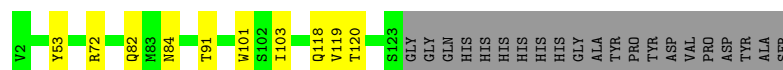


• Molecule 3: Nanosota-EB2

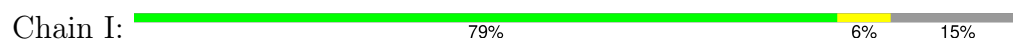
Chain G:  80% 6% 15%



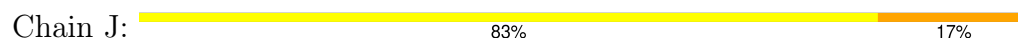
- Molecule 3: Nanosota-EB2



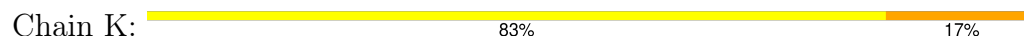
- Molecule 3: Nanosota-EB2



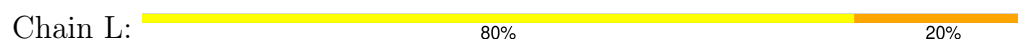
- Molecule 4: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 4: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	62575	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.526	Depositor
Minimum map value	-0.219	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.05	Depositor
Map size ($\text{\AA}$)	398.39984, 398.39984, 398.39984	wwPDB
Map dimensions	450, 450, 450	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.885333, 0.885333, 0.885333	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.29	0/2802	0.56	0/3809
1	B	0.33	0/2816	0.58	0/3829
1	C	0.32	0/2718	0.59	0/3695
2	D	0.29	0/945	0.66	1/1282 (0.1%)
2	E	0.26	0/945	0.60	1/1282 (0.1%)
2	F	0.27	0/945	0.67	1/1282 (0.1%)
3	G	0.17	0/948	0.46	0/1284
3	H	0.25	0/948	0.58	0/1284
3	I	0.19	0/948	0.50	0/1284
All	All	0.29	0/14015	0.58	3/19031 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	105	PRO	N-CA-C	6.93	119.16	110.70
2	D	105	PRO	N-CA-C	6.60	118.75	110.70
2	E	105	PRO	N-CA-C	6.04	118.07	110.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2734	0	2659	26	0
1	B	2748	0	2673	29	0
1	C	2651	0	2578	32	0
2	D	927	0	901	9	0
2	E	927	0	901	8	0
2	F	927	0	901	8	0
3	G	929	0	879	5	0
3	H	929	0	879	7	0
3	I	929	0	879	7	0
4	J	72	0	61	2	0
4	K	72	0	61	1	0
5	L	61	0	52	1	0
6	A	14	0	13	0	0
6	B	14	0	13	1	0
6	C	14	0	13	1	0
All	All	13948	0	13463	112	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (112) 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:198:PRO:HD2	1:A:547:LEU:H	1.61	0.65
2:E:15:GLY:H	2:E:86:LEU:HB3	1.61	0.64
1:A:85:ARG:NH1	1:A:178:GLU:OE2	2.32	0.62
3:G:99:ARG:HH11	3:G:100:PRO:HD2	1.65	0.61
1:C:46:SER:HB2	1:C:50:LYS:HB2	1.82	0.61
1:A:603:ILE:HD12	1:B:612:PRO:HA	1.83	0.60
2:D:15:GLY:H	2:D:86:LEU:HB3	1.67	0.59
1:A:263:SER:HG	2:E:29:SER:HG	1.51	0.59
3:I:52:ASN:O	3:I:72:ARG:NH1	2.35	0.59
1:C:236:VAL:HG11	1:C:260:ILE:HD11	1.84	0.58
1:C:516:HIS:NE2	1:C:545:GLU:OE2	2.38	0.57
1:C:529:LEU:HB3	1:C:532:ILE:HD12	1.87	0.57
1:A:507:ALA:HB1	3:H:103:ILE:HG23	1.89	0.54
1:B:506:ASN:O	3:I:57:ARG:NH2	2.41	0.54
1:A:164:ARG:NH1	1:C:575:ALA:O	2.40	0.54
1:A:46:SER:HB2	1:A:50:LYS:HB2	1.90	0.53
2:F:37:TYR:HE1	2:F:47:LEU:HD13	1.72	0.53
2:F:15:GLY:H	2:F:86:LEU:HB3	1.73	0.53
1:B:70:LEU:HG	1:B:179:GLY:HA2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:3:GLN:NE2	2:E:25:SER:OG	2.42	0.53
1:C:88:PHE:HB3	1:C:149:GLY:H	1.74	0.53
1:A:89:ARG:NH1	1:B:534:TYR:O	2.42	0.52
1:A:238:ASN:O	1:A:272:LYS:NZ	2.41	0.52
1:B:163:ASP:OD1	1:B:543:TYR:OH	2.26	0.52
1:C:198:PRO:HD2	1:C:547:LEU:H	1.74	0.52
2:E:72:ARG:NH2	2:E:74:ASN:OD1	2.40	0.52
1:A:240:THR:HG21	1:A:260:ILE:HG21	1.92	0.51
2:D:20:LEU:HD13	2:D:81:LEU:HD22	1.92	0.51
1:B:114:LYS:NZ	1:B:120:GLU:OE2	2.44	0.50
1:B:195:SER:O	1:B:196:SER:C	2.54	0.50
1:A:76:ALA:O	1:A:85:ARG:NH2	2.45	0.50
3:I:83:MET:HB3	3:I:86:LEU:HD21	1.94	0.49
3:I:40:ALA:HB3	3:I:43:LYS:HB2	1.93	0.49
1:A:197:HIS:HA	1:A:546:GLY:HA2	1.95	0.49
2:D:83:MET:HE1	2:D:86:LEU:HD13	1.95	0.49
1:A:56:LYS:NZ	1:C:599:GLY:O	2.45	0.49
1:A:163:ASP:OD1	1:A:543:TYR:OH	2.29	0.48
1:B:191:LYS:O	1:B:195:SER:HB2	2.13	0.48
1:B:128:GLY:HA3	1:B:580:ARG:HH12	1.78	0.48
3:G:40:ALA:HB3	3:G:43:LYS:HB2	1.96	0.48
3:H:101:TRP:CD1	4:J:2:NAG:H4	2.49	0.48
1:B:37:VAL:HG12	1:B:185:ILE:HB	1.96	0.47
1:C:49:ASP:OD1	1:C:595:GLN:NE2	2.47	0.47
1:C:526:ALA:O	1:C:530:ALA:HB3	2.14	0.47
1:A:88:PHE:HB3	1:A:149:GLY:H	1.79	0.47
2:D:20:LEU:HB2	2:D:81:LEU:HB3	1.96	0.47
3:H:82:GLN:NE2	3:H:84:ASN:OD1	2.42	0.47
1:B:575:ALA:O	1:C:164:ARG:NH1	2.48	0.47
1:B:198:PRO:HD2	1:B:547:LEU:H	1.79	0.47
1:C:51:LEU:O	1:C:596:ARG:NH2	2.41	0.47
1:B:218:ILE:HA	6:B:701:NAG:H81	1.97	0.47
1:A:126:PRO:HD2	1:A:129:ILE:HG13	1.97	0.46
1:A:37:VAL:HG12	1:A:185:ILE:HB	1.96	0.46
1:B:563:ASN:OD1	5:L:1:NAG:N2	2.48	0.46
1:C:122:LEU:O	1:C:172:ARG:NH1	2.48	0.46
1:A:127:ASP:OD1	1:A:127:ASP:N	2.47	0.46
1:B:183:PHE:HB3	1:B:585:LEU:HD21	1.98	0.46
1:C:163:ASP:OD1	1:C:543:TYR:OH	2.27	0.46
1:C:505:VAL:H	3:G:105:ASN:HD21	1.62	0.46
3:H:91:THR:HG23	3:H:120:THR:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:526:ALA:HB1	1:C:567:GLN:HE22	1.80	0.46
1:B:275:TRP:HE3	2:D:102:VAL:HG22	1.80	0.46
1:A:196:SER:O	1:A:197:HIS:C	2.58	0.45
1:C:275:TRP:HE3	2:F:102:VAL:HG22	1.81	0.45
2:E:86:LEU:HD21	2:E:120:VAL:HG22	1.99	0.45
1:C:141:VAL:HG22	1:C:220:TYR:HB2	1.99	0.45
2:F:4:LEU:HD21	2:F:98:ALA:HB2	1.98	0.45
1:B:56:LYS:O	1:B:62:GLN:NE2	2.48	0.44
2:E:105:PRO:N	2:E:106:PRO:HD2	2.32	0.44
1:C:64:ARG:HD2	1:C:186:LEU:HD11	1.98	0.44
1:C:111:LEU:HD12	1:C:170:ILE:HG21	1.99	0.44
2:D:20:LEU:HD22	2:D:81:LEU:HD13	2.00	0.44
2:F:105:PRO:N	2:F:106:PRO:HD2	2.33	0.44
1:A:263:SER:OG	2:E:29:SER:OG	2.25	0.43
2:F:104:PRO:HG2	2:F:106:PRO:HG2	2.00	0.43
1:B:38:ILE:HB	1:B:187:PRO:HD3	1.99	0.43
2:F:18:LEU:O	2:F:82:GLN:NE2	2.52	0.43
3:H:53:TYR:HA	3:H:72:ARG:HH12	1.83	0.43
1:A:521:GLN:HE22	1:A:524:GLY:HA3	1.83	0.43
1:B:193:PHE:HZ	1:B:554:LEU:HD21	1.83	0.43
1:A:506:ASN:ND2	1:A:511:CYS:SG	2.92	0.43
1:B:122:LEU:HB2	1:B:172:ARG:HB2	2.00	0.43
1:A:602:HIS:O	1:A:603:ILE:HG13	2.18	0.42
1:B:63:LEU:HD23	1:B:185:ILE:HG12	2.02	0.42
1:B:505:VAL:H	3:I:105:ASN:ND2	2.18	0.42
1:C:52:VAL:HG13	1:C:54:ARG:H	1.85	0.42
2:D:37:TYR:HE1	2:D:47:LEU:HD13	1.83	0.42
3:I:76:LYS:HB3	3:I:76:LYS:HE2	1.81	0.42
2:E:104:PRO:HG2	2:E:106:PRO:HG2	2.02	0.42
3:G:91:THR:HG23	3:G:120:THR:HA	2.02	0.42
2:D:105:PRO:O	2:D:106:PRO:C	2.62	0.42
3:G:34:MET:HE2	3:G:34:MET:HB3	1.91	0.42
3:H:118:GLN:NE2	3:H:119:VAL:O	2.52	0.42
1:B:59:SER:OG	1:B:61:ASN:OD1	2.30	0.41
1:C:218:ILE:HA	6:C:701:NAG:H81	2.02	0.41
1:C:563:ASN:OD1	4:K:1:NAG:N2	2.52	0.41
1:A:79:VAL:HG21	1:A:141:VAL:HG21	2.02	0.41
1:B:247:ARG:NH2	1:B:278:ASN:O	2.54	0.41
2:D:35:GLY:HA2	2:D:50:ALA:HA	2.03	0.41
1:A:609:CYS:SG	1:C:600:THR:OG1	2.70	0.41
3:H:101:TRP:HD1	4:J:2:NAG:H4	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:ASP:OD1	1:B:283:THR:N	2.54	0.41
1:C:273:LEU:HD22	2:F:105:PRO:HG2	2.02	0.41
1:B:572:PHE:HA	1:C:522:ASP:OD1	2.21	0.40
1:C:115:LYS:HE2	1:C:147:CYS:HB2	2.03	0.40
1:C:127:ASP:N	1:C:127:ASP:OD1	2.51	0.40
1:B:79:VAL:HB	1:B:80:PRO:HD3	2.03	0.40
1:B:89:ARG:NH2	1:C:536:GLY:O	2.55	0.40
1:C:162:TYR:CE1	1:C:176:PHE:HB3	2.56	0.40
1:B:505:VAL:H	3:I:105:ASN:HD21	1.67	0.40
1:C:238:ASN:O	1:C:272:LYS:NZ	2.43	0.40
1:C:70:LEU:HD23	1:C:75:VAL:HG21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	343/509 (67%)	320 (93%)	21 (6%)	2 (1%)	21	50
1	B	345/509 (68%)	328 (95%)	17 (5%)	0	100	100
1	C	330/509 (65%)	313 (95%)	16 (5%)	1 (0%)	36	64
2	D	120/143 (84%)	108 (90%)	11 (9%)	1 (1%)	16	43
2	E	120/143 (84%)	114 (95%)	5 (4%)	1 (1%)	16	43
2	F	120/143 (84%)	111 (92%)	8 (7%)	1 (1%)	16	43
3	G	120/143 (84%)	119 (99%)	1 (1%)	0	100	100
3	H	120/143 (84%)	115 (96%)	5 (4%)	0	100	100
3	I	120/143 (84%)	116 (97%)	4 (3%)	0	100	100
All	All	1738/2385 (73%)	1644 (95%)	88 (5%)	6 (0%)	37	64

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	603	ILE
1	C	527	ILE
1	A	281	ILE
2	D	55	VAL
2	E	55	VAL
2	F	55	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	293/428 (68%)	291 (99%)	2 (1%)	76	91
1	B	295/428 (69%)	293 (99%)	2 (1%)	76	91
1	C	283/428 (66%)	282 (100%)	1 (0%)	84	94
2	D	97/113 (86%)	97 (100%)	0	100	100
2	E	97/113 (86%)	97 (100%)	0	100	100
2	F	97/113 (86%)	97 (100%)	0	100	100
3	G	95/111 (86%)	95 (100%)	0	100	100
3	H	95/111 (86%)	95 (100%)	0	100	100
3	I	95/111 (86%)	95 (100%)	0	100	100
All	All	1447/1956 (74%)	1442 (100%)	5 (0%)	84	95

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	196	SER
1	A	197	HIS
1	B	527	ILE
1	B	529	LEU
1	C	521	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	110	ASN
1	B	197	HIS
1	B	268	ASN
1	B	278	ASN
1	B	506	ASN
1	B	595	GLN
1	C	61	ASN
1	C	506	ASN
1	C	508	GLN
1	C	514	ASN
1	C	521	GLN
1	C	551	GLN
2	D	97	ASN
2	E	3	GLN
2	E	77	ASN
2	E	117	GLN
2	F	77	ASN
2	F	82	GLN
3	G	13	GLN
3	G	39	GLN

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](#)

17 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
4	NAG	J	1	4,1	14,14,15	0.83	0	17,19,21	1.92	5 (29%)
4	NAG	J	2	4	14,14,15	0.75	0	17,19,21	1.05	1 (5%)
4	BMA	J	3	4	11,11,12	0.88	0	15,15,17	2.74	6 (40%)
4	MAN	J	4	4	11,11,12	0.69	0	15,15,17	3.14	3 (20%)
4	MAN	J	5	4	11,11,12	0.72	0	15,15,17	1.34	1 (6%)
4	MAN	J	6	4	11,11,12	0.57	0	15,15,17	2.08	1 (6%)
4	NAG	K	1	4,1	14,14,15	0.78	0	17,19,21	1.62	5 (29%)
4	NAG	K	2	4	14,14,15	0.77	0	17,19,21	1.25	2 (11%)
4	BMA	K	3	4	11,11,12	0.74	0	15,15,17	3.24	6 (40%)
4	MAN	K	4	4	11,11,12	0.78	0	15,15,17	1.65	2 (13%)
4	MAN	K	5	4	11,11,12	0.73	0	15,15,17	1.11	1 (6%)
4	MAN	K	6	4	11,11,12	0.73	0	15,15,17	1.46	1 (6%)
5	NAG	L	1	5,1	14,14,15	0.83	0	17,19,21	1.77	5 (29%)
5	NAG	L	2	5	14,14,15	0.75	0	17,19,21	1.20	1 (5%)
5	BMA	L	3	5	11,11,12	0.75	0	15,15,17	3.14	5 (33%)
5	MAN	L	4	5	11,11,12	0.70	0	15,15,17	1.22	1 (6%)
5	MAN	L	5	5	11,11,12	0.89	1 (9%)	15,15,17	0.95	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
4	NAG	J	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	J	2	4	-	2/6/23/26	0/1/1/1
4	BMA	J	3	4	-	2/2/19/22	0/1/1/1
4	MAN	J	4	4	-	0/2/19/22	0/1/1/1
4	MAN	J	5	4	-	0/2/19/22	0/1/1/1
4	MAN	J	6	4	-	2/2/19/22	0/1/1/1
4	NAG	K	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	K	2	4	-	0/6/23/26	0/1/1/1
4	BMA	K	3	4	-	2/2/19/22	0/1/1/1
4	MAN	K	4	4	-	2/2/19/22	0/1/1/1
4	MAN	K	5	4	-	0/2/19/22	0/1/1/1
4	MAN	K	6	4	-	0/2/19/22	0/1/1/1
5	NAG	L	1	5,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	L	2	5	-	2/6/23/26	0/1/1/1
5	BMA	L	3	5	-	1/2/19/22	0/1/1/1
5	MAN	L	4	5	-	0/2/19/22	1/1/1/1
5	MAN	L	5	5	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	5	MAN	O5-C1	-2.22	1.40	1.43

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	K	3	BMA	C1-O5-C5	10.69	126.52	112.19
5	L	3	BMA	C1-O5-C5	10.56	126.34	112.19
4	J	4	MAN	C1-O5-C5	9.78	125.29	112.19
4	J	3	BMA	C1-O5-C5	7.94	122.83	112.19
4	J	6	MAN	C1-O5-C5	7.46	122.19	112.19
4	J	4	MAN	C1-C2-C3	5.19	117.19	109.64
4	K	6	MAN	C1-O5-C5	4.53	118.26	112.19
4	J	1	NAG	C2-N2-C7	4.24	128.59	122.90
4	K	3	BMA	O3-C3-C4	3.96	119.72	110.38
4	J	5	MAN	C1-O5-C5	3.87	117.38	112.19
4	K	4	MAN	C1-O5-C5	3.72	117.18	112.19
5	L	4	MAN	C1-O5-C5	3.65	117.08	112.19
5	L	1	NAG	O5-C1-C2	-3.65	105.64	111.29
4	J	3	BMA	C2-C3-C4	3.44	116.92	110.86
4	J	1	NAG	O5-C1-C2	-3.44	105.97	111.29
4	J	3	BMA	C3-C4-C5	3.38	116.36	110.23
4	J	1	NAG	O3-C3-C2	-3.36	102.41	109.40
4	K	1	NAG	C1-O5-C5	3.26	116.56	112.19
5	L	1	NAG	C2-N2-C7	3.07	127.02	122.90
5	L	2	NAG	C1-O5-C5	3.06	116.29	112.19
4	K	1	NAG	O5-C1-C2	-3.06	106.56	111.29
4	K	5	MAN	C1-O5-C5	3.02	116.24	112.19
4	K	4	MAN	C1-C2-C3	-3.00	105.28	109.64
5	L	1	NAG	C1-O5-C5	2.92	116.10	112.19
5	L	3	BMA	C3-C4-C5	2.88	115.45	110.23
4	K	2	NAG	C1-O5-C5	2.83	115.98	112.19
5	L	1	NAG	O4-C4-C5	-2.79	102.45	109.32
4	J	1	NAG	C1-O5-C5	2.77	115.90	112.19
5	L	3	BMA	O5-C5-C4	2.73	117.47	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	3	BMA	O3-C3-C2	-2.70	104.54	110.05
4	K	1	NAG	C2-N2-C7	2.67	126.48	122.90
4	K	3	BMA	O5-C5-C4	2.62	117.19	110.83
4	J	4	MAN	C2-C3-C4	2.51	115.28	110.86
4	J	3	BMA	O4-C4-C3	-2.44	104.63	110.38
4	J	2	NAG	C1-O5-C5	2.38	115.37	112.19
4	J	1	NAG	O3-C3-C4	2.37	115.95	110.38
4	J	3	BMA	O5-C5-C4	2.36	116.57	110.83
4	K	2	NAG	C2-N2-C7	2.30	125.98	122.90
5	L	3	BMA	O4-C4-C3	-2.27	105.03	110.38
4	K	1	NAG	O3-C3-C4	2.18	115.50	110.38
5	L	1	NAG	O3-C3-C2	-2.17	104.89	109.40
5	L	3	BMA	O3-C3-C2	-2.08	105.81	110.05
4	K	3	BMA	O4-C4-C3	-2.08	105.47	110.38
4	K	3	BMA	O2-C2-C3	2.05	114.41	110.15
4	K	1	NAG	C1-C2-N2	2.05	113.66	110.43
4	K	3	BMA	C3-C4-C5	2.03	113.91	110.23

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	K	4	MAN	C4-C5-C6-O6
4	J	3	BMA	C4-C5-C6-O6
4	K	3	BMA	O5-C5-C6-O6
5	L	2	NAG	O5-C5-C6-O6
4	K	4	MAN	O5-C5-C6-O6
4	J	3	BMA	O5-C5-C6-O6
5	L	1	NAG	O5-C5-C6-O6
4	J	2	NAG	O5-C5-C6-O6
4	K	3	BMA	C4-C5-C6-O6
4	J	2	NAG	C4-C5-C6-O6
5	L	1	NAG	C4-C5-C6-O6
4	J	1	NAG	C8-C7-N2-C2
4	J	1	NAG	O7-C7-N2-C2
5	L	2	NAG	C4-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
5	L	3	BMA	O5-C5-C6-O6
4	J	6	MAN	C4-C5-C6-O6
4	J	6	MAN	O5-C5-C6-O6

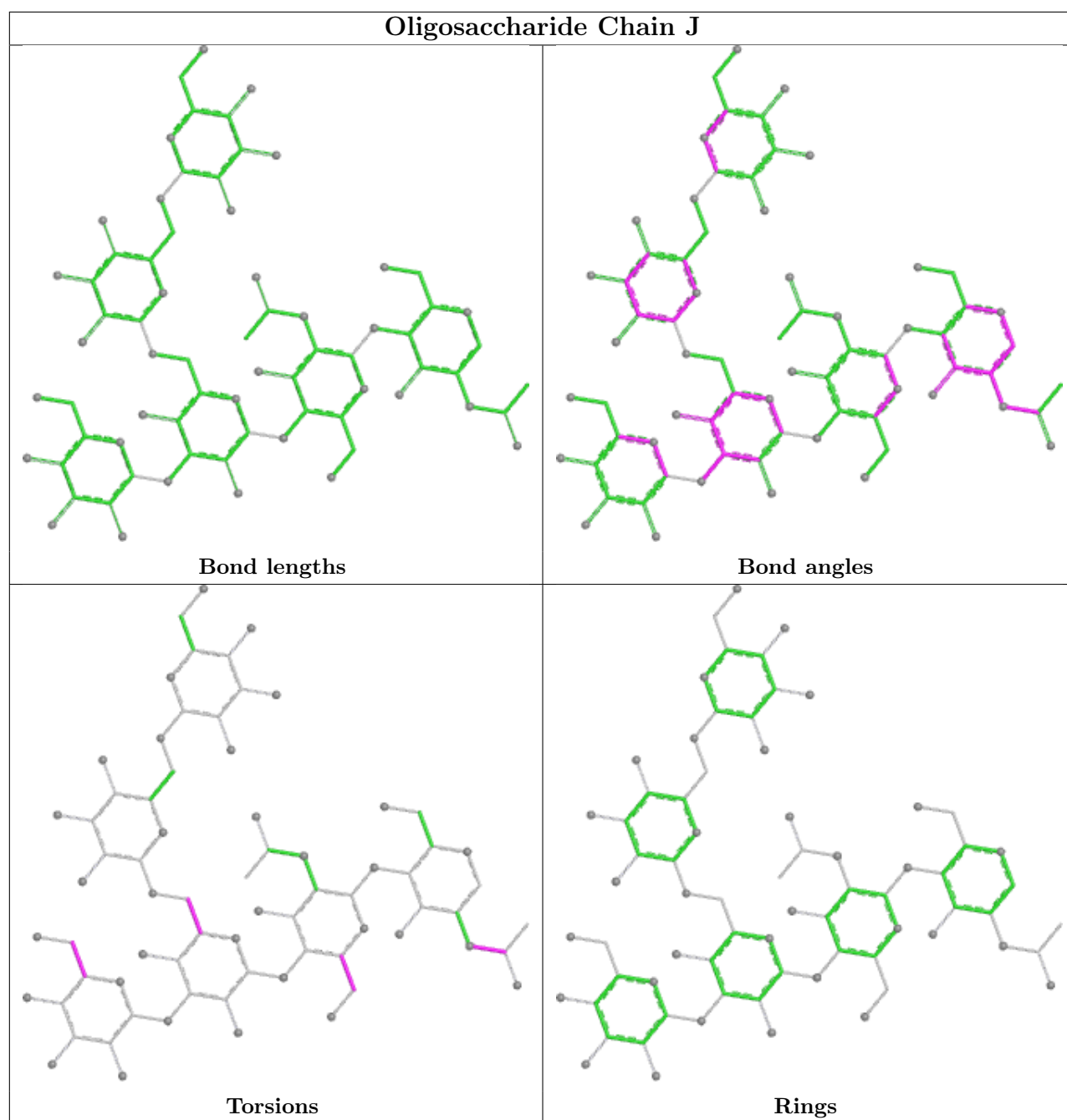
All (1) ring outliers are listed below:

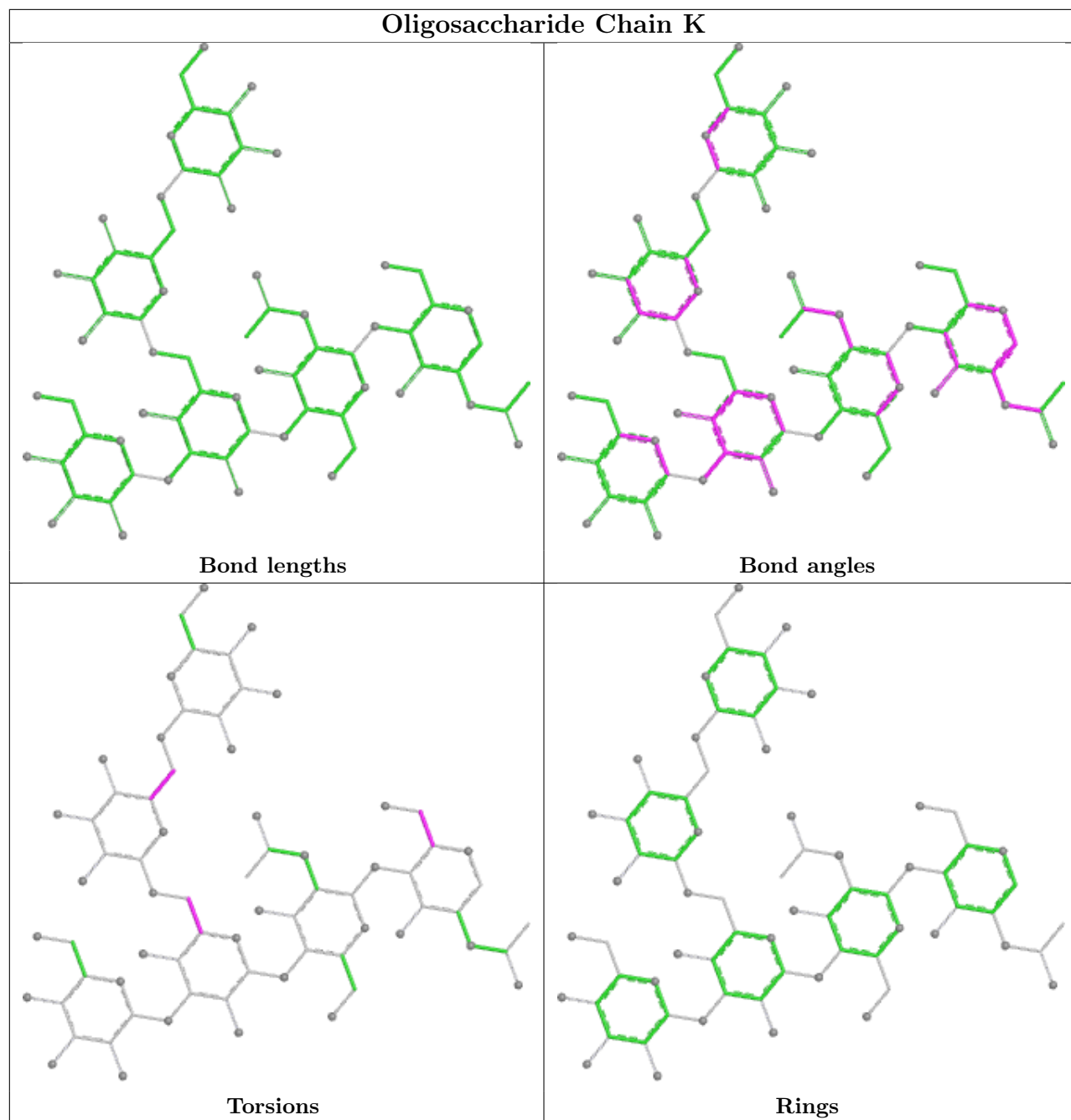
Mol	Chain	Res	Type	Atoms
5	L	4	MAN	C1-C2-C3-C4-C5-O5

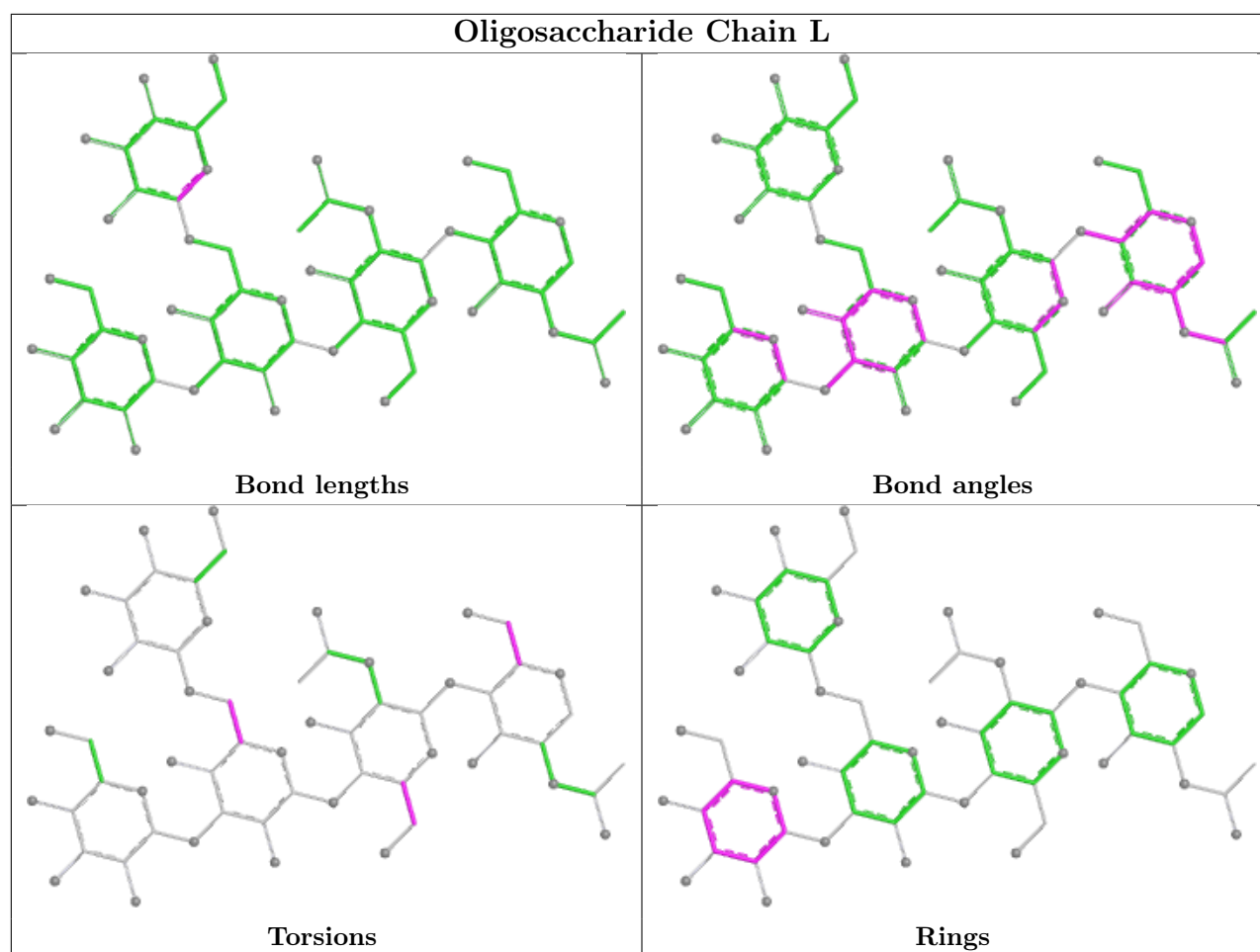
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	1	NAG	1	0
4	J	2	NAG	2	0
4	K	1	NAG	1	0

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	$\# Z > 2$
6	NAG	B	701	1	14,14,15	0.90	1 (7%)	17,19,21	1.82	2 (11%)
6	NAG	C	701	1	14,14,15	0.76	1 (7%)	17,19,21	2.15	1 (5%)
6	NAG	A	701	1	14,14,15	0.65	0	17,19,21	1.83	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
6	NAG	B	701	1	-	0/6/23/26	0/1/1/1
6	NAG	C	701	1	-	0/6/23/26	0/1/1/1
6	NAG	A	701	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	701	NAG	C1-C2	2.77	1.56	1.52
6	C	701	NAG	C1-C2	2.13	1.55	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	701	NAG	C1-O5-C5	7.79	122.63	112.19
6	B	701	NAG	C1-O5-C5	6.65	121.09	112.19
6	A	701	NAG	C1-O5-C5	5.06	118.96	112.19
6	A	701	NAG	C2-N2-C7	3.02	126.95	122.90
6	A	701	NAG	O3-C3-C2	-2.29	104.64	109.40
6	B	701	NAG	C3-C4-C5	-2.11	106.41	110.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	701	NAG	1	0
6	C	701	NAG	1	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.

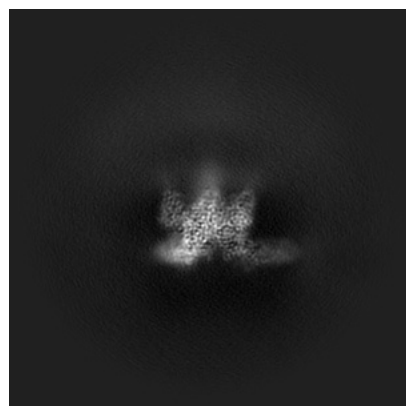
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-71328. These allow visual inspection of the internal detail of the map and identification of artifacts.

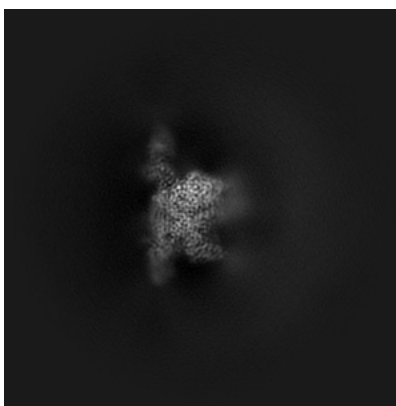
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

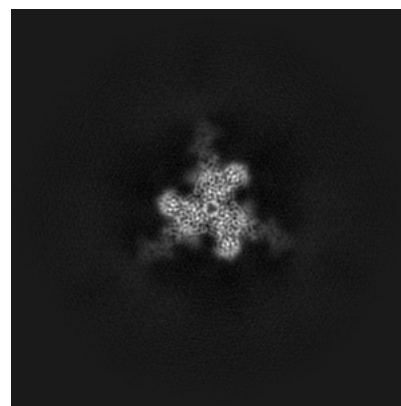
6.1.1 Primary map



X

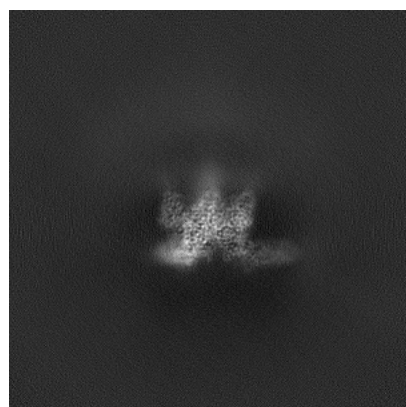


Y

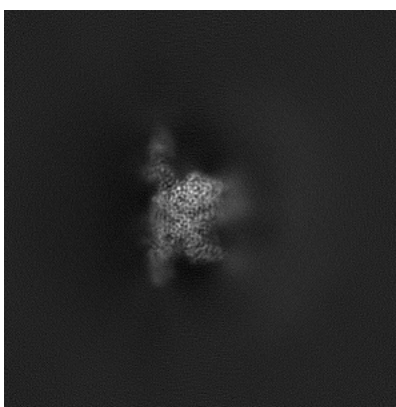


Z

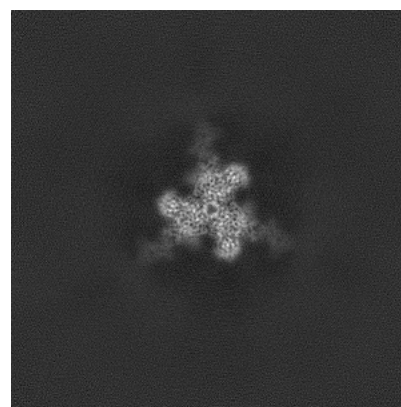
6.1.2 Raw map



X



Y

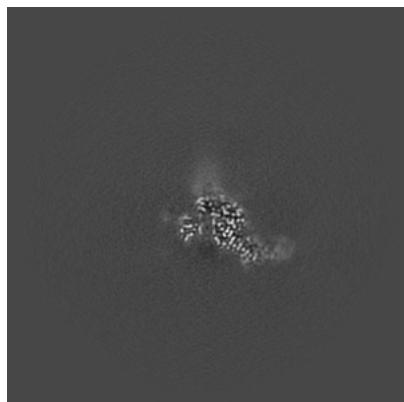


Z

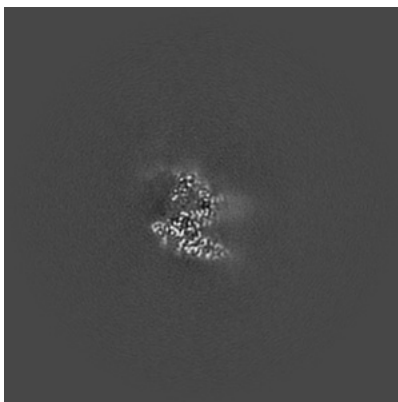
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

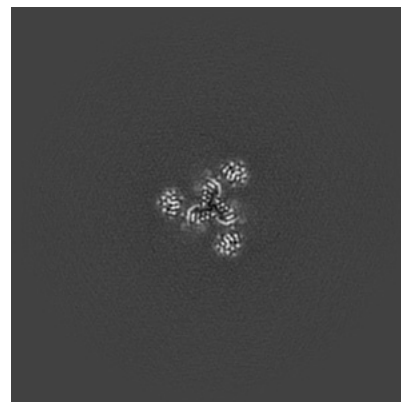
6.2.1 Primary map



X Index: 225

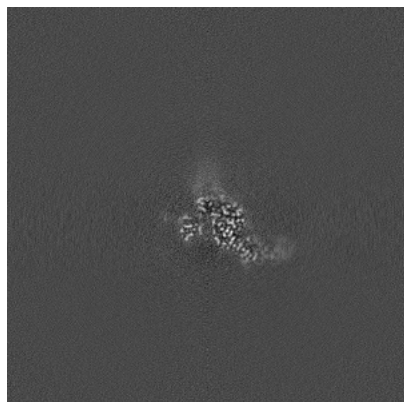


Y Index: 225



Z Index: 225

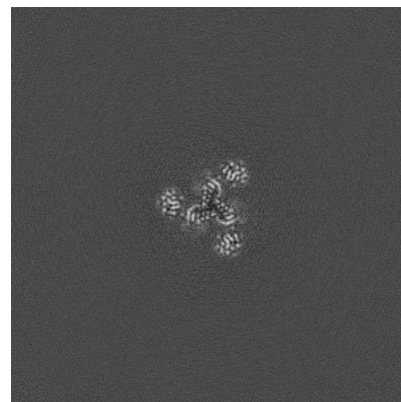
6.2.2 Raw map



X Index: 225



Y Index: 225

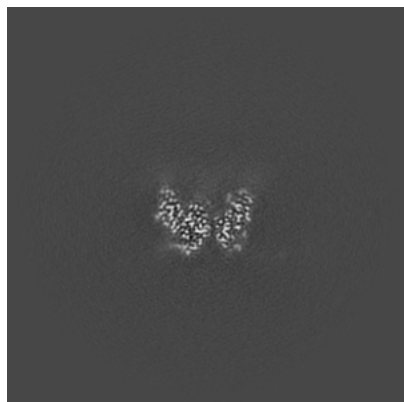


Z Index: 225

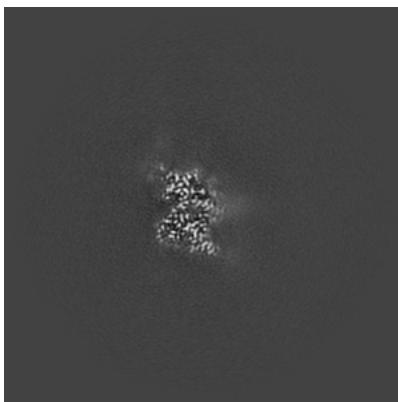
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

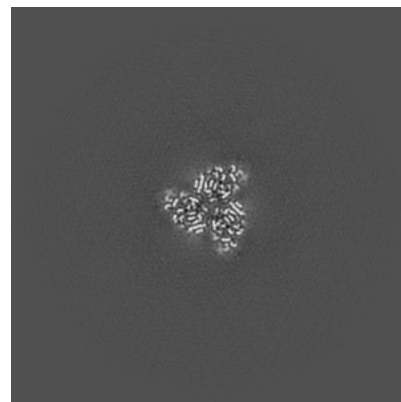
6.3.1 Primary map



X Index: 244

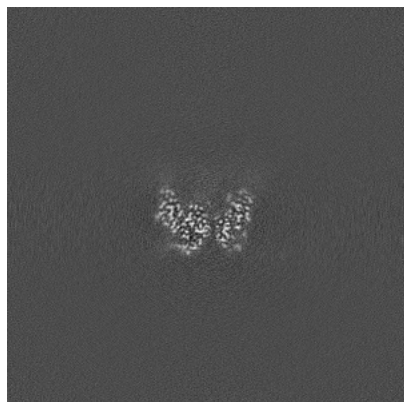


Y Index: 217

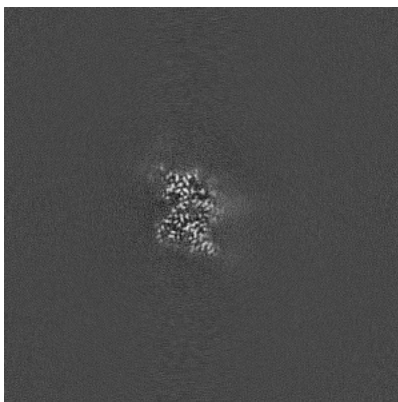


Z Index: 207

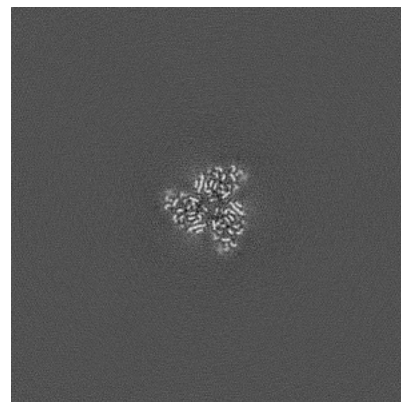
6.3.2 Raw map



X Index: 244



Y Index: 217

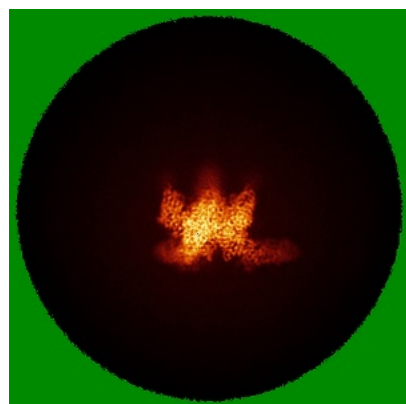


Z Index: 207

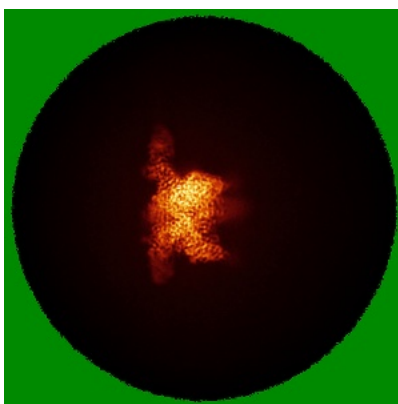
The images above show the largest variance slices of the map in three orthogonal directions.

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

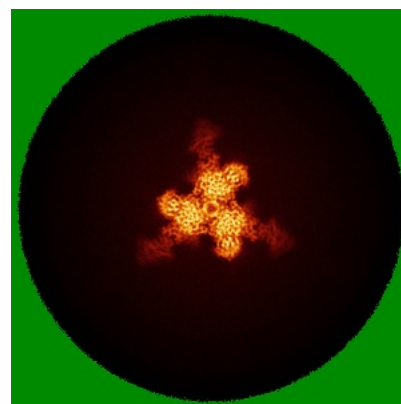
6.4.1 Primary map



X

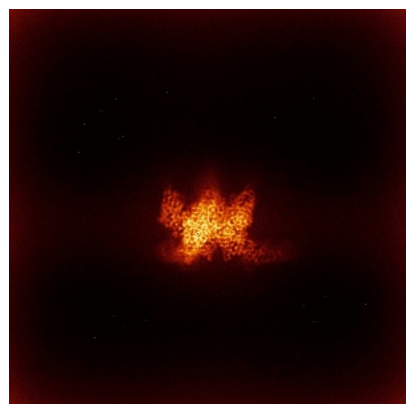


Y

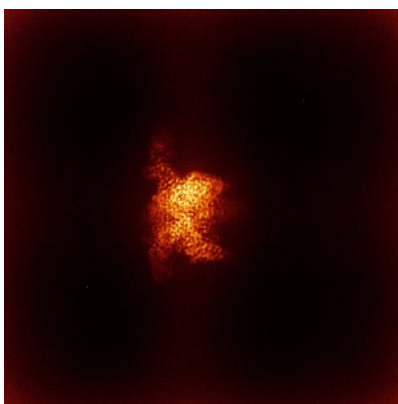


Z

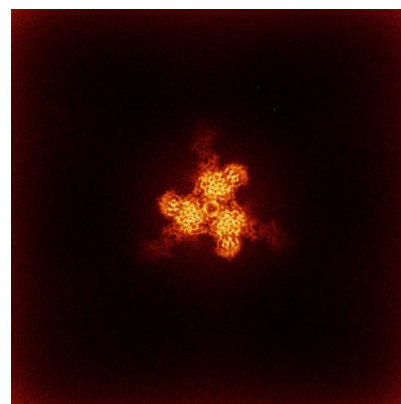
6.4.2 Raw map



X



Y

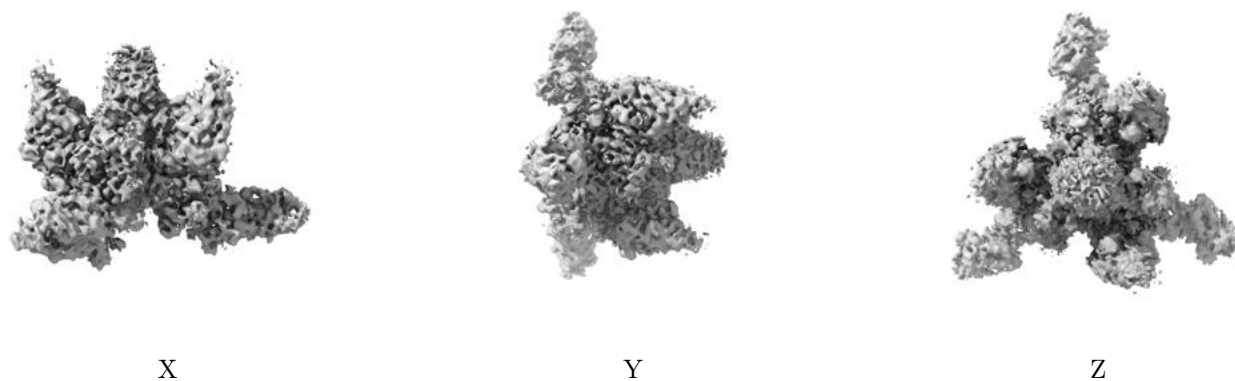


Z

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.

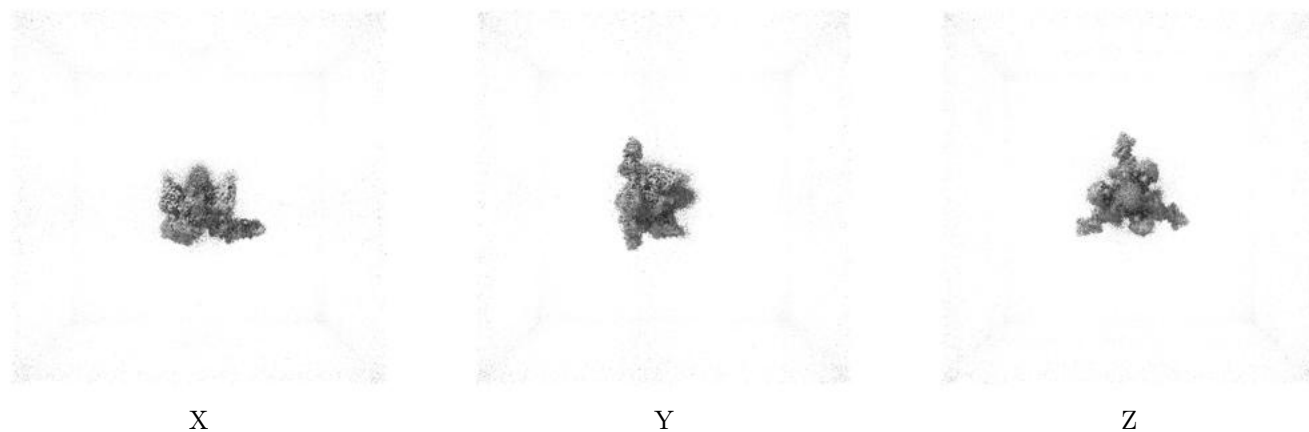
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.05. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

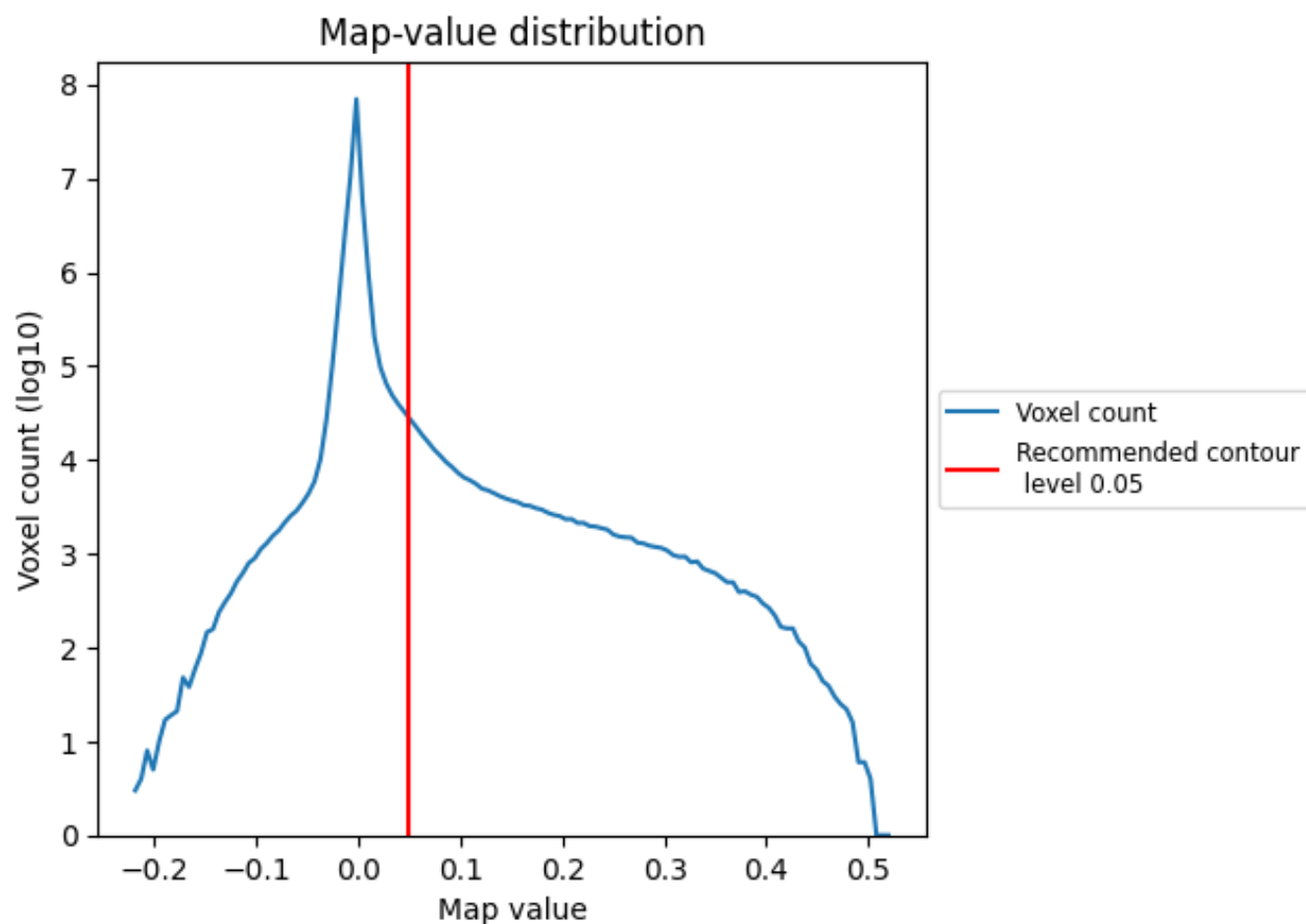
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

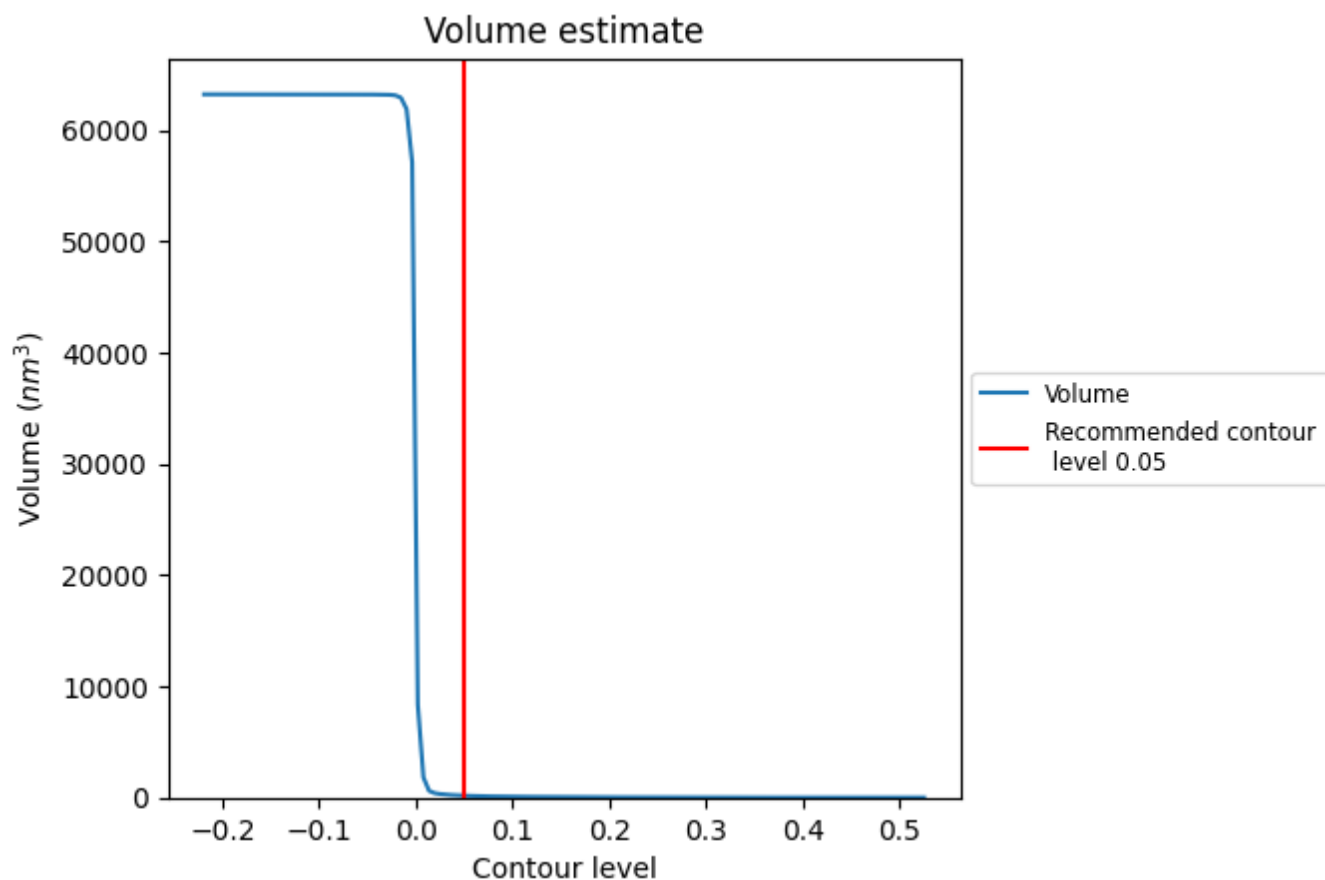
This section contains the results of statistical analysis of the map.

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.

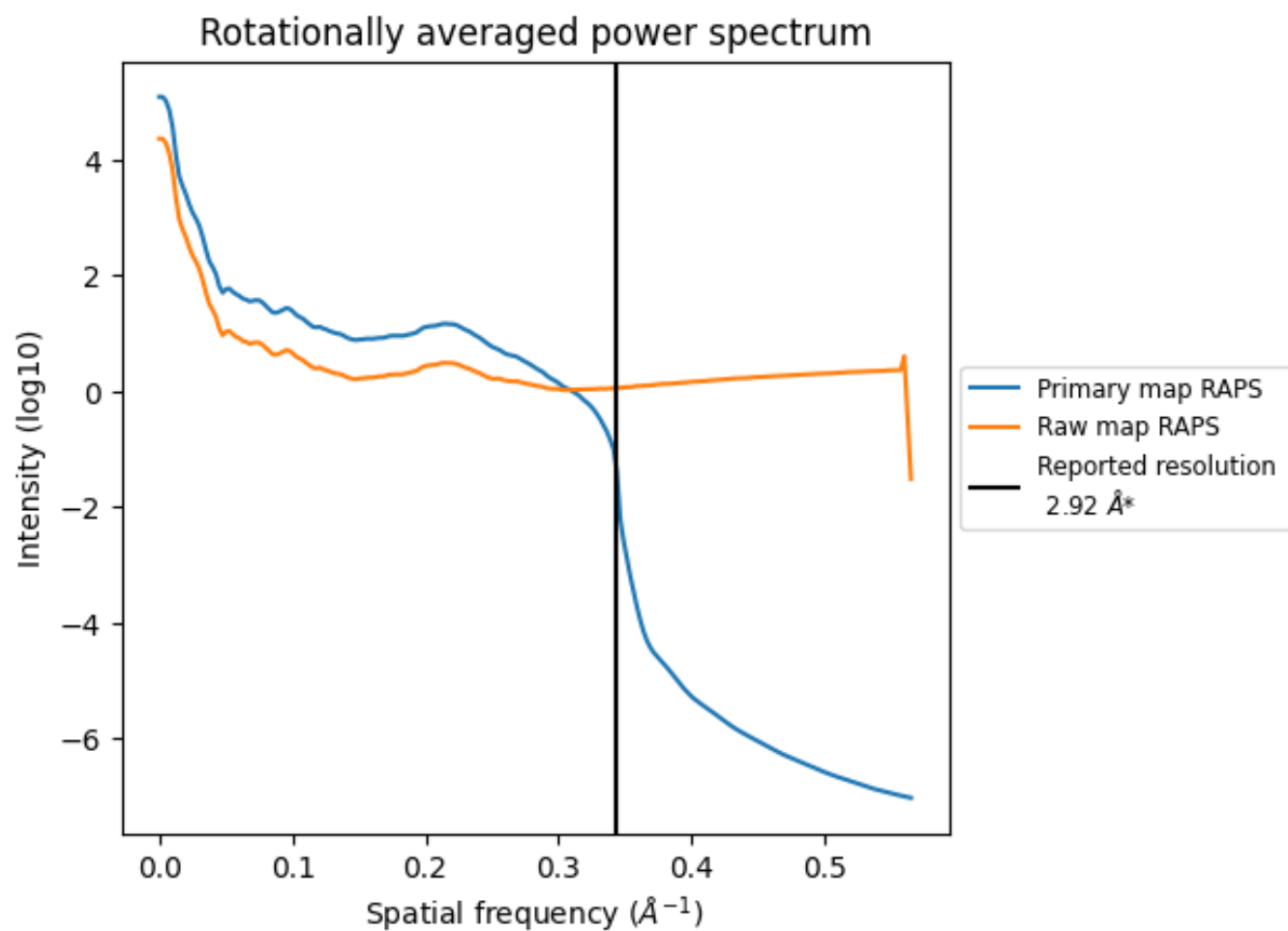
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 171 nm³; this corresponds to an approximate mass of 154 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 ⓘ

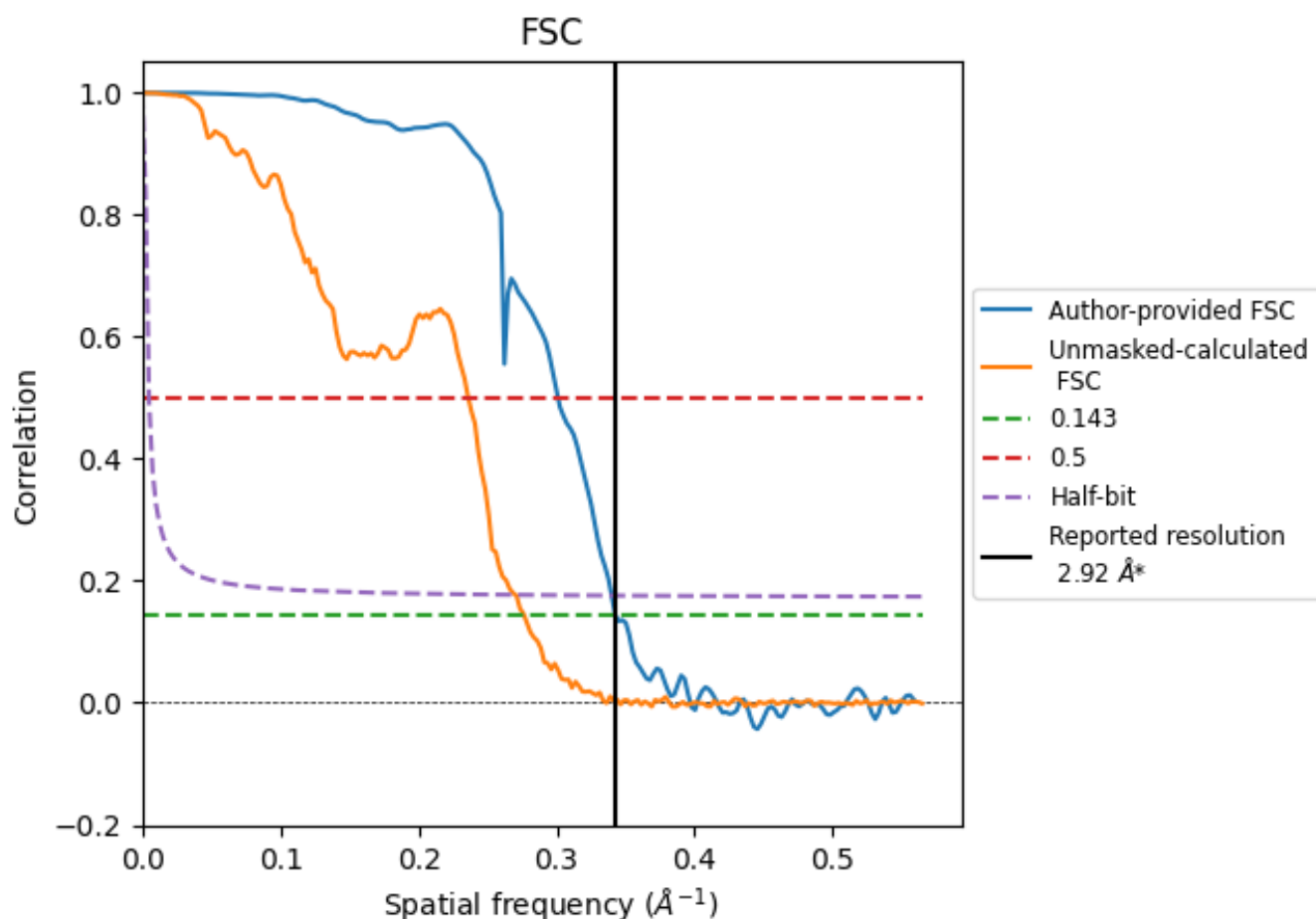


*Reported resolution corresponds to spatial frequency of 0.342 Å⁻¹

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 0.342 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

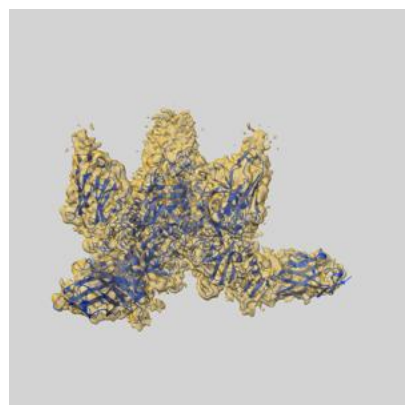
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.92	-	-
Author-provided FSC curve	2.92	3.32	2.94
Unmasked-calculated*	3.61	4.23	3.70

*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 3.61 differs from the reported value 2.92 by more than 10 %

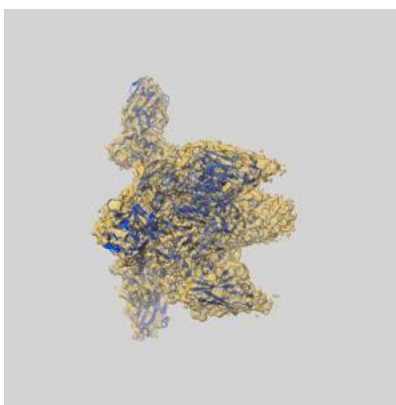
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-71328 and PDB model 9P6X. Per-residue inclusion information can be found in section 3 on page 11.

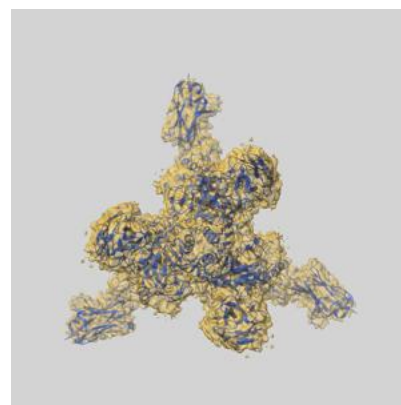
9.1 Map-model overlay [i](#)



X



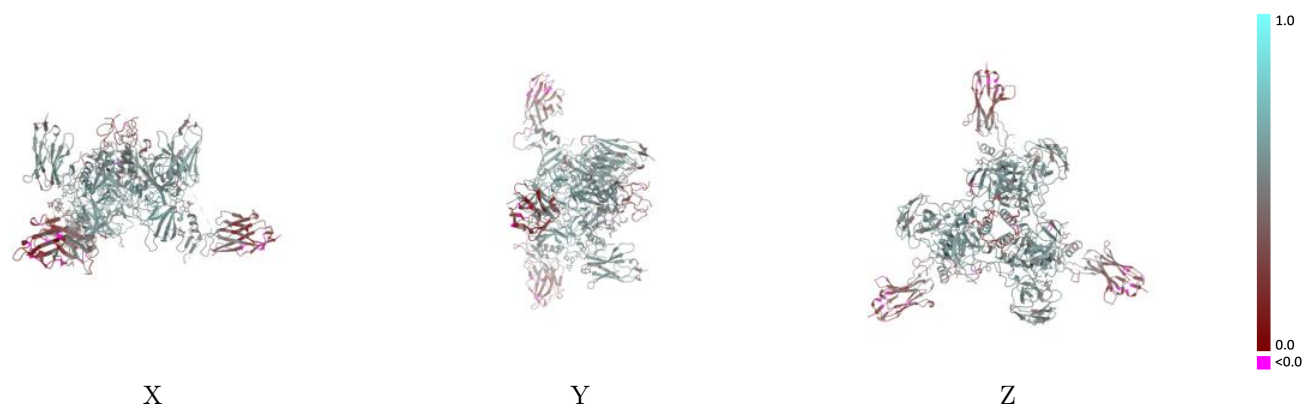
Y



Z

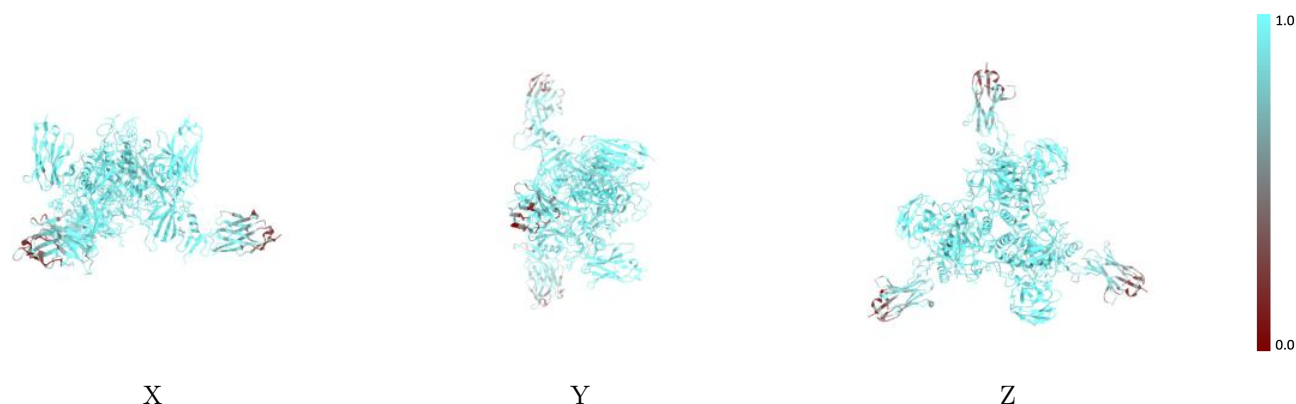
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



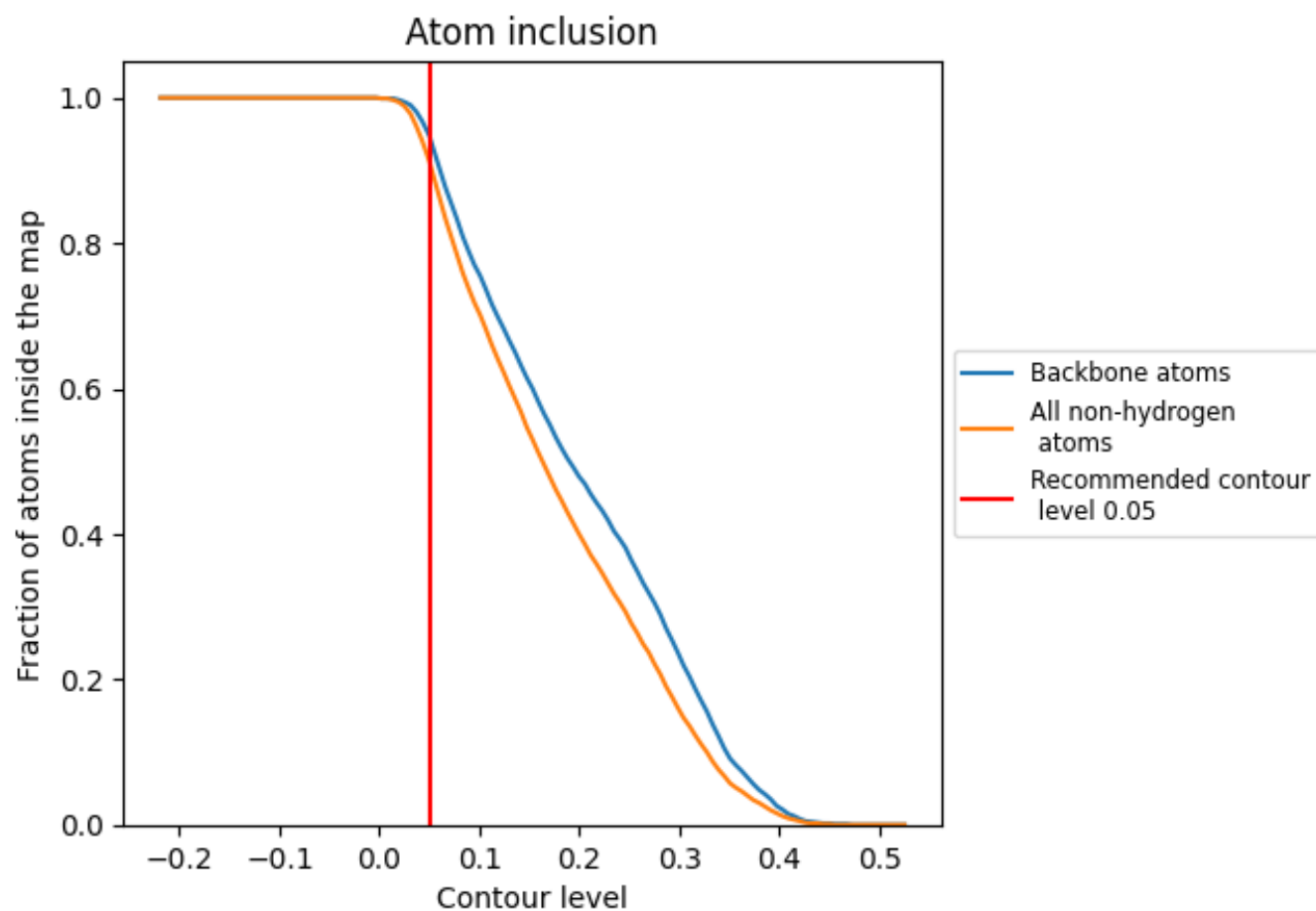
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.05).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 91% 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 (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9130	0.4730
A	0.9620	0.5200
B	0.9560	0.5130
C	0.9680	0.5250
D	0.6950	0.2800
E	0.6930	0.2780
F	0.7030	0.2880
G	0.9750	0.5310
H	0.9760	0.5160
I	0.9830	0.5320
J	0.9860	0.4800
K	0.9440	0.4850
L	1.0000	0.5590

