



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 08:37 AM UTC

PDB ID : 7K37 / pdb_00007k37
EMDB ID : EMD-22652
Title : Structure of full-length influenza HA with a head-binding antibody at pH 7.8
Authors : Gui, M.; Xiang, Y.; Gao, J.
Deposited on : 2020-09-10
Resolution : 2.80 Å (reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

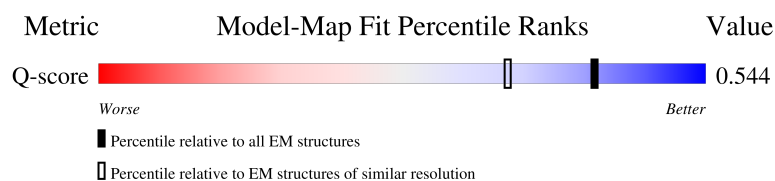
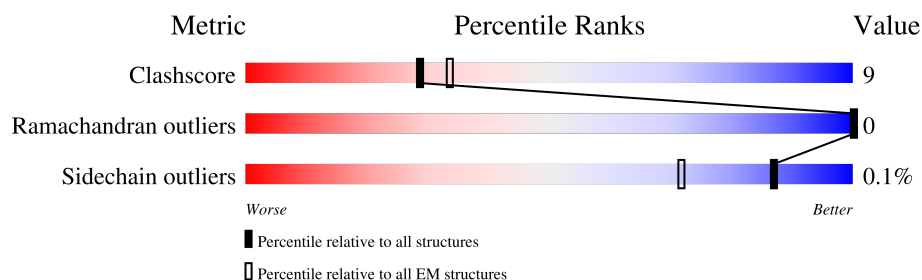
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.80 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.



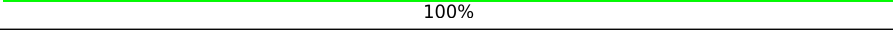
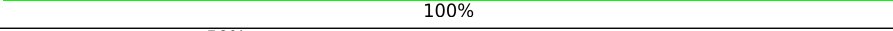
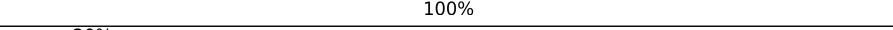
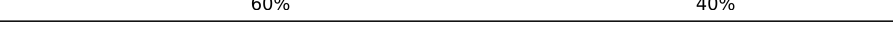
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11806 (2.30 - 3.30)

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	345	 81% 11% 8%
1	C	345	 80% 12% 8%
1	E	345	 80% 12% 8%
2	B	254	 55% 13% 32%

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Mol	Chain	Length	Quality of chain
2	D	254	
2	F	254	
3	G	219	
3	I	219	
3	K	219	
4	H	235	
4	J	235	
4	L	235	
5	M	2	
5	P	2	
5	S	2	
6	N	5	
6	Q	5	
6	T	5	
7	O	8	
7	R	8	
7	U	8	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 21561 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	318	Total	C	N	O	S	0	0
			2454	1537	431	473	13		
1	C	318	Total	C	N	O	S	0	0
			2454	1537	431	473	13		
1	E	318	Total	C	N	O	S	0	0
			2454	1537	431	473	13		

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	172	Total	C	N	O	S	0	0
			1398	867	246	279	6		
2	D	172	Total	C	N	O	S	0	0
			1398	867	246	279	6		
2	F	172	Total	C	N	O	S	0	0
			1398	867	246	279	6		

There are 99 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	222	SER	-	expression tag	UNP C7RWD9
B	223	MET	-	expression tag	UNP C7RWD9
B	224	GLY	-	expression tag	UNP C7RWD9
B	225	TRP	-	expression tag	UNP C7RWD9
B	226	SER	-	expression tag	UNP C7RWD9
B	227	HIS	-	expression tag	UNP C7RWD9
B	228	PRO	-	expression tag	UNP C7RWD9
B	229	GLN	-	expression tag	UNP C7RWD9
B	230	PHE	-	expression tag	UNP C7RWD9
B	231	GLU	-	expression tag	UNP C7RWD9
B	232	LYS	-	expression tag	UNP C7RWD9
B	233	GLY	-	expression tag	UNP C7RWD9
B	234	GLY	-	expression tag	UNP C7RWD9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	235	GLY	-	expression tag	UNP C7RWD9
B	236	ALA	-	expression tag	UNP C7RWD9
B	237	ARG	-	expression tag	UNP C7RWD9
B	238	GLY	-	expression tag	UNP C7RWD9
B	239	GLY	-	expression tag	UNP C7RWD9
B	240	SER	-	expression tag	UNP C7RWD9
B	241	GLY	-	expression tag	UNP C7RWD9
B	242	GLY	-	expression tag	UNP C7RWD9
B	243	GLY	-	expression tag	UNP C7RWD9
B	244	SER	-	expression tag	UNP C7RWD9
B	245	TRP	-	expression tag	UNP C7RWD9
B	246	SER	-	expression tag	UNP C7RWD9
B	247	HIS	-	expression tag	UNP C7RWD9
B	248	PRO	-	expression tag	UNP C7RWD9
B	249	GLN	-	expression tag	UNP C7RWD9
B	250	PHE	-	expression tag	UNP C7RWD9
B	251	GLU	-	expression tag	UNP C7RWD9
B	252	LYS	-	expression tag	UNP C7RWD9
B	253	GLY	-	expression tag	UNP C7RWD9
B	254	PHE	-	expression tag	UNP C7RWD9
D	222	SER	-	expression tag	UNP C7RWD9
D	223	MET	-	expression tag	UNP C7RWD9
D	224	GLY	-	expression tag	UNP C7RWD9
D	225	TRP	-	expression tag	UNP C7RWD9
D	226	SER	-	expression tag	UNP C7RWD9
D	227	HIS	-	expression tag	UNP C7RWD9
D	228	PRO	-	expression tag	UNP C7RWD9
D	229	GLN	-	expression tag	UNP C7RWD9
D	230	PHE	-	expression tag	UNP C7RWD9
D	231	GLU	-	expression tag	UNP C7RWD9
D	232	LYS	-	expression tag	UNP C7RWD9
D	233	GLY	-	expression tag	UNP C7RWD9
D	234	GLY	-	expression tag	UNP C7RWD9
D	235	GLY	-	expression tag	UNP C7RWD9
D	236	ALA	-	expression tag	UNP C7RWD9
D	237	ARG	-	expression tag	UNP C7RWD9
D	238	GLY	-	expression tag	UNP C7RWD9
D	239	GLY	-	expression tag	UNP C7RWD9
D	240	SER	-	expression tag	UNP C7RWD9
D	241	GLY	-	expression tag	UNP C7RWD9
D	242	GLY	-	expression tag	UNP C7RWD9
D	243	GLY	-	expression tag	UNP C7RWD9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	244	SER	-	expression tag	UNP C7RWD9
D	245	TRP	-	expression tag	UNP C7RWD9
D	246	SER	-	expression tag	UNP C7RWD9
D	247	HIS	-	expression tag	UNP C7RWD9
D	248	PRO	-	expression tag	UNP C7RWD9
D	249	GLN	-	expression tag	UNP C7RWD9
D	250	PHE	-	expression tag	UNP C7RWD9
D	251	GLU	-	expression tag	UNP C7RWD9
D	252	LYS	-	expression tag	UNP C7RWD9
D	253	GLY	-	expression tag	UNP C7RWD9
D	254	PHE	-	expression tag	UNP C7RWD9
F	222	SER	-	expression tag	UNP C7RWD9
F	223	MET	-	expression tag	UNP C7RWD9
F	224	GLY	-	expression tag	UNP C7RWD9
F	225	TRP	-	expression tag	UNP C7RWD9
F	226	SER	-	expression tag	UNP C7RWD9
F	227	HIS	-	expression tag	UNP C7RWD9
F	228	PRO	-	expression tag	UNP C7RWD9
F	229	GLN	-	expression tag	UNP C7RWD9
F	230	PHE	-	expression tag	UNP C7RWD9
F	231	GLU	-	expression tag	UNP C7RWD9
F	232	LYS	-	expression tag	UNP C7RWD9
F	233	GLY	-	expression tag	UNP C7RWD9
F	234	GLY	-	expression tag	UNP C7RWD9
F	235	GLY	-	expression tag	UNP C7RWD9
F	236	ALA	-	expression tag	UNP C7RWD9
F	237	ARG	-	expression tag	UNP C7RWD9
F	238	GLY	-	expression tag	UNP C7RWD9
F	239	GLY	-	expression tag	UNP C7RWD9
F	240	SER	-	expression tag	UNP C7RWD9
F	241	GLY	-	expression tag	UNP C7RWD9
F	242	GLY	-	expression tag	UNP C7RWD9
F	243	GLY	-	expression tag	UNP C7RWD9
F	244	SER	-	expression tag	UNP C7RWD9
F	245	TRP	-	expression tag	UNP C7RWD9
F	246	SER	-	expression tag	UNP C7RWD9
F	247	HIS	-	expression tag	UNP C7RWD9
F	248	PRO	-	expression tag	UNP C7RWD9
F	249	GLN	-	expression tag	UNP C7RWD9
F	250	PHE	-	expression tag	UNP C7RWD9
F	251	GLU	-	expression tag	UNP C7RWD9
F	252	LYS	-	expression tag	UNP C7RWD9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	253	GLY	-	expression tag	UNP C7RWD9
F	254	PHE	-	expression tag	UNP C7RWD9

- Molecule 3 is a protein called antibody Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	202	Total	C	N	O	S	2	0
			1502	932	258	308	4		
3	I	202	Total	C	N	O	S	2	0
			1502	932	258	308	4		
3	K	202	Total	C	N	O	S	2	0
			1502	932	258	308	4		

- Molecule 4 is a protein called antibody Fab heavy chain.

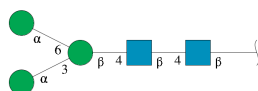
Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	212	Total	C	N	O	S	0	0
			1608	1022	269	307	10		
4	J	212	Total	C	N	O	S	0	0
			1608	1022	269	307	10		
4	L	212	Total	C	N	O	S	0	0
			1608	1022	269	307	10		

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



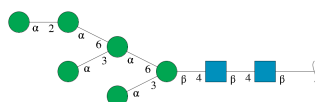
Mol	Chain	Residues	Atoms				AltConf	Trace
5	M	2	Total	C	N	O	0	0
			28	16	2	10		
5	P	2	Total	C	N	O	0	0
			28	16	2	10		
5	S	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 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
6	N	5	Total	C	N	O	0	0
			61	34	2	25		
6	Q	5	Total	C	N	O	0	0
			61	34	2	25		
6	T	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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
7	O	8	Total	C	N	O	0	0
			94	52	2	40		
7	R	8	Total	C	N	O	0	0
			94	52	2	40		
7	U	8	Total	C	N	O	0	0
			94	52	2	40		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



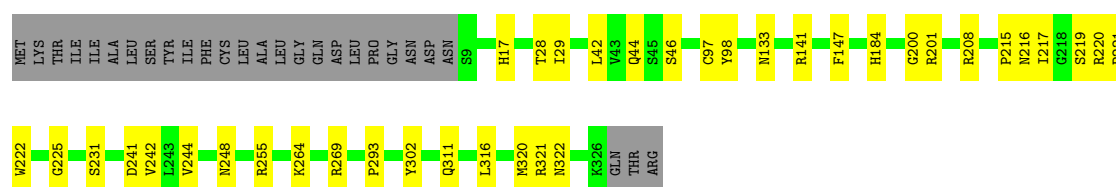
Mol	Chain	Residues	Atoms				AltConf
8	A	1	Total	C	N	O	0
			14	8	1	5	
8	A	1	Total	C	N	O	0
			14	8	1	5	
8	B	1	Total	C	N	O	0
			14	8	1	5	
8	C	1	Total	C	N	O	0
			14	8	1	5	
8	C	1	Total	C	N	O	0
			14	8	1	5	
8	D	1	Total	C	N	O	0
			14	8	1	5	
8	E	1	Total	C	N	O	0
			14	8	1	5	
8	E	1	Total	C	N	O	0
			14	8	1	5	
8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

3 Residue-property plots [i](#)

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

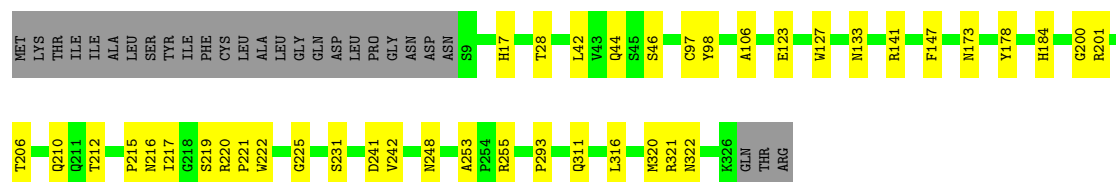
• Molecule 1: Hemagglutinin

Chain A: 



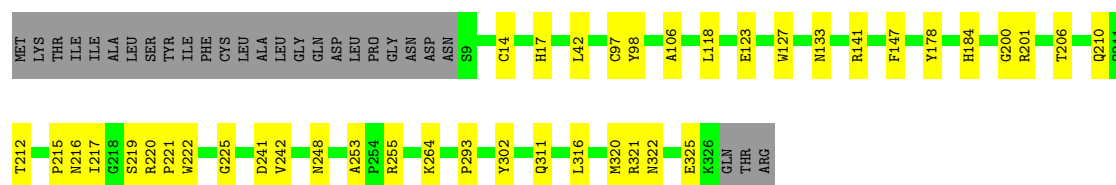
• Molecule 1: Hemagglutinin

Chain C: 



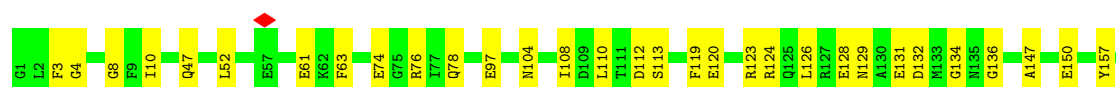
• Molecule 1: Hemagglutinin

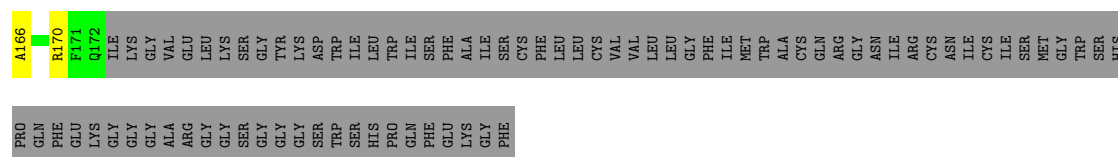
Chain E: 



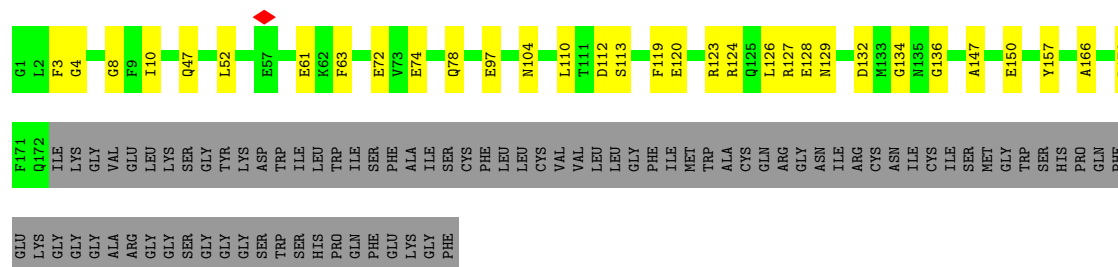
• Molecule 2: Hemagglutinin

Chain B: 

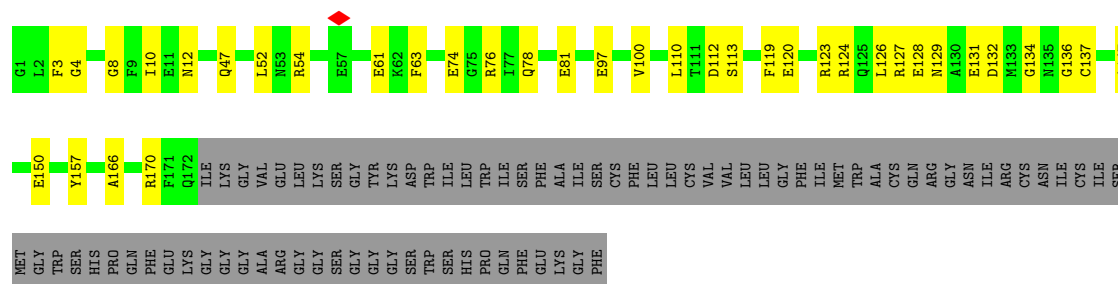




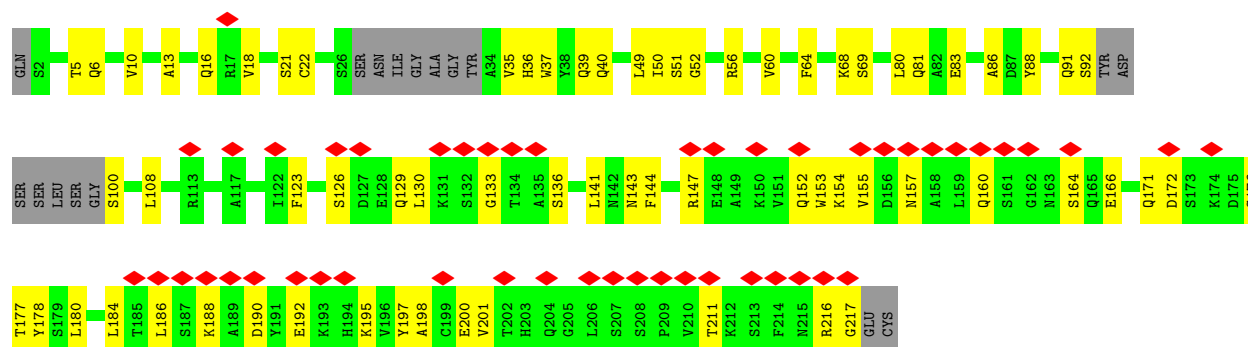
• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin

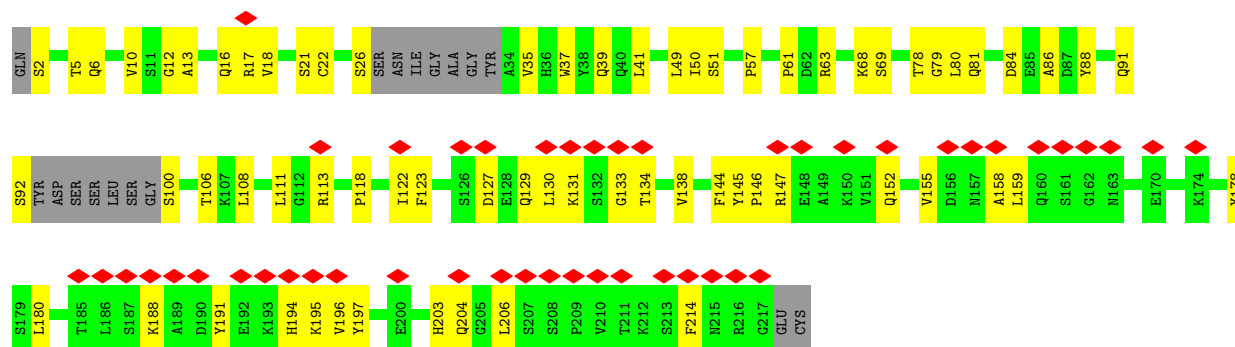


• Molecule 3: antibody Fab light chain

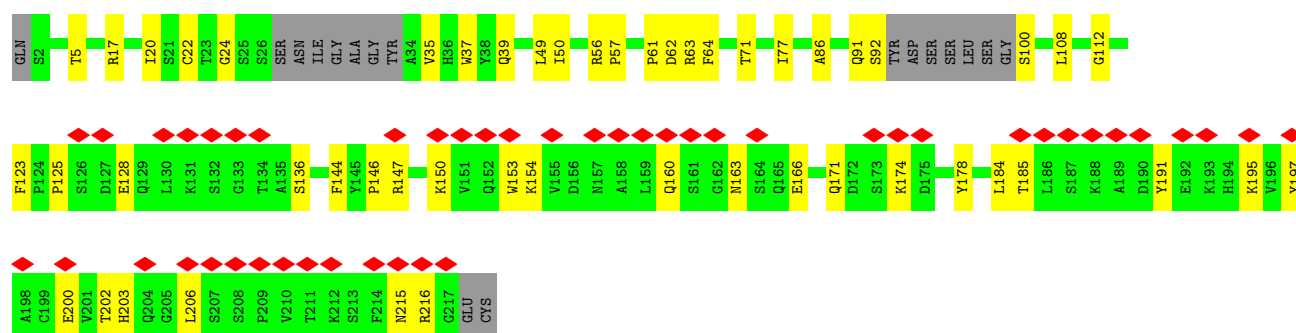


• Molecule 3: antibody Fab light chain

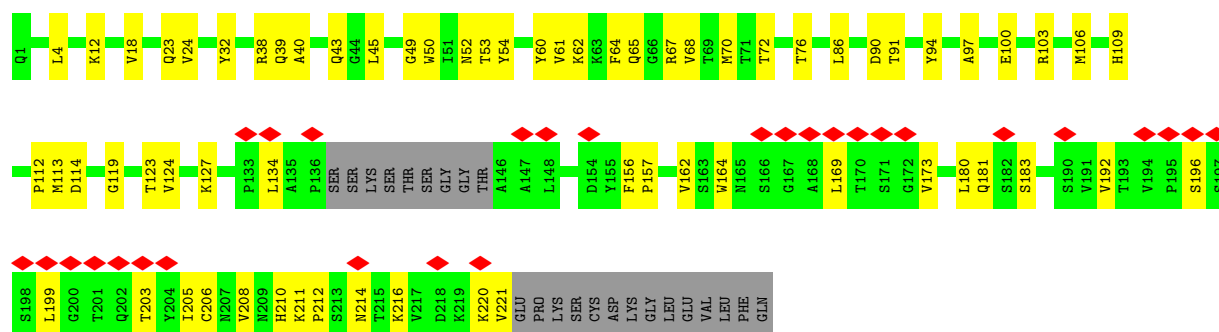




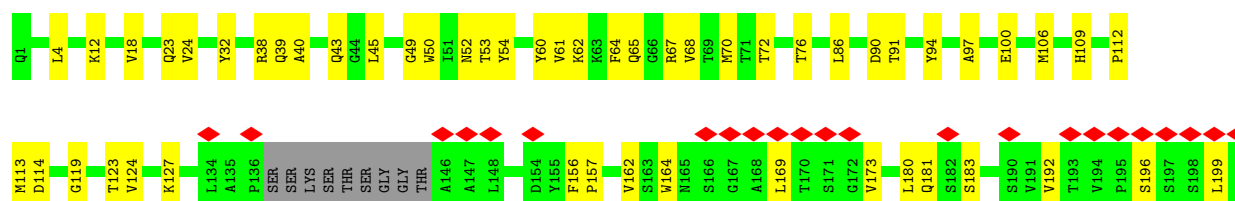
• Molecule 3: antibody Fab light chain

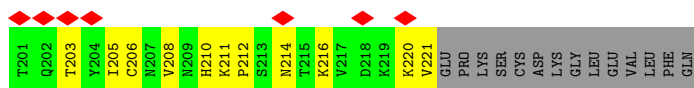


• Molecule 4: antibody Fab heavy chain

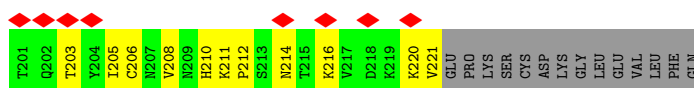
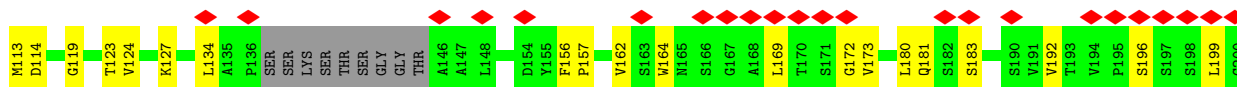


• Molecule 4: antibody Fab heavy chain





- Molecule 4: antibody Fab heavy chain



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

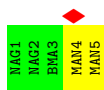


- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

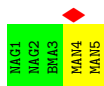


- Molecule 6: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

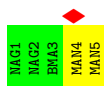




- Molecule 6: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 6: 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



- Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]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



MAG1
MAG2
EM/A3
MAN4
MAN5
MAN6
MAN7
MAN8

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	359922	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 ($e^-/\text{\AA}^2$)	50	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	0.607	Depositor
Minimum map value	-0.306	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	399.0, 399.0, 399.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.33, 1.33, 1.33	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: NAG, BMA, MAN

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	0.26	0/2510	0.34	0/3420
1	C	0.25	0/2510	0.33	0/3420
1	E	0.26	0/2510	0.33	0/3420
2	B	0.24	0/1422	0.36	0/1911
2	D	0.24	0/1422	0.36	0/1911
2	F	0.25	0/1422	0.36	0/1911
3	G	0.17	0/1531	0.41	0/2079
3	I	0.17	0/1531	0.43	0/2079
3	K	0.16	0/1531	0.40	0/2079
4	H	0.21	0/1648	0.42	0/2246
4	J	0.21	0/1648	0.41	0/2246
4	L	0.21	0/1648	0.41	0/2246
All	All	0.23	0/21333	0.38	0/28968

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	2454	0	2403	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2454	0	2403	31	0
1	E	2454	0	2403	35	0
2	B	1398	0	1316	22	0
2	D	1398	0	1316	20	0
2	F	1398	0	1316	25	0
3	G	1502	0	1452	51	0
3	I	1502	0	1452	53	0
3	K	1502	0	1452	36	0
4	H	1608	0	1584	47	0
4	J	1608	0	1584	44	0
4	L	1608	0	1584	44	0
5	M	28	0	25	0	0
5	P	28	0	25	0	0
5	S	28	0	25	0	0
6	N	61	0	52	0	0
6	Q	61	0	52	0	0
6	T	61	0	52	0	0
7	O	94	0	79	1	0
7	R	94	0	79	0	0
7	U	94	0	79	0	0
8	A	28	0	26	0	0
8	B	14	0	13	1	0
8	C	28	0	26	0	0
8	D	14	0	13	1	0
8	E	28	0	26	0	0
8	F	14	0	13	1	0
All	All	21561	0	20850	400	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (400) 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)
3:I:118:PRO:HB3	3:I:144:PHE:HB3	1.65	0.79
3:I:17:ARG:HH11	3:I:79:GLY:H	1.34	0.76
2:B:74:GLU:OE1	2:B:78:GLN:NE2	2.20	0.75
2:D:120:GLU:OE1	2:D:123:ARG:NH1	2.20	0.75
1:A:293:PRO:HG3	2:B:52:LEU:HD21	1.68	0.75
2:F:120:GLU:OE1	2:F:123:ARG:NH1	2.20	0.75
2:B:120:GLU:OE1	2:B:123:ARG:NH1	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:74:GLU:OE1	2:F:78:GLN:NE2	2.20	0.74
2:D:74:GLU:OE1	2:D:78:GLN:NE2	2.20	0.74
3:I:147:ARG:HB3	3:I:178:TYR:HD2	1.53	0.73
3:G:37:TRP:HB2	3:G:50:ILE:HB	1.69	0.73
3:I:63:ARG:HH12	3:I:81:GLN:HG3	1.53	0.71
3:I:17:ARG:HE	3:I:78:THR:HA	1.55	0.70
4:L:67:ARG:NH2	4:L:90:ASP:OD2	2.25	0.70
3:I:155:VAL:HG23	3:I:158:ALA:HB3	1.73	0.70
4:H:67:ARG:NH2	4:H:90:ASP:OD2	2.25	0.70
4:J:181:GLN:NE2	4:J:183:SER:OG	2.22	0.70
3:I:146:PRO:HD2	3:I:203:HIS:CE1	2.27	0.69
4:H:181:GLN:NE2	4:H:183:SER:OG	2.22	0.69
4:J:67:ARG:NH2	4:J:90:ASP:OD2	2.25	0.69
1:E:293:PRO:HG3	2:F:52:LEU:HD21	1.75	0.69
3:G:129:GLN:OE1	3:G:136:SER:OG	2.11	0.68
1:C:201:ARG:NH2	1:E:217:ILE:O	2.27	0.68
4:L:181:GLN:NE2	4:L:183:SER:OG	2.22	0.68
4:L:214:ASN:OD1	4:L:216:LYS:NZ	2.27	0.68
1:A:311:GLN:NE2	2:B:97:GLU:OE1	2.26	0.67
3:K:17:ARG:HA	3:K:77:ILE:O	1.94	0.67
4:H:214:ASN:OD1	4:H:216:LYS:NZ	2.27	0.67
4:J:214:ASN:OD1	4:J:216:LYS:NZ	2.27	0.67
4:L:12:LYS:HB2	4:L:124:VAL:HG22	1.78	0.66
4:H:12:LYS:HB2	4:H:124:VAL:HG22	1.78	0.65
3:G:166:GLU:HB3	3:G:180:LEU:HD11	1.78	0.65
4:H:97:ALA:HB1	4:H:113:MET:HB3	1.79	0.65
1:C:212:THR:HB	1:E:216:ASN:HD22	1.61	0.65
4:J:97:ALA:HB1	4:J:113:MET:HB3	1.79	0.65
4:L:97:ALA:HB1	4:L:113:MET:HB3	1.79	0.65
4:J:12:LYS:HB2	4:J:124:VAL:HG22	1.78	0.64
3:K:160:GLN:HE21	3:K:184:LEU:HD21	1.61	0.64
3:G:133:GLY:HA2	3:G:188:LYS:HB2	1.78	0.64
4:H:169:LEU:HD13	4:H:192:VAL:HG11	1.80	0.64
4:J:169:LEU:HD13	4:J:192:VAL:HG11	1.80	0.64
3:I:41:LEU:HD23	3:I:86:ALA:HB2	1.77	0.64
4:L:169:LEU:HD13	4:L:192:VAL:HG11	1.80	0.64
1:C:293:PRO:HG3	2:D:52:LEU:HD21	1.78	0.64
4:J:4:LEU:HD22	4:J:24:VAL:HG12	1.80	0.64
4:H:4:LEU:HD22	4:H:24:VAL:HG12	1.80	0.64
2:B:124:ARG:HD3	2:D:134:GLY:HA2	1.80	0.63
3:I:17:ARG:NH1	3:I:79:GLY:H	1.96	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:133:ASN:OD1	1:E:255:ARG:NH2	2.31	0.62
3:I:144:PHE:HB2	3:I:203:HIS:CE1	2.34	0.62
3:I:130:LEU:HA	3:I:188:LYS:HZ2	1.63	0.62
4:L:4:LEU:HD22	4:L:24:VAL:HG12	1.80	0.62
1:E:216:ASN:OD1	1:E:220:ARG:NH2	2.34	0.61
1:C:216:ASN:OD1	1:C:220:ARG:NH2	2.33	0.61
4:H:210:HIS:CD2	4:H:212:PRO:HD2	2.36	0.61
3:I:130:LEU:HD13	3:I:188:LYS:HZ3	1.66	0.61
4:L:210:HIS:CD2	4:L:212:PRO:HD2	2.36	0.61
3:K:112:GLY:H	3:K:171:GLN:HE21	1.47	0.61
1:A:216:ASN:OD1	1:A:220:ARG:NH2	2.34	0.61
4:J:210:HIS:CD2	4:J:212:PRO:HD2	2.36	0.60
4:H:49:GLY:HA3	4:H:70:MET:HE1	1.84	0.60
3:K:24:GLY:O	3:K:71:THR:OG1	2.19	0.60
4:L:49:GLY:HA3	4:L:70:MET:HE1	1.84	0.60
3:I:147:ARG:HB3	3:I:178:TYR:CD2	2.36	0.59
4:J:49:GLY:HA3	4:J:70:MET:HE1	1.84	0.59
3:I:92:SER:O	3:I:100:SER:OG	2.21	0.59
3:K:92:SER:O	3:K:100:SER:OG	2.20	0.59
1:C:133:ASN:OD1	1:C:255:ARG:NH2	2.31	0.59
3:G:18:VAL:HG11	3:G:108:LEU:HD11	1.83	0.59
1:E:311:GLN:NE2	2:F:97:GLU:OE1	2.36	0.58
4:L:205:ILE:HA	4:L:220:LYS:HA	1.85	0.58
4:H:53:THR:HG22	4:H:72:THR:HG21	1.86	0.58
3:I:2:SER:N	3:I:26:SER:HG	2.01	0.58
4:J:205:ILE:HA	4:J:220:LYS:HA	1.85	0.58
1:A:133:ASN:OD1	1:A:255:ARG:NH2	2.31	0.58
4:J:53:THR:HG22	4:J:72:THR:HG21	1.86	0.58
3:I:129:GLN:NE2	3:I:134:THR:O	2.37	0.57
3:K:163:ASN:ND2	3:K:185:THR:O	2.21	0.57
4:H:205:ILE:HA	4:H:220:LYS:HA	1.85	0.57
3:I:63:ARG:NH2	3:I:84:ASP:OD1	2.37	0.57
4:J:23:GLN:NE2	4:J:76:THR:O	2.32	0.57
3:K:37:TRP:HB2	3:K:50:ILE:HG12	1.86	0.57
3:K:150:LYS:O	3:K:202:THR:OG1	2.19	0.57
1:C:212:THR:CB	1:E:216:ASN:HD22	2.17	0.57
4:L:53:THR:HG22	4:L:72:THR:HG21	1.86	0.57
1:A:208:ARG:HH22	2:D:72:GLU:CD	2.13	0.57
3:G:154:LYS:HB2	3:G:198:ALA:HB3	1.87	0.57
2:D:124:ARG:HD3	2:F:134:GLY:HA2	1.87	0.57
3:I:122:ILE:HG21	3:I:214:PHE:CD1	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:146:PRO:HD2	3:K:203:HIS:CE1	2.41	0.56
3:K:195:LYS:O	3:K:216:ARG:NH1	2.38	0.56
3:I:152:GLN:HB3	3:I:159:LEU:HD11	1.86	0.56
3:G:92:SER:O	3:G:100:SER:OG	2.23	0.56
4:L:53:THR:O	4:L:54:TYR:HB3	2.06	0.56
3:I:145:TYR:HB3	3:I:146:PRO:HD3	1.87	0.56
1:C:123:GLU:OE2	1:C:178:TYR:OH	2.20	0.56
3:I:13:ALA:O	3:I:16:GLN:NE2	2.39	0.55
4:J:53:THR:O	4:J:54:TYR:HB3	2.06	0.55
3:I:113:ARG:HA	3:I:145:TYR:OH	2.06	0.55
4:H:53:THR:O	4:H:54:TYR:HB3	2.06	0.55
3:I:155:VAL:HG12	3:I:197:TYR:CD1	2.41	0.55
1:A:216:ASN:HD22	1:E:212:THR:HB	1.71	0.55
4:H:54:TYR:O	7:O:5:MAN:O6	2.17	0.55
1:C:106:ALA:HB3	2:F:76:ARG:HD2	1.88	0.54
8:F:301:NAG:H83	8:F:301:NAG:H3	1.90	0.54
8:D:301:NAG:H83	8:D:301:NAG:H3	1.90	0.54
3:I:195:LYS:HG3	3:I:196:VAL:HG23	1.89	0.54
2:F:47:GLN:HB3	2:F:110:LEU:HD11	1.90	0.53
3:G:49:LEU:HD21	3:G:88:TYR:HE2	1.74	0.53
2:B:128:GLU:O	2:B:170:ARG:NH1	2.41	0.53
8:B:301:NAG:H83	8:B:301:NAG:H3	1.90	0.53
1:C:311:GLN:NE2	2:D:97:GLU:OE1	2.41	0.53
3:I:39:GLN:HB3	3:I:49:LEU:CD2	2.39	0.53
4:J:60:TYR:CE1	4:J:70:MET:HE3	2.44	0.53
4:L:60:TYR:CE1	4:L:70:MET:HE3	2.44	0.53
2:D:127:ARG:NH2	2:F:131:GLU:OE1	2.31	0.53
3:G:68:LYS:HG2	3:G:69:SER:O	2.09	0.53
4:H:60:TYR:CE1	4:H:70:MET:HE3	2.44	0.53
1:A:97:CYS:SG	1:A:98:TYR:N	2.82	0.53
3:G:13:ALA:N	3:G:16:GLN:HE22	2.07	0.52
1:C:97:CYS:SG	1:C:98:TYR:N	2.82	0.52
2:B:131:GLU:OE1	2:F:127:ARG:NH2	2.30	0.52
3:G:130:LEU:O	3:G:188:LYS:HE3	2.09	0.52
1:C:42:LEU:HD11	1:C:316:LEU:HB2	1.92	0.52
2:D:47:GLN:HB3	2:D:110:LEU:HD11	1.90	0.52
2:D:128:GLU:O	2:D:170:ARG:NH1	2.41	0.52
3:G:216:ARG:NH2	3:G:217:GLY:O	2.41	0.52
4:H:210:HIS:HD2	4:H:212:PRO:HD2	1.75	0.52
4:J:40:ALA:HB3	4:J:43:GLN:HB2	1.92	0.52
3:G:155:VAL:HG22	3:G:197:TYR:CE2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:61:PRO:HB2	3:I:63:ARG:HG2	1.92	0.52
2:D:3:PHE:HB2	2:D:112:ASP:OD1	2.10	0.52
4:H:40:ALA:HB3	4:H:43:GLN:HB2	1.92	0.52
3:I:191:TYR:O	3:I:197:TYR:OH	2.25	0.52
3:K:39:GLN:HB3	3:K:49:LEU:CD2	2.39	0.52
1:A:201:ARG:NH2	1:C:217:ILE:O	2.43	0.51
2:B:3:PHE:HB2	2:B:112:ASP:OD1	2.10	0.51
4:L:40:ALA:HB3	4:L:43:GLN:HB2	1.92	0.51
2:B:47:GLN:HB3	2:B:110:LEU:HD11	1.91	0.51
1:E:42:LEU:HD11	1:E:316:LEU:HB2	1.92	0.51
3:G:172:ASP:HB2	3:G:176:SER:H	1.75	0.51
3:I:130:LEU:HA	3:I:188:LYS:NZ	2.26	0.51
3:K:154:LYS:HD3	3:K:200:GLU:HB2	1.93	0.51
4:L:94:TYR:O	4:L:119:GLY:HA2	2.11	0.51
3:G:13:ALA:O	3:G:16:GLN:NE2	2.44	0.51
4:J:94:TYR:O	4:J:119:GLY:HA2	2.11	0.51
1:A:42:LEU:HD11	1:A:316:LEU:HB2	1.92	0.51
2:F:3:PHE:HB2	2:F:112:ASP:OD1	2.10	0.51
1:C:28:THR:HG22	2:D:104:ASN:HB3	1.92	0.51
3:G:143:ASN:HD22	3:G:177:THR:HG21	1.76	0.51
1:C:222:TRP:CZ2	1:C:225:GLY:HA2	2.46	0.51
3:G:10:VAL:HG11	3:G:18:VAL:HG13	1.93	0.51
4:H:157:PRO:HD2	4:H:210:HIS:CE1	2.46	0.51
4:L:106:MET:HE3	4:L:109:HIS:NE2	2.26	0.51
4:L:210:HIS:HD2	4:L:212:PRO:HD2	1.75	0.51
2:F:128:GLU:O	2:F:170:ARG:NH1	2.41	0.50
3:I:39:GLN:HB3	3:I:49:LEU:HD21	1.93	0.50
4:J:86:LEU:HD13	4:J:124:VAL:HG21	1.93	0.50
3:K:5:THR:O	3:K:22:CYS:HA	2.11	0.50
1:E:141:ARG:NH1	1:E:147:PHE:O	2.40	0.50
1:E:97:CYS:SG	1:E:98:TYR:N	2.82	0.50
2:F:3:PHE:CD2	2:F:113:SER:HA	2.46	0.50
4:H:86:LEU:HD13	4:H:124:VAL:HG21	1.93	0.50
4:J:157:PRO:HD2	4:J:210:HIS:CE1	2.46	0.50
4:L:157:PRO:HD2	4:L:210:HIS:CE1	2.46	0.50
4:J:210:HIS:HD2	4:J:212:PRO:HD2	1.75	0.50
3:K:49:LEU:O	3:K:57:PRO:HD2	2.10	0.50
2:B:3:PHE:CD2	2:B:113:SER:HA	2.46	0.50
1:E:222:TRP:CZ2	1:E:225:GLY:HA2	2.47	0.50
3:G:141:LEU:HD21	3:G:201:VAL:HB	1.94	0.50
3:K:203:HIS:HB3	3:K:206:LEU:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:3:PHE:CD2	2:D:113:SER:HA	2.46	0.50
4:H:106:MET:HE3	4:H:109:HIS:NE2	2.26	0.50
3:K:144:PHE:HB2	3:K:203:HIS:CE1	2.47	0.50
4:J:106:MET:HE3	4:J:109:HIS:NE2	2.26	0.49
1:A:241:ASP:OD1	1:A:242:VAL:N	2.40	0.49
4:H:94:TYR:O	4:H:119:GLY:HA2	2.11	0.49
3:K:191:TYR:O	3:K:197:TYR:OH	2.30	0.49
1:A:217:ILE:O	1:E:201:ARG:NH2	2.45	0.49
3:I:146:PRO:HG2	3:I:204:GLN:HB2	1.93	0.49
4:L:86:LEU:HD13	4:L:124:VAL:HG21	1.93	0.49
1:A:222:TRP:CZ2	1:A:225:GLY:HA2	2.47	0.49
3:G:126:SER:O	3:G:130:LEU:HG	2.12	0.49
3:G:147:ARG:HB2	3:G:178:TYR:HE2	1.77	0.49
4:H:114:ASP:OD1	4:H:114:ASP:N	2.45	0.49
3:I:10:VAL:HG11	3:I:18:VAL:HG13	1.94	0.49
3:K:39:GLN:HB3	3:K:49:LEU:HD21	1.95	0.49
4:L:23:GLN:NE2	4:L:76:THR:O	2.33	0.49
1:E:123:GLU:OE2	1:E:178:TYR:OH	2.20	0.49
3:G:160:GLN:HE22	3:G:164:SER:HB3	1.78	0.49
1:A:141:ARG:NH1	1:A:147:PHE:O	2.40	0.49
3:G:155:VAL:HG23	3:G:160:GLN:HB2	1.95	0.49
3:I:197:TYR:HB2	3:I:214:PHE:CE1	2.47	0.49
3:G:81:GLN:NE2	3:G:83:GLU:OE2	2.44	0.48
1:E:325:GLU:HG3	2:F:12:ASN:HD21	1.78	0.48
4:L:114:ASP:N	4:L:114:ASP:OD1	2.45	0.48
2:B:134:GLY:HA2	2:F:124:ARG:HD3	1.94	0.48
3:G:35:VAL:HA	3:G:91:GLN:O	2.12	0.48
4:H:211:LYS:HB2	4:H:212:PRO:HD3	1.96	0.48
3:K:37:TRP:CD1	3:K:50:ILE:HD11	2.48	0.48
1:C:210:GLN:HE22	1:E:184:HIS:CD2	2.32	0.48
4:J:43:GLN:OE1	4:J:180:LEU:HD21	2.14	0.48
4:L:43:GLN:OE1	4:L:180:LEU:HD21	2.14	0.48
3:I:203:HIS:CD2	3:I:206:LEU:H	2.32	0.48
1:C:210:GLN:OE1	1:E:220:ARG:NE	2.46	0.48
3:G:49:LEU:HD21	3:G:88:TYR:CE2	2.48	0.48
3:K:174:LYS:NZ	4:L:172:GLY:H	2.11	0.48
4:L:127:LYS:HB3	4:L:156:PHE:H	1.79	0.48
1:C:241:ASP:OD1	1:C:242:VAL:N	2.40	0.48
3:I:63:ARG:HD2	3:I:78:THR:O	2.14	0.48
4:J:127:LYS:HB3	4:J:156:PHE:H	1.79	0.48
3:G:166:GLU:OE1	3:G:166:GLU:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:TRP:CE2	1:A:225:GLY:HA2	2.50	0.47
1:C:206:THR:C	1:E:221:PRO:HG2	2.39	0.47
4:H:23:GLN:NE2	4:H:76:THR:O	2.33	0.47
3:K:20:ILE:HD11	3:K:108:LEU:HD11	1.96	0.47
3:I:37:TRP:HB2	3:I:50:ILE:HG12	1.95	0.47
2:F:10:ILE:HD13	2:F:136:GLY:HA3	1.96	0.47
4:H:43:GLN:OE1	4:H:180:LEU:HD21	2.14	0.47
4:H:61:VAL:HB	4:H:64:PHE:HD2	1.80	0.47
4:J:211:LYS:HB2	4:J:212:PRO:HD3	1.96	0.47
2:D:10:ILE:HD13	2:D:136:GLY:HA3	1.96	0.47
1:A:28:THR:HG21	2:B:108:ILE:HD12	1.94	0.47
1:C:141:ARG:NH1	1:C:147:PHE:O	2.40	0.47
4:L:211:LYS:HB2	4:L:212:PRO:HD3	1.96	0.47
2:B:147:ALA:O	2:B:150:GLU:HG3	2.15	0.47
2:D:147:ALA:O	2:D:150:GLU:HG3	2.15	0.47
1:E:222:TRP:CE2	1:E:225:GLY:HA2	2.49	0.47
2:F:147:ALA:O	2:F:150:GLU:HG3	2.15	0.47
3:G:171:GLN:HG2	3:G:178:TYR:CE1	2.50	0.47
4:H:205:ILE:HG13	4:H:220:LYS:HB3	1.96	0.47
3:G:36:HIS:CG	4:H:112:PRO:HB3	2.49	0.47
4:H:164:TRP:CZ3	4:H:206:CYS:HB3	2.50	0.47
1:A:221:PRO:HG2	1:E:206:THR:C	2.40	0.47
2:B:10:ILE:HD13	2:B:136:GLY:HA3	1.96	0.47
4:J:61:VAL:HB	4:J:64:PHE:HD2	1.80	0.47
3:G:18:VAL:HG21	3:G:80:LEU:HD13	1.95	0.46
4:H:38:ARG:NH2	4:H:64:PHE:HE1	2.13	0.46
1:C:222:TRP:CE2	1:C:225:GLY:HA2	2.49	0.46
4:J:164:TRP:CZ3	4:J:206:CYS:HB3	2.50	0.46
3:K:61:PRO:HB2	3:K:63:ARG:HG2	1.96	0.46
4:L:61:VAL:HB	4:L:64:PHE:HD2	1.80	0.46
3:K:215:ASN:OD1	3:K:216:ARG:N	2.48	0.46
4:L:157:PRO:HD2	4:L:210:HIS:HE1	1.81	0.46
4:L:164:TRP:CZ3	4:L:206:CYS:HB3	2.50	0.46
3:G:123:PHE:CD2	4:H:134:LEU:HB3	2.51	0.46
4:J:38:ARG:NH2	4:J:64:PHE:HE1	2.13	0.46
4:L:38:ARG:NH2	4:L:64:PHE:HE1	2.13	0.46
4:J:114:ASP:OD1	4:J:114:ASP:N	2.45	0.46
1:A:184:HIS:CD2	1:E:210:GLN:HE22	2.33	0.46
3:G:51:SER:OG	3:G:52:GLY:N	2.49	0.46
1:C:212:THR:HB	1:E:216:ASN:ND2	2.30	0.46
4:H:127:LYS:HB3	4:H:156:PHE:H	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:192:GLU:HA	3:G:195:LYS:HE3	1.98	0.46
3:I:5:THR:O	3:I:22:CYS:HA	2.16	0.46
3:K:153:TRP:HB3	3:K:160:GLN:NE2	2.31	0.46
3:K:35:VAL:HA	3:K:91:GLN:O	2.16	0.45
3:K:147:ARG:HB2	3:K:178:TYR:CD2	2.50	0.45
1:E:241:ASP:OD1	1:E:242:VAL:N	2.40	0.45
3:I:123:PHE:HB2	3:I:138:VAL:HG13	1.98	0.45
3:I:37:TRP:CD1	3:I:50:ILE:HD11	2.51	0.45
3:I:80:LEU:HD11	3:I:108:LEU:HD21	1.98	0.45
4:J:203:THR:OG1	4:J:221:VAL:HB	2.17	0.45
4:J:205:ILE:HG13	4:J:220:LYS:HB3	1.96	0.45
1:C:201:ARG:HD3	1:C:248:ASN:OD1	2.17	0.45
4:L:203:THR:OG1	4:L:221:VAL:HB	2.17	0.45
3:I:51:SER:HB3	4:J:112:PRO:HG3	1.99	0.45
4:L:205:ILE:HG13	4:L:220:LYS:HB3	1.96	0.45
4:H:162:VAL:HG22	4:H:208:VAL:HG22	1.99	0.45
1:A:28:THR:HG22	2:B:104:ASN:HB3	1.99	0.45
1:A:201:ARG:HD3	1:A:248:ASN:OD1	2.17	0.45
3:G:144:PHE:HE1	3:G:180:LEU:N	2.15	0.45
4:H:157:PRO:HD2	4:H:210:HIS:HE1	1.81	0.45
4:J:157:PRO:HD2	4:J:210:HIS:HE1	1.81	0.45
1:E:201:ARG:HD3	1:E:248:ASN:OD1	2.17	0.45
1:E:201:ARG:N	1:E:248:ASN:OD1	2.47	0.45
3:G:49:LEU:HD12	3:G:50:ILE:HG12	1.98	0.45
3:G:200:GLU:HG3	3:G:211:THR:HA	1.98	0.45
4:H:203:THR:OG1	4:H:221:VAL:HB	2.17	0.45
3:K:56:ARG:NH2	3:K:62:ASP:OD1	2.50	0.45
1:C:44:GLN:NE2	1:C:46:SER:O	2.42	0.44
3:K:39:GLN:HG3	3:K:86:ALA:HB3	1.99	0.44
1:A:244:VAL:HG23	1:C:221:PRO:HG3	1.99	0.44
1:A:264:LYS:O	1:A:302:TYR:OH	2.32	0.44
3:I:6:GLN:HA	3:I:21:SER:O	2.18	0.44
1:A:29:ILE:O	2:F:54:ARG:HD2	2.17	0.44
2:D:129:ASN:HA	2:D:166:ALA:HB1	2.00	0.44
3:G:195:LYS:HE2	3:G:195:LYS:HA	1.99	0.44
1:A:44:GLN:NE2	1:A:46:SER:O	2.42	0.44
4:J:196:SER:HA	4:J:199:LEU:HD23	2.00	0.44
3:K:128:GLU:OE1	3:K:128:GLU:N	2.37	0.44
2:F:129:ASN:HA	2:F:166:ALA:HB1	2.00	0.44
4:L:62:LYS:HA	4:L:65:GLN:HG2	2.00	0.44
3:K:61:PRO:HD2	3:K:64:PHE:CE2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:144:PHE:HE2	3:K:147:ARG:HA	1.82	0.44
3:G:40:GLN:O	3:G:86:ALA:HB1	2.17	0.43
3:G:152:GLN:OE1	3:G:153:TRP:N	2.51	0.43
4:H:196:SER:HA	4:H:199:LEU:HD23	2.00	0.43
2:B:4:GLY:O	2:B:8:GLY:HA3	2.18	0.43
4:H:91:THR:HG23	4:H:123:THR:HA	2.01	0.43
3:I:194:HIS:HB2	3:I:197:TYR:CE1	2.53	0.43
3:K:123:PHE:CD2	4:L:134:LEU:HB3	2.53	0.43
4:L:162:VAL:HG22	4:L:208:VAL:HG22	1.99	0.43
3:G:5:THR:O	3:G:22:CYS:HA	2.19	0.43
3:I:17:ARG:NE	3:I:78:THR:HA	2.30	0.43
2:B:129:ASN:HA	2:B:166:ALA:HB1	2.00	0.43
1:E:321:ARG:NH1	1:E:322:ASN:O	2.52	0.43
4:J:162:VAL:HG22	4:J:208:VAL:HG22	1.99	0.43
2:D:4:GLY:O	2:D:8:GLY:HA3	2.18	0.43
2:F:4:GLY:O	2:F:8:GLY:HA3	2.18	0.43
3:G:39:GLN:HG3	3:G:86:ALA:HB3	2.01	0.43
3:G:56:ARG:HD3	3:G:64:PHE:O	2.19	0.43
3:G:136:SER:HA	3:G:184:LEU:O	2.18	0.43
1:C:201:ARG:N	1:C:248:ASN:OD1	2.47	0.43
4:H:156:PHE:HB3	4:H:157:PRO:HD3	2.01	0.43
2:F:61:GLU:HG2	2:F:63:PHE:CZ	2.54	0.43
3:G:22:CYS:HB2	3:G:37:TRP:CH2	2.54	0.43
4:J:91:THR:HG23	4:J:123:THR:HA	2.01	0.43
3:G:16:GLN:H	3:G:80:LEU:HD23	1.83	0.42
4:L:196:SER:HA	4:L:199:LEU:HD23	2.00	0.42
1:E:118:LEU:HD23	1:E:118:LEU:HA	1.92	0.42
3:I:144:PHE:CZ	3:I:180:LEU:HB2	2.54	0.42
1:E:264:LYS:O	1:E:302:TYR:OH	2.33	0.42
3:I:127:ASP:O	3:I:131:LYS:HG2	2.19	0.42
1:C:173:ASN:OD1	1:C:173:ASN:N	2.52	0.42
1:C:321:ARG:NH1	1:C:322:ASN:O	2.52	0.42
1:E:42:LEU:HD12	2:F:100:VAL:HG22	2.01	0.42
2:F:119:PHE:HZ	2:F:132:ASP:OD1	2.03	0.42
3:I:35:VAL:HA	3:I:91:GLN:O	2.19	0.42
3:I:49:LEU:O	3:I:57:PRO:HD2	2.20	0.42
4:J:62:LYS:HD3	4:J:65:GLN:HG3	2.02	0.42
3:K:160:GLN:OE1	3:K:160:GLN:N	2.53	0.42
4:L:18:VAL:HB	4:L:86:LEU:HD11	2.02	0.42
4:L:91:THR:HG23	4:L:123:THR:HA	2.01	0.42
1:C:200:GLY:O	1:C:215:PRO:HD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:119:PHE:HZ	2:D:132:ASP:OD1	2.03	0.42
1:E:200:GLY:O	1:E:215:PRO:HD2	2.19	0.42
3:I:133:GLY:C	3:I:188:LYS:HE3	2.44	0.42
4:J:18:VAL:HB	4:J:86:LEU:HD11	2.02	0.42
4:J:62:LYS:HA	4:J:65:GLN:HG2	2.00	0.42
3:G:154:LYS:HE3	3:G:157:ASN:HA	2.01	0.42
4:H:62:LYS:HA	4:H:65:GLN:HG2	2.00	0.42
3:I:6:GLN:HG2	3:I:106:THR:OG1	2.20	0.42
4:L:50:TRP:CH2	4:L:52:ASN:HB2	2.55	0.42
1:A:321:ARG:NH1	1:A:322:ASN:O	2.52	0.42
3:G:86:ALA:H	3:G:108:LEU:HB2	1.84	0.42
4:J:50:TRP:CH2	4:J:52:ASN:HB2	2.55	0.42
1:A:200:GLY:O	1:A:215:PRO:HD2	2.19	0.42
2:B:76:ARG:NH2	2:F:81:GLU:OE1	2.49	0.42
3:G:186:LEU:HB3	3:G:190:ASP:OD1	2.20	0.42
4:H:173:VAL:HG12	4:H:192:VAL:HB	2.02	0.42
3:I:194:HIS:HB2	3:I:197:TYR:HE1	1.85	0.42
4:J:32:TYR:CD2	4:J:100:GLU:HA	2.55	0.42
4:J:156:PHE:HB3	4:J:157:PRO:HD3	2.01	0.42
3:K:125:PRO:HB3	3:K:136:SER:O	2.20	0.42
3:K:166:GLU:OE2	3:K:166:GLU:N	2.53	0.42
1:A:320:MET:HE2	1:A:320:MET:HB3	1.86	0.42
2:B:119:PHE:HZ	2:B:132:ASP:OD1	2.03	0.42
1:E:320:MET:HE2	1:E:320:MET:HB3	1.86	0.42
4:H:50:TRP:CH2	4:H:52:ASN:HB2	2.55	0.42
1:A:184:HIS:NE2	1:A:231:SER:OG	2.51	0.41
2:D:61:GLU:HG2	2:D:63:PHE:CZ	2.54	0.41
4:L:32:TYR:CD2	4:L:100:GLU:HA	2.55	0.41
3:G:6:GLN:HA	3:G:21:SER:O	2.19	0.41
3:I:68:LYS:HG2	3:I:69:SER:O	2.20	0.41
2:B:61:GLU:HG2	2:B:63:PHE:CZ	2.54	0.41
4:L:156:PHE:HB3	4:L:157:PRO:HD3	2.01	0.41
4:J:39:GLN:HB2	4:J:45:LEU:HD23	2.03	0.41
4:L:173:VAL:HG12	4:L:192:VAL:HB	2.02	0.41
2:B:126:LEU:HD23	2:B:157:TYR:CE2	2.56	0.41
1:C:127:TRP:CZ2	1:C:253:ALA:HB1	2.56	0.41
1:E:14:CYS:HA	2:F:137:CYS:HA	2.03	0.41
4:H:62:LYS:HD3	4:H:65:GLN:HG3	2.02	0.41
4:H:18:VAL:HB	4:H:86:LEU:HD11	2.02	0.41
1:C:17:HIS:HB2	1:C:320:MET:SD	2.61	0.41
3:G:200:GLU:N	3:G:200:GLU:OE1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:146:PRO:HD2	3:K:203:HIS:HE1	1.83	0.41
1:C:184:HIS:NE2	1:C:231:SER:OG	2.51	0.41
1:E:127:TRP:CZ2	1:E:253:ALA:HB1	2.56	0.41
2:F:126:LEU:HD23	2:F:157:TYR:CE2	2.56	0.41
3:G:155:VAL:HG13	3:G:197:TYR:CZ	2.56	0.41
4:H:49:GLY:CA	4:H:70:MET:HE1	2.51	0.41
4:H:60:TYR:CE1	4:H:68:VAL:HG12	2.56	0.41
4:J:49:GLY:CA	4:J:70:MET:HE1	2.51	0.41
4:J:60:TYR:CE1	4:J:68:VAL:HG12	2.56	0.41
4:J:173:VAL:HG12	4:J:192:VAL:HB	2.02	0.41
4:L:62:LYS:HD3	4:L:65:GLN:HG3	2.02	0.41
2:D:126:LEU:HD23	2:D:157:TYR:CE2	2.56	0.41
1:E:17:HIS:HB2	1:E:320:MET:SD	2.61	0.41
4:L:81:MET:HE2	4:L:81:MET:HB3	1.98	0.41
4:H:32:TYR:CD2	4:H:100:GLU:HA	2.55	0.40
3:I:12:GLY:O	3:I:111:LEU:N	2.54	0.40
1:A:17:HIS:HB2	1:A:320:MET:SD	2.61	0.40
4:L:39:GLN:HB2	4:L:45:LEU:HD23	2.03	0.40
4:H:39:GLN:HB2	4:H:45:LEU:HD23	2.03	0.40
1:A:201:ARG:N	1:A:248:ASN:OD1	2.47	0.40
1:A:269:ARG:HB3	4:H:103:ARG:HD2	2.04	0.40
2:B:76:ARG:HD2	1:E:106:ALA:HB3	2.03	0.40
3:G:49:LEU:HA	3:G:60:VAL:HG21	2.02	0.40
3:I:39:GLN:HB2	3:I:88:TYR:CE2	2.57	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	316/345 (92%)	312 (99%)	4 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	316/345 (92%)	312 (99%)	4 (1%)	0	100	100
1	E	316/345 (92%)	312 (99%)	4 (1%)	0	100	100
2	B	170/254 (67%)	157 (92%)	13 (8%)	0	100	100
2	D	170/254 (67%)	157 (92%)	13 (8%)	0	100	100
2	F	170/254 (67%)	157 (92%)	13 (8%)	0	100	100
3	G	196/219 (90%)	186 (95%)	10 (5%)	0	100	100
3	I	196/219 (90%)	187 (95%)	9 (5%)	0	100	100
3	K	196/219 (90%)	187 (95%)	9 (5%)	0	100	100
4	H	208/235 (88%)	194 (93%)	14 (7%)	0	100	100
4	J	208/235 (88%)	194 (93%)	14 (7%)	0	100	100
4	L	208/235 (88%)	194 (93%)	14 (7%)	0	100	100
All	All	2670/3159 (84%)	2549 (96%)	121 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	280/303 (92%)	279 (100%)	1 (0%)	84	94
1	C	280/303 (92%)	279 (100%)	1 (0%)	84	94
1	E	280/303 (92%)	279 (100%)	1 (0%)	84	94
2	B	147/212 (69%)	147 (100%)	0	100	100
2	D	147/212 (69%)	147 (100%)	0	100	100
2	F	147/212 (69%)	147 (100%)	0	100	100
3	G	171/184 (93%)	171 (100%)	0	100	100
3	I	171/184 (93%)	171 (100%)	0	100	100
3	K	171/184 (93%)	171 (100%)	0	100	100
4	H	180/200 (90%)	180 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	J	180/200 (90%)	180 (100%)	0	100	100
4	L	180/200 (90%)	180 (100%)	0	100	100
All	All	2334/2697 (86%)	2331 (100%)	3 (0%)	87	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	SER
1	C	219	SER
1	E	219	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	191	GLN
1	A	211	GLN
1	C	191	GLN
1	C	211	GLN
1	E	17	HIS
1	E	191	GLN
1	E	211	GLN
3	G	160	GLN
3	I	6	GLN
3	I	157	ASN
3	I	163	ASN
3	I	203	HIS
3	K	203	HIS

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 ⓘ

45 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	$\# Z > 2$
5	NAG	M	1	5,1	14,14,15	0.41	0	17,19,21	0.47	0
5	NAG	M	2	5	14,14,15	0.21	0	17,19,21	0.45	0
6	NAG	N	1	1,6	14,14,15	0.44	0	17,19,21	0.48	0
6	NAG	N	2	6	14,14,15	0.29	0	17,19,21	0.48	0
6	BMA	N	3	6	11,11,12	0.49	0	15,15,17	0.84	0
6	MAN	N	4	6	11,11,12	0.55	0	15,15,17	0.94	2 (13%)
6	MAN	N	5	6	11,11,12	0.65	0	15,15,17	0.88	1 (6%)
7	NAG	O	1	1,7	14,14,15	0.37	0	17,19,21	0.40	0
7	NAG	O	2	7	14,14,15	0.16	0	17,19,21	0.70	0
7	BMA	O	3	7	11,11,12	0.59	0	15,15,17	0.97	1 (6%)
7	MAN	O	4	7	11,11,12	0.68	0	15,15,17	1.73	5 (33%)
7	MAN	O	5	7	11,11,12	0.52	0	15,15,17	1.05	2 (13%)
7	MAN	O	6	7	11,11,12	0.51	0	15,15,17	0.97	2 (13%)
7	MAN	O	7	7	11,11,12	1.06	1 (9%)	15,15,17	0.84	1 (6%)
7	MAN	O	8	7	11,11,12	0.63	0	15,15,17	0.86	1 (6%)
5	NAG	P	1	5,1	14,14,15	0.41	0	17,19,21	0.47	0
5	NAG	P	2	5	14,14,15	0.21	0	17,19,21	0.45	0
6	NAG	Q	1	1,6	14,14,15	0.43	0	17,19,21	0.47	0
6	NAG	Q	2	6	14,14,15	0.29	0	17,19,21	0.48	0
6	BMA	Q	3	6	11,11,12	0.48	0	15,15,17	0.84	0
6	MAN	Q	4	6	11,11,12	0.55	0	15,15,17	0.94	2 (13%)
6	MAN	Q	5	6	11,11,12	0.64	0	15,15,17	0.88	1 (6%)
7	NAG	R	1	1,7	14,14,15	0.37	0	17,19,21	0.40	0
7	NAG	R	2	7	14,14,15	0.15	0	17,19,21	0.70	0
7	BMA	R	3	7	11,11,12	0.59	0	15,15,17	0.97	1 (6%)
7	MAN	R	4	7	11,11,12	0.69	0	15,15,17	1.73	5 (33%)
7	MAN	R	5	7	11,11,12	0.50	0	15,15,17	1.05	2 (13%)
7	MAN	R	6	7	11,11,12	0.50	0	15,15,17	0.97	2 (13%)
7	MAN	R	7	7	11,11,12	1.04	1 (9%)	15,15,17	0.84	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MAN	R	8	7	11,11,12	0.63	0	15,15,17	0.86	1 (6%)
5	NAG	S	1	5,1	14,14,15	0.41	0	17,19,21	0.47	0
5	NAG	S	2	5	14,14,15	0.21	0	17,19,21	0.45	0
6	NAG	T	1	1,6	14,14,15	0.43	0	17,19,21	0.47	0
6	NAG	T	2	6	14,14,15	0.29	0	17,19,21	0.49	0
6	BMA	T	3	6	11,11,12	0.49	0	15,15,17	0.84	0
6	MAN	T	4	6	11,11,12	0.56	0	15,15,17	0.94	2 (13%)
6	MAN	T	5	6	11,11,12	0.65	0	15,15,17	0.88	1 (6%)
7	NAG	U	1	1,7	14,14,15	0.37	0	17,19,21	0.40	0
7	NAG	U	2	7	14,14,15	0.15	0	17,19,21	0.70	0
7	BMA	U	3	7	11,11,12	0.60	0	15,15,17	0.96	0
7	MAN	U	4	7	11,11,12	0.68	0	15,15,17	1.73	5 (33%)
7	MAN	U	5	7	11,11,12	0.51	0	15,15,17	1.05	2 (13%)
7	MAN	U	6	7	11,11,12	0.50	0	15,15,17	0.98	2 (13%)
7	MAN	U	7	7	11,11,12	1.04	1 (9%)	15,15,17	0.83	1 (6%)
7	MAN	U	8	7	11,11,12	0.64	0	15,15,17	0.87	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	M	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	M	2	5	-	2/6/23/26	0/1/1/1
6	NAG	N	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	N	2	6	-	0/6/23/26	0/1/1/1
6	BMA	N	3	6	-	0/2/19/22	0/1/1/1
6	MAN	N	4	6	-	2/2/19/22	0/1/1/1
6	MAN	N	5	6	-	0/2/19/22	0/1/1/1
7	NAG	O	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	O	2	7	-	2/6/23/26	0/1/1/1
7	BMA	O	3	7	-	0/2/19/22	0/1/1/1
7	MAN	O	4	7	-	0/2/19/22	0/1/1/1
7	MAN	O	5	7	-	0/2/19/22	0/1/1/1
7	MAN	O	6	7	-	0/2/19/22	0/1/1/1
7	MAN	O	7	7	-	2/2/19/22	0/1/1/1
7	MAN	O	8	7	-	0/2/19/22	0/1/1/1
5	NAG	P	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	P	2	5	-	2/6/23/26	0/1/1/1
6	NAG	Q	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	Q	2	6	-	0/6/23/26	0/1/1/1
6	BMA	Q	3	6	-	0/2/19/22	0/1/1/1
6	MAN	Q	4	6	-	2/2/19/22	0/1/1/1
6	MAN	Q	5	6	-	0/2/19/22	0/1/1/1
7	NAG	R	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	R	2	7	-	2/6/23/26	0/1/1/1
7	BMA	R	3	7	-	0/2/19/22	0/1/1/1
7	MAN	R	4	7	-	0/2/19/22	0/1/1/1
7	MAN	R	5	7	-	0/2/19/22	0/1/1/1
7	MAN	R	6	7	-	0/2/19/22	0/1/1/1
7	MAN	R	7	7	-	2/2/19/22	0/1/1/1
7	MAN	R	8	7	-	0/2/19/22	0/1/1/1
5	NAG	S	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	S	2	5	-	2/6/23/26	0/1/1/1
6	NAG	T	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	T	2	6	-	0/6/23/26	0/1/1/1
6	BMA	T	3	6	-	0/2/19/22	0/1/1/1
6	MAN	T	4	6	-	2/2/19/22	0/1/1/1
6	MAN	T	5	6	-	0/2/19/22	0/1/1/1
7	NAG	U	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	U	2	7	-	2/6/23/26	0/1/1/1
7	BMA	U	3	7	-	0/2/19/22	0/1/1/1
7	MAN	U	4	7	-	0/2/19/22	0/1/1/1
7	MAN	U	5	7	-	0/2/19/22	0/1/1/1
7	MAN	U	6	7	-	0/2/19/22	0/1/1/1
7	MAN	U	7	7	-	2/2/19/22	0/1/1/1
7	MAN	U	8	7	-	0/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	O	7	MAN	O5-C1	-3.13	1.38	1.43
7	R	7	MAN	O5-C1	-3.10	1.38	1.43
7	U	7	MAN	O5-C1	-3.08	1.38	1.43

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	U	4	MAN	O3-C3-C2	3.10	116.38	110.05
7	O	4	MAN	O3-C3-C2	3.10	116.37	110.05
7	R	4	MAN	O3-C3-C2	3.09	116.36	110.05
7	U	4	MAN	C3-C4-C5	-3.01	104.77	110.23
7	O	4	MAN	C3-C4-C5	-3.00	104.79	110.23
7	R	4	MAN	C3-C4-C5	-2.99	104.81	110.23
7	R	4	MAN	C1-C2-C3	-2.89	105.44	109.64
7	U	4	MAN	C1-C2-C3	-2.89	105.44	109.64
7	O	4	MAN	C1-C2-C3	-2.88	105.45	109.64
7	R	5	MAN	C1-O5-C5	2.66	115.75	112.19
7	O	5	MAN	C1-O5-C5	2.63	115.70	112.19
7	U	5	MAN	C1-O5-C5	2.61	115.69	112.19
7	U	6	MAN	C1-O5-C5	2.50	115.54	112.19
7	R	6	MAN	C1-O5-C5	2.49	115.53	112.19
7	O	7	MAN	O2-C2-C3	-2.48	105.02	110.15
7	R	7	MAN	O2-C2-C3	-2.47	105.03	110.15
7	O	6	MAN	C1-O5-C5	2.47	115.50	112.19
7	R	5	MAN	O2-C2-C3	-2.47	105.04	110.15
7	U	5	MAN	O2-C2-C3	-2.46	105.05	110.15
7	O	5	MAN	O2-C2-C3	-2.46	105.05	110.15
7	U	7	MAN	O2-C2-C3	-2.45	105.08	110.15
7	O	6	MAN	O2-C2-C3	-2.35	105.29	110.15
7	U	6	MAN	O2-C2-C3	-2.35	105.29	110.15
6	Q	4	MAN	C1-O5-C5	2.33	115.31	112.19
7	R	4	MAN	O3-C3-C4	2.32	115.85	110.38
6	T	4	MAN	C1-O5-C5	2.32	115.30	112.19
7	O	4	MAN	O3-C3-C4	2.31	115.83	110.38
7	R	6	MAN	O2-C2-C3	-2.31	105.36	110.15
7	U	4	MAN	O3-C3-C4	2.31	115.82	110.38
6	N	4	MAN	C1-O5-C5	2.30	115.27	112.19
7	U	4	MAN	C1-O5-C5	2.29	115.25	112.19
7	O	4	MAN	C1-O5-C5	2.27	115.23	112.19
7	R	4	MAN	C1-O5-C5	2.27	115.23	112.19
6	N	4	MAN	O2-C2-C3	-2.19	105.61	110.15
6	T	4	MAN	O2-C2-C3	-2.19	105.61	110.15
6	Q	4	MAN	O2-C2-C3	-2.17	105.65	110.15
7	U	8	MAN	O2-C2-C3	-2.17	105.67	110.15
7	O	8	MAN	O2-C2-C3	-2.15	105.69	110.15
7	R	8	MAN	O2-C2-C3	-2.15	105.70	110.15
6	T	5	MAN	O2-C2-C3	-2.13	105.75	110.15
6	N	5	MAN	O2-C2-C3	-2.13	105.75	110.15
6	Q	5	MAN	O2-C2-C3	-2.12	105.76	110.15
7	O	3	BMA	O2-C2-C3	-2.02	105.97	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	R	3	BMA	O2-C2-C3	-2.01	105.98	110.15

There are no chirality outliers.

All (30) torsion outliers are listed below:

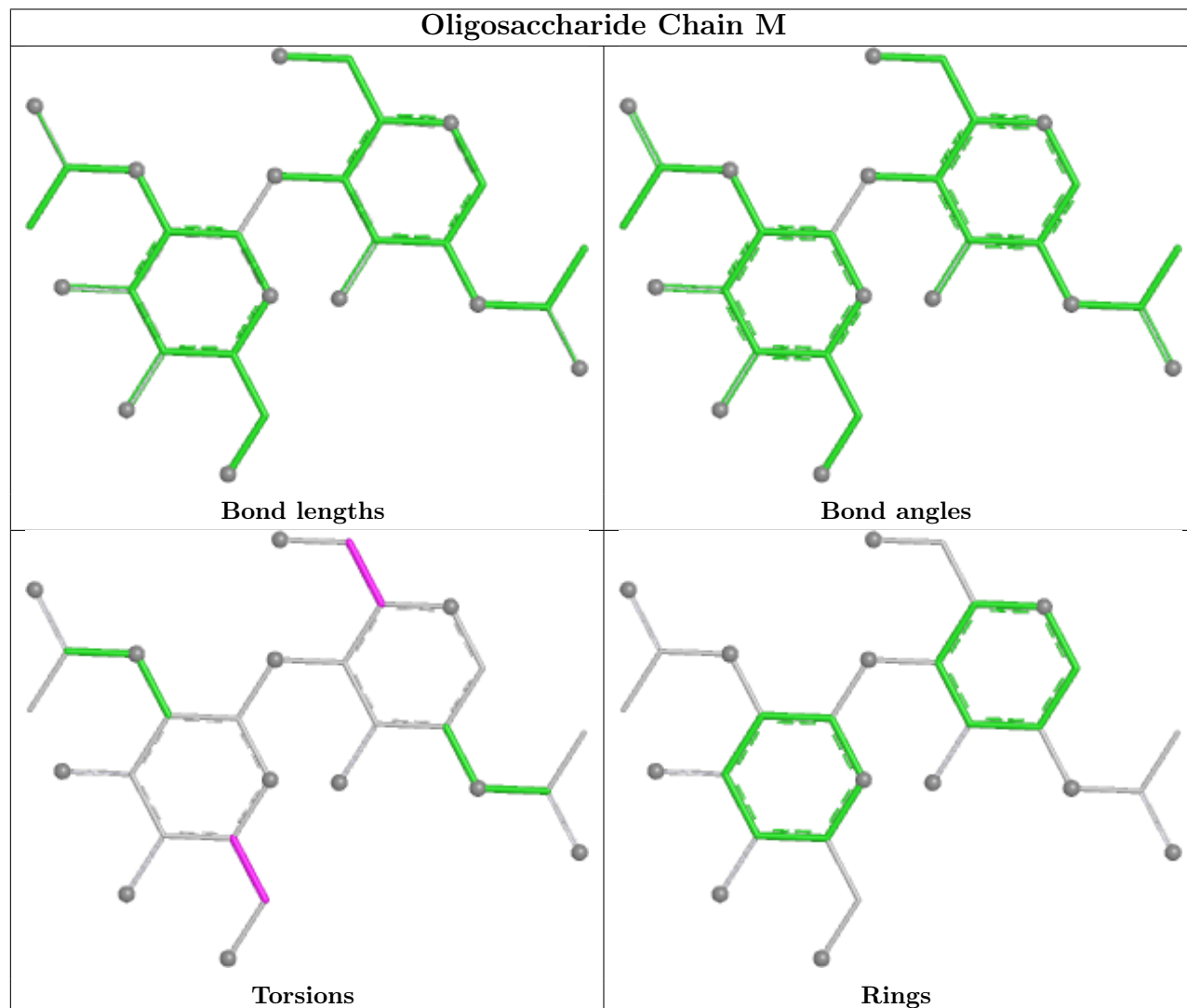
Mol	Chain	Res	Type	Atoms
7	O	7	MAN	O5-C5-C6-O6
7	R	7	MAN	O5-C5-C6-O6
7	U	7	MAN	O5-C5-C6-O6
7	O	7	MAN	C4-C5-C6-O6
7	R	7	MAN	C4-C5-C6-O6
7	U	7	MAN	C4-C5-C6-O6
5	P	2	NAG	O5-C5-C6-O6
5	M	2	NAG	O5-C5-C6-O6
5	S	2	NAG	O5-C5-C6-O6
6	N	4	MAN	O5-C5-C6-O6
6	Q	4	MAN	O5-C5-C6-O6
6	T	4	MAN	O5-C5-C6-O6
5	M	2	NAG	C4-C5-C6-O6
5	P	2	NAG	C4-C5-C6-O6
5	S	2	NAG	C4-C5-C6-O6
5	M	1	NAG	C4-C5-C6-O6
5	P	1	NAG	C4-C5-C6-O6
5	S	1	NAG	C4-C5-C6-O6
5	M	1	NAG	O5-C5-C6-O6
5	P	1	NAG	O5-C5-C6-O6
5	S	1	NAG	O5-C5-C6-O6
7	O	2	NAG	C3-C2-N2-C7
7	R	2	NAG	C3-C2-N2-C7
7	U	2	NAG	C3-C2-N2-C7
7	O	2	NAG	C1-C2-N2-C7
7	R	2	NAG	C1-C2-N2-C7
7	U	2	NAG	C1-C2-N2-C7
6	N	4	MAN	C4-C5-C6-O6
6	T	4	MAN	C4-C5-C6-O6
6	Q	4	MAN	C4-C5-C6-O6

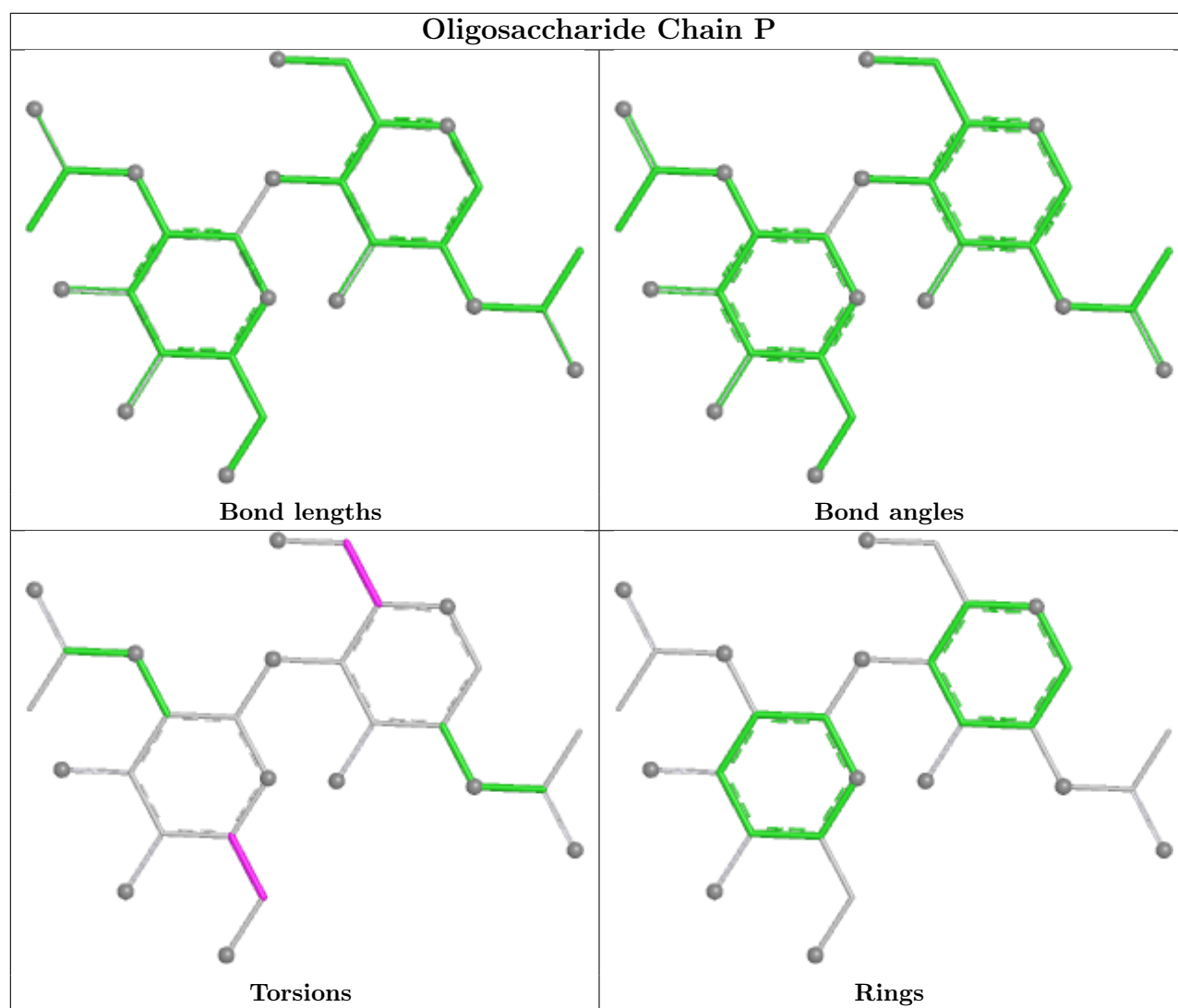
There are no ring outliers.

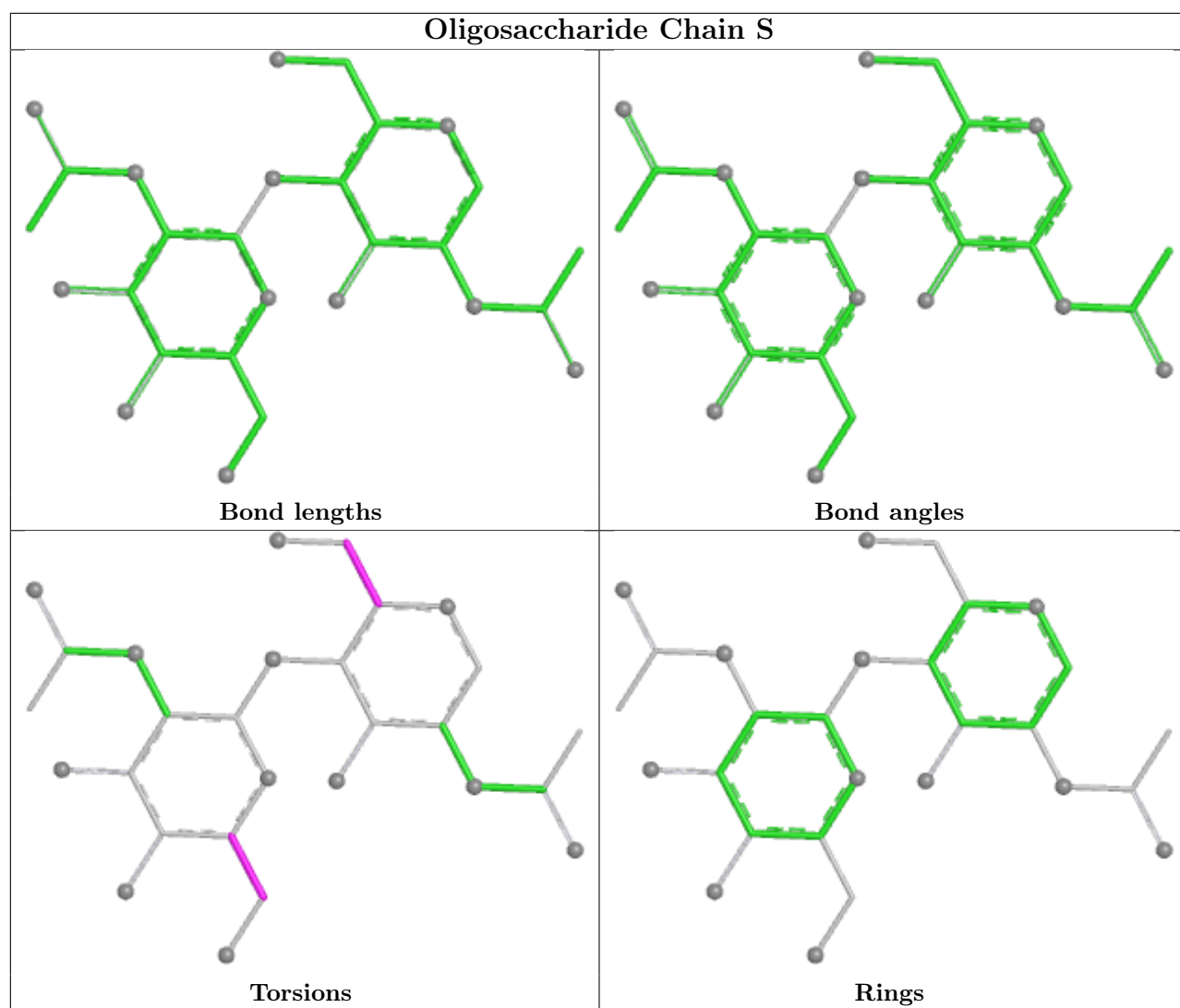
1 monomer is involved in 1 short contact:

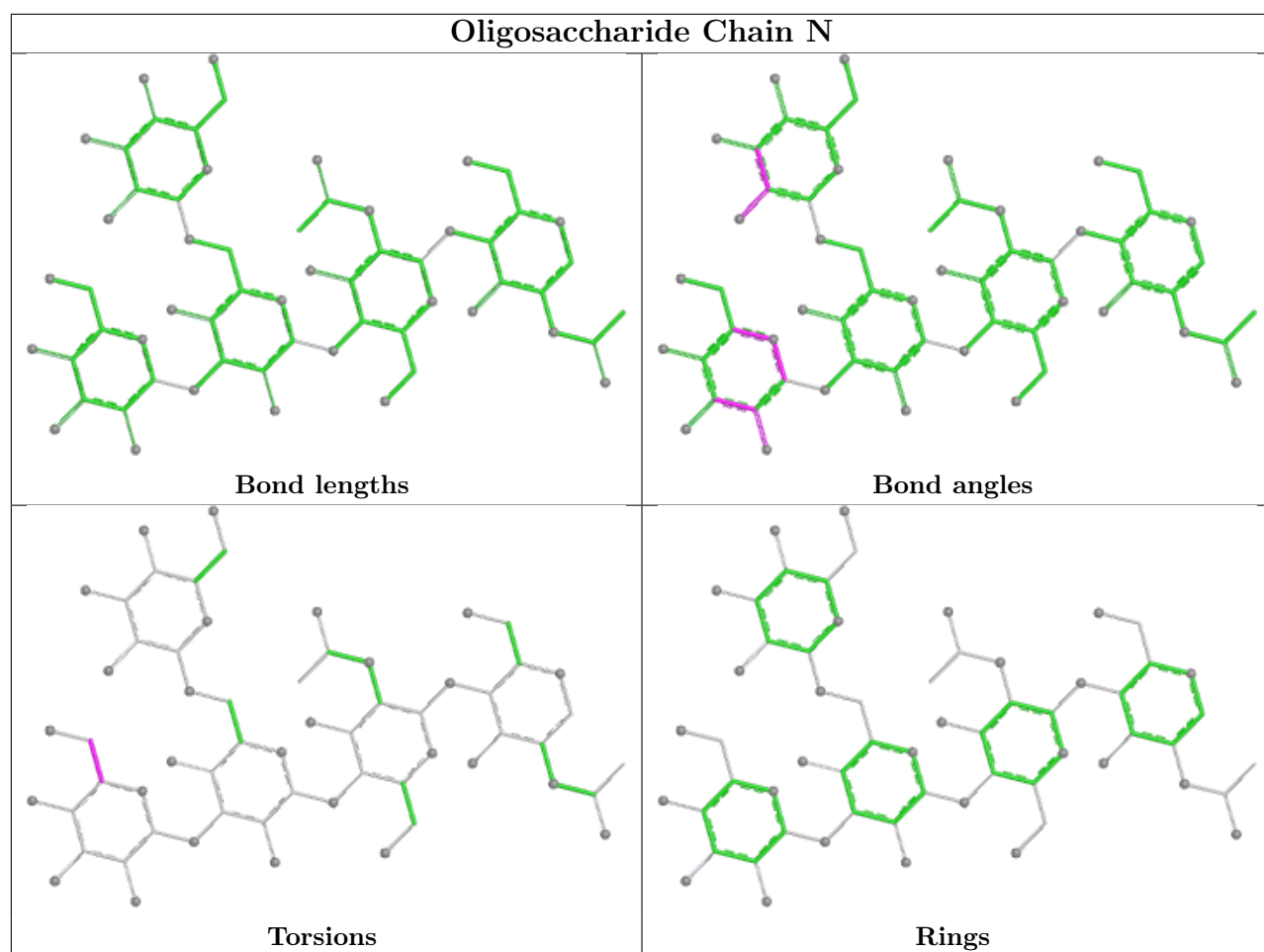
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	O	5	MAN	1	0

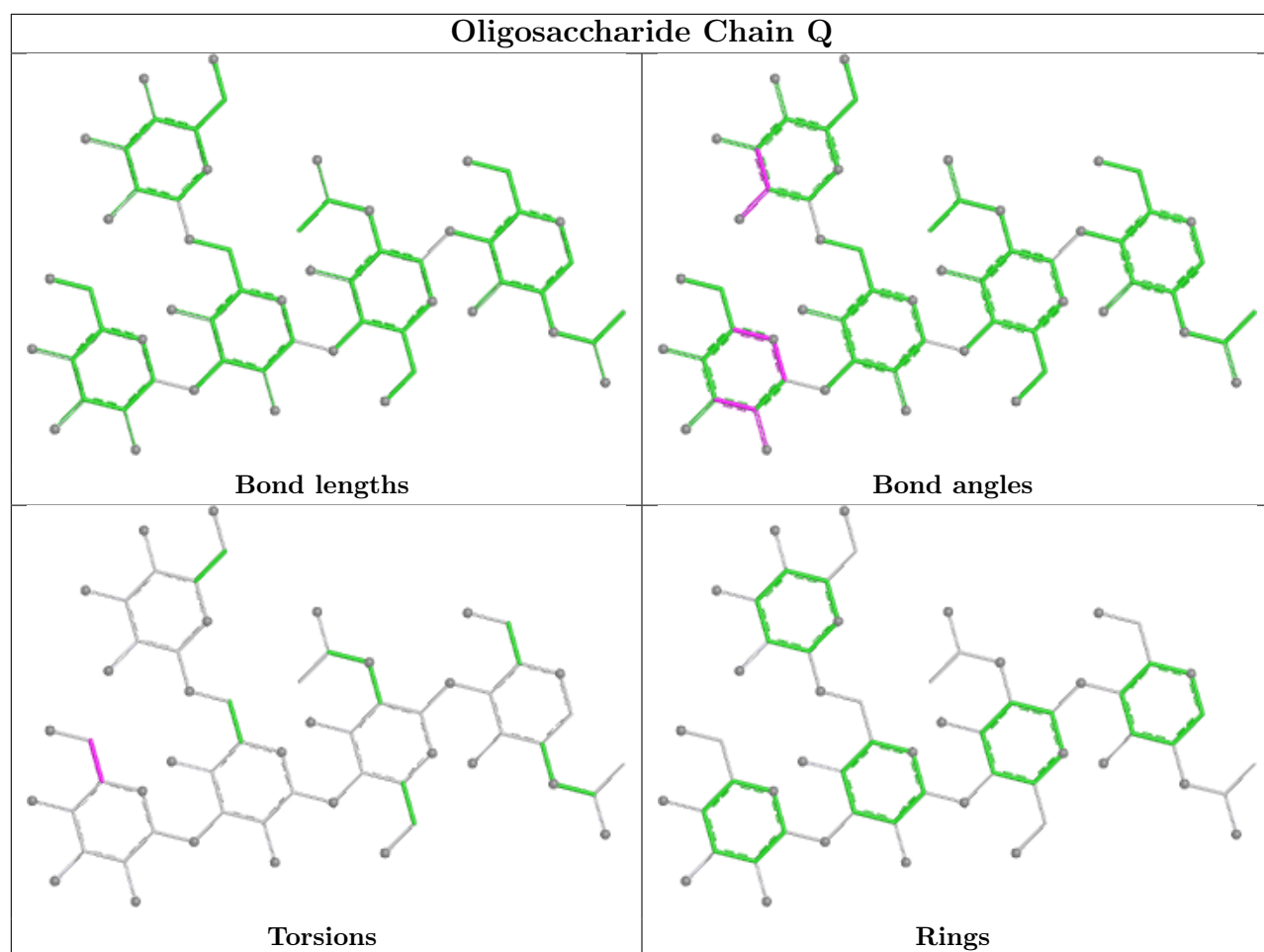
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

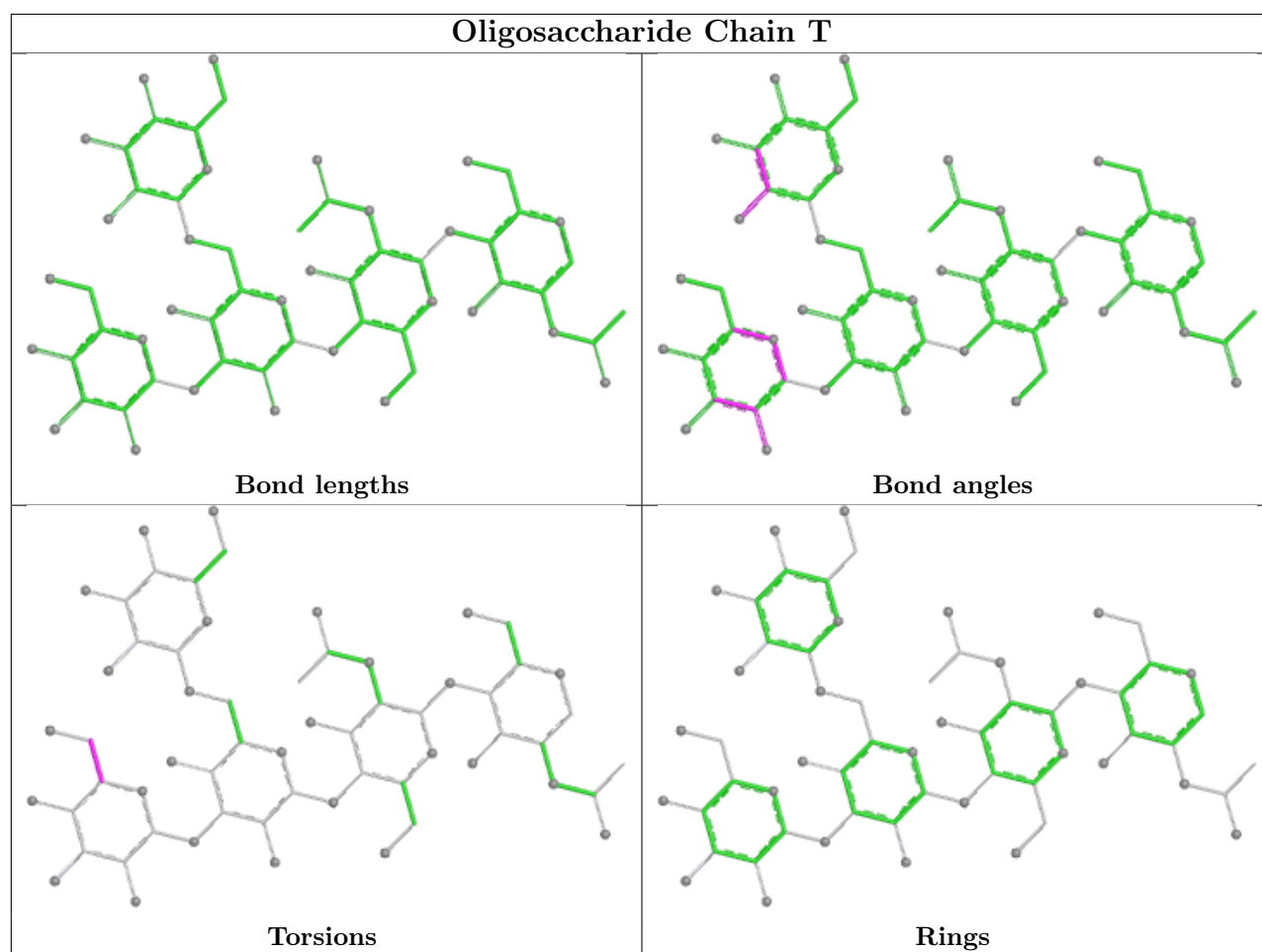


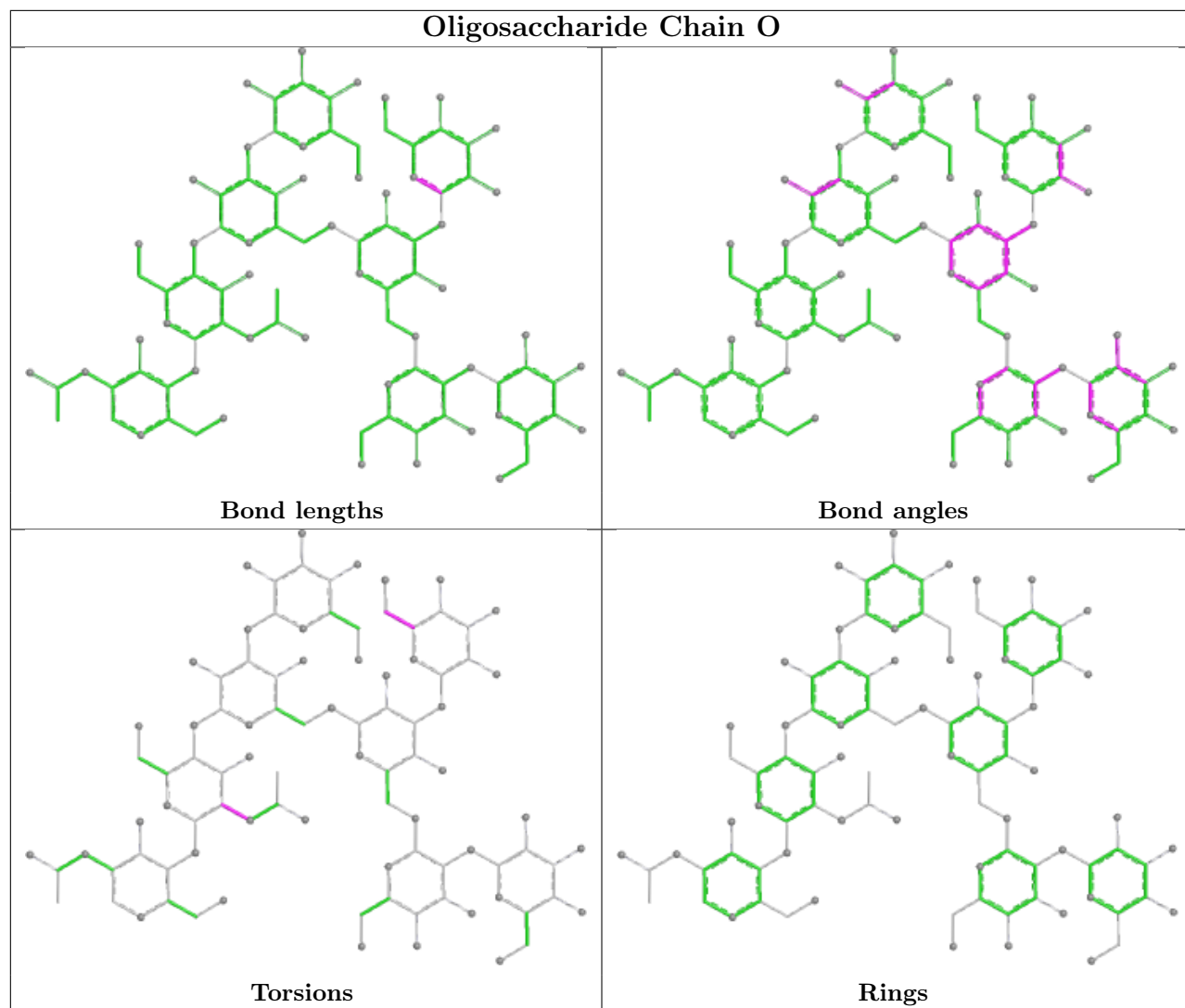


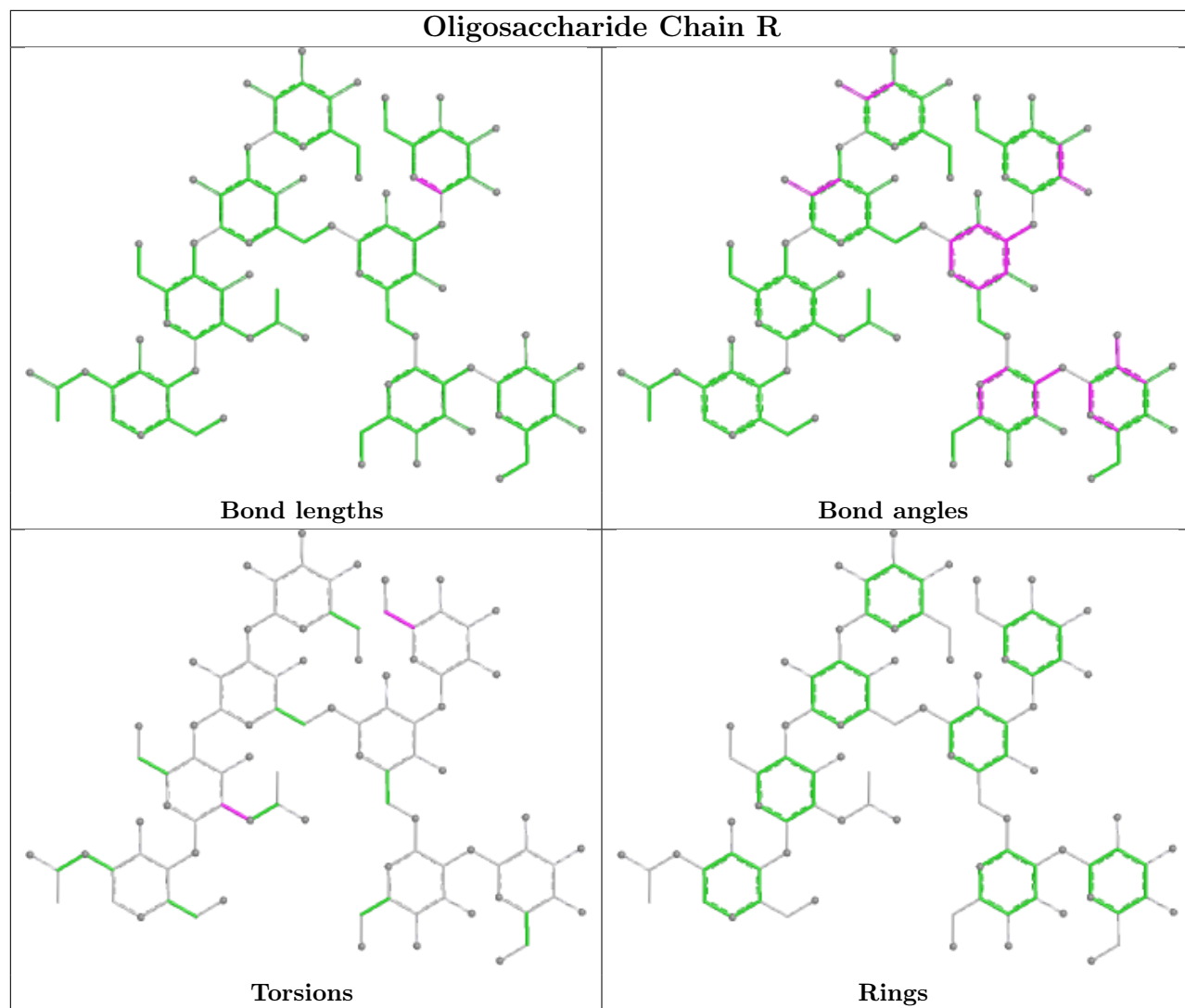


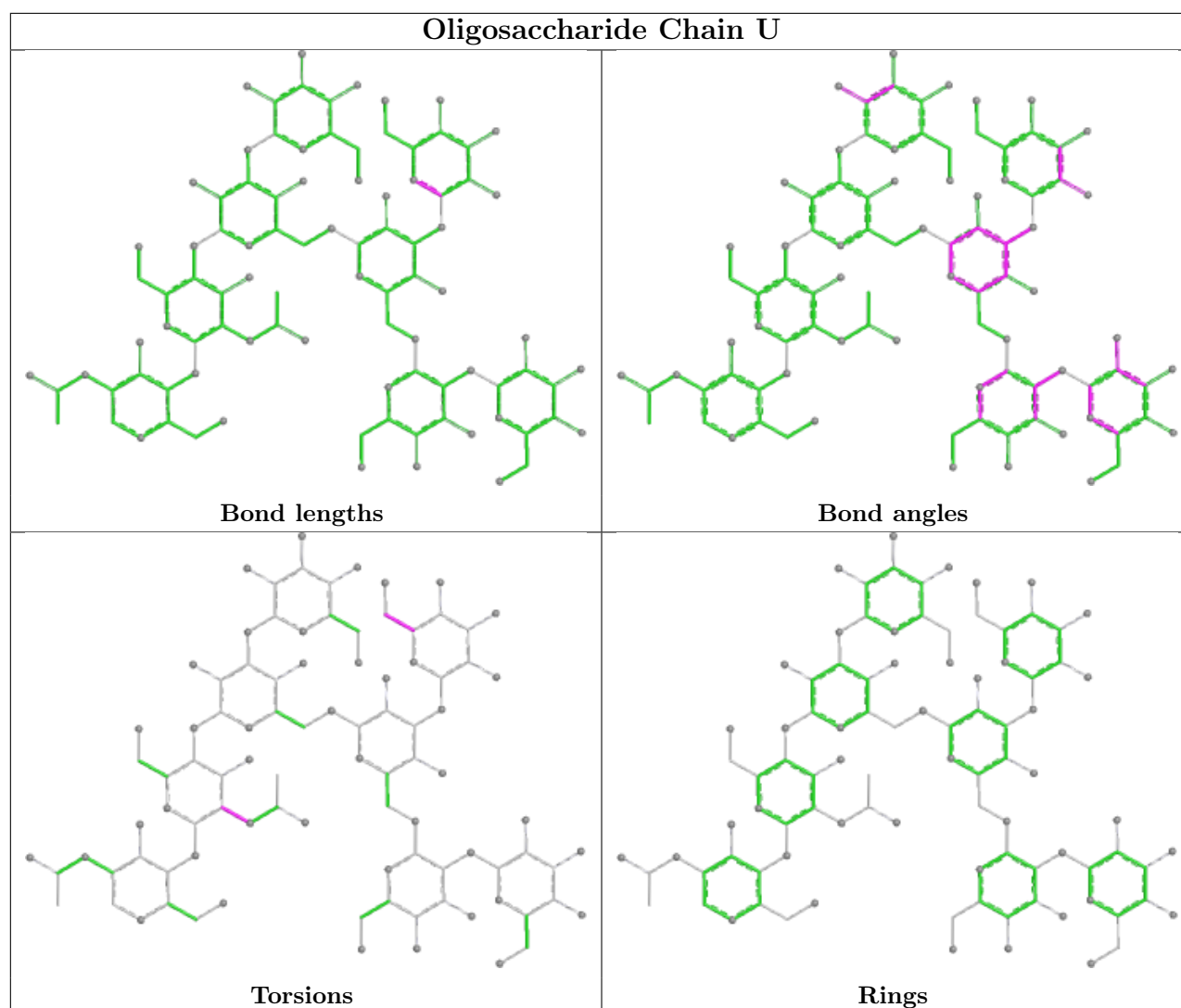












5.6 Ligand geometry [i](#)

9 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	$\# Z > 2$
8	NAG	C	402	1	14,14,15	0.18	0	17,19,21	0.50	0
8	NAG	A	401	1	14,14,15	0.45	0	17,19,21	0.52	0
8	NAG	B	301	2	14,14,15	0.36	0	17,19,21	1.34	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	C	401	1	14,14,15	0.44	0	17,19,21	0.52	0
8	NAG	A	402	1	14,14,15	0.18	0	17,19,21	0.48	0
8	NAG	E	401	1	14,14,15	0.45	0	17,19,21	0.52	0
8	NAG	D	301	2	14,14,15	0.38	0	17,19,21	1.35	2 (11%)
8	NAG	F	301	2	14,14,15	0.37	0	17,19,21	1.35	2 (11%)
8	NAG	E	402	1	14,14,15	0.18	0	17,19,21	0.50	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
8	NAG	C	402	1	-	2/6/23/26	0/1/1/1
8	NAG	A	401	1	-	4/6/23/26	0/1/1/1
8	NAG	B	301	2	-	6/6/23/26	0/1/1/1
8	NAG	C	401	1	-	4/6/23/26	0/1/1/1
8	NAG	A	402	1	-	2/6/23/26	0/1/1/1
8	NAG	E	401	1	-	4/6/23/26	0/1/1/1
8	NAG	D	301	2	-	6/6/23/26	0/1/1/1
8	NAG	F	301	2	-	6/6/23/26	0/1/1/1
8	NAG	E	402	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	F	301	NAG	C2-N2-C7	4.61	129.08	122.90
8	D	301	NAG	C2-N2-C7	4.61	129.08	122.90
8	B	301	NAG	C2-N2-C7	4.59	129.06	122.90
8	B	301	NAG	C1-C2-N2	2.25	113.97	110.43
8	D	301	NAG	C1-C2-N2	2.24	113.96	110.43
8	F	301	NAG	C1-C2-N2	2.22	113.92	110.43

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	401	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
8	C	401	NAG	O5-C5-C6-O6
8	E	401	NAG	O5-C5-C6-O6
8	A	402	NAG	O5-C5-C6-O6
8	C	402	NAG	O5-C5-C6-O6
8	E	402	NAG	O5-C5-C6-O6
8	A	401	NAG	C4-C5-C6-O6
8	C	401	NAG	C4-C5-C6-O6
8	E	401	NAG	C4-C5-C6-O6
8	B	301	NAG	C8-C7-N2-C2
8	B	301	NAG	O7-C7-N2-C2
8	D	301	NAG	C8-C7-N2-C2
8	D	301	NAG	O7-C7-N2-C2
8	F	301	NAG	C8-C7-N2-C2
8	F	301	NAG	O7-C7-N2-C2
8	A	402	NAG	C4-C5-C6-O6
8	C	402	NAG	C4-C5-C6-O6
8	E	402	NAG	C4-C5-C6-O6
8	F	301	NAG	C4-C5-C6-O6
8	D	301	NAG	C4-C5-C6-O6
8	B	301	NAG	C4-C5-C6-O6
8	F	301	NAG	O5-C5-C6-O6
8	D	301	NAG	O5-C5-C6-O6
8	B	301	NAG	O5-C5-C6-O6
8	A	401	NAG	C3-C2-N2-C7
8	C	401	NAG	C3-C2-N2-C7
8	E	401	NAG	C3-C2-N2-C7
8	A	401	NAG	C1-C2-N2-C7
8	B	301	NAG	C1-C2-N2-C7
8	C	401	NAG	C1-C2-N2-C7
8	D	301	NAG	C1-C2-N2-C7
8	E	401	NAG	C1-C2-N2-C7
8	F	301	NAG	C1-C2-N2-C7
8	B	301	NAG	C3-C2-N2-C7
8	D	301	NAG	C3-C2-N2-C7
8	F	301	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	301	NAG	1	0
8	D	301	NAG	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	F	301	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.

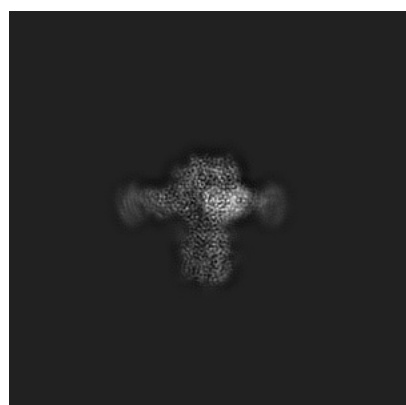
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22652. These allow visual inspection of the internal detail of the map and identification of artifacts.

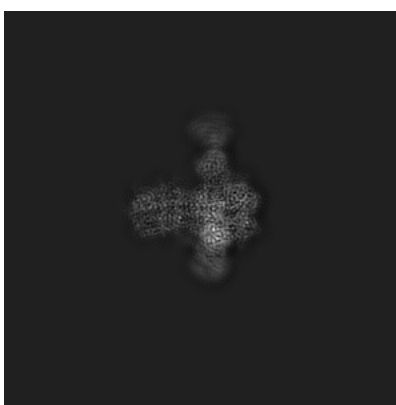
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

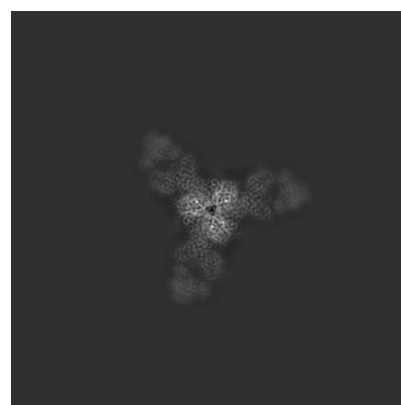
6.1.1 Primary map



X



Y



Z

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

6.2 Central slices [i](#)

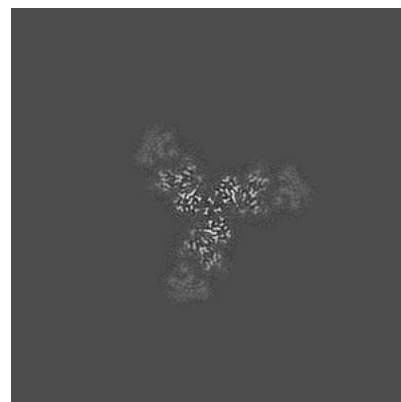
6.2.1 Primary map



X Index: 150



Y Index: 150

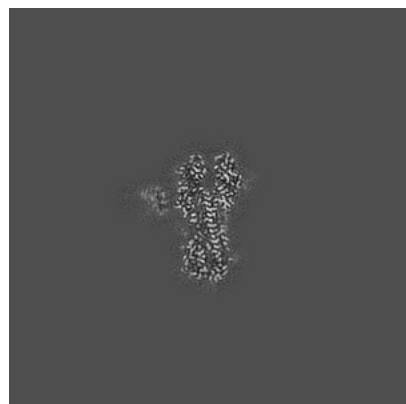


Z Index: 150

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

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 154



Y Index: 156

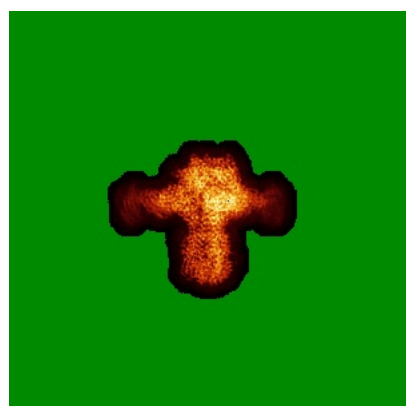


Z Index: 157

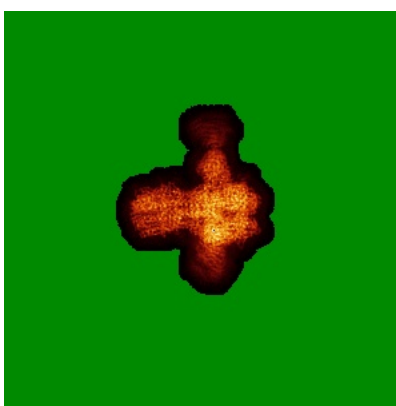
The images above show the largest variance slices of the map in three orthogonal directions.

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

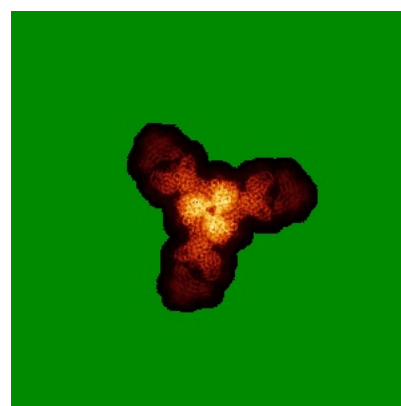
6.4.1 Primary map



X



Y

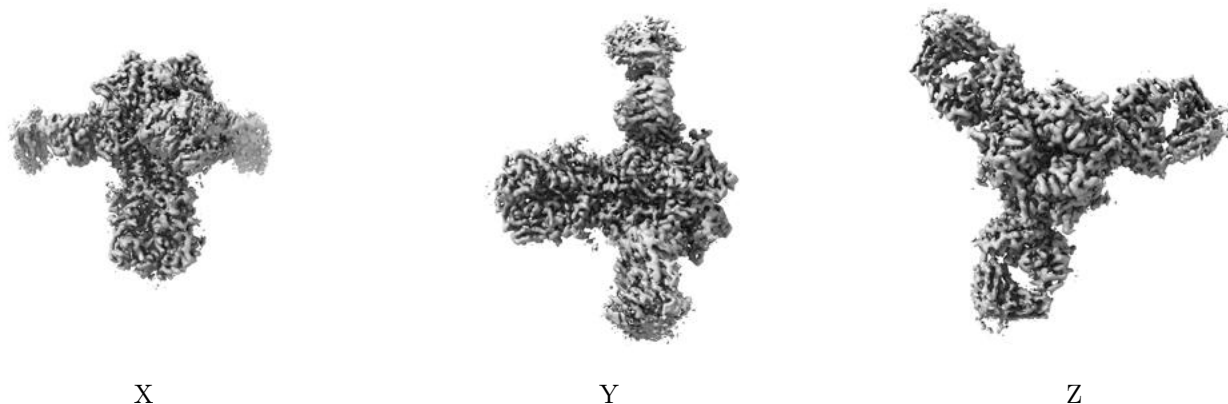


Z

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

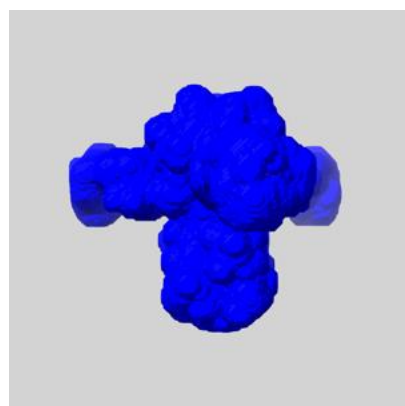
6.6 Mask visualisation [i](#)

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

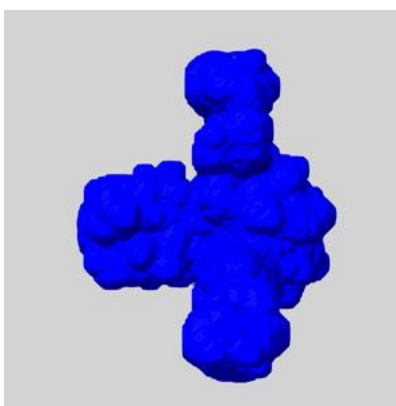
A mask typically either:

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

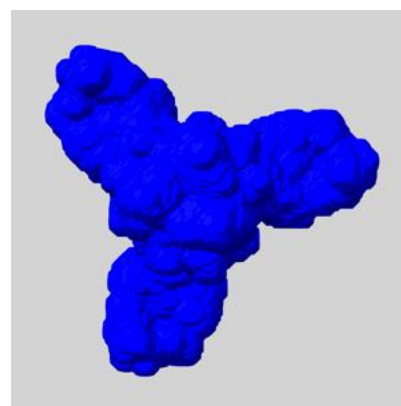
6.6.1 emd_22652_msk_1.map [i](#)



X



Y

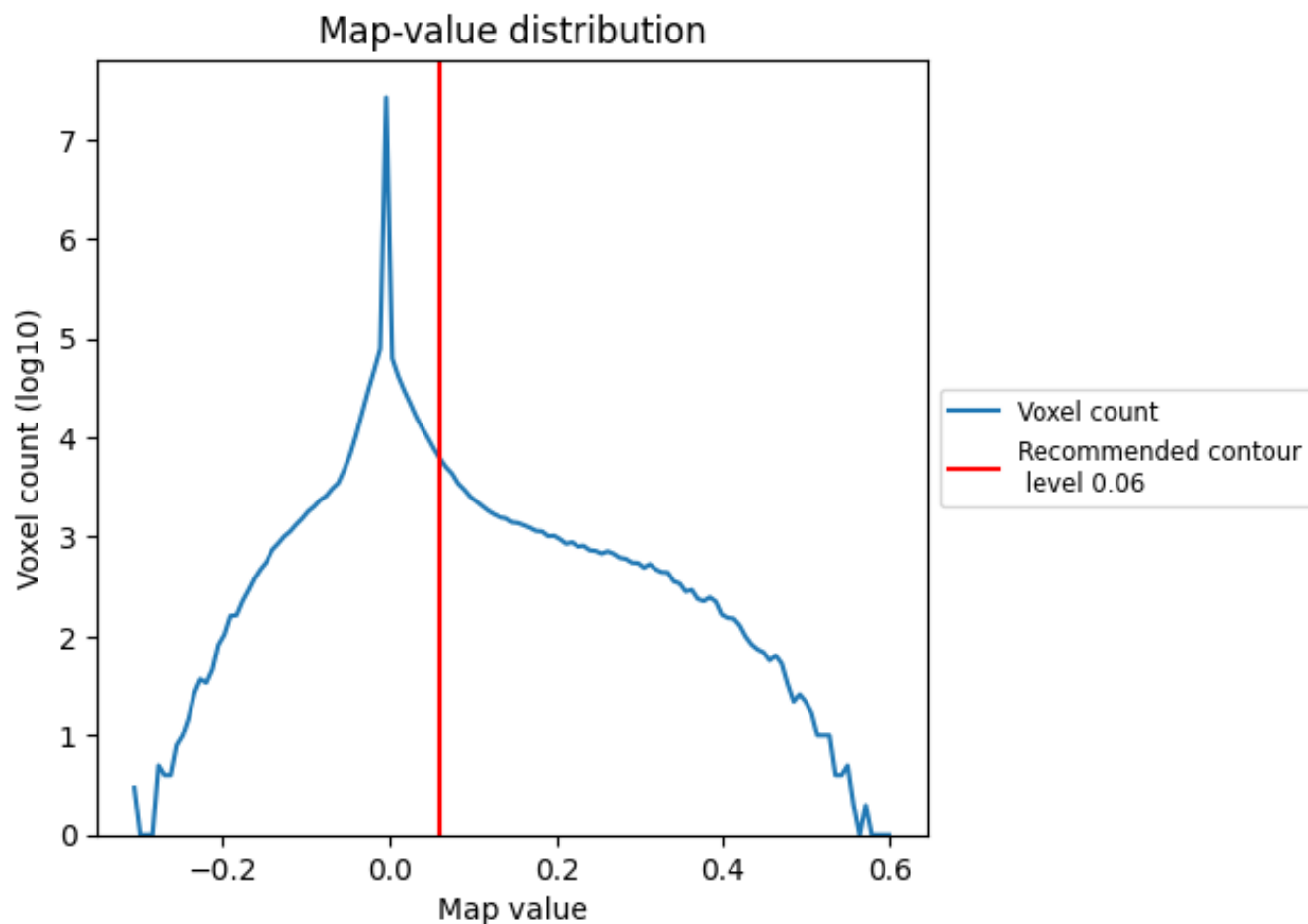


Z

7 Map analysis [i](#)

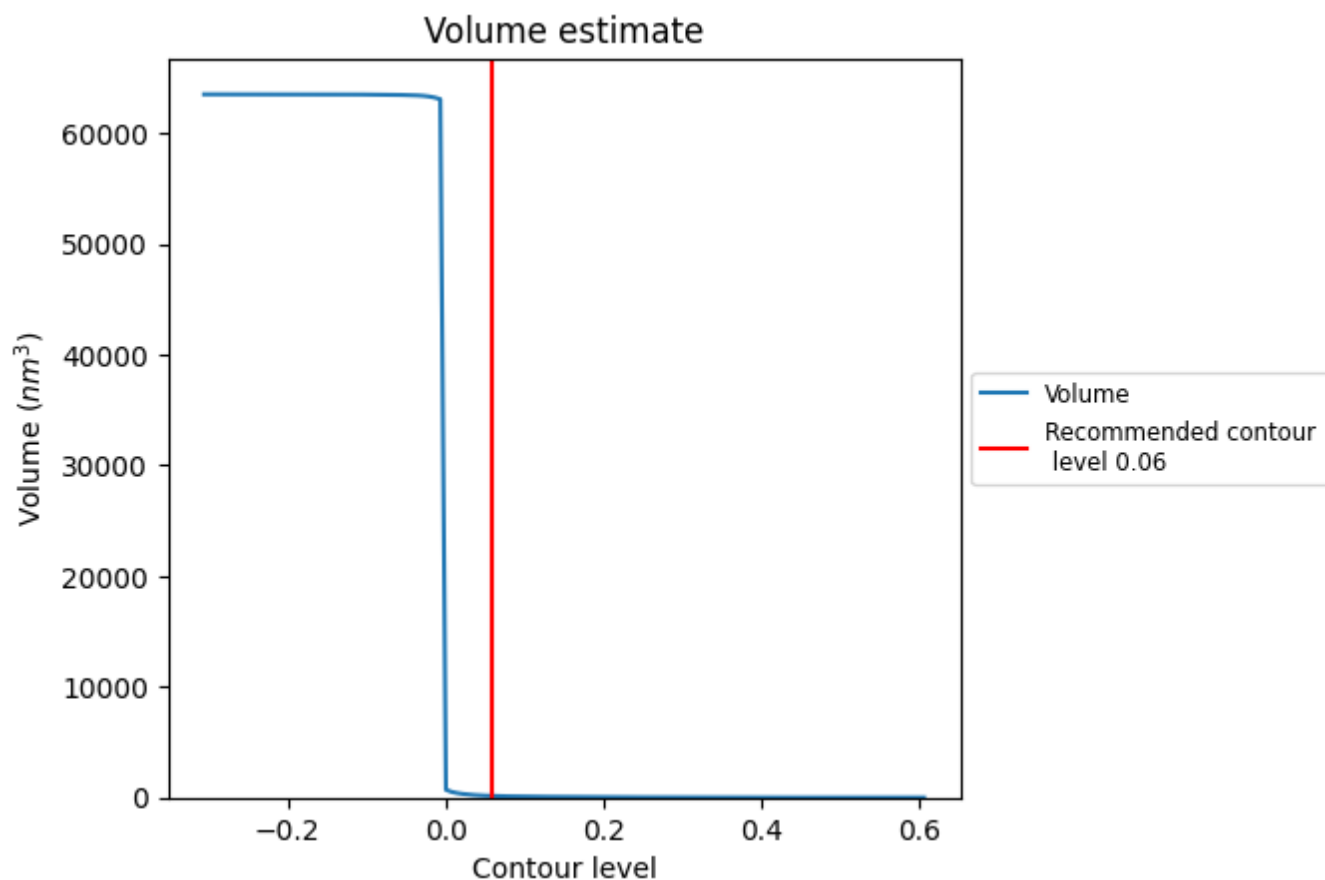
This section contains the results of statistical analysis of the map.

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.

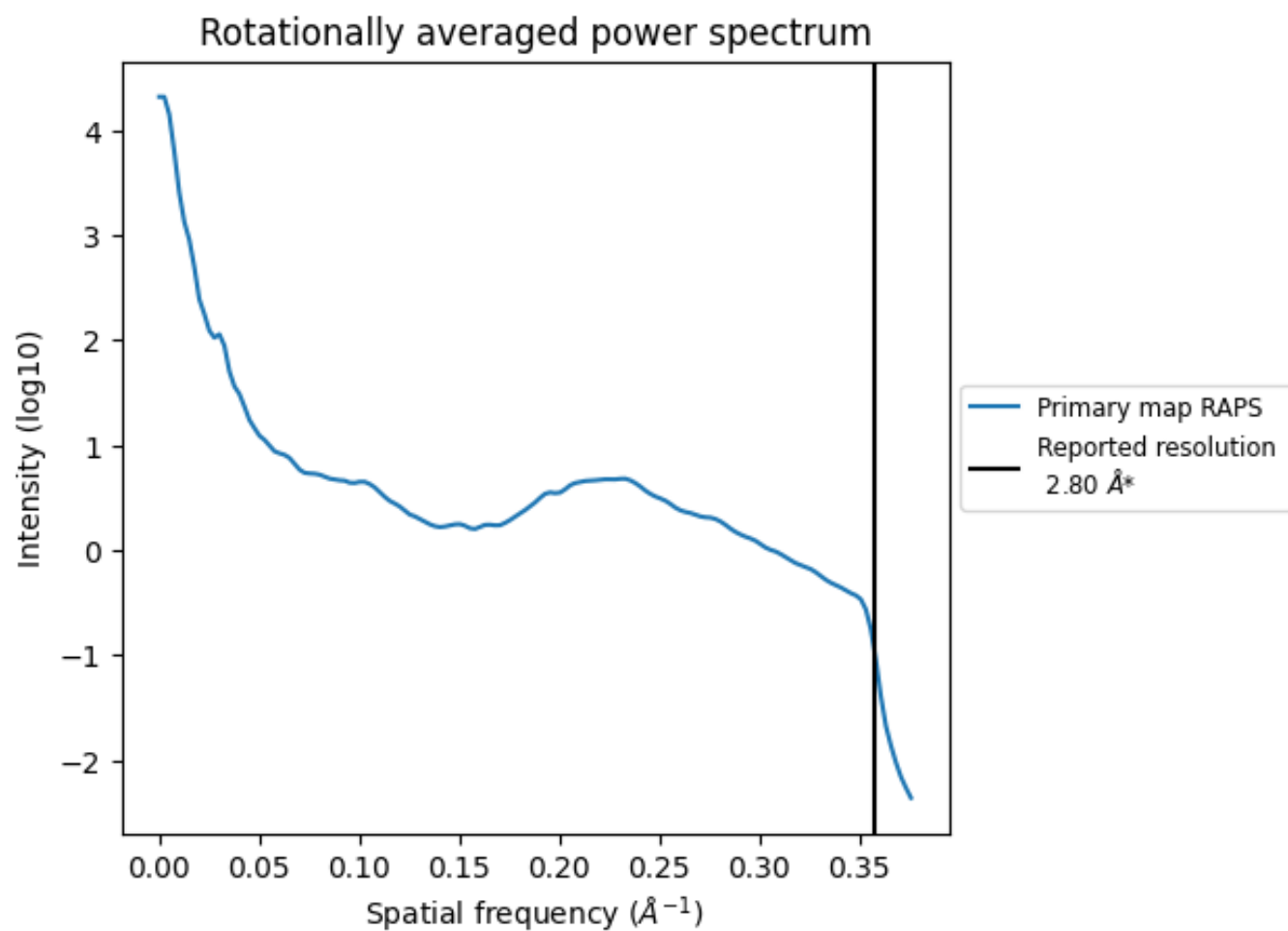
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 145 nm³; this corresponds to an approximate mass of 131 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 ⓘ

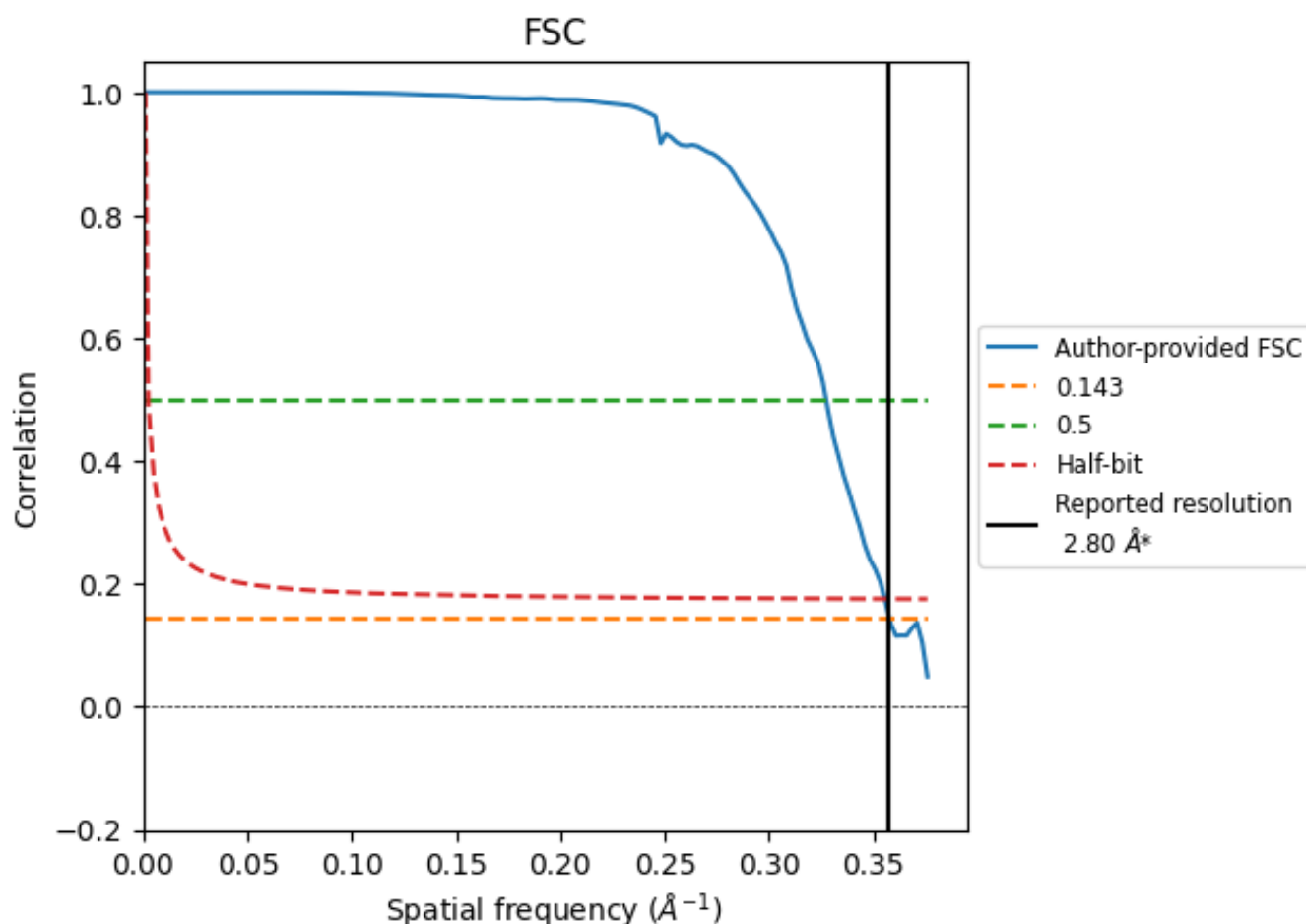


*Reported resolution corresponds to spatial frequency of 0.357 Å⁻¹

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 0.357 Å⁻¹

8.2 Resolution estimates [i](#)

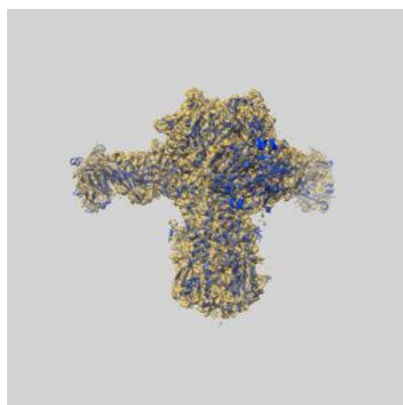
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.80	-	-
Author-provided FSC curve	2.79	3.06	2.81
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

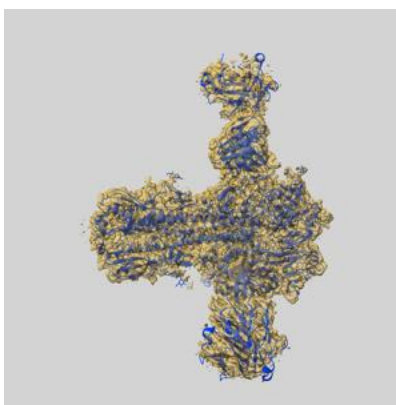
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-22652 and PDB model 7K37. Per-residue inclusion information can be found in section [3](#) on page [10](#).

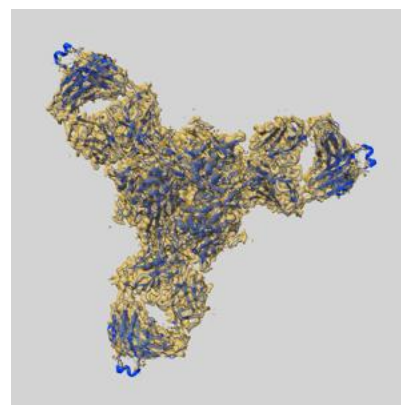
9.1 Map-model overlay [i](#)



X



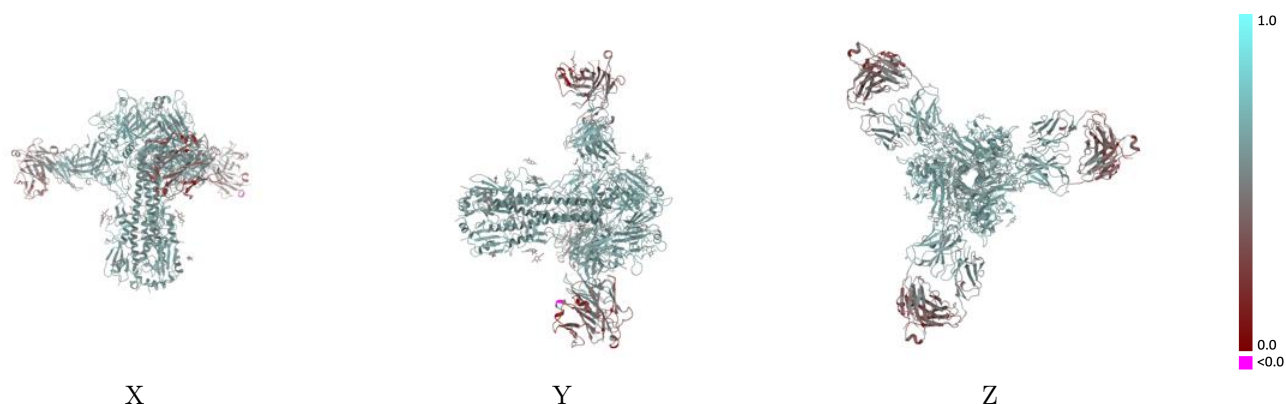
Y



Z

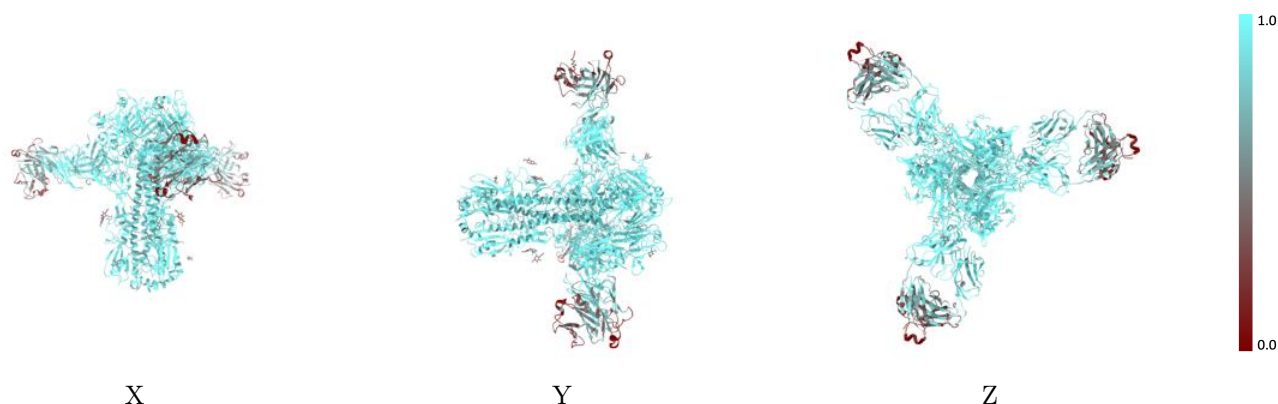
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



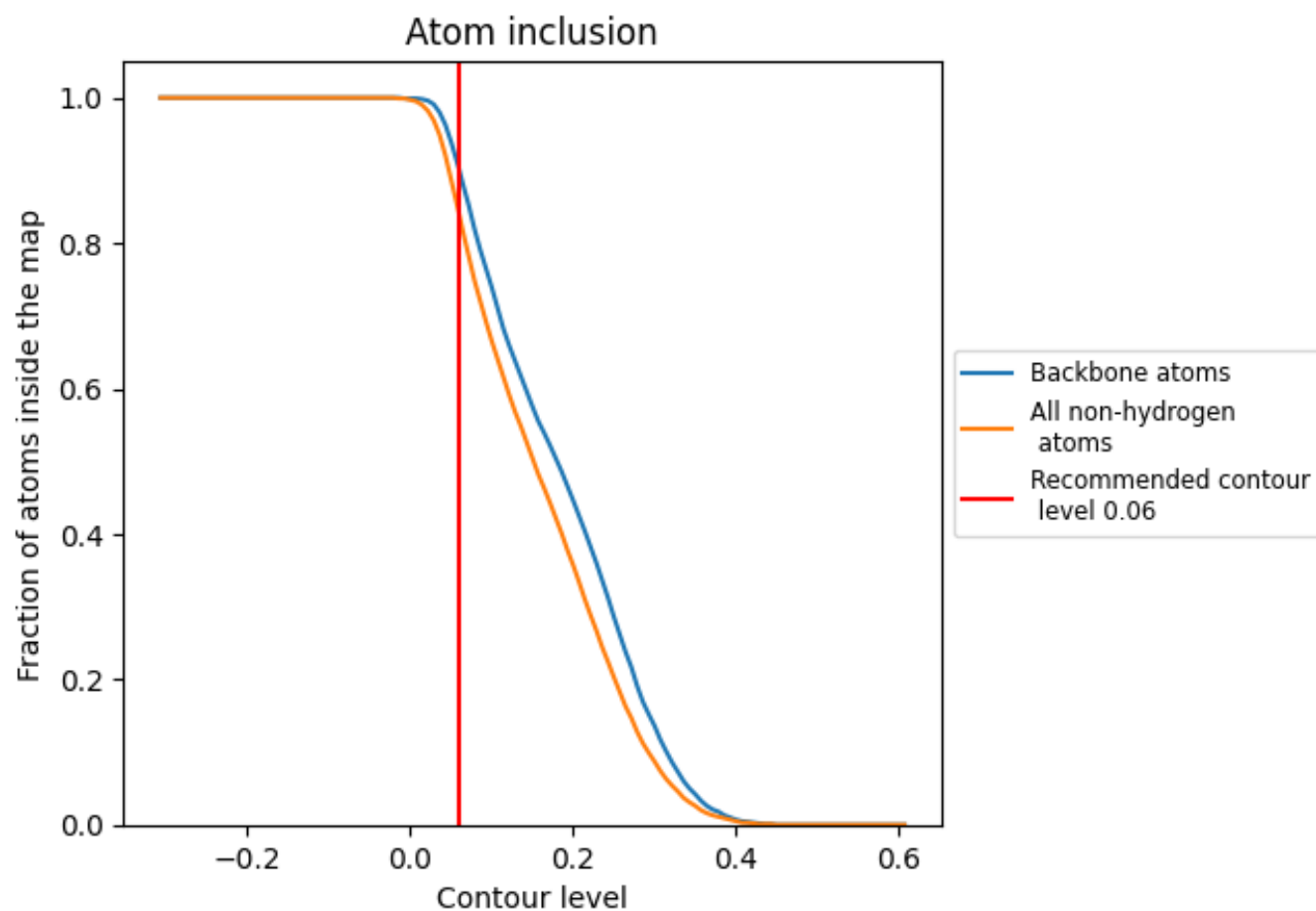
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.06).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 84% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8420	 0.5440
A	 0.9610	 0.6130
B	 0.9090	 0.5840
C	 0.9660	 0.6130
D	 0.9050	 0.5840
E	 0.9680	 0.6150
F	 0.9060	 0.5830
G	 0.6420	 0.4190
H	 0.7890	 0.5220
I	 0.6480	 0.4200
J	 0.7900	 0.5200
K	 0.6420	 0.4240
L	 0.7740	 0.5180
M	 0.5000	 0.4040
N	 0.8360	 0.5470
O	 0.8720	 0.5500
P	 0.4290	 0.4050
Q	 0.8530	 0.5610
R	 0.8620	 0.5270
S	 0.4640	 0.4100
T	 0.8530	 0.5500
U	 0.8830	 0.5400

