



Full wwPDB EM Validation Report ⓘ

May 11, 2026 – 12:26 PM JST

PDB ID : 9X5W / pdb_00009x5w
EMDB ID : EMD-66599
Title : B/Brisbane/60/2008 HA in complex with BP-1A
Authors : Nguyen, V.H.T.; Ma, C.
Deposited on : 2025-10-14
Resolution : 2.76 Å (reported)
Based on initial model : 6FYW

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 : 1.8.5 (274361), CSD as541be (2020)
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

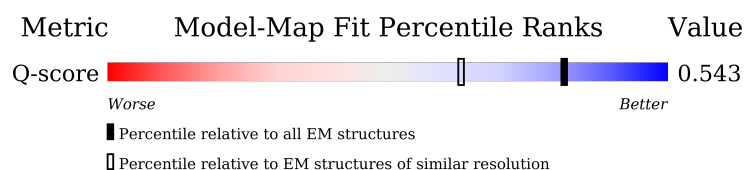
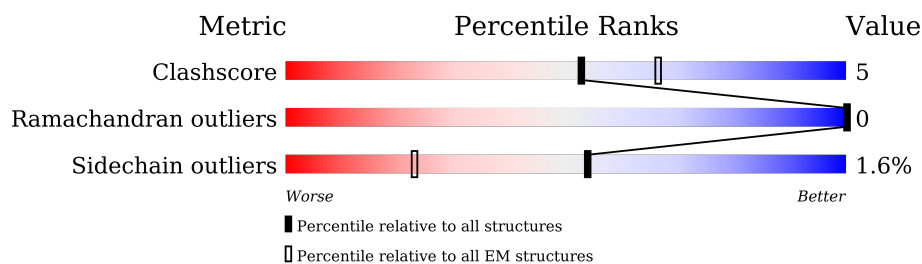
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.76 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.




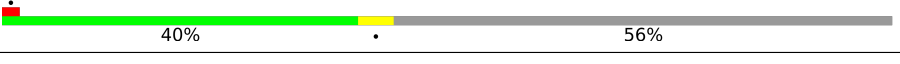



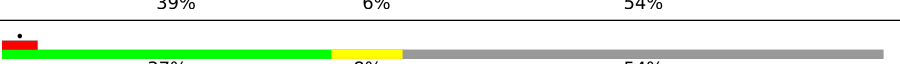
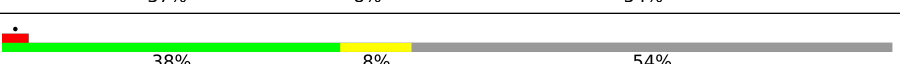
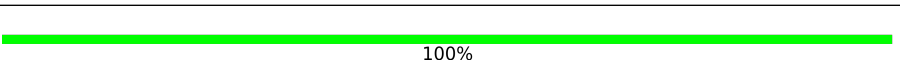
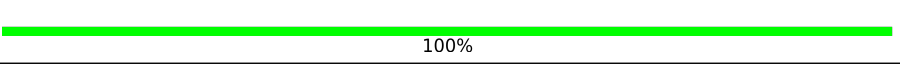
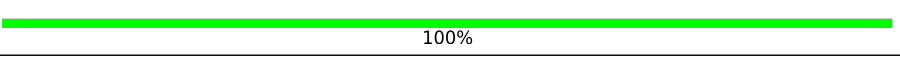
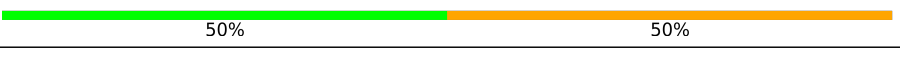
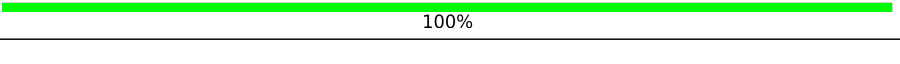
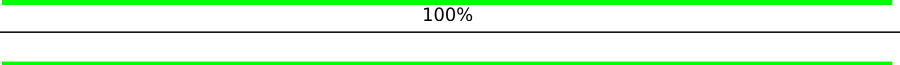
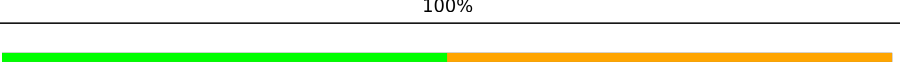
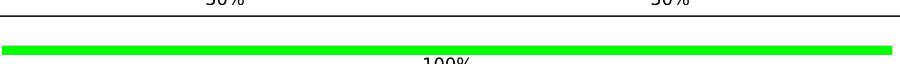
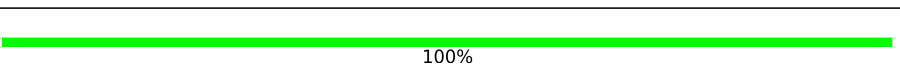
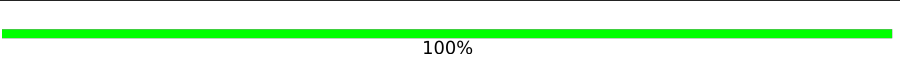



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10642 (2.26 - 3.26)

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	362	 79% 14% 7%
1	C	362	 79% 14% 7%
1	E	362	 78% 14% 7%
2	B	235	 39% 5% 56%

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Mol	Chain	Length	Quality of chain
2	D	235	
2	F	235	
3	G	474	
3	H	474	
3	J	474	
4	I	233	
4	K	233	
4	L	233	
5	M	2	
5	N	2	
5	O	2	
5	P	2	
5	Q	2	
5	R	2	
5	S	2	
5	T	2	
5	U	2	
5	V	2	
5	W	2	
5	X	2	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 15978 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 HA1 chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	336	Total	C	N	O	S	3	0
			2570	1612	458	486	14		
1	C	336	Total	C	N	O	S	3	0
			2570	1612	458	486	14		
1	E	336	Total	C	N	O	S	3	0
			2570	1612	458	486	14		

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	103	Total	C	N	O	S	0	0
			807	502	137	166	2		
2	D	103	Total	C	N	O	S	0	0
			807	502	137	166	2		
2	F	103	Total	C	N	O	S	0	0
			807	502	137	166	2		

There are 141 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	536	ASP	-	expression tag	UNP C0LT38
B	537	ILE	-	expression tag	UNP C0LT38
B	538	ARG	-	expression tag	UNP C0LT38
B	539	SER	-	expression tag	UNP C0LT38
B	540	LEU	-	expression tag	UNP C0LT38
B	541	VAL	-	expression tag	UNP C0LT38
B	542	PRO	-	expression tag	UNP C0LT38
B	543	ARG	-	expression tag	UNP C0LT38
B	544	GLY	-	expression tag	UNP C0LT38
B	545	SER	-	expression tag	UNP C0LT38
B	546	PRO	-	expression tag	UNP C0LT38
B	547	GLY	-	expression tag	UNP C0LT38
B	548	SER	-	expression tag	UNP C0LT38

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Chain	Residue	Modelled	Actual	Comment	Reference
B	549	GLY	-	expression tag	UNP C0LT38
B	550	TYR	-	expression tag	UNP C0LT38
B	551	ILE	-	expression tag	UNP C0LT38
B	552	PRO	-	expression tag	UNP C0LT38
B	553	GLU	-	expression tag	UNP C0LT38
B	554	ALA	-	expression tag	UNP C0LT38
B	555	PRO	-	expression tag	UNP C0LT38
B	556	ARG	-	expression tag	UNP C0LT38
B	557	ASP	-	expression tag	UNP C0LT38
B	558	GLY	-	expression tag	UNP C0LT38
B	559	GLN	-	expression tag	UNP C0LT38
B	560	ALA	-	expression tag	UNP C0LT38
B	561	TYR	-	expression tag	UNP C0LT38
B	562	VAL	-	expression tag	UNP C0LT38
B	563	ARG	-	expression tag	UNP C0LT38
B	564	LYS	-	expression tag	UNP C0LT38
B	565	ASP	-	expression tag	UNP C0LT38
B	566	GLY	-	expression tag	UNP C0LT38
B	567	GLU	-	expression tag	UNP C0LT38
B	568	TRP	-	expression tag	UNP C0LT38
B	569	VAL	-	expression tag	UNP C0LT38
B	570	LEU	-	expression tag	UNP C0LT38
B	571	LEU	-	expression tag	UNP C0LT38
B	572	SER	-	expression tag	UNP C0LT38
B	573	THR	-	expression tag	UNP C0LT38
B	574	PHE	-	expression tag	UNP C0LT38
B	575	LEU	-	expression tag	UNP C0LT38
B	576	GLY	-	expression tag	UNP C0LT38
B	577	HIS	-	expression tag	UNP C0LT38
B	578	HIS	-	expression tag	UNP C0LT38
B	579	HIS	-	expression tag	UNP C0LT38
B	580	HIS	-	expression tag	UNP C0LT38
B	581	HIS	-	expression tag	UNP C0LT38
B	582	HIS	-	expression tag	UNP C0LT38
D	536	ASP	-	expression tag	UNP C0LT38
D	537	ILE	-	expression tag	UNP C0LT38
D	538	ARG	-	expression tag	UNP C0LT38
D	539	SER	-	expression tag	UNP C0LT38
D	540	LEU	-	expression tag	UNP C0LT38
D	541	VAL	-	expression tag	UNP C0LT38
D	542	PRO	-	expression tag	UNP C0LT38
D	543	ARG	-	expression tag	UNP C0LT38

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Chain	Residue	Modelled	Actual	Comment	Reference
D	544	GLY	-	expression tag	UNP C0LT38
D	545	SER	-	expression tag	UNP C0LT38
D	546	PRO	-	expression tag	UNP C0LT38
D	547	GLY	-	expression tag	UNP C0LT38
D	548	SER	-	expression tag	UNP C0LT38
D	549	GLY	-	expression tag	UNP C0LT38
D	550	TYR	-	expression tag	UNP C0LT38
D	551	ILE	-	expression tag	UNP C0LT38
D	552	PRO	-	expression tag	UNP C0LT38
D	553	GLU	-	expression tag	UNP C0LT38
D	554	ALA	-	expression tag	UNP C0LT38
D	555	PRO	-	expression tag	UNP C0LT38
D	556	ARG	-	expression tag	UNP C0LT38
D	557	ASP	-	expression tag	UNP C0LT38
D	558	GLY	-	expression tag	UNP C0LT38
D	559	GLN	-	expression tag	UNP C0LT38
D	560	ALA	-	expression tag	UNP C0LT38
D	561	TYR	-	expression tag	UNP C0LT38
D	562	VAL	-	expression tag	UNP C0LT38
D	563	ARG	-	expression tag	UNP C0LT38
D	564	LYS	-	expression tag	UNP C0LT38
D	565	ASP	-	expression tag	UNP C0LT38
D	566	GLY	-	expression tag	UNP C0LT38
D	567	GLU	-	expression tag	UNP C0LT38
D	568	TRP	-	expression tag	UNP C0LT38
D	569	VAL	-	expression tag	UNP C0LT38
D	570	LEU	-	expression tag	UNP C0LT38
D	571	LEU	-	expression tag	UNP C0LT38
D	572	SER	-	expression tag	UNP C0LT38
D	573	THR	-	expression tag	UNP C0LT38
D	574	PHE	-	expression tag	UNP C0LT38
D	575	LEU	-	expression tag	UNP C0LT38
D	576	GLY	-	expression tag	UNP C0LT38
D	577	HIS	-	expression tag	UNP C0LT38
D	578	HIS	-	expression tag	UNP C0LT38
D	579	HIS	-	expression tag	UNP C0LT38
D	580	HIS	-	expression tag	UNP C0LT38
D	581	HIS	-	expression tag	UNP C0LT38
D	582	HIS	-	expression tag	UNP C0LT38
F	536	ASP	-	expression tag	UNP C0LT38
F	537	ILE	-	expression tag	UNP C0LT38
F	538	ARG	-	expression tag	UNP C0LT38

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Chain	Residue	Modelled	Actual	Comment	Reference
F	539	SER	-	expression tag	UNP C0LT38
F	540	LEU	-	expression tag	UNP C0LT38
F	541	VAL	-	expression tag	UNP C0LT38
F	542	PRO	-	expression tag	UNP C0LT38
F	543	ARG	-	expression tag	UNP C0LT38
F	544	GLY	-	expression tag	UNP C0LT38
F	545	SER	-	expression tag	UNP C0LT38
F	546	PRO	-	expression tag	UNP C0LT38
F	547	GLY	-	expression tag	UNP C0LT38
F	548	SER	-	expression tag	UNP C0LT38
F	549	GLY	-	expression tag	UNP C0LT38
F	550	TYR	-	expression tag	UNP C0LT38
F	551	ILE	-	expression tag	UNP C0LT38
F	552	PRO	-	expression tag	UNP C0LT38
F	553	GLU	-	expression tag	UNP C0LT38
F	554	ALA	-	expression tag	UNP C0LT38
F	555	PRO	-	expression tag	UNP C0LT38
F	556	ARG	-	expression tag	UNP C0LT38
F	557	ASP	-	expression tag	UNP C0LT38
F	558	GLY	-	expression tag	UNP C0LT38
F	559	GLN	-	expression tag	UNP C0LT38
F	560	ALA	-	expression tag	UNP C0LT38
F	561	TYR	-	expression tag	UNP C0LT38
F	562	VAL	-	expression tag	UNP C0LT38
F	563	ARG	-	expression tag	UNP C0LT38
F	564	LYS	-	expression tag	UNP C0LT38
F	565	ASP	-	expression tag	UNP C0LT38
F	566	GLY	-	expression tag	UNP C0LT38
F	567	GLU	-	expression tag	UNP C0LT38
F	568	TRP	-	expression tag	UNP C0LT38
F	569	VAL	-	expression tag	UNP C0LT38
F	570	LEU	-	expression tag	UNP C0LT38
F	571	LEU	-	expression tag	UNP C0LT38
F	572	SER	-	expression tag	UNP C0LT38
F	573	THR	-	expression tag	UNP C0LT38
F	574	PHE	-	expression tag	UNP C0LT38
F	575	LEU	-	expression tag	UNP C0LT38
F	576	GLY	-	expression tag	UNP C0LT38
F	577	HIS	-	expression tag	UNP C0LT38
F	578	HIS	-	expression tag	UNP C0LT38
F	579	HIS	-	expression tag	UNP C0LT38
F	580	HIS	-	expression tag	UNP C0LT38

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Chain	Residue	Modelled	Actual	Comment	Reference
F	581	HIS	-	expression tag	UNP C0LT38
F	582	HIS	-	expression tag	UNP C0LT38

- Molecule 3 is a protein called BP-1A heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	125	Total	C	N	O	S	0	0
			988	623	167	194	4		
3	G	125	Total	C	N	O	S	0	0
			988	623	167	194	4		
3	J	125	Total	C	N	O	S	0	0
			988	623	167	194	4		

- Molecule 4 is a protein called BP-1A light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	L	107	Total	C	N	O	S	0	0
			807	503	141	160	3		
4	I	107	Total	C	N	O	S	0	0
			807	503	141	160	3		
4	K	107	Total	C	N	O	S	0	0
			807	503	141	160	3		

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



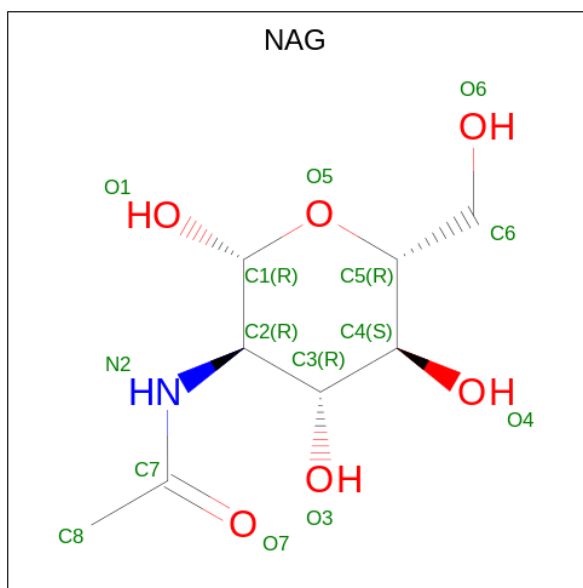
Mol	Chain	Residues	Atoms				AltConf	Trace
5	M	2	Total	C	N	O	0	0
			28	16	2	10		
5	N	2	Total	C	N	O	0	0
			28	16	2	10		
5	O	2	Total	C	N	O	0	0
			28	16	2	10		
5	P	2	Total	C	N	O	0	0
			28	16	2	10		
5	Q	2	Total	C	N	O	0	0
			28	16	2	10		

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Mol	Chain	Residues	Atoms				AltConf	Trace
5	R	2	Total	C	N	O	0	0
			28	16	2	10		
5	S	2	Total	C	N	O	0	0
			28	16	2	10		
5	T	2	Total	C	N	O	0	0
			28	16	2	10		
5	U	2	Total	C	N	O	0	0
			28	16	2	10		
5	V	2	Total	C	N	O	0	0
			28	16	2	10		
5	W	2	Total	C	N	O	0	0
			28	16	2	10		
5	X	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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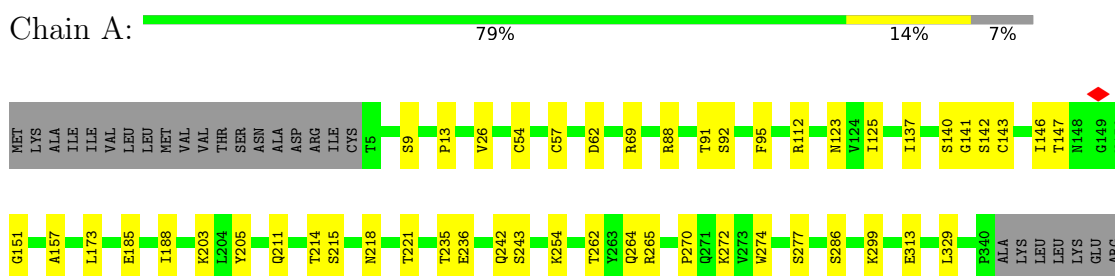
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Mol	Chain	Residues	Atoms				AltConf
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	

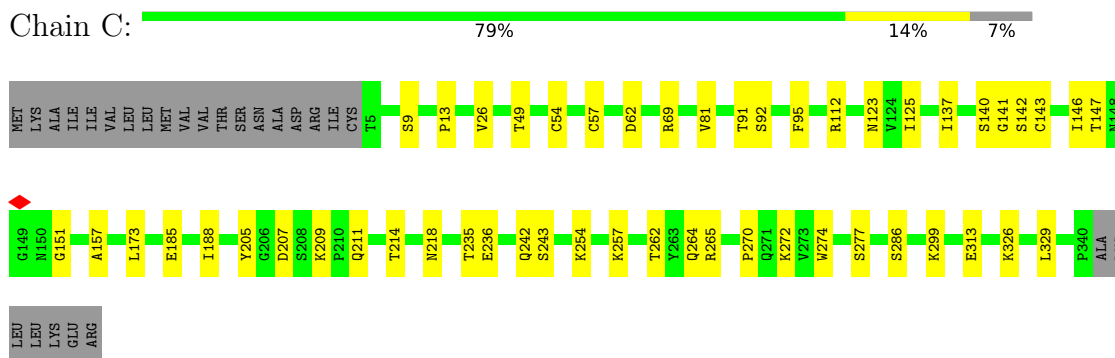
3 Residue-property plots

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

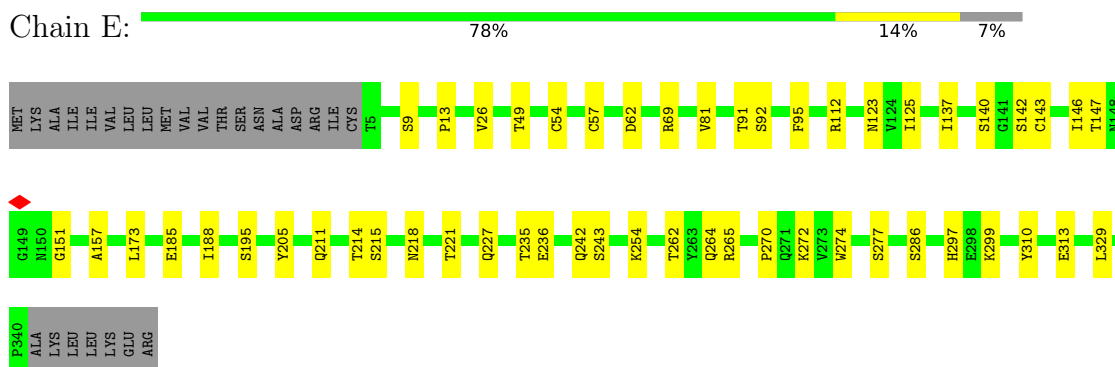
- Molecule 1: Hemagglutinin HA1 chain



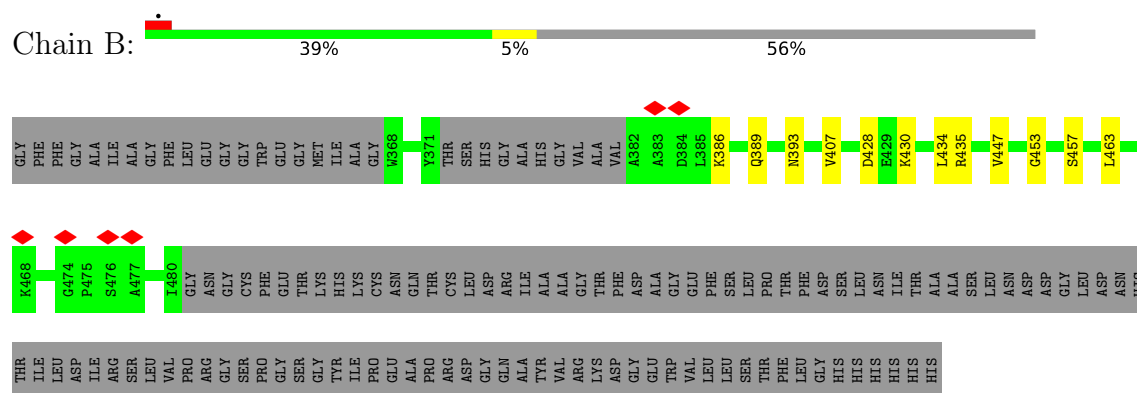
- Molecule 1: Hemagglutinin HA1 chain



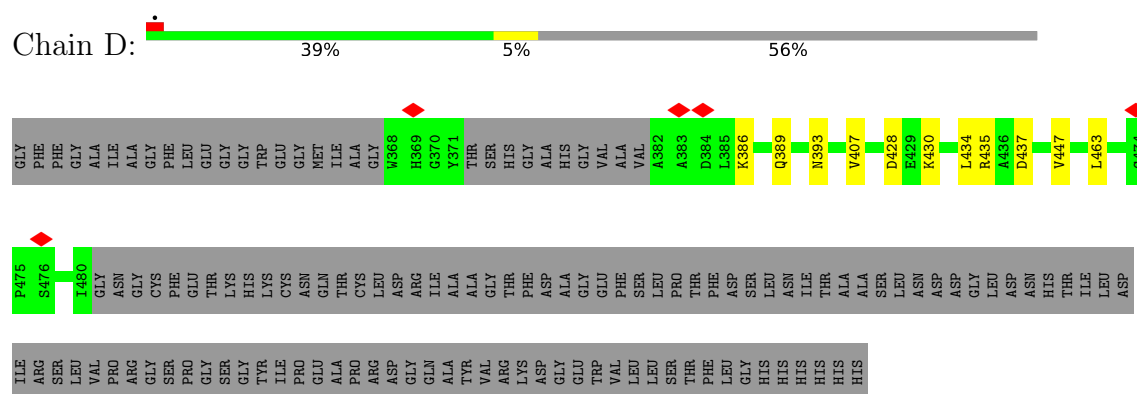
- Molecule 1: Hemagglutinin HA1 chain



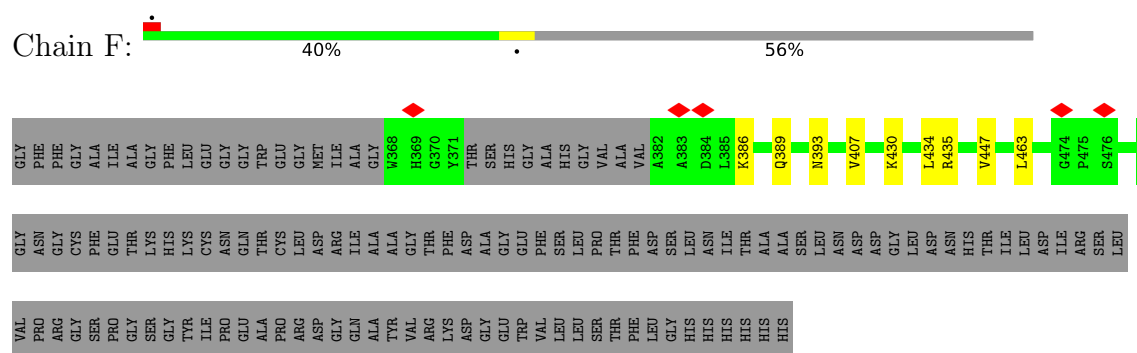
- Molecule 2: Hemagglutinin HA2 chain



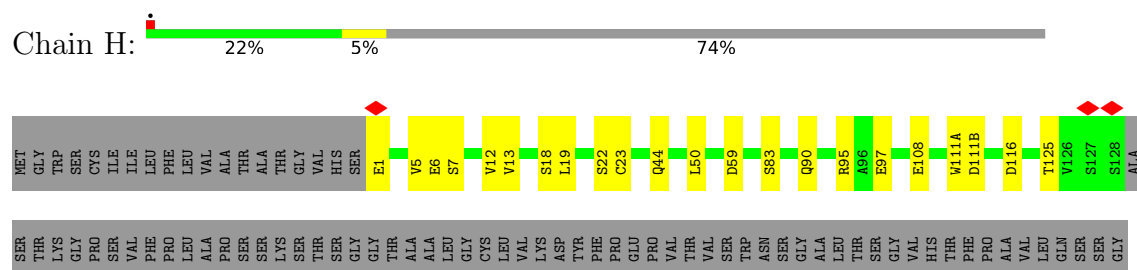
- Molecule 2: Hemagglutinin HA2 chain



- Molecule 2: Hemagglutinin HA2 chain

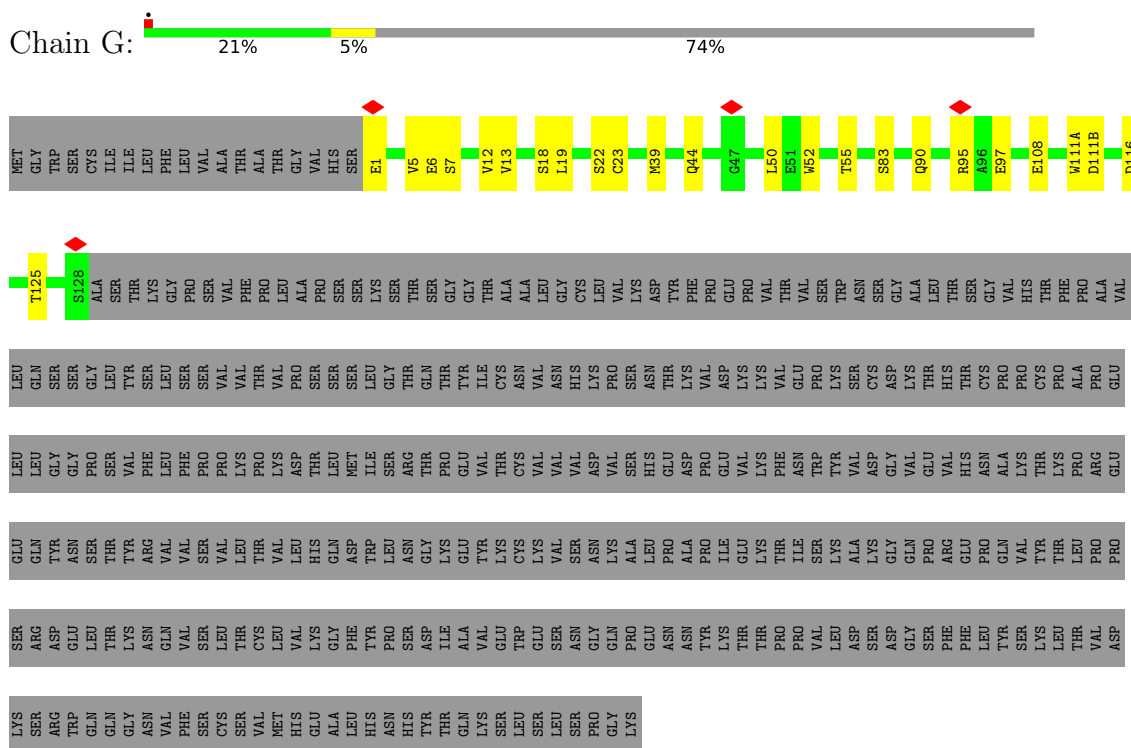


- Molecule 3: BP-1A heavy chain

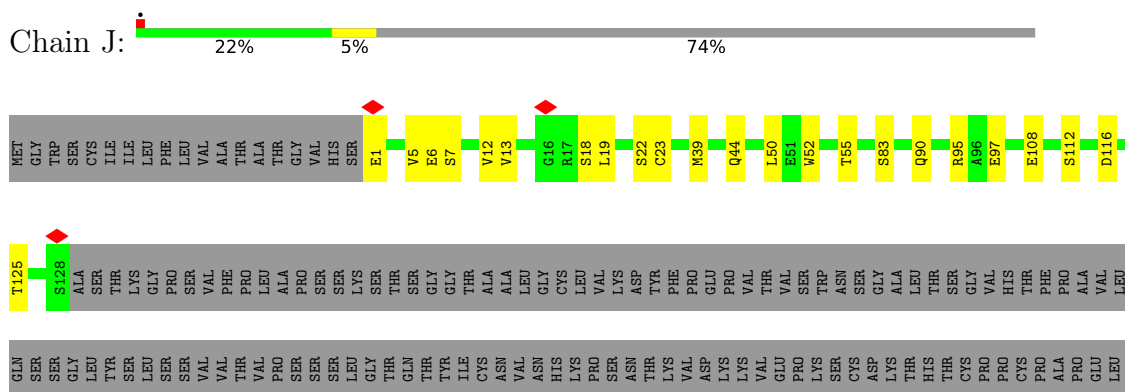


[illegible]

- Molecule 3: BP-1A heavy chain




- Molecule 3: BP-1A heavy chain




LEU	GLY	ASP	GLY	ASN	PRO	GLN	THR	THR	VAL	PHE	LEU	PRO	PRO	LEU	PRO	LYS	THR	ASP	THR	GLN	VAL	THR	THR	LEU	PRO	ARG	GLU	SER
GLN	TYR	ASN	SER	GLN	THR	THR	VAL	ARG	VAL	PHE	VAL	VAL	VAL	LEU	THR	VAL	LEU	GLN	ASP	THR	GLN	VAL	THR	LEU	PRO	ARG	GLU	SER
ARG	ASP	GLU	THR	THR	LYS	ASN	GLN	VAL	ASN	GLN	VAL	VAL	VAL	LEU	THR	VAL	VAL	GLY	THR	THR	VAL	VAL	THR	LEU	PRO	ARG	GLU	SER
SER	ARG	TRP	GLN	GLN	GLY	ASN	VAL	PHE	VAL	PHE	SER	CYS	SER	VAL	MET	HIS	GLY	ALA	LEU	HIS	THR	PRO	ASN	HIS	THR	TYR	GLY	LYS

• Molecule 4: BP-1A light chain

Chain L:  38% 8% 54%


LYS	ASP	SER	THR	TYR	SER	LEU	VAL	PHE	LEU	SER	THR	LEU	SER	ASP	GLY	GLN	LEU	VAL	GLY	LYS	THR	VAL	TYR	ALA	CYS	GLU	VAL	THR	HIS	GLN	GLY	LEU	SER	SER	PRO	VAL	THR	LYS	THR	PHE	ASN	ASN	VAL	THR	PRO	ASP	GLY	GLU	CYS									
THR	VAL	ALA	ALA	PRO	SER	VAL	PHE	ILE	PHE	PRO	PRO	SER	ASP	GLU	GLN	LEU	VAL	LYS	SER	GLY	THR	ALA	SER	VAL	VAL	CYS	LEU	LEU	ASN	ASN	PHE	TYR	PRO	ARG	GLU	ALA	LYS	VAL	GLN	THR	LYS	VAL	ASP	ASN	ALA	LEU	GLN	SER	GLY	ASN	SER	VAL	THR	THR	GLU	GLN	ASP	SER
MET	GLY	TRP	SER	CYS	ILE	ILE	PHE	LEU	VAL	ALA	ALA	THR	THR	THR	GLY	HIS	SER	E1	T5	Q6	S7	P8	A9	T10	V11	G16	E17	R18	C23	R24	S37	Q43	L52	L53	I54	S55	R67	A68	S69	G70	I71	Q95	S96	E97	D98	F99	I126	K127										

• Molecule 4: BP-1A light chain

Chain I:  39% 6% 54%

SER	SER	LEU	LEU	SER	THR	THR	LEU	LEU	SER	LYS	ALA	ASP	THR	GLY	GLN	VAL	THR	VAL	E1	Q6	S7	P8	A9	T10	V11	G16	E17	R18	S22	C23	R24	S37	G47	R67	A68	S69	S99	G70	I71	Q95	S96	E97	D98	F99	I126	K127	ARG	THR	THR	VAL	ALA	ALA
SER	VAL	PHE	ILE	SER	PRO	PRO	LEU	SER	ASP	GLY	GLN	LEU	LYS	SER	GLY	THR	ALA	VAL	VAL	VAL	LEU	ASN	PHE	TTR	PRO	ARG	GLY	ALA	LYS	VAL	GLN	THR	VAL	ASP	ASN	ALA	LEU	GLN	SER	GLY	ASN	SER	THR	GLY	THR	THR	THR	THR	VAL	ALA	ALA	
SER	LEU	SER	THR	THR	LEU	LEU	SER	LEU	SER	LYS	ALA	ASP	THR	GLY	GLN	VAL	THR	VAL	E1	Q6	S7	P8	A9	T10	V11	G16	E17	R18	S22	C23	R24	S37	G47	R67	A68	S69	S99	G70	I71	Q95	S96	E97	D98	F99	I126	K127	ARG	THR	THR	VAL	ALA	ALA

• Molecule 4: BP-1A light chain

Chain K:  37% 8% 54%

LYS	ASP	SER	THR	TYR	SER	LEU	SER	THR	LEU	SER	THR	LEU	SER	ASP	GLY	GLN	VAL	GLY	E1	Q6	S7	P8	A9	T10	V11	G16	E17	R18	S22	C23	R24	S37	Q43	L52	L53	ASN	I54	S55	R67	I71	Q95	S96	E97	D98	F99	R116	I126	K127	ARG									
THR	VAL	ALA	ALA	PRO	SER	VAL	PHE	ILE	PHE	PRO	PRO	THR	SER	ASP	GLU	GLN	LEU	VAL	GLY	SER	SER	GLY	THR	ALA	VAL	VAL	LEU	LEU	ASN	ASN	PHE	TYR	PRO	ARG	GLU	ALA	LYS	VAL	GLN	TRP	LYS	VAL	ASP	ASP	ASN	GLY	LEU	GLN	SER	GLN	GLU	SER	VAL	THR	GLU	GLN	ASP	SER

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

Chain M:  100%



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

Chain N:  100%



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

Chain O:  100%



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

Chain P:  50% 50%



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

Chain Q:  100%



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

Chain R:  100%



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

Chain S:  100%



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

Chain T:  50% 50%



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

Chain U:  100%



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

Chain V:  100%



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

Chain W:  100%



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

Chain X:  50% 50%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	343055	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 ($e^-/\text{\AA}^2$)	46.40	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	4.855	Depositor
Minimum map value	-3.841	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.075	Depositor
Recommended contour level	0.227	Depositor
Map size (\AA)	318.72, 318.72, 318.72	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.83, 0.83, 0.83	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

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	0.25	0/2630	0.48	0/3574
1	C	0.25	0/2630	0.48	0/3574
1	E	0.25	0/2630	0.48	0/3574
2	B	0.24	0/812	0.57	0/1092
2	D	0.24	0/812	0.57	0/1092
2	F	0.24	0/812	0.57	0/1092
3	G	0.30	0/1013	0.73	5/1372 (0.4%)
3	H	0.30	0/1013	0.73	5/1372 (0.4%)
3	J	0.30	0/1013	0.73	5/1372 (0.4%)
4	I	0.25	0/824	0.70	3/1117 (0.3%)
4	K	0.25	0/824	0.70	3/1117 (0.3%)
4	L	0.25	0/824	0.70	3/1117 (0.3%)
All	All	0.26	0/15837	0.58	24/21465 (0.1%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	97	GLU	CA-CB-CG	9.64	133.38	114.10
4	K	97	GLU	CA-CB-CG	9.62	133.33	114.10
4	I	97	GLU	CA-CB-CG	9.60	133.30	114.10
3	J	97	GLU	CA-CB-CG	8.74	131.58	114.10
3	G	97	GLU	CA-CB-CG	8.73	131.57	114.10
3	H	97	GLU	CA-CB-CG	8.72	131.54	114.10
4	L	95	GLN	CA-CB-CG	6.70	127.50	114.10
4	I	95	GLN	CA-CB-CG	6.67	127.45	114.10
4	K	95	GLN	CA-CB-CG	6.67	127.43	114.10
3	H	95	ARG	CG-CD-NE	6.33	125.93	112.00
3	J	95	ARG	CG-CD-NE	6.31	125.88	112.00
3	G	95	ARG	CG-CD-NE	6.28	125.81	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	1	GLU	CA-CB-CG	6.26	126.63	114.10
3	G	1	GLU	CA-CB-CG	6.25	126.59	114.10
3	H	1	GLU	CA-CB-CG	6.24	126.59	114.10
4	K	97	GLU	CB-CG-CD	6.13	123.02	112.60
4	L	97	GLU	CB-CG-CD	6.13	123.02	112.60
4	I	97	GLU	CB-CG-CD	6.13	123.01	112.60
3	G	1	GLU	CB-CG-CD	5.43	121.83	112.60
3	J	1	GLU	CB-CG-CD	5.43	121.83	112.60
3	H	1	GLU	CB-CG-CD	5.42	121.81	112.60
3	H	95	ARG	CD-NE-CZ	5.28	131.79	124.40
3	G	95	ARG	CD-NE-CZ	5.23	131.72	124.40
3	J	95	ARG	CD-NE-CZ	5.20	131.68	124.40

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	2570	0	2562	30	0
1	C	2570	0	2562	31	0
1	E	2570	0	2562	29	0
2	B	807	0	818	8	0
2	D	807	0	818	7	0
2	F	807	0	818	6	0
3	G	988	0	928	13	0
3	H	988	0	928	12	0
3	J	988	0	928	12	0
4	I	807	0	790	9	0
4	K	807	0	790	12	0
4	L	807	0	790	11	0
5	M	28	0	25	0	0
5	N	28	0	25	0	0
5	O	28	0	25	0	0
5	P	28	0	25	1	0
5	Q	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	R	28	0	25	0	0
5	S	28	0	25	0	0
5	T	28	0	25	1	0
5	U	28	0	25	0	0
5	V	28	0	25	0	0
5	W	28	0	25	0	0
5	X	28	0	25	1	0
6	A	42	0	39	0	0
6	C	42	0	39	0	0
6	E	42	0	39	0	0
All	All	15978	0	15711	167	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 5.

All (167) 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:235:THR:HG22	1:A:236:GLU:HG3	1.41	1.03
1:C:235:THR:HG22	1:C:236:GLU:HG3	1.40	0.99
1:E:235:THR:HG22	1:E:236:GLU:HG3	1.41	0.99
3:H:6:GLU:HG2	3:H:23:CYS:CB	2.12	0.80
3:G:6:GLU:HG2	3:G:23:CYS:CB	2.12	0.80
3:J:6:GLU:HG2	3:J:23:CYS:CB	2.11	0.80
3:J:6:GLU:HG2	3:J:23:CYS:HB2	1.65	0.78
3:G:6:GLU:HG2	3:G:23:CYS:HB2	1.65	0.77
4:K:7:SER:O	4:K:22:SER:OG	2.03	0.77
3:H:6:GLU:HG2	3:H:23:CYS:HB2	1.65	0.76
4:I:7:SER:O	4:I:22:SER:OG	2.02	0.74
4:K:126:ILE:HG22	4:K:127:LYS:N	2.08	0.68
4:I:126:ILE:HG22	4:I:127:LYS:N	2.08	0.67
1:A:235:THR:HG22	1:A:236:GLU:CG	2.23	0.67
4:L:126:ILE:HG22	4:L:127:LYS:N	2.08	0.66
3:J:108:GLU:HA	3:J:116:ASP:OD2	1.99	0.63
1:C:235:THR:HG22	1:C:236:GLU:CG	2.23	0.62
1:E:235:THR:HG22	1:E:236:GLU:CG	2.23	0.62
3:H:108:GLU:HA	3:H:116:ASP:OD2	1.99	0.62
3:G:108:GLU:HA	3:G:116:ASP:OD2	1.99	0.62
1:E:54:CYS:HB3	1:E:57:CYS:H	1.65	0.60
2:B:386:LYS:O	2:B:386:LYS:HG2	2.01	0.60
1:C:54:CYS:HB3	1:C:57:CYS:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:CYS:HB3	1:A:57:CYS:H	1.65	0.60
2:F:386:LYS:O	2:F:386:LYS:HG2	2.01	0.60
2:D:386:LYS:HG2	2:D:386:LYS:O	2.01	0.59
1:C:185:GLU:HG2	1:C:277:SER:HB3	1.85	0.59
1:E:185:GLU:HG2	1:E:277:SER:HB3	1.85	0.59
1:A:185:GLU:HG2	1:A:277:SER:HB3	1.85	0.58
2:D:407:VAL:HB	2:D:435:ARG:HH22	1.69	0.58
2:F:407:VAL:HB	2:F:435:ARG:HH22	1.69	0.58
1:A:146:ILE:HG13	1:A:147:THR:HG23	1.86	0.58
4:L:126:ILE:CG2	4:L:127:LYS:N	2.67	0.58
1:E:243:SER:HA	5:X:1:NAG:H62	1.86	0.58
1:C:146:ILE:HG13	1:C:147:THR:HG23	1.86	0.57
4:I:126:ILE:CG2	4:I:127:LYS:N	2.67	0.57
1:A:141:GLY:HA3	3:H:111(B):ASP:HA	1.86	0.57
1:A:243:SER:HA	5:P:1:NAG:H62	1.86	0.57
2:B:407:VAL:HB	2:B:435:ARG:HH22	1.69	0.57
1:E:146:ILE:HG13	1:E:147:THR:HG23	1.86	0.57
4:K:126:ILE:CG2	4:K:127:LYS:N	2.67	0.57
1:C:218:ASN:HD21	1:C:254:LYS:H	1.53	0.56
1:A:218:ASN:HD21	1:A:254:LYS:H	1.53	0.56
1:C:243:SER:HA	5:T:1:NAG:H62	1.87	0.56
1:E:218:ASN:HD21	1:E:254:LYS:H	1.53	0.55
3:J:18:SER:OG	3:J:90:GLN:NE2	2.40	0.55
1:E:235:THR:CG2	1:E:236:GLU:HG3	2.27	0.55
3:H:18:SER:OG	3:H:90:GLN:NE2	2.40	0.55
3:G:18:SER:OG	3:G:90:GLN:NE2	2.40	0.55
1:A:62:ASP:OD2	1:A:112:ARG:NH2	2.38	0.53
1:C:137:ILE:HD12	1:C:157:ALA:HB1	1.90	0.53
1:C:141:GLY:HA3	3:G:111(B):ASP:HA	1.91	0.53
1:C:235:THR:CG2	1:C:236:GLU:HG3	2.27	0.53
1:A:137:ILE:HD12	1:A:157:ALA:HB1	1.90	0.53
1:E:137:ILE:HD12	1:E:157:ALA:HB1	1.90	0.52
3:J:112:SER:O	4:K:116:ARG:NH2	2.42	0.52
1:A:235:THR:CG2	1:A:236:GLU:HG3	2.27	0.51
1:C:62:ASP:OD2	1:C:112:ARG:NH2	2.38	0.51
4:L:67:ARG:NH1	4:L:71:ILE:O	2.44	0.51
4:I:67:ARG:NH1	4:I:71:ILE:O	2.44	0.51
4:I:96:SER:HA	4:I:99:PHE:HD2	1.77	0.50
4:I:8:PRO:HG2	4:I:11:VAL:HG22	1.93	0.50
4:K:96:SER:HA	4:K:99:PHE:HD2	1.77	0.50
4:K:67:ARG:NH1	4:K:71:ILE:O	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:GLY:N	3:H:111(A):TRP:O	2.34	0.50
4:L:8:PRO:HG2	4:L:11:VAL:HG22	1.93	0.50
2:B:428:ASP:OD1	2:F:430:LYS:NZ	2.40	0.50
4:K:8:PRO:HG2	4:K:11:VAL:HG22	1.93	0.49
4:L:96:SER:HA	4:L:99:PHE:HD2	1.77	0.49
1:E:62:ASP:OD2	1:E:112:ARG:NH2	2.38	0.49
3:H:6:GLU:HG2	3:H:23:CYS:HA	1.96	0.48
1:A:329:LEU:HD13	2:B:447:VAL:HG21	1.96	0.47
1:E:142:SER:OG	1:E:242:GLN:NE2	2.47	0.47
4:K:6:GLN:HG3	4:K:23:CYS:HB2	1.97	0.47
1:A:142:SER:OG	1:A:242:GLN:NE2	2.47	0.47
2:B:430:LYS:NZ	2:D:428:ASP:OD1	2.42	0.46
1:E:329:LEU:HD13	2:F:447:VAL:HG21	1.97	0.46
1:A:13:PRO:HA	1:A:26:VAL:O	2.16	0.46
4:I:127:LYS:O	4:I:127:LYS:HG2	2.16	0.46
4:K:127:LYS:O	4:K:127:LYS:HG2	2.16	0.46
4:L:6:GLN:HG3	4:L:23:CYS:HB2	1.97	0.46
4:I:6:GLN:HG3	4:I:23:CYS:HB2	1.97	0.46
3:J:6:GLU:HG2	3:J:23:CYS:HA	1.96	0.46
3:G:6:GLU:HG2	3:G:23:CYS:HA	1.96	0.46
3:J:6:GLU:HG2	3:J:23:CYS:CA	2.46	0.46
3:H:6:GLU:HG2	3:H:23:CYS:CA	2.46	0.46
2:D:430:LYS:HE2	2:D:434:LEU:HD11	1.98	0.46
2:F:430:LYS:HE2	2:F:434:LEU:HD11	1.98	0.45
4:L:127:LYS:HG2	4:L:127:LYS:O	2.16	0.45
1:C:13:PRO:HA	1:C:26:VAL:O	2.16	0.45
1:C:329:LEU:HD13	2:D:447:VAL:HG21	1.98	0.45
1:E:13:PRO:HA	1:E:26:VAL:O	2.16	0.45
3:G:6:GLU:HG2	3:G:23:CYS:CA	2.45	0.45
1:C:95:PHE:N	1:C:143:CYS:SG	2.90	0.44
3:G:12:VAL:HG22	3:G:125:THR:HB	1.99	0.44
1:A:95:PHE:N	1:A:143:CYS:SG	2.90	0.44
2:B:430:LYS:HE2	2:B:434:LEU:HD11	1.98	0.44
2:B:389:GLN:O	2:B:393:ASN:ND2	2.51	0.44
1:C:142:SER:OG	1:C:242:GLN:NE2	2.47	0.44
4:I:126:ILE:CG2	4:I:127:LYS:H	2.31	0.44
1:C:69:ARG:NH1	1:C:92:SER:OG	2.51	0.43
1:C:299:LYS:HG3	1:C:313:GLU:HG3	2.00	0.43
3:J:12:VAL:HG22	3:J:125:THR:HB	1.99	0.43
1:E:123:ASN:OD1	1:E:272:LYS:NZ	2.47	0.43
1:A:299:LYS:HG3	1:A:313:GLU:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:95:PHE:N	1:E:143:CYS:SG	2.90	0.43
3:H:12:VAL:HG22	3:H:125:THR:HB	1.99	0.43
1:E:211:GLN:HB3	1:E:264:GLN:HB2	2.01	0.43
3:J:39:MET:HE3	3:J:39:MET:HB3	1.89	0.43
1:A:69:ARG:NH1	1:A:92:SER:OG	2.52	0.43
1:A:211:GLN:HB3	1:A:264:GLN:HB2	2.01	0.43
1:C:123:ASN:OD1	1:C:272:LYS:NZ	2.47	0.43
2:D:389:GLN:O	2:D:393:ASN:ND2	2.51	0.43
1:E:205:TYR:HB3	1:E:265:ARG:HB3	2.01	0.43
2:F:389:GLN:O	2:F:393:ASN:ND2	2.51	0.43
1:C:211:GLN:HB3	1:C:264:GLN:HB2	2.01	0.42
1:A:91:THR:OG1	1:A:236:GLU:O	2.37	0.42
1:C:140:SER:N	1:C:151:GLY:O	2.52	0.42
1:C:141:GLY:N	3:G:111(A):TRP:O	2.40	0.42
1:A:205:TYR:HB3	1:A:265:ARG:HB3	2.01	0.42
2:B:453:GLY:O	2:B:457:SER:OG	2.36	0.42
1:C:188:ILE:HB	1:C:274:TRP:HB2	2.01	0.42
1:C:207:ASP:OD2	1:C:209:LYS:NZ	2.44	0.42
1:E:140:SER:N	1:E:151:GLY:O	2.52	0.42
1:E:173:LEU:O	1:E:262:THR:HA	2.20	0.42
1:A:173:LEU:O	1:A:262:THR:HA	2.20	0.42
1:C:205:TYR:HB3	1:C:265:ARG:HB3	2.01	0.42
1:E:214:THR:HB	1:E:262:THR:HB	2.01	0.42
4:L:126:ILE:CG2	4:L:127:LYS:H	2.31	0.42
1:C:214:THR:HB	1:C:262:THR:HB	2.01	0.42
1:E:69:ARG:NH1	1:E:92:SER:OG	2.52	0.42
1:E:299:LYS:HG3	1:E:313:GLU:HG3	2.00	0.42
1:C:173:LEU:O	1:C:262:THR:HA	2.20	0.42
3:J:13:VAL:HG21	3:J:19:LEU:HD22	2.02	0.42
1:A:188:ILE:HB	1:A:274:TRP:HB2	2.01	0.42
4:L:43:GLN:HB2	4:L:53:LEU:HD11	2.02	0.42
1:E:188:ILE:HB	1:E:274:TRP:HB2	2.01	0.42
3:H:44:GLN:HB2	3:H:50:LEU:HD23	2.02	0.42
3:G:52:TRP:NE1	3:G:55:THR:OG1	2.43	0.42
3:J:44:GLN:HB2	3:J:50:LEU:HD23	2.02	0.42
1:A:140:SER:N	1:A:151:GLY:O	2.52	0.41
4:K:126:ILE:CG2	4:K:127:LYS:H	2.31	0.41
1:A:214:THR:HB	1:A:262:THR:HB	2.01	0.41
1:C:91:THR:OG1	1:C:236:GLU:O	2.37	0.41
3:H:13:VAL:HG21	3:H:19:LEU:HD22	2.02	0.41
1:E:91:THR:OG1	1:E:236:GLU:O	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:39:MET:HE3	3:G:39:MET:HB3	1.89	0.41
4:K:43:GLN:HB2	4:K:53:LEU:HD11	2.02	0.41
1:E:49:THR:HG23	1:E:81:VAL:HG12	2.03	0.41
1:E:297:HIS:NE2	1:E:310:TYR:OH	2.49	0.41
3:G:13:VAL:HG21	3:G:19:LEU:HD22	2.02	0.41
1:A:125:ILE:HD12	1:A:270:PRO:HG2	2.03	0.41
1:C:49:THR:HG23	1:C:81:VAL:HG12	2.03	0.41
1:C:326:LYS:NZ	2:D:437:ASP:OD1	2.54	0.41
1:E:125:ILE:HD12	1:E:270:PRO:HG2	2.03	0.41
3:G:44:GLN:HB2	3:G:50:LEU:HD23	2.02	0.41
3:J:52:TRP:NE1	3:J:55:THR:OG1	2.43	0.41
1:A:88:ARG:HD3	1:C:257:LYS:HE3	2.03	0.41
1:A:203:LYS:HD3	3:H:59:ASP:OD1	2.21	0.41
1:E:215:SER:O	1:E:221:THR:HA	2.21	0.41
1:A:215:SER:O	1:A:221:THR:HA	2.21	0.40
4:L:52:LEU:HD21	4:L:55:SER:HB3	2.04	0.40
1:C:125:ILE:HD12	1:C:270:PRO:HG2	2.03	0.40
4:K:52:LEU:HD21	4:K:55:SER:HB3	2.04	0.40
1:E:195:SER:OG	1:E:227:GLN:O	2.35	0.40
4:L:5:THR:OG1	4:L:24:ARG:HG2	2.22	0.40
1:A:123:ASN:OD1	1:A:272:LYS:NZ	2.47	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	337/362 (93%)	328 (97%)	9 (3%)	0	100	100
1	C	337/362 (93%)	328 (97%)	9 (3%)	0	100	100
1	E	337/362 (93%)	328 (97%)	9 (3%)	0	100	100
2	B	99/235 (42%)	96 (97%)	3 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	99/235 (42%)	96 (97%)	3 (3%)	0	100	100
2	F	99/235 (42%)	96 (97%)	3 (3%)	0	100	100
3	G	123/474 (26%)	113 (92%)	10 (8%)	0	100	100
3	H	123/474 (26%)	113 (92%)	10 (8%)	0	100	100
3	J	123/474 (26%)	113 (92%)	10 (8%)	0	100	100
4	I	105/233 (45%)	98 (93%)	7 (7%)	0	100	100
4	K	105/233 (45%)	98 (93%)	7 (7%)	0	100	100
4	L	105/233 (45%)	98 (93%)	7 (7%)	0	100	100
All	All	1992/3912 (51%)	1905 (96%)	87 (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	287/307 (94%)	285 (99%)	2 (1%)	76	86
1	C	287/307 (94%)	285 (99%)	2 (1%)	76	86
1	E	287/307 (94%)	285 (99%)	2 (1%)	76	86
2	B	91/191 (48%)	90 (99%)	1 (1%)	65	79
2	D	91/191 (48%)	90 (99%)	1 (1%)	65	79
2	F	91/191 (48%)	90 (99%)	1 (1%)	65	79
3	G	106/420 (25%)	102 (96%)	4 (4%)	29	52
3	H	106/420 (25%)	102 (96%)	4 (4%)	29	52
3	J	106/420 (25%)	102 (96%)	4 (4%)	29	52
4	I	88/199 (44%)	86 (98%)	2 (2%)	44	66
4	K	88/199 (44%)	86 (98%)	2 (2%)	44	66
4	L	88/199 (44%)	86 (98%)	2 (2%)	44	66
All	All	1716/3351 (51%)	1689 (98%)	27 (2%)	54	73

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	286	SER
2	B	463	LEU
1	C	9	SER
1	C	286	SER
2	D	463	LEU
1	E	9	SER
1	E	286	SER
2	F	463	LEU
3	H	5	VAL
3	H	7	SER
3	H	22	SER
3	H	83	SER
4	L	37	SER
4	L	96	SER
3	G	5	VAL
3	G	7	SER
3	G	22	SER
3	G	83	SER
4	I	37	SER
4	I	96	SER
3	J	5	VAL
3	J	7	SER
3	J	22	SER
3	J	83	SER
4	K	37	SER
4	K	96	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	HIS
1	A	122	HIS
1	A	171	ASN
1	A	200	GLN
1	A	218	ASN
1	A	234	GLN
1	A	253	GLN
1	A	264	GLN
2	B	421	HIS
2	B	442	GLN

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Mol	Chain	Res	Type
2	B	456	ASN
1	C	99	HIS
1	C	122	HIS
1	C	171	ASN
1	C	200	GLN
1	C	218	ASN
1	C	234	GLN
1	C	253	GLN
1	C	264	GLN
2	D	369	HIS
2	D	393	ASN
2	D	421	HIS
2	D	442	GLN
2	D	456	ASN
1	E	99	HIS
1	E	122	HIS
1	E	171	ASN
1	E	200	GLN
1	E	218	ASN
1	E	234	GLN
1	E	253	GLN
1	E	264	GLN
2	F	442	GLN
2	F	456	ASN
3	H	90	GLN
3	G	90	GLN
3	J	90	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

24 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	1,5	14,14,15	0.33	0	17,19,21	0.68	0
5	NAG	M	2	5	14,14,15	0.29	0	17,19,21	0.42	0
5	NAG	N	1	1,5	14,14,15	0.32	0	17,19,21	0.47	0
5	NAG	N	2	5	14,14,15	0.34	0	17,19,21	0.47	0
5	NAG	O	1	1,5	14,14,15	0.29	0	17,19,21	0.58	0
5	NAG	O	2	5	14,14,15	0.40	0	17,19,21	0.51	0
5	NAG	P	1	1,5	14,14,15	0.88	1 (7%)	17,19,21	1.03	1 (5%)
5	NAG	P	2	5	14,14,15	0.46	0	17,19,21	0.44	0
5	NAG	Q	1	1,5	14,14,15	0.32	0	17,19,21	0.68	0
5	NAG	Q	2	5	14,14,15	0.30	0	17,19,21	0.41	0
5	NAG	R	1	1,5	14,14,15	0.31	0	17,19,21	0.47	0
5	NAG	R	2	5	14,14,15	0.33	0	17,19,21	0.47	0
5	NAG	S	1	1,5	14,14,15	0.29	0	17,19,21	0.58	0
5	NAG	S	2	5	14,14,15	0.39	0	17,19,21	0.52	0
5	NAG	T	1	1,5	14,14,15	0.87	1 (7%)	17,19,21	1.02	1 (5%)
5	NAG	T	2	5	14,14,15	0.46	0	17,19,21	0.43	0
5	NAG	U	1	1,5	14,14,15	0.34	0	17,19,21	0.68	0
5	NAG	U	2	5	14,14,15	0.29	0	17,19,21	0.41	0
5	NAG	V	1	1,5	14,14,15	0.32	0	17,19,21	0.47	0
5	NAG	V	2	5	14,14,15	0.33	0	17,19,21	0.48	0
5	NAG	W	1	1,5	14,14,15	0.29	0	17,19,21	0.58	0
5	NAG	W	2	5	14,14,15	0.41	0	17,19,21	0.51	0
5	NAG	X	1	1,5	14,14,15	0.88	1 (7%)	17,19,21	1.02	1 (5%)
5	NAG	X	2	5	14,14,15	0.46	0	17,19,21	0.44	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
5	NAG	M	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	M	2	5	-	0/6/23/26	0/1/1/1
5	NAG	N	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	N	2	5	-	0/6/23/26	0/1/1/1
5	NAG	O	1	1,5	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	O	2	5	-	2/6/23/26	0/1/1/1
5	NAG	P	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	P	2	5	-	0/6/23/26	0/1/1/1
5	NAG	Q	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	0/6/23/26	0/1/1/1
5	NAG	R	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	R	2	5	-	0/6/23/26	0/1/1/1
5	NAG	S	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	S	2	5	-	2/6/23/26	0/1/1/1
5	NAG	T	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	T	2	5	-	0/6/23/26	0/1/1/1
5	NAG	U	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	U	2	5	-	0/6/23/26	0/1/1/1
5	NAG	V	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	V	2	5	-	0/6/23/26	0/1/1/1
5	NAG	W	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	W	2	5	-	2/6/23/26	0/1/1/1
5	NAG	X	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	X	2	5	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	P	1	NAG	C1-C2	2.59	1.56	1.52
5	X	1	NAG	C1-C2	2.56	1.56	1.52
5	T	1	NAG	C1-C2	2.56	1.56	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	P	1	NAG	C1-O5-C5	3.44	116.86	112.19
5	X	1	NAG	C1-O5-C5	3.43	116.84	112.19
5	T	1	NAG	C1-O5-C5	3.42	116.83	112.19

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	M	1	NAG	O5-C5-C6-O6
5	Q	1	NAG	O5-C5-C6-O6

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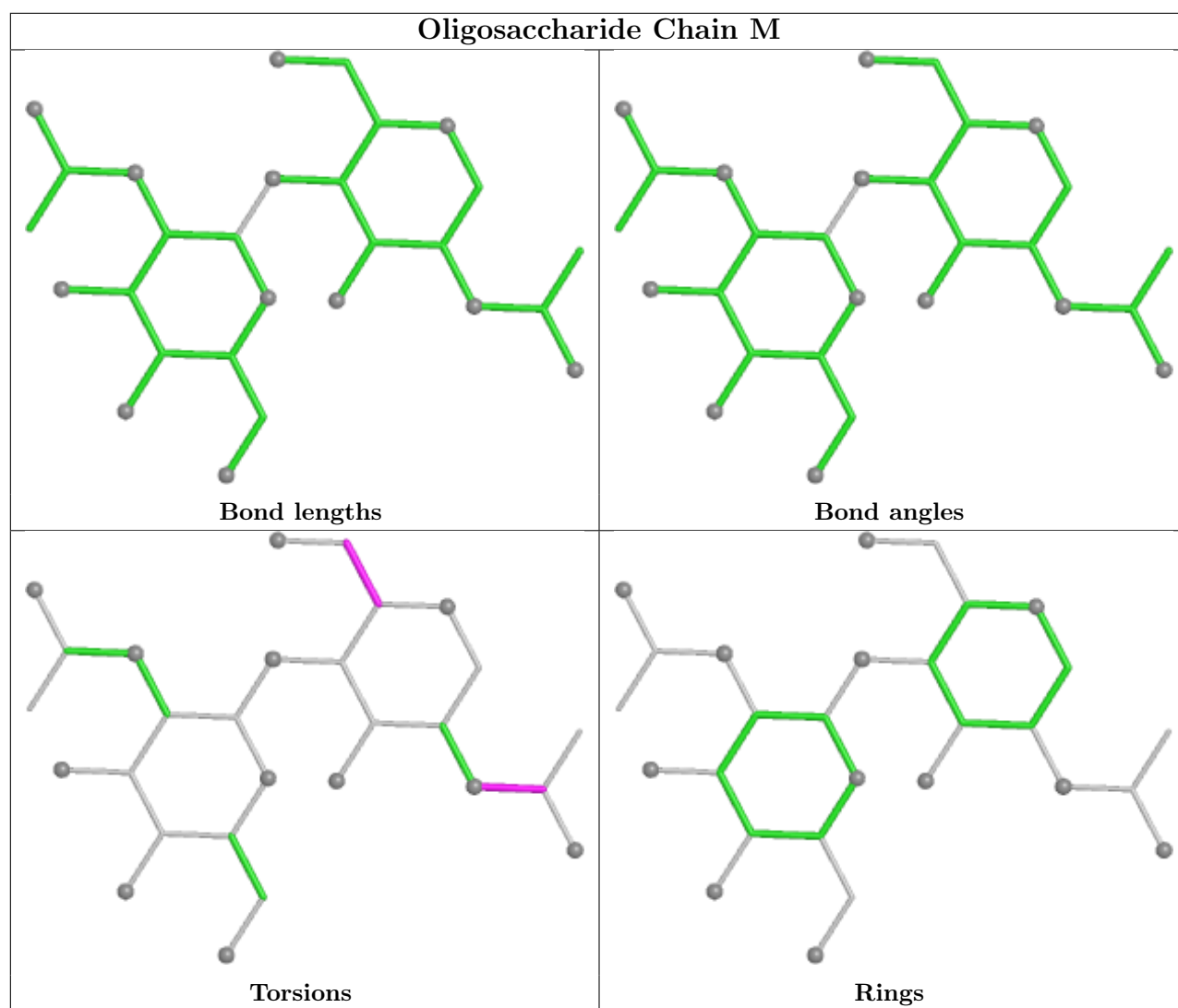
Mol	Chain	Res	Type	Atoms
5	U	1	NAG	O5-C5-C6-O6
5	M	1	NAG	C8-C7-N2-C2
5	M	1	NAG	O7-C7-N2-C2
5	Q	1	NAG	C8-C7-N2-C2
5	Q	1	NAG	O7-C7-N2-C2
5	U	1	NAG	C8-C7-N2-C2
5	U	1	NAG	O7-C7-N2-C2
5	M	1	NAG	C4-C5-C6-O6
5	Q	1	NAG	C4-C5-C6-O6
5	U	1	NAG	C4-C5-C6-O6
5	S	2	NAG	C4-C5-C6-O6
5	O	2	NAG	C4-C5-C6-O6
5	W	2	NAG	C4-C5-C6-O6
5	S	2	NAG	O5-C5-C6-O6
5	W	2	NAG	O5-C5-C6-O6
5	O	2	NAG	O5-C5-C6-O6

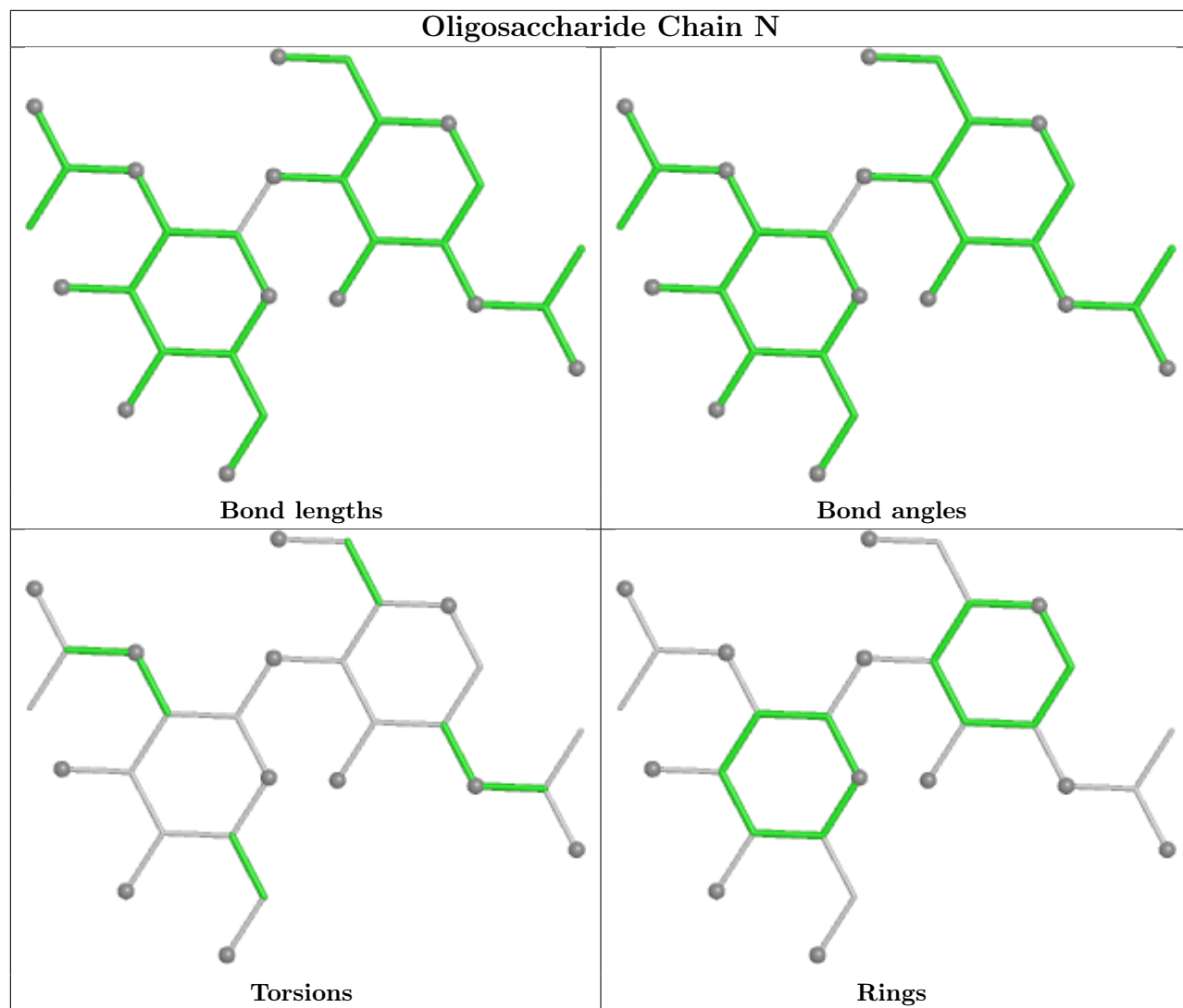
There are no ring outliers.

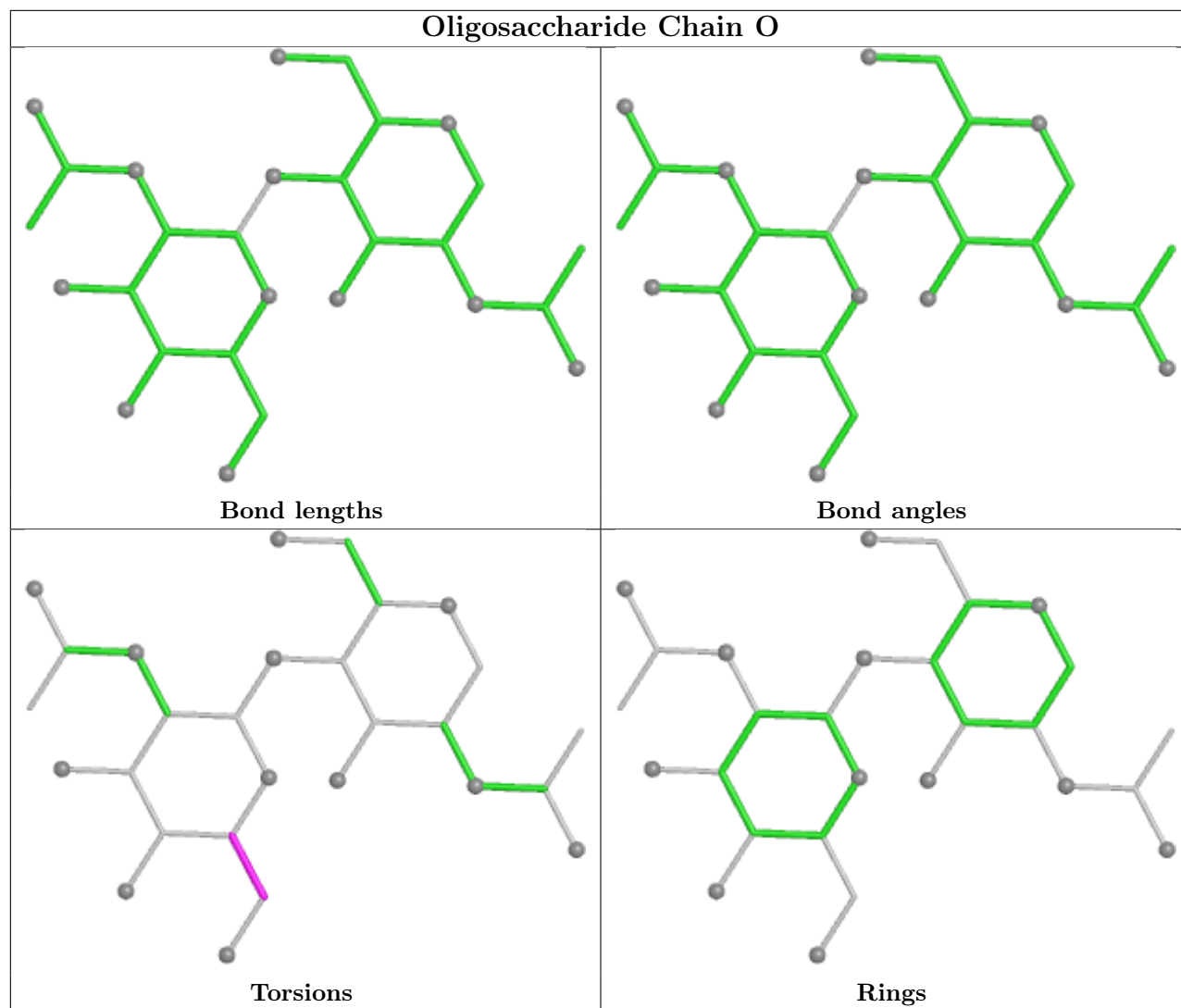
3 monomers are involved in 3 short contacts:

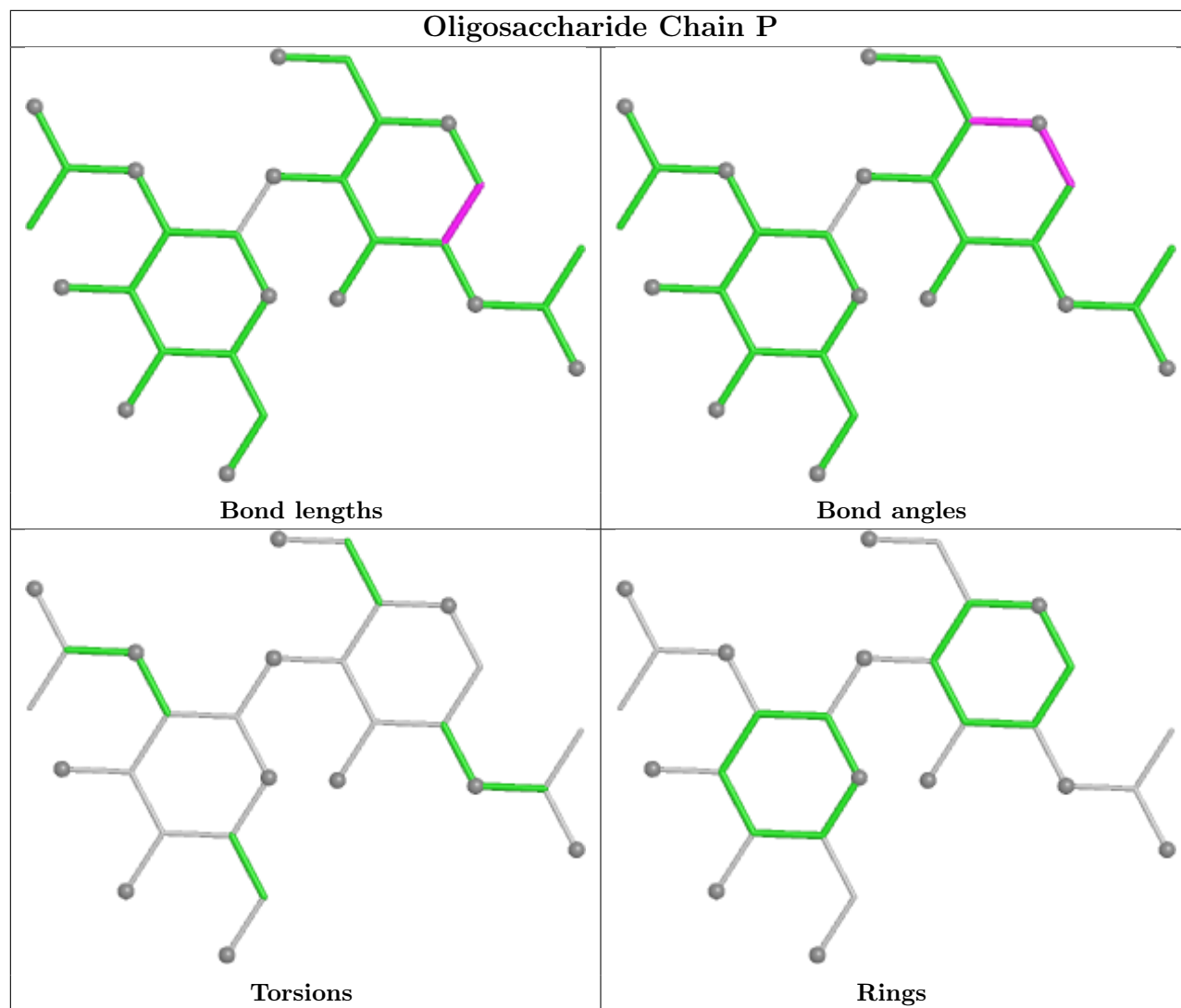
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	T	1	NAG	1	0
5	X	1	NAG	1	0
5	P	1	NAG	1	0

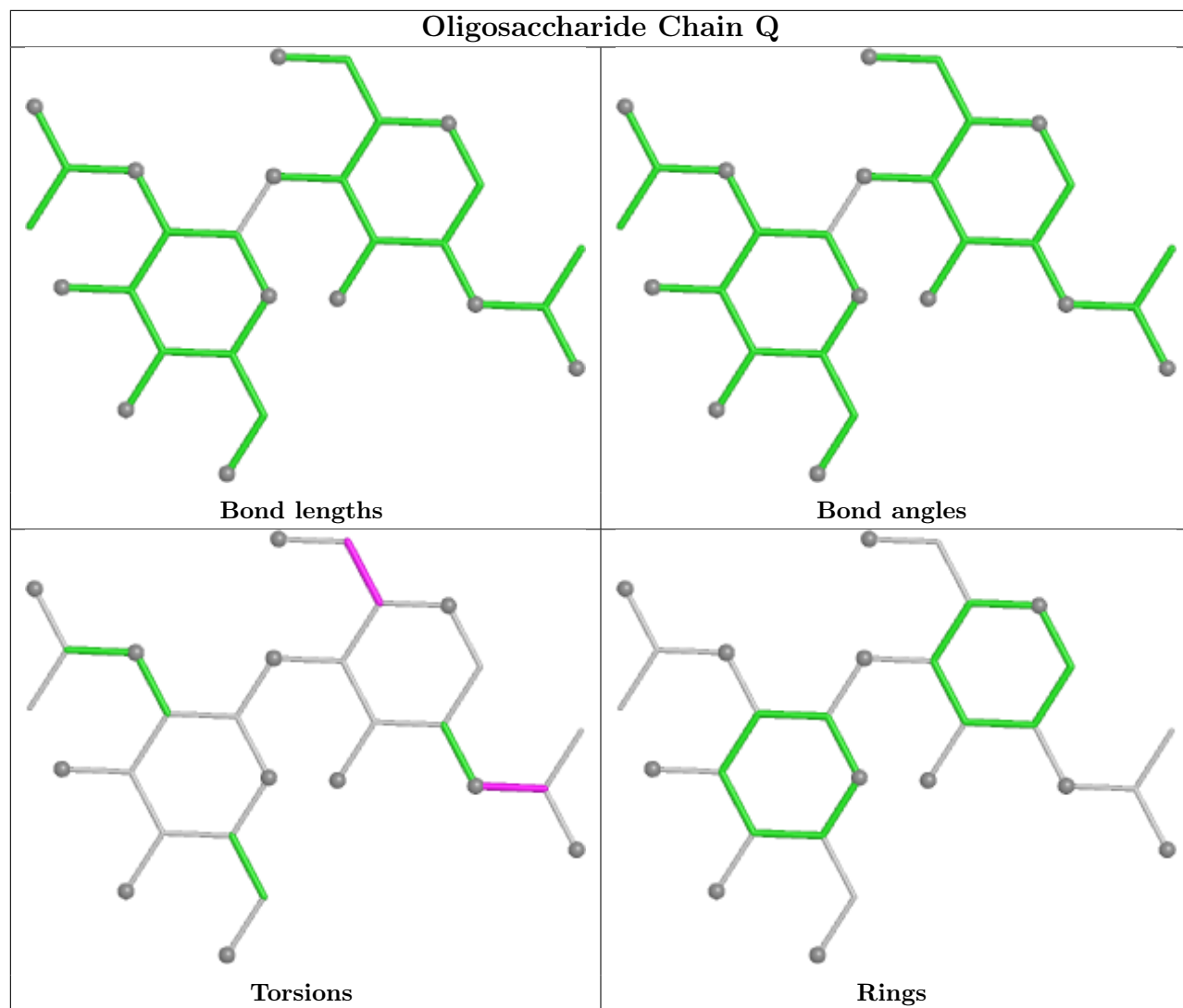
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

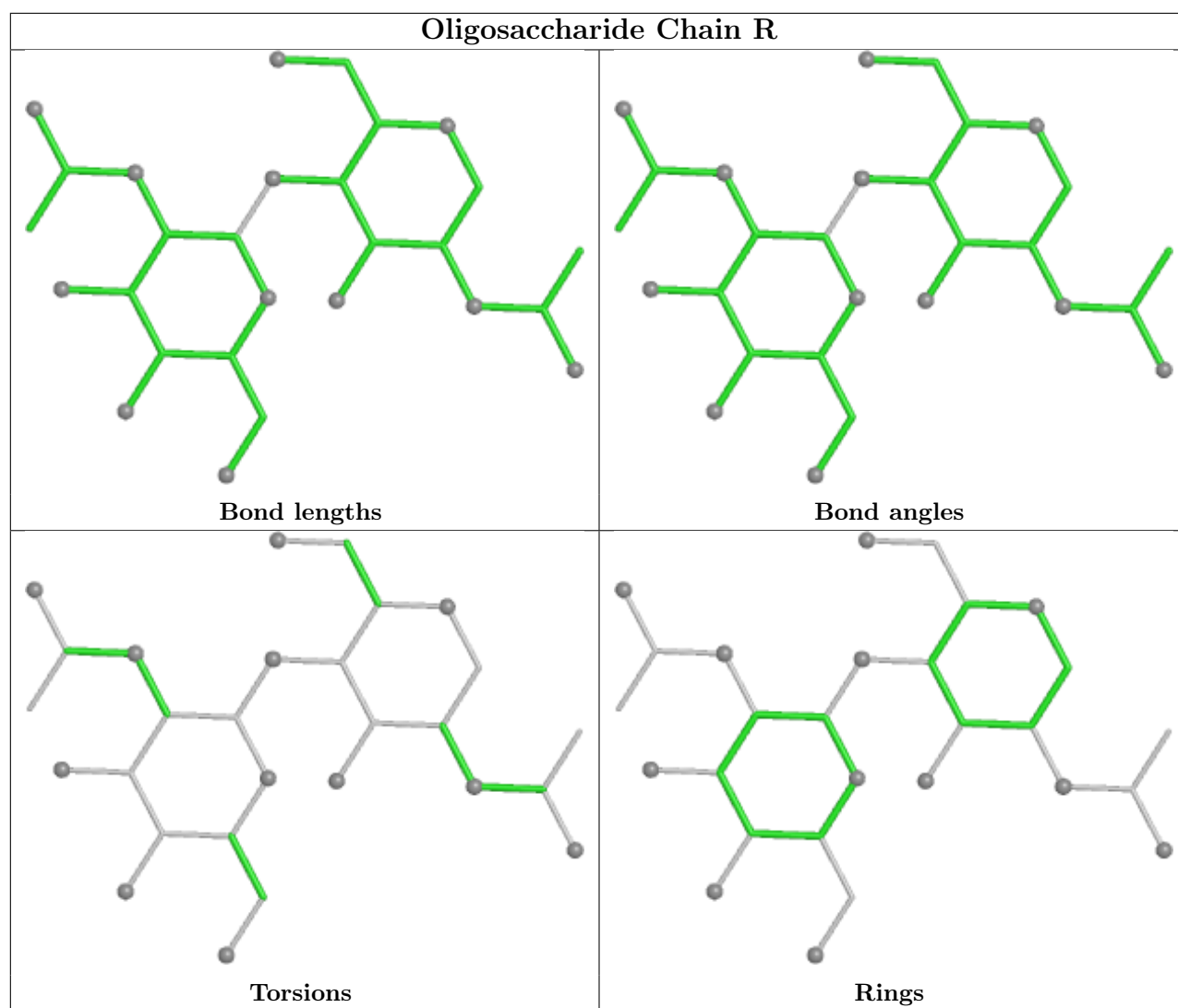


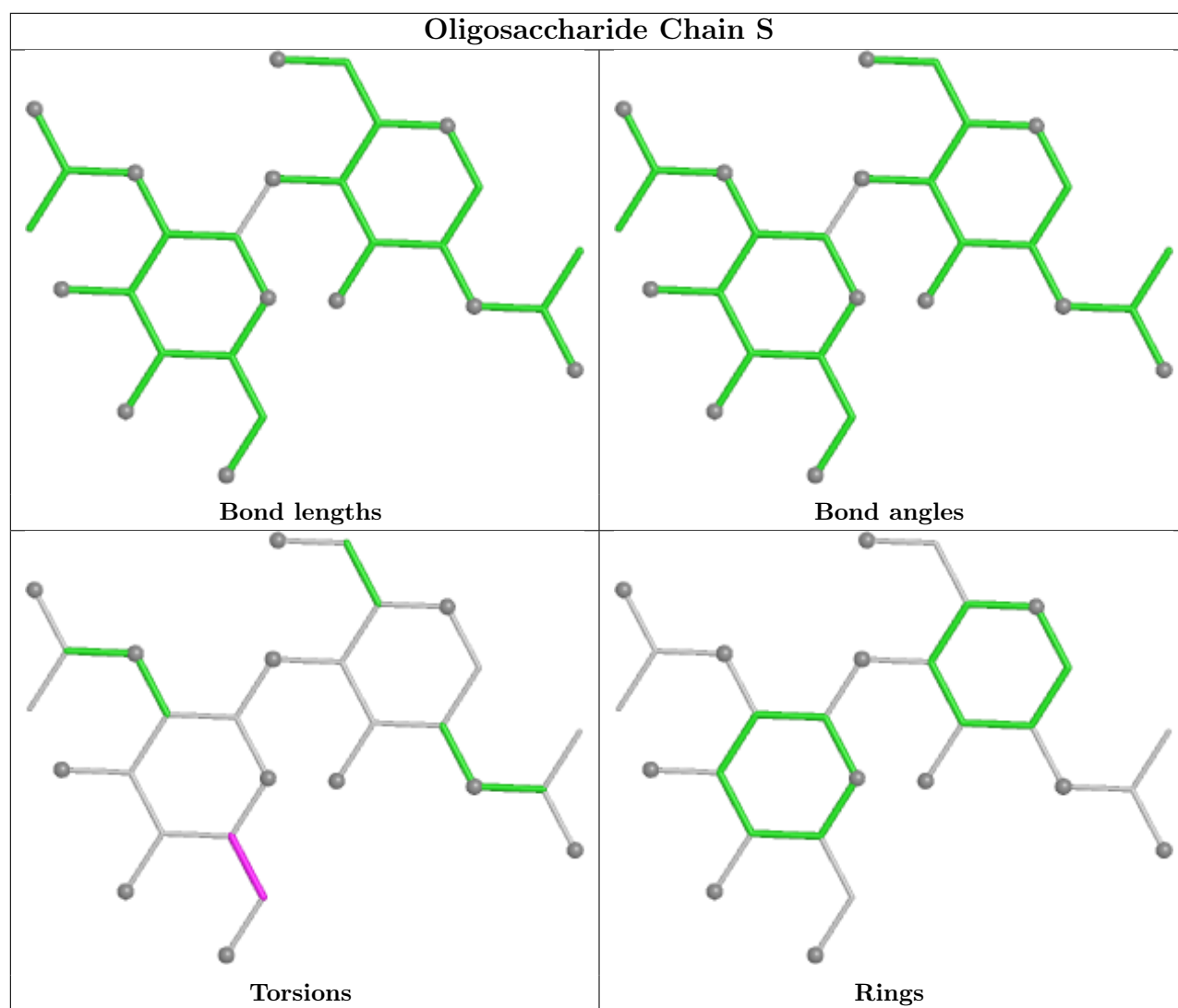


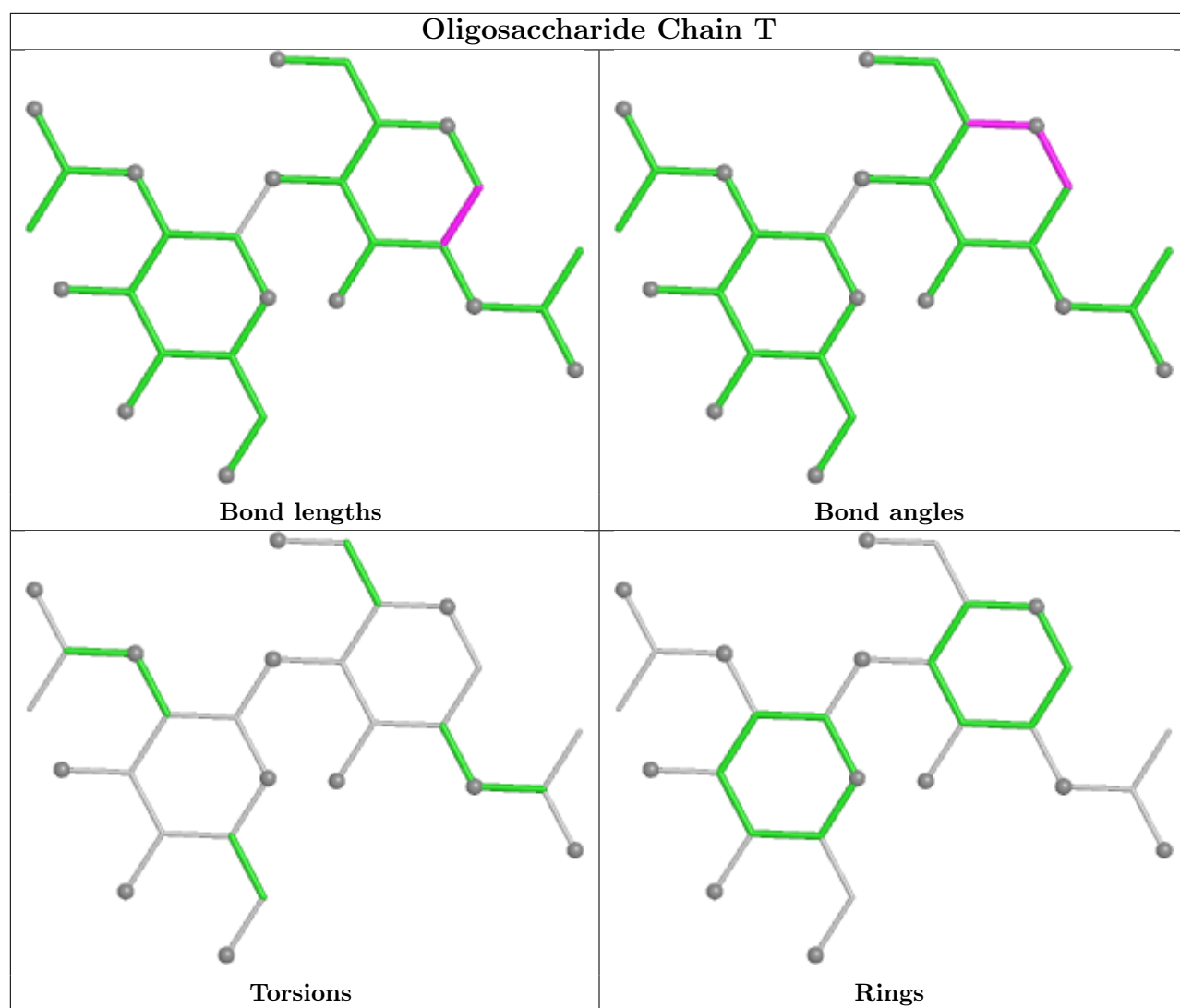


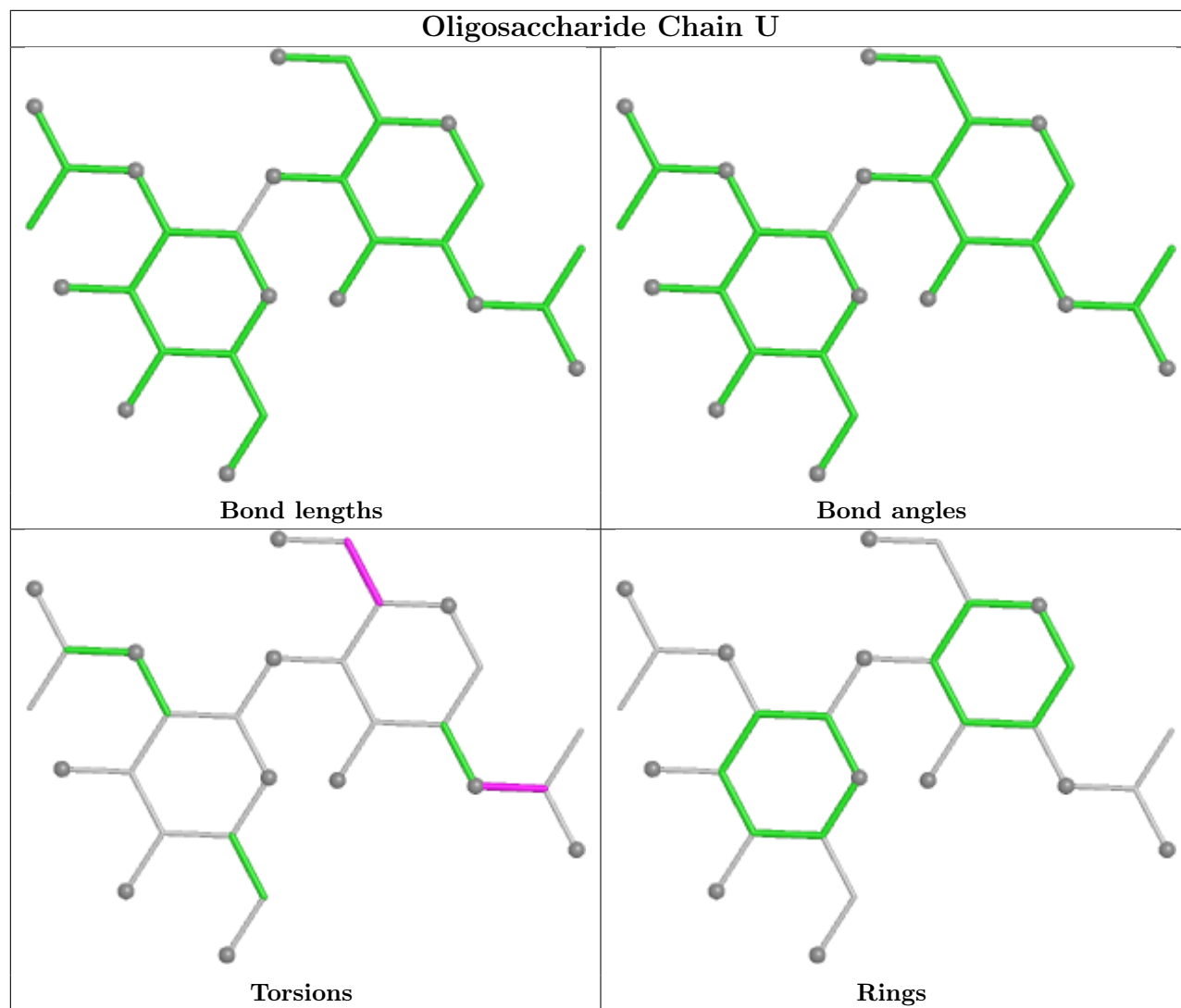


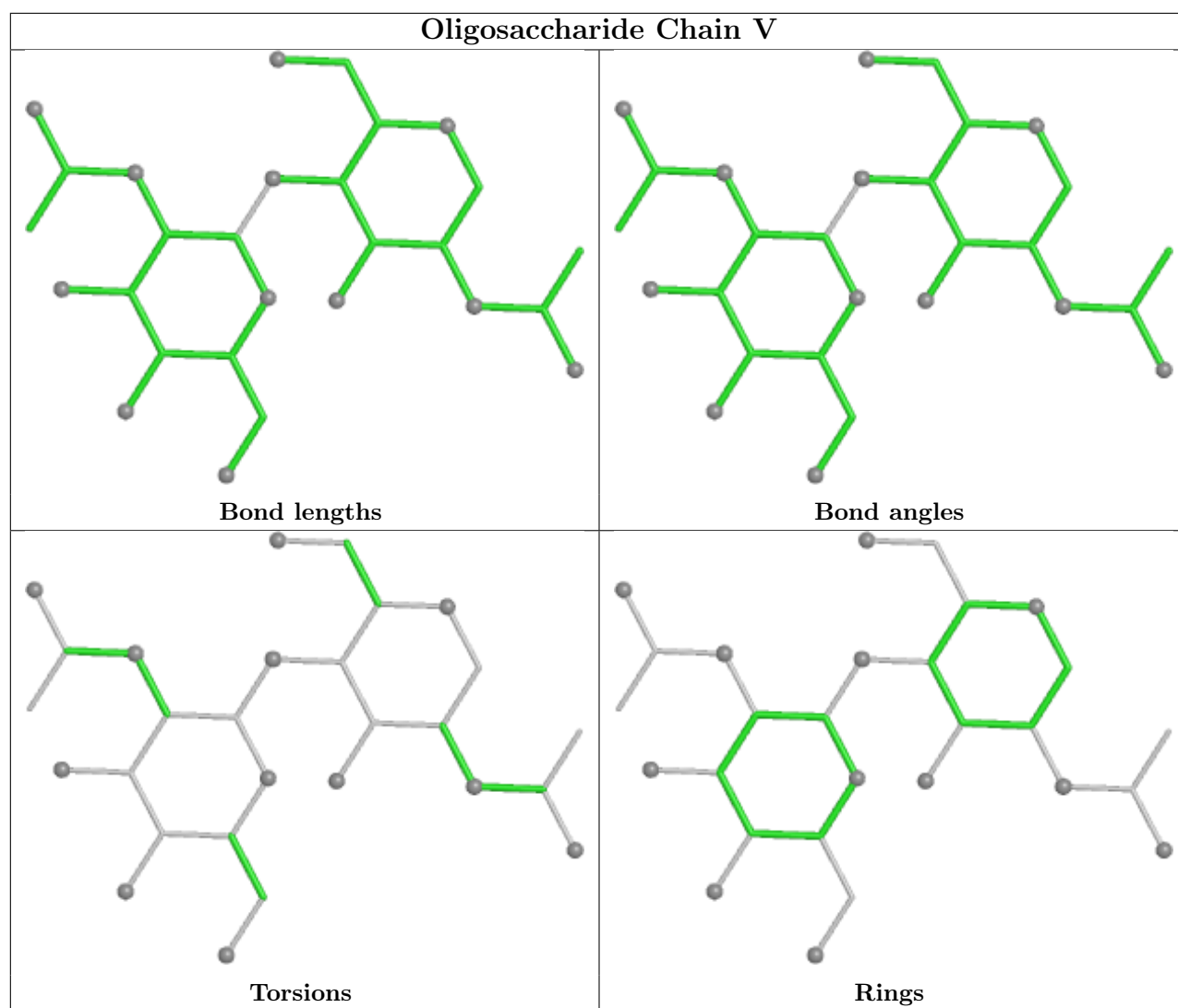


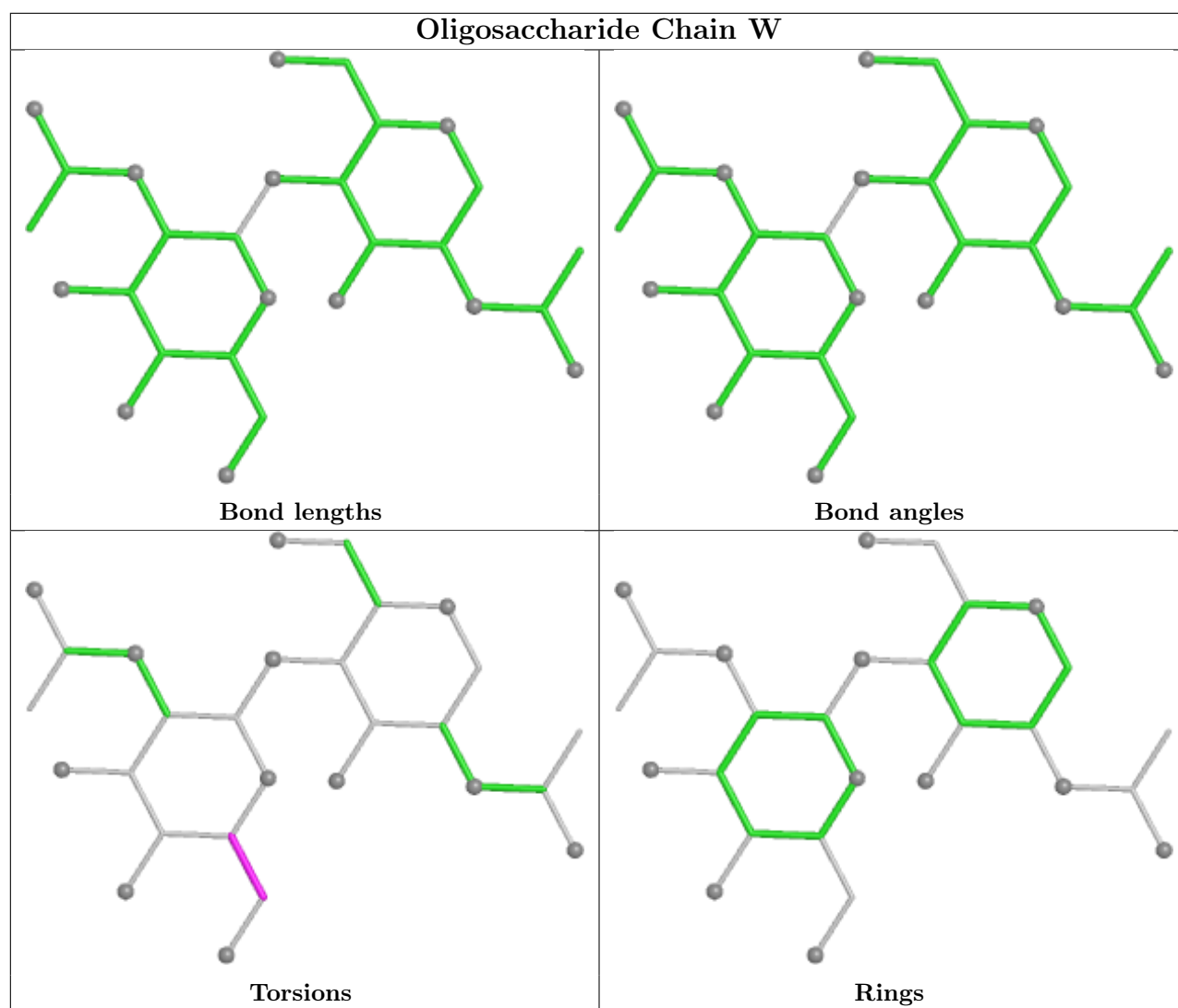


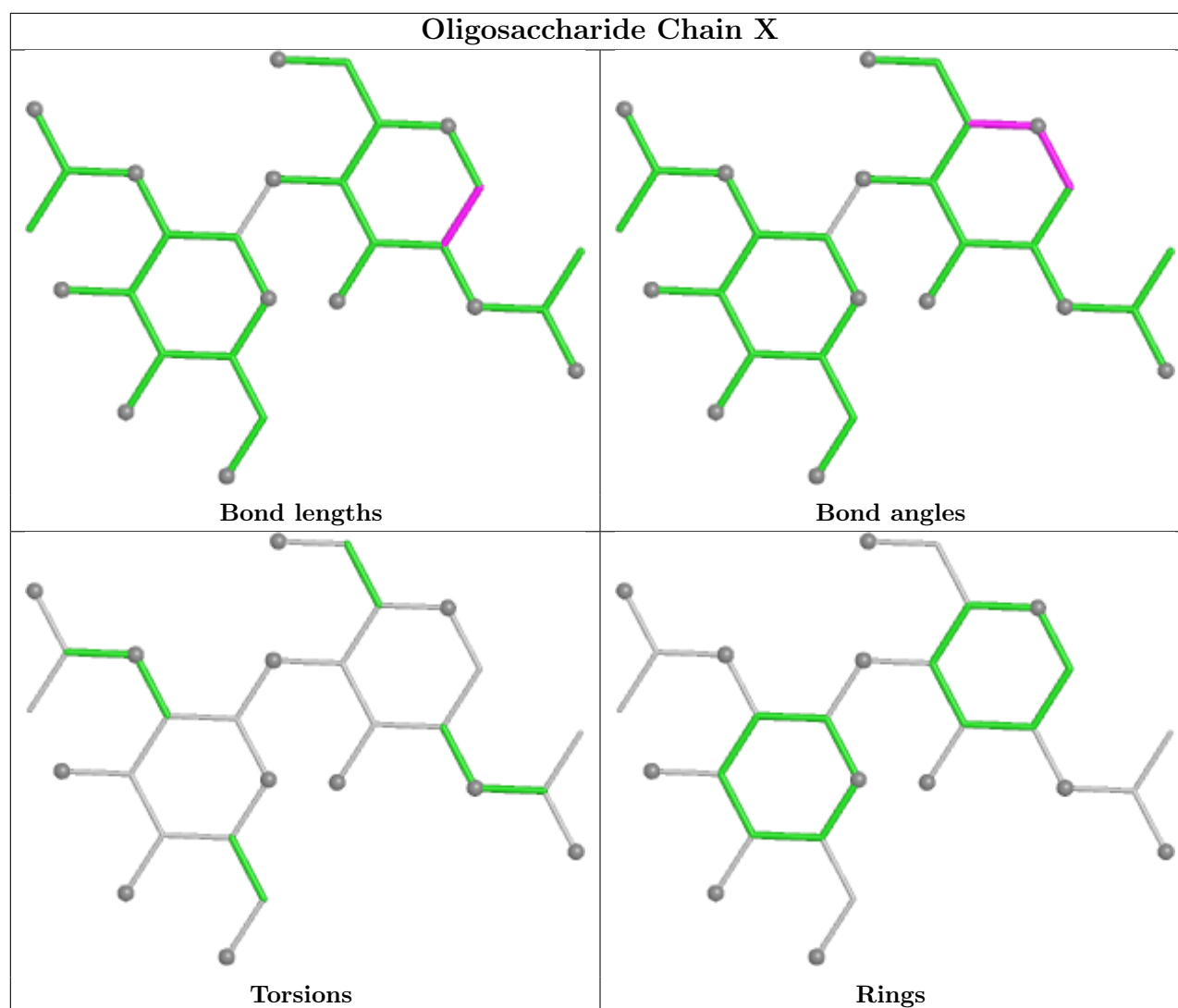












5.6 Ligand geometry [i](#)

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	$\# Z > 2$
6	NAG	C	401	1	14,14,15	0.37	0	17,19,21	0.56	0
6	NAG	E	403	1	14,14,15	0.24	0	17,19,21	0.60	0
6	NAG	C	402	1	14,14,15	0.21	0	17,19,21	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	403	1	14,14,15	0.23	0	17,19,21	0.60	0
6	NAG	C	403	1	14,14,15	0.24	0	17,19,21	0.60	0
6	NAG	E	402	1	14,14,15	0.19	0	17,19,21	0.37	0
6	NAG	E	401	1	14,14,15	0.37	0	17,19,21	0.55	0
6	NAG	A	401	1	14,14,15	0.37	0	17,19,21	0.56	0
6	NAG	A	402	1	14,14,15	0.19	0	17,19,21	0.37	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
6	NAG	C	401	1	-	2/6/23/26	0/1/1/1
6	NAG	E	403	1	-	0/6/23/26	0/1/1/1
6	NAG	C	402	1	-	2/6/23/26	0/1/1/1
6	NAG	A	403	1	-	0/6/23/26	0/1/1/1
6	NAG	C	403	1	-	0/6/23/26	0/1/1/1
6	NAG	E	402	1	-	2/6/23/26	0/1/1/1
6	NAG	E	401	1	-	2/6/23/26	0/1/1/1
6	NAG	A	401	1	-	2/6/23/26	0/1/1/1
6	NAG	A	402	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	402	NAG	O5-C5-C6-O6
6	E	402	NAG	O5-C5-C6-O6
6	C	402	NAG	O5-C5-C6-O6
6	A	402	NAG	C4-C5-C6-O6
6	C	402	NAG	C4-C5-C6-O6
6	E	402	NAG	C4-C5-C6-O6
6	A	401	NAG	O5-C5-C6-O6
6	C	401	NAG	O5-C5-C6-O6
6	E	401	NAG	O5-C5-C6-O6
6	A	401	NAG	C4-C5-C6-O6

Continued on next page...

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Mol	Chain	Res	Type	Atoms
6	C	401	NAG	C4-C5-C6-O6
6	E	401	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

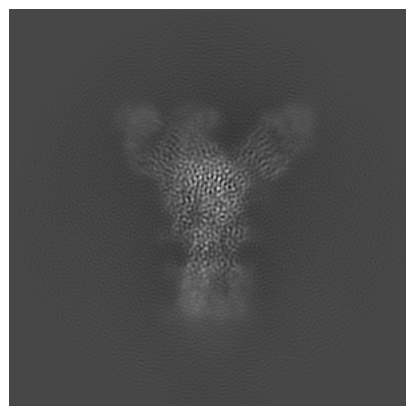
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-66599. These allow visual inspection of the internal detail of the map and identification of artifacts.

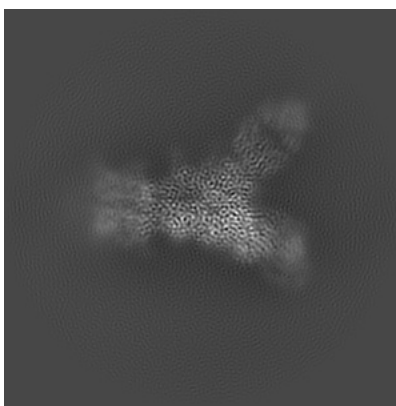
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

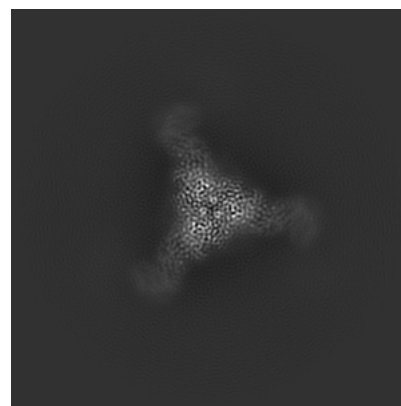
6.1.1 Primary map



X



Y

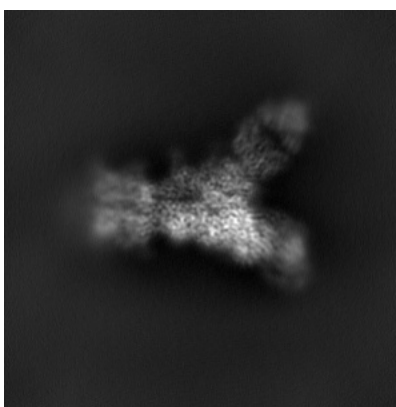


Z

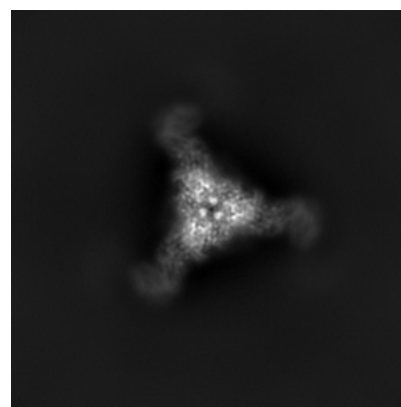
6.1.2 Raw map



X



Y

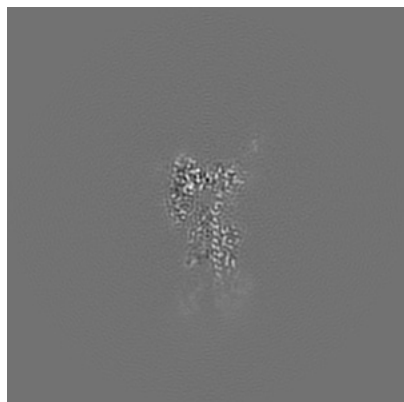


Z

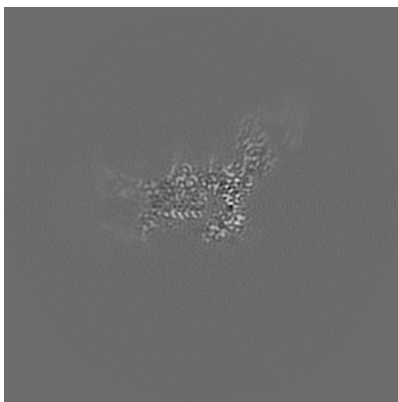
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

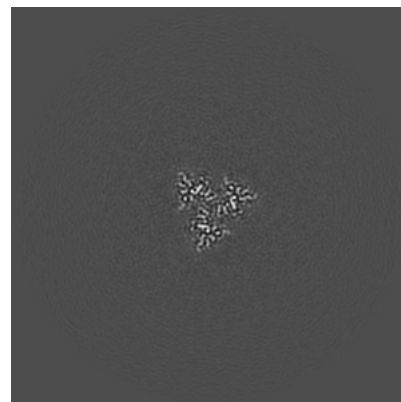
6.2.1 Primary map



X Index: 192

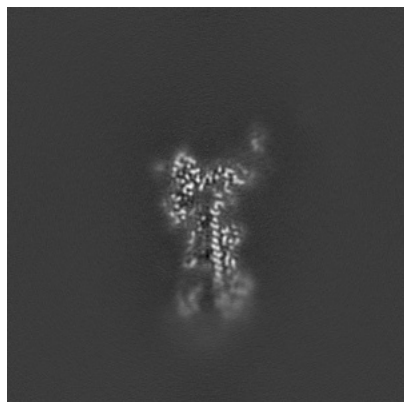


Y Index: 192

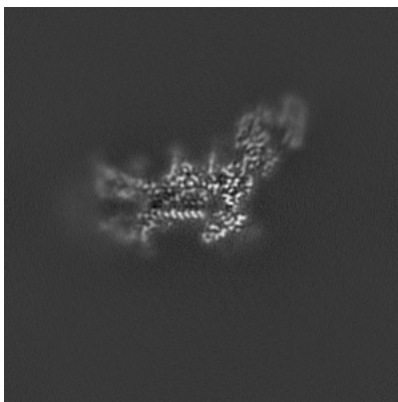


Z Index: 192

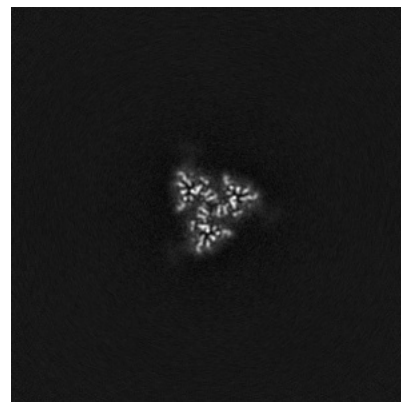
6.2.2 Raw map



X Index: 192



Y Index: 192

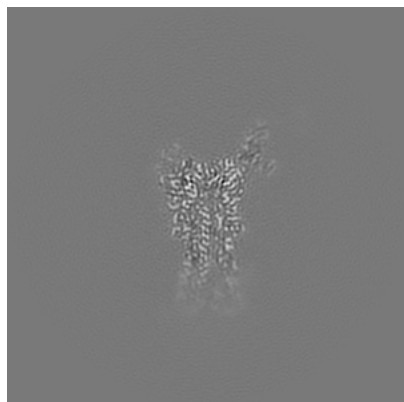


Z Index: 192

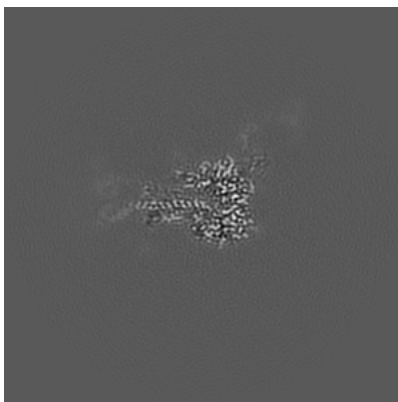
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

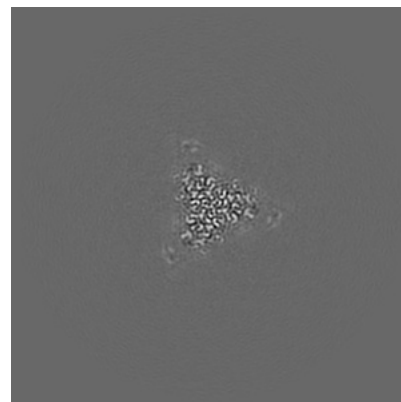
6.3.1 Primary map



X Index: 183

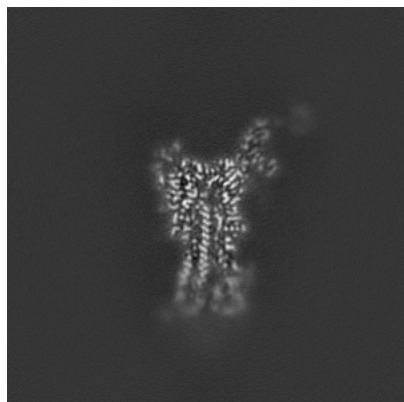


Y Index: 201

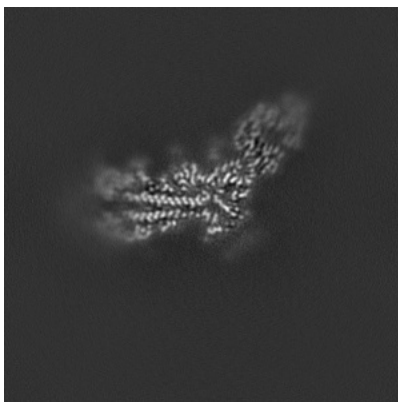


Z Index: 221

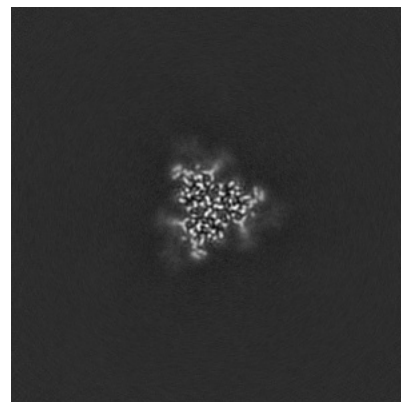
6.3.2 Raw map



X Index: 183



Y Index: 187



Z Index: 215

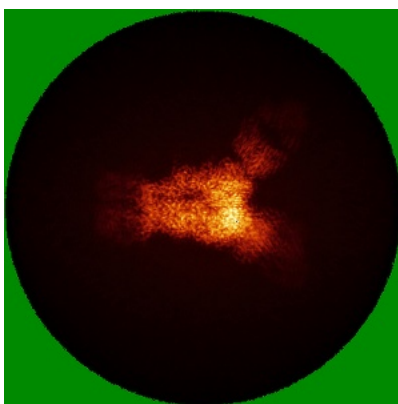
The images above show the largest variance slices of the map in three orthogonal directions.

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

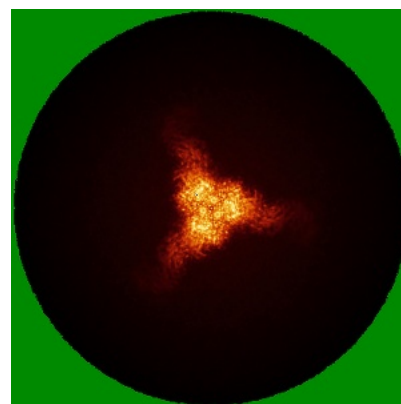
6.4.1 Primary map



X

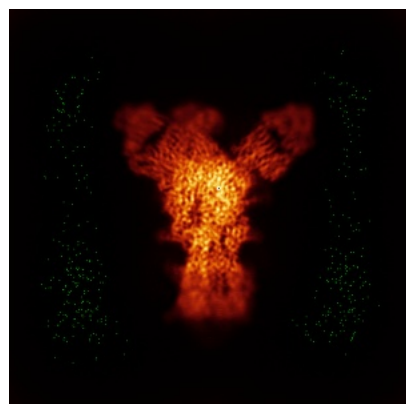


Y

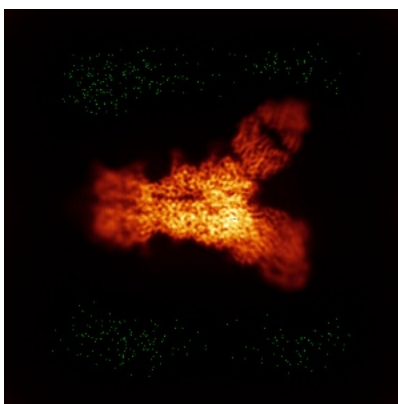


Z

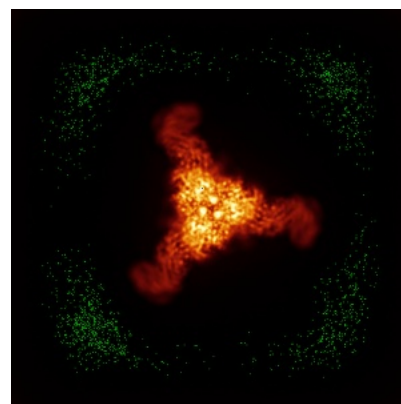
6.4.2 Raw map



X



Y

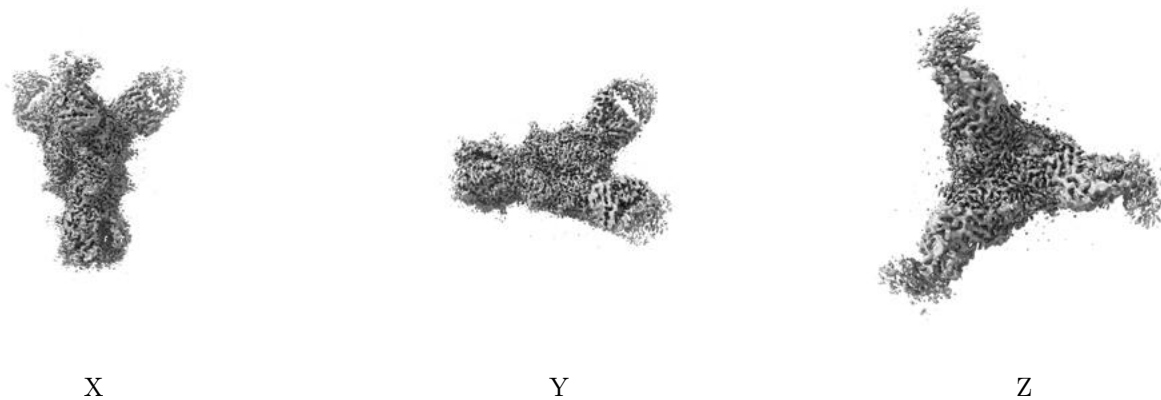


Z

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.

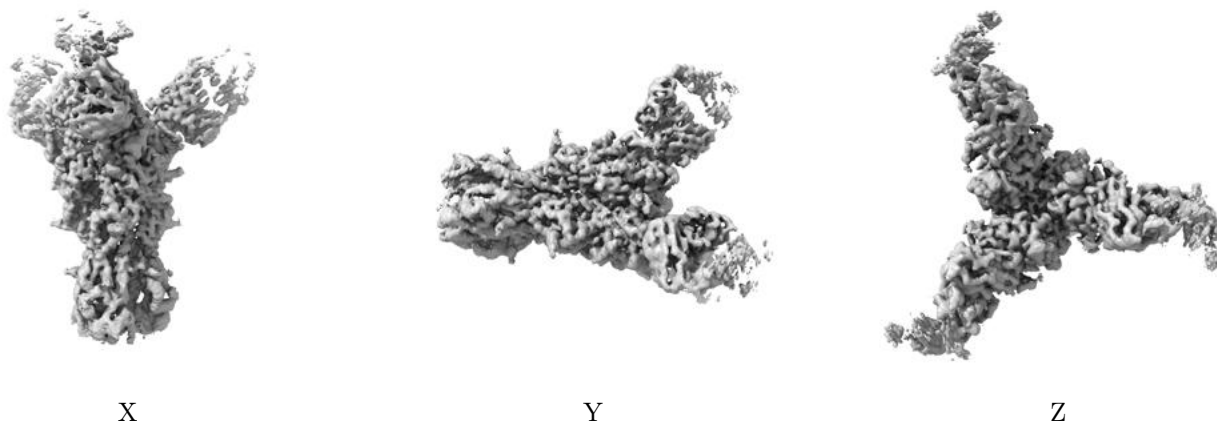
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.227. 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.

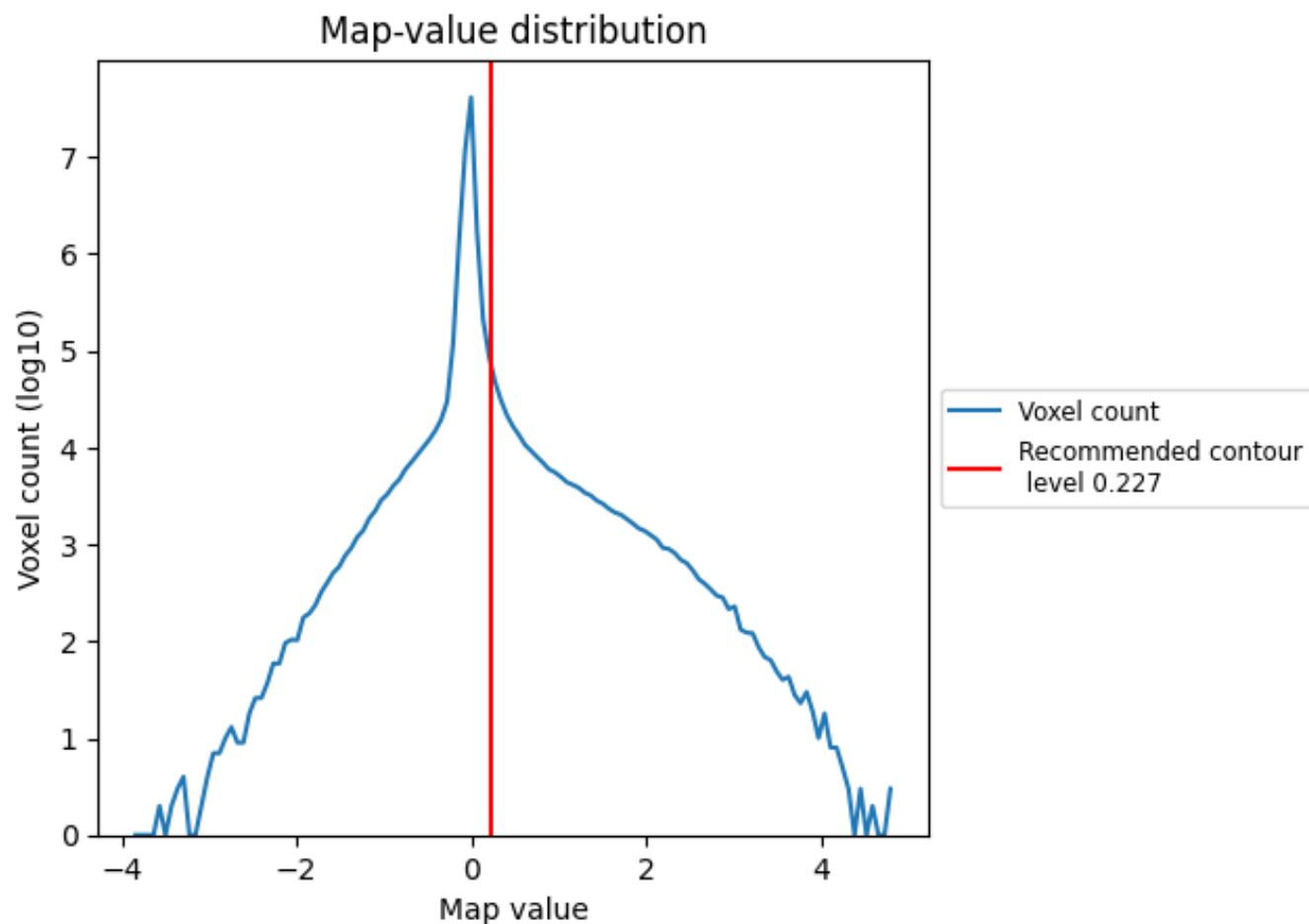
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

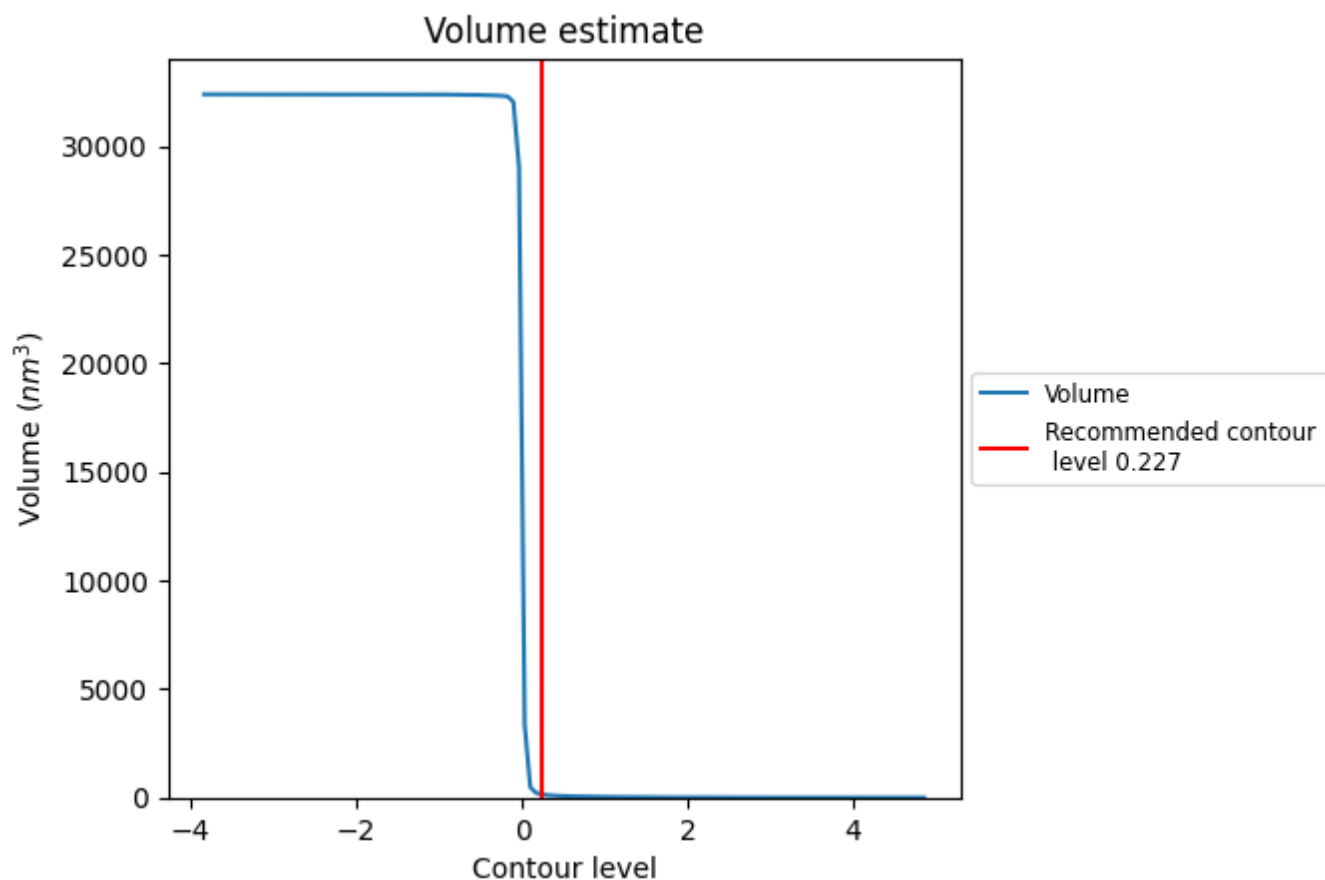
This section contains the results of statistical analysis of the map.

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.

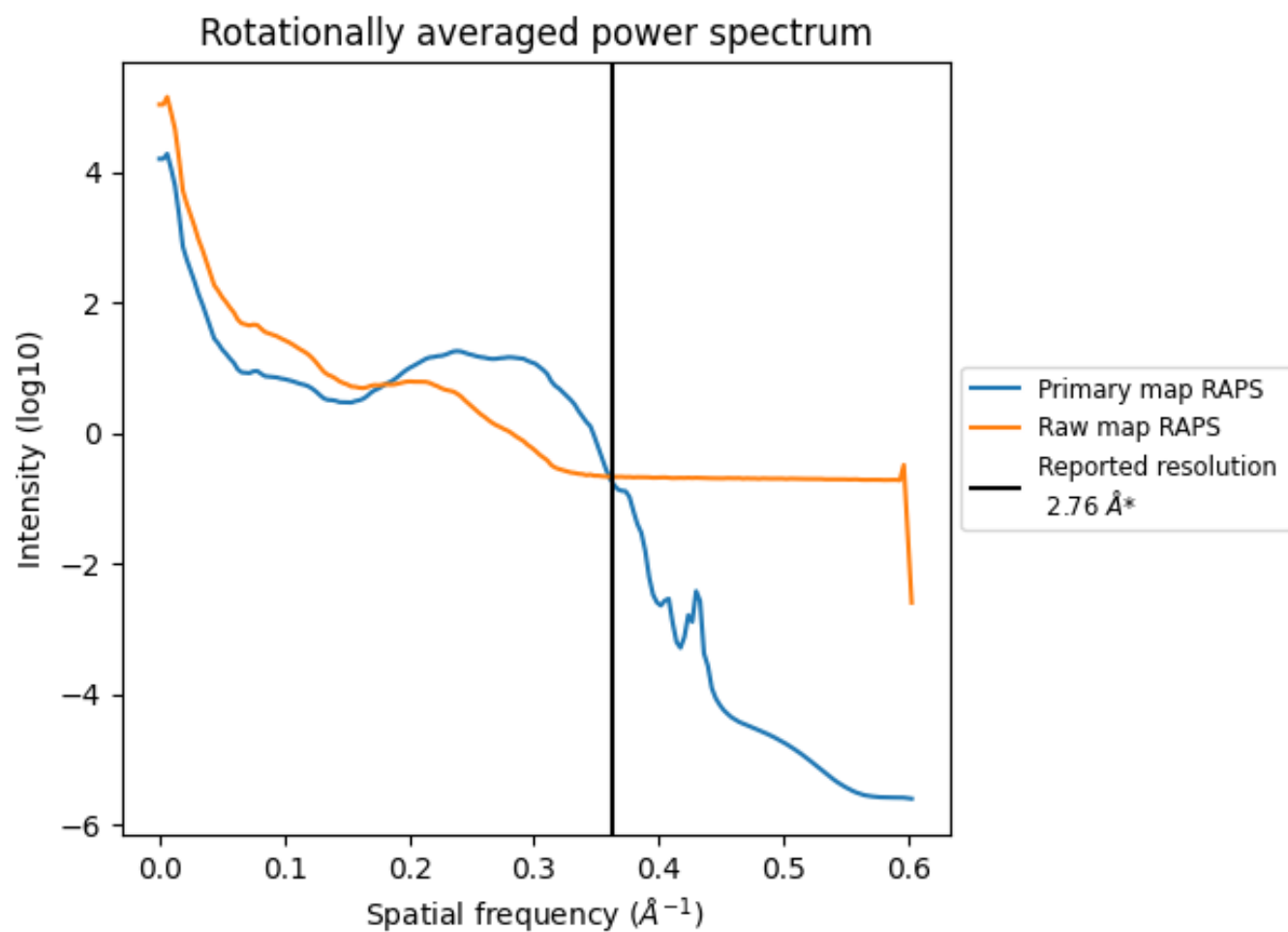
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 158 nm³; this corresponds to an approximate mass of 143 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 ⓘ

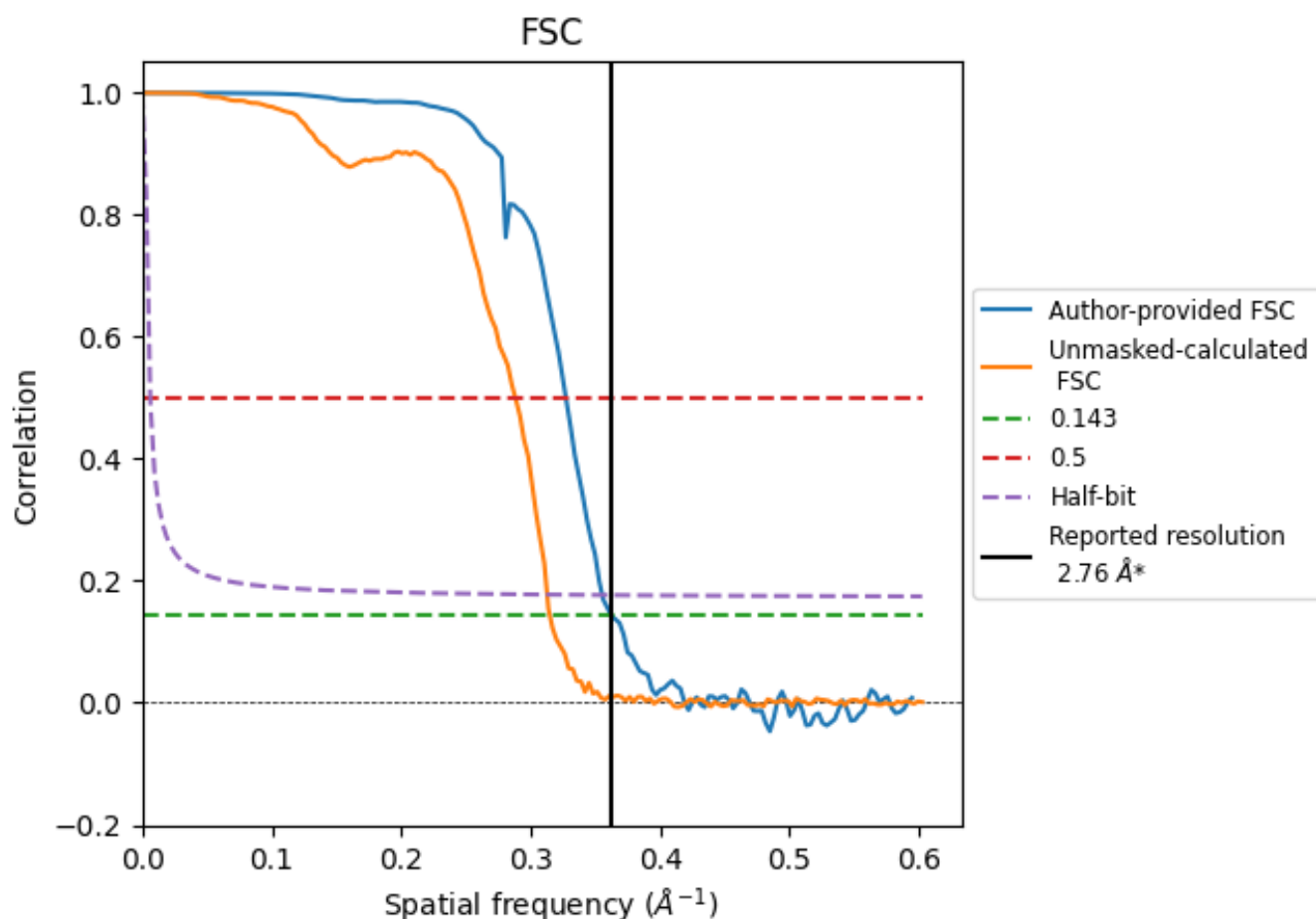


*Reported resolution corresponds to spatial frequency of 0.362 Å⁻¹

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 0.362 \AA^{-1}

8.2 Resolution estimates [i](#)

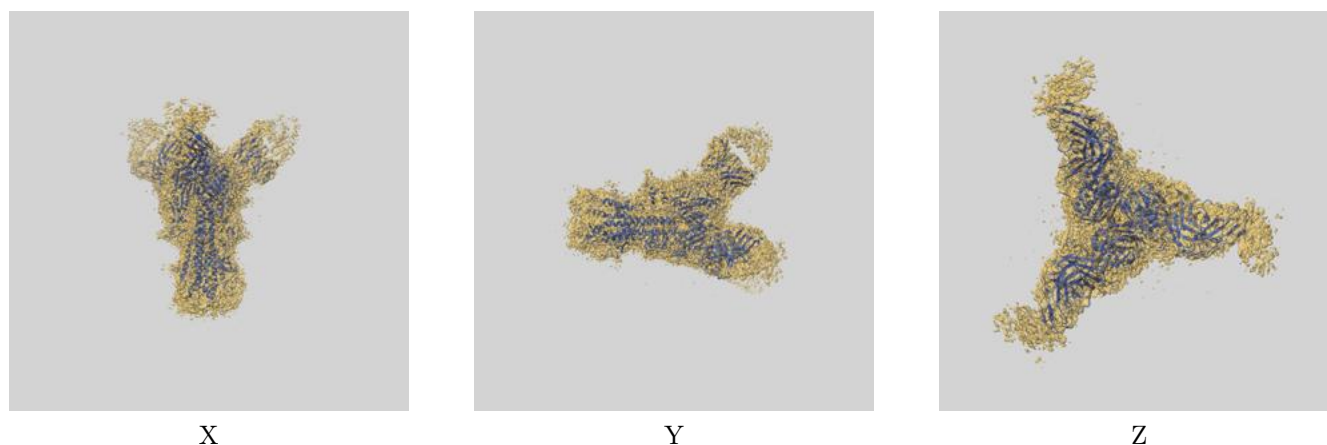
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.76	-	-
Author-provided FSC curve	2.76	3.06	2.81
Unmasked-calculated*	3.18	3.47	3.20

*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 3.18 differs from the reported value 2.76 by more than 10 %

9 Map-model fit [i](#)

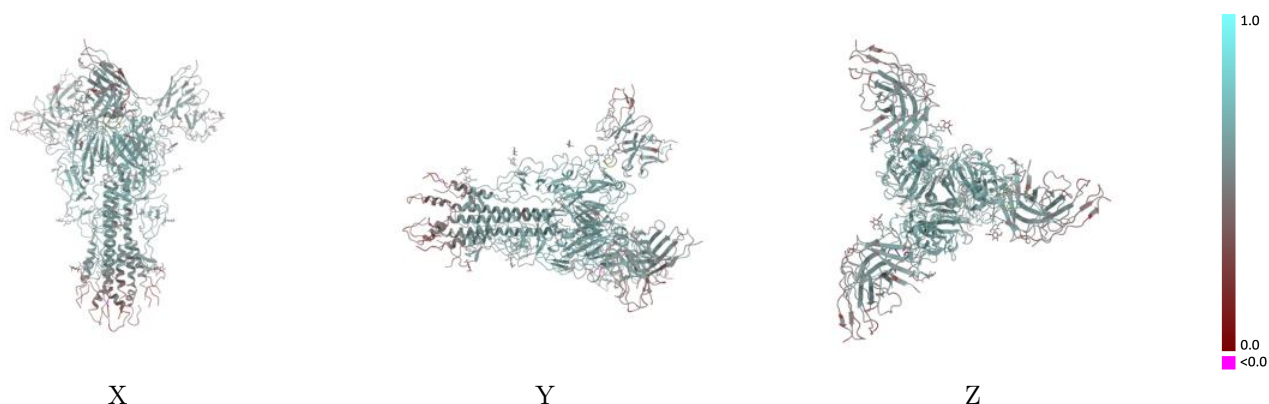
This section contains information regarding the fit between EMDB map EMD-66599 and PDB model 9X5W. Per-residue inclusion information can be found in [section 3](#) on [page 11](#).

9.1 Map-model overlay [i](#)



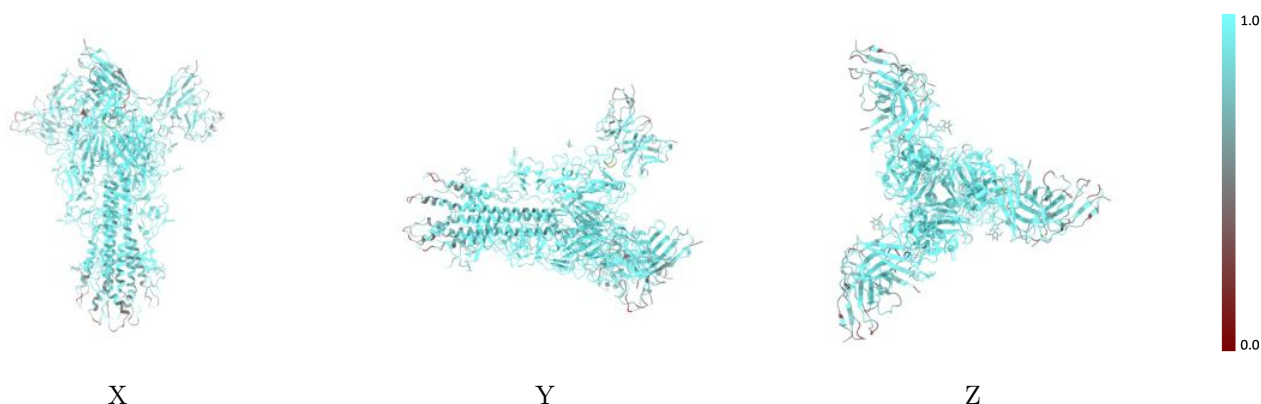
The images above show the 3D surface view of the map at the recommended contour level 0.227 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



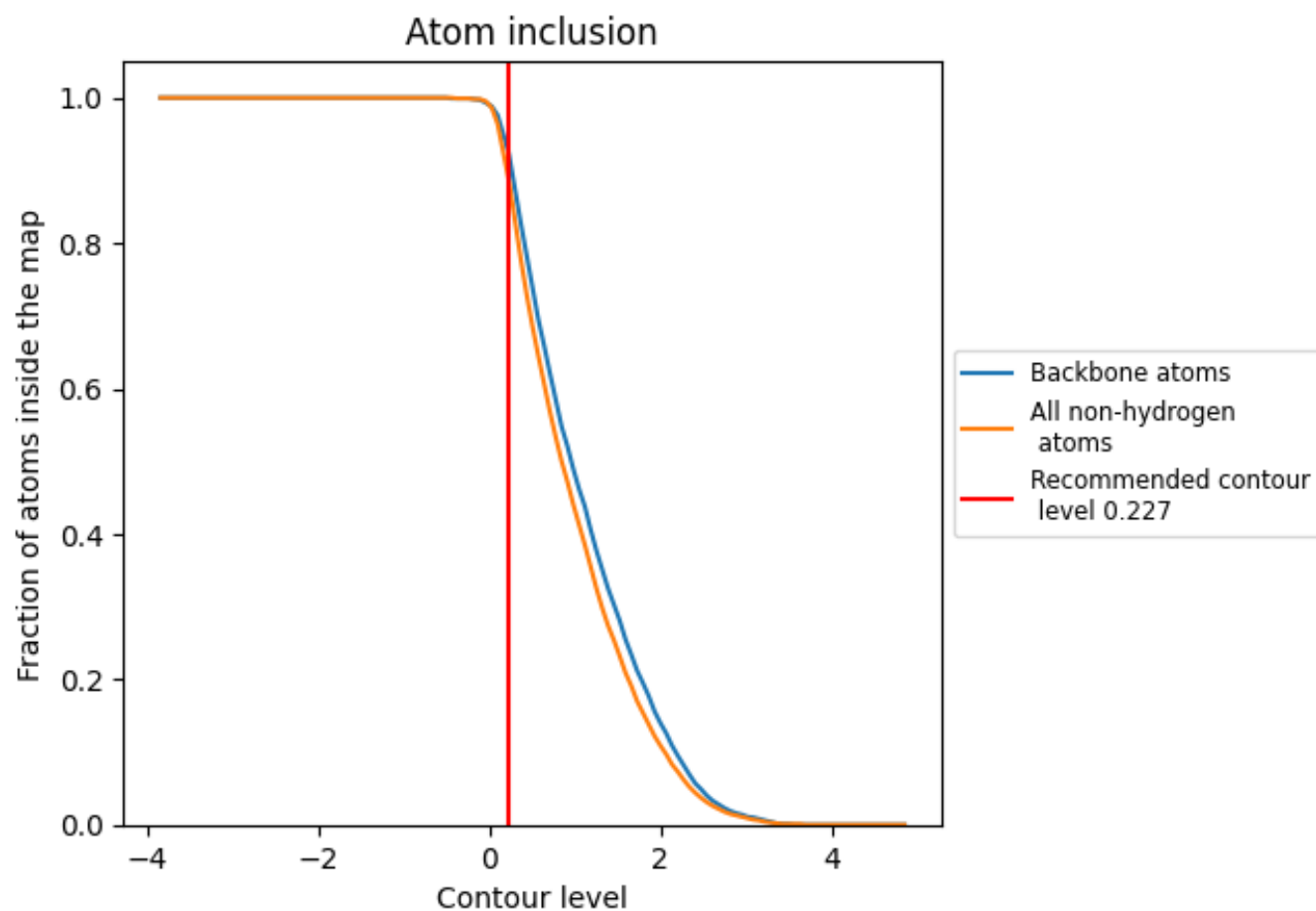
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.227).



















































9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 89% 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 (0.227) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8890	 0.5430
A	 0.9640	 0.5960
B	 0.8400	 0.4920
C	 0.9650	 0.5930
D	 0.8360	 0.4900
E	 0.9630	 0.5950
F	 0.8410	 0.4920
G	 0.8450	 0.5220
H	 0.8510	 0.5250
I	 0.7550	 0.4560
J	 0.8360	 0.5210
K	 0.7410	 0.4530
L	 0.7580	 0.4620
M	 0.9640	 0.5350
N	 0.9290	 0.4520
O	 0.8210	 0.4740
P	 0.8930	 0.5510
Q	 0.9640	 0.5440
R	 0.8930	 0.4460
S	 0.8210	 0.4710
T	 0.8930	 0.5490
U	 0.9640	 0.5420
V	 0.9290	 0.4400
W	 0.8210	 0.4820
X	 0.8930	 0.5520

