



Full wwPDB EM Validation Report ⓘ

Nov 30, 2022 – 12:29 AM JST

PDB ID : 7XIY
EMDB ID : EMD-33212
Title : SARS-CoV-2 Omicron BA.3 variant spike
Authors : Wang, X.; Wang, L.
Deposited on : 2022-04-14
Resolution : 3.07 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.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

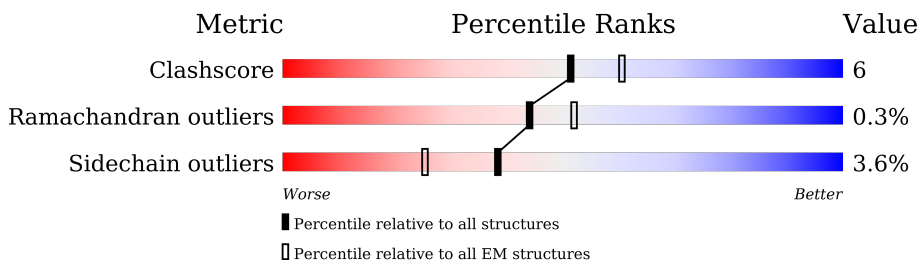
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.07 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	1267	 21% 70% 16% 14%
1	B	1267	 21% 70% 15% 14%
1	C	1267	 21% 71% 14% 14%
2	D	2	 100%
2	E	2	 100%
2	F	2	 100%
2	G	2	 100%
2	H	2	 100%

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Mol	Chain	Length	Quality of chain
2	I	2	 100%
2	J	2	 100%
2	K	2	 100%
2	L	2	 100%
2	M	2	 100%
2	N	2	 100%
2	O	2	 100%

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 26089 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	1085	8477	5423	1410	1607	37	0	0
1	A	1085	8477	5423	1410	1607	37	0	0
1	C	1085	8477	5423	1410	1607	37	0	0

There are 123 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	67	VAL	ALA	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	95	ILE	THR	variant	UNP P0DTC2
B	142	ASP	GLY	variant	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	?	-	ASN	deletion	UNP P0DTC2
B	212	ILE	LEU	variant	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	446	SER	GLY	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	484	ALA	GLU	variant	UNP P0DTC2
B	493	ARG	GLN	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	505	HIS	TYR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	HIS	PRO	variant	UNP P0DTC2
B	683	ALA	ARG	variant	UNP P0DTC2
B	685	ALA	ARG	variant	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	variant	UNP P0DTC2
B	892	PRO	ALA	variant	UNP P0DTC2
B	899	PRO	ALA	variant	UNP P0DTC2
B	942	PRO	ALA	variant	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
A	67	VAL	ALA	variant	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	95	ILE	THR	variant	UNP P0DTC2
A	142	ASP	GLY	variant	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	?	-	ASN	deletion	UNP P0DTC2
A	212	ILE	LEU	variant	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2
A	493	ARG	GLN	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	614	GLY	ASP	variant	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	HIS	PRO	variant	UNP P0DTC2
A	683	ALA	ARG	variant	UNP P0DTC2
A	685	ALA	ARG	variant	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	variant	UNP P0DTC2
A	892	PRO	ALA	variant	UNP P0DTC2
A	899	PRO	ALA	variant	UNP P0DTC2
A	942	PRO	ALA	variant	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	67	VAL	ALA	variant	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	95	ILE	THR	variant	UNP P0DTC2
C	142	ASP	GLY	variant	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	?	-	ASN	deletion	UNP P0DTC2
C	212	ILE	LEU	variant	UNP P0DTC2
C	339	ASP	GLY	variant	UNP P0DTC2
C	371	PHE	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	405	ASN	ASP	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	446	SER	GLY	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2
C	484	ALA	GLU	variant	UNP P0DTC2
C	493	ARG	GLN	variant	UNP P0DTC2
C	498	ARG	GLN	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	505	HIS	TYR	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	681	HIS	PRO	variant	UNP P0DTC2
C	683	ALA	ARG	variant	UNP P0DTC2
C	685	ALA	ARG	variant	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	817	PRO	PHE	variant	UNP P0DTC2
C	892	PRO	ALA	variant	UNP P0DTC2
C	899	PRO	ALA	variant	UNP P0DTC2
C	942	PRO	ALA	variant	UNP P0DTC2
C	954	HIS	GLN	variant	UNP P0DTC2
C	969	LYS	ASN	variant	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2

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



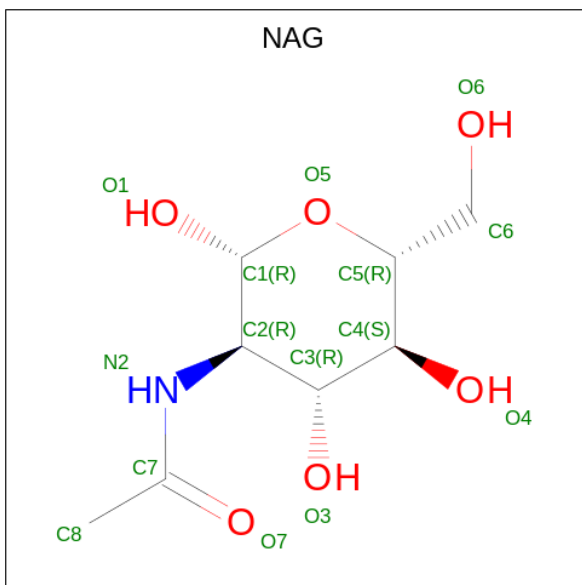
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
2	D	2	28	16	2	10	0	0
2	E	2	28	16	2	10	0	0
2	F	2	28	16	2	10	0	0
2	G	2	28	16	2	10	0	0
2	H	2	28	16	2	10	0	0
2	I	2	28	16	2	10	0	0
2	J	2	28	16	2	10	0	0
2	K	2	28	16	2	10	0	0
2	L	2	28	16	2	10	0	0

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

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
3	B	1	Total	C	N	O	0
			112	64	8	40	
3	B	1	Total	C	N	O	0
			112	64	8	40	
3	B	1	Total	C	N	O	0
			112	64	8	40	
3	B	1	Total	C	N	O	0
			112	64	8	40	
3	B	1	Total	C	N	O	0
			112	64	8	40	
3	B	1	Total	C	N	O	0
			112	64	8	40	
3	B	1	Total	C	N	O	0
			112	64	8	40	

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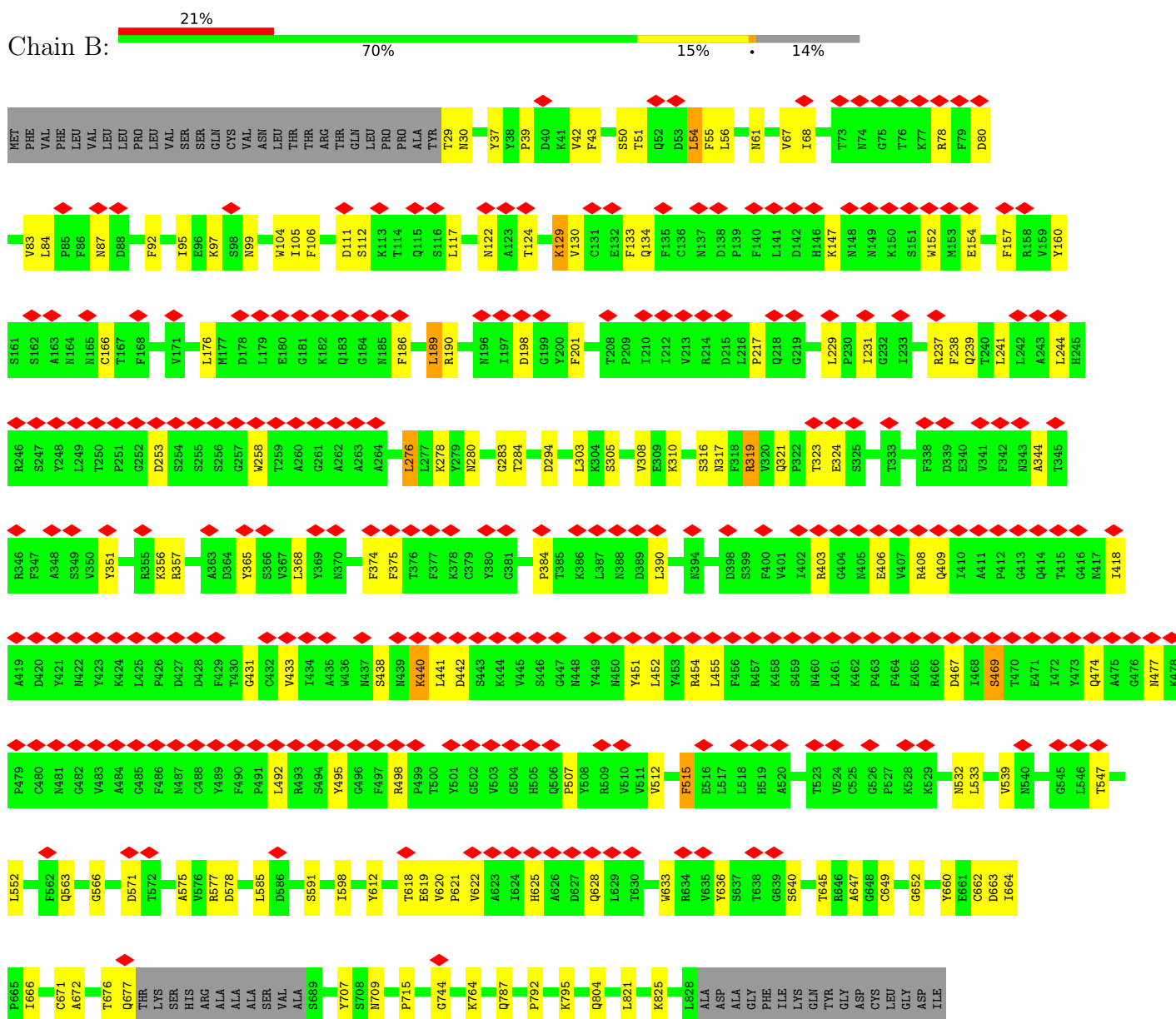
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Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	A	1	Total 98	C 56	N 7	O 35	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0
3	C	1	Total 112	C 64	N 8	O 40	0

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: Spike glycoprotein



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

Chain I:  100%



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

Chain J:  100%



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

Chain K:  100%



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

Chain L:  100%



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

Chain M:  100%



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

Chain N:  100%



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

Chain O:  100%

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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	185916	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K2 BASE (4k x 4k)	Depositor
Maximum map value	3.223	Depositor
Minimum map value	-1.505	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.081	Depositor
Recommended contour level	0.4	Depositor
Map size (\AA)	332.8, 332.8, 332.8	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.04, 1.04, 1.04	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	0.27	0/8684	0.57	9/11829 (0.1%)
1	B	0.34	2/8684 (0.0%)	0.60	7/11829 (0.1%)
1	C	0.29	1/8684 (0.0%)	0.56	3/11829 (0.0%)
All	All	0.30	3/26052 (0.0%)	0.58	19/35487 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	384	PRO	CG-CD	-16.45	0.96	1.50
1	C	373	PRO	CG-CD	-7.25	1.26	1.50
1	B	384	PRO	N-CD	7.17	1.57	1.47

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	384	PRO	N-CD-CG	-16.42	78.57	103.20
1	B	217	PRO	CA-N-CD	-11.86	94.90	111.50
1	C	373	PRO	CA-N-CD	-11.50	95.40	111.50
1	A	384	PRO	CA-N-CD	-11.25	95.75	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	384	PRO	CA-N-CD	-10.88	96.27	111.50
1	B	384	PRO	CA-CB-CG	-8.79	87.30	104.00
1	C	373	PRO	N-CD-CG	-8.65	90.22	103.20
1	A	110	LEU	CA-CB-CG	8.14	134.02	115.30
1	B	1118	ASP	CB-CG-OD1	6.93	124.54	118.30
1	B	217	PRO	N-CD-CG	-5.83	94.46	103.20
1	A	110	LEU	CB-CG-CD2	5.69	120.67	111.00
1	A	384	PRO	N-CD-CG	-5.60	94.80	103.20
1	A	53	ASP	CB-CG-OD1	5.51	123.26	118.30
1	A	117	LEU	CA-CB-CG	5.40	127.73	115.30
1	A	303	LEU	CA-CB-CG	5.33	127.55	115.30
1	B	54	LEU	CA-CB-CG	5.29	127.47	115.30
1	A	367	VAL	CG1-CB-CG2	-5.20	102.58	110.90
1	C	1118	ASP	CB-CG-OD1	5.10	122.89	118.30
1	A	118	LEU	CA-CB-CG	5.07	126.97	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	620	VAL	Peptide
1	B	620	VAL	Peptide
1	C	620	VAL	Peptide

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	8477	0	8258	113	0
1	B	8477	0	8257	111	0
1	C	8477	0	8257	102	0
2	D	28	0	25	0	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
2	I	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	28	0	25	0	0
2	K	28	0	25	0	0
2	L	28	0	25	0	0
2	M	28	0	25	0	0
2	N	28	0	25	0	0
2	O	28	0	25	0	0
3	A	98	0	91	0	0
3	B	112	0	104	0	0
3	C	112	0	104	0	0
All	All	26089	0	25371	297	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 6.

All (297) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:547:THR:OG1	1:A:978:ASN:HB3	1.57	1.03
1:A:748:GLU:OE2	1:A:981:LEU:HD13	1.68	0.94
1:A:748:GLU:OE2	1:A:981:LEU:CD1	2.31	0.78
1:B:1051:SER:HG	1:B:1064:HIS:HD1	1.33	0.77
1:B:547:THR:OG1	1:A:978:ASN:CB	2.38	0.71
1:B:973:ILE:HD12	1:B:983:ARG:HH21	1.57	0.68
1:C:324:GLU:HB2	1:C:539:VAL:HG12	1.76	0.68
1:A:1051:SER:HG	1:A:1064:HIS:HD1	1.42	0.67
1:C:332:ILE:O	1:C:528:LYS:NZ	2.28	0.65
1:B:324:GLU:HB2	1:B:539:VAL:HG12	1.79	0.64
1:A:324:GLU:HB2	1:A:539:VAL:HG12	1.79	0.63
1:B:201:PHE:HB3	1:B:229:LEU:HB3	1.80	0.63
1:B:825:LYS:HE3	1:B:939:SER:HA	1.81	0.63
1:A:504:GLY:HA3	1:C:373:PRO:HB2	1.83	0.60
1:B:106:PHE:HB2	1:B:117:LEU:HB2	1.83	0.60
1:B:39:PRO:HG2	1:B:51:THR:HG21	1.84	0.59
1:C:403:ARG:HA	1:C:507:PRO:HA	1.85	0.58
1:C:316:SER:OG	1:C:317:ASN:N	2.36	0.58
1:B:50:SER:HB2	1:B:276:LEU:HD12	1.84	0.58
1:B:129:LYS:HD3	1:B:166:CYS:HB3	1.86	0.58
1:A:337:PRO:HD2	1:A:358:ILE:HG23	1.87	0.57
1:C:605:SER:OG	1:C:606:ASN:N	2.37	0.57
1:B:976:VAL:HG12	1:B:979:ASP:H	1.69	0.57
1:C:742:ILE:O	1:C:1000:ARG:NH1	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:ILE:HB	1:A:395:VAL:HB	1.87	0.57
1:C:383:SER:HB3	1:C:386:LYS:HB2	1.87	0.56
1:B:357:ARG:HH21	1:A:230:PRO:HG2	1.71	0.56
1:A:317:ASN:HA	1:A:594:GLY:HA2	1.87	0.56
1:A:752:LEU:HD21	1:A:990:GLU:HB2	1.86	0.56
1:A:360:ASN:H	1:A:523:THR:HB	1.71	0.56
1:A:804:GLN:NE2	1:A:935:GLN:OE1	2.39	0.55
1:B:792:PRO:O	1:B:795:LYS:NZ	2.39	0.55
1:B:563:GLN:O	1:B:577:ARG:NH1	2.40	0.55
1:A:191:GLU:HB2	1:A:223:LEU:HD21	1.89	0.55
1:A:640:SER:HB3	1:A:652:GLY:HA2	1.89	0.55
1:A:777:ASN:OD1	1:A:1019:ARG:NH1	2.40	0.55
1:B:374:PHE:H	1:C:405:ASN:HD21	1.53	0.54
1:A:405:ASN:O	1:A:409:GLN:NE2	2.41	0.54
1:B:566:GLY:HA3	1:B:575:ALA:HB3	1.89	0.54
1:A:53:ASP:OD1	1:A:54:LEU:N	2.41	0.54
1:A:383:SER:HB3	1:A:386:LYS:HB2	1.90	0.54
1:C:29:THR:OG1	1:C:30:ASN:N	2.40	0.54
1:B:95:ILE:HD13	1:B:189:LEU:HG	1.90	0.54
1:A:474:GLN:NE2	1:A:477:ASN:O	2.41	0.54
1:B:474:GLN:NE2	1:B:477:ASN:O	2.41	0.54
1:C:460:ASN:O	1:C:462:LYS:NZ	2.40	0.54
1:A:287:ASP:OD1	1:A:288:ALA:N	2.41	0.54
1:B:984:LEU:HD12	1:B:988:GLU:HG3	1.90	0.53
1:C:467:ASP:N	1:C:467:ASP:OD1	2.41	0.53
1:C:676:THR:OG1	1:C:677:GLN:N	2.41	0.53
1:C:253:ASP:N	1:C:253:ASP:OD1	2.40	0.53
1:A:853:GLN:HB2	1:A:858:LEU:HB2	1.91	0.53
1:C:474:GLN:NE2	1:C:477:ASN:O	2.41	0.53
1:B:280:ASN:HD21	1:B:284:THR:HB	1.74	0.53
1:B:1107:ARG:HH22	1:A:896:ILE:HD11	1.74	0.53
1:A:316:SER:OG	1:A:317:ASN:N	2.42	0.53
1:B:804:GLN:NE2	1:B:935:GLN:OE1	2.39	0.52
1:A:676:THR:OG1	1:A:677:GLN:N	2.41	0.52
1:C:571:ASP:OD1	1:C:571:ASP:N	2.39	0.52
1:C:566:GLY:HA3	1:C:575:ALA:HB3	1.92	0.52
1:A:547:THR:O	1:C:978:ASN:ND2	2.43	0.52
1:C:941:THR:HB	1:C:944:ALA:HB2	1.91	0.52
1:B:452:LEU:HD12	1:B:492:LEU:HB3	1.92	0.52
1:A:50:SER:HB2	1:A:276:LEU:HD12	1.92	0.51
1:B:1093:GLY:O	1:A:904:TYR:OH	2.29	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:403:ARG:HA	1:B:507:PRO:HA	1.93	0.51
1:B:1035:GLY:HA3	1:C:1040:VAL:HG21	1.92	0.51
1:A:280:ASN:HD21	1:A:284:THR:HB	1.76	0.50
1:C:884:SER:O	1:C:887:THR:OG1	2.30	0.50
1:B:310:LYS:NZ	1:B:663:ASP:OD2	2.41	0.50
1:B:147:LYS:O	1:B:147:LYS:NZ	2.42	0.50
1:B:612:TYR:HB2	1:B:649:CYS:HB3	1.94	0.50
1:A:403:ARG:HA	1:A:507:PRO:HA	1.94	0.50
1:C:193:VAL:HB	1:C:204:TYR:HB2	1.93	0.50
1:B:552:LEU:HD12	1:B:585:LEU:HB3	1.94	0.49
1:B:1040:VAL:HG21	1:A:1035:GLY:HA3	1.92	0.49
1:A:715:PRO:HA	1:A:1072:GLU:HA	1.94	0.49
1:B:406:GLU:HA	1:B:409:GLN:HB2	1.93	0.49
1:B:467:ASP:OD1	1:B:467:ASP:N	2.45	0.49
1:C:152:TRP:NE1	1:C:154:GLU:OE2	2.45	0.49
1:C:473:TYR:HD1	1:C:474:GLN:H	1.59	0.49
1:B:316:SER:OG	1:B:317:ASN:N	2.44	0.49
1:A:977:LEU:HD22	1:A:1000:ARG:HH12	1.78	0.49
1:B:941:THR:HB	1:B:944:ALA:HB2	1.93	0.49
1:A:375:PHE:HB3	1:A:434:ILE:HG23	1.95	0.49
1:B:433:VAL:HA	1:B:512:VAL:HA	1.93	0.49
1:A:707:TYR:HB3	1:C:792:PRO:HG2	1.93	0.49
1:A:985:ASP:HB3	1:C:413:GLY:HA3	1.95	0.49
1:C:402:ILE:HD11	1:C:406:GLU:HB2	1.95	0.48
1:A:61:ASN:N	1:A:61:ASN:OD1	2.46	0.48
1:B:29:THR:OG1	1:B:30:ASN:N	2.45	0.48
1:A:555:SER:HB2	1:A:557:LYS:HG2	1.94	0.48
1:C:804:GLN:NE2	1:C:935:GLN:OE1	2.41	0.48
1:A:83:VAL:HG21	1:A:237:ARG:HH21	1.77	0.48
1:A:916:LEU:HD12	1:A:923:ILE:HD13	1.94	0.48
1:C:115:GLN:HG2	1:C:132:GLU:HG3	1.96	0.48
1:C:37:TYR:OH	1:C:54:LEU:O	2.31	0.48
1:C:339:ASP:OD1	1:C:339:ASP:N	2.47	0.48
1:A:617:CYS:HB2	1:A:649:CYS:HB2	1.61	0.48
1:C:244:LEU:HG	1:C:247:SER:HB2	1.96	0.48
1:C:645:THR:HG22	1:C:647:ALA:H	1.79	0.48
1:C:1051:SER:OG	1:C:1064:HIS:ND1	2.43	0.48
1:C:360:ASN:H	1:C:523:THR:HB	1.78	0.48
1:A:378:LYS:HE3	1:A:408:ARG:HA	1.95	0.48
1:C:984:LEU:HD23	1:C:988:GLU:HG3	1.96	0.47
1:A:213:VAL:O	1:A:214:ARG:NE	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:ARG:NE	1:A:459:SER:O	2.47	0.47
1:A:571:ASP:OD1	1:A:571:ASP:N	2.42	0.47
1:B:253:ASP:OD1	1:B:253:ASP:N	2.48	0.47
1:B:37:TYR:OH	1:B:54:LEU:O	2.29	0.47
1:B:676:THR:OG1	1:B:677:GLN:N	2.47	0.47
1:A:662:CYS:HB2	1:A:671:CYS:HB3	1.85	0.47
1:C:792:PRO:O	1:C:795:LYS:NZ	2.48	0.47
1:B:104:TRP:HB3	1:B:238:PHE:HE1	1.80	0.47
1:A:169:GLU:HG3	1:A:171:VAL:HG23	1.96	0.47
1:A:519:HIS:HB3	1:A:520:ALA:H	1.44	0.47
1:A:54:LEU:HD12	1:A:270:LEU:HB3	1.97	0.47
1:A:707:TYR:HB2	1:C:883:THR:HG23	1.96	0.47
1:C:367:VAL:HG23	1:C:370:ASN:HD21	1.79	0.47
1:B:67:VAL:HG22	1:B:80:ASP:HB3	1.96	0.46
1:B:571:ASP:OD1	1:B:571:ASP:N	2.39	0.46
1:B:666:ILE:HD11	1:B:672:ALA:HB2	1.98	0.46
1:B:764:LYS:NZ	1:C:314:GLN:OE1	2.48	0.46
1:B:927:PHE:HZ	1:B:1052:PHE:HE2	1.62	0.46
1:C:147:LYS:O	1:C:147:LYS:NZ	2.43	0.46
1:B:598:ILE:HG23	1:B:664:ILE:HG21	1.96	0.46
1:A:37:TYR:HB3	1:A:223:LEU:HB2	1.98	0.46
1:A:452:LEU:HD22	1:A:494:SER:HB3	1.98	0.46
1:A:743:CYS:HA	1:A:977:LEU:HD21	1.97	0.46
1:C:1048:HIS:HA	1:C:1066:THR:HG22	1.97	0.46
1:B:440:LYS:HD2	1:B:441:LEU:HG	1.98	0.46
1:A:744:GLY:H	1:A:977:LEU:HD23	1.79	0.46
1:A:986:PRO:HA	1:A:989:ALA:HB3	1.97	0.46
1:C:280:ASN:HD21	1:C:284:THR:HB	1.81	0.46
1:B:303:LEU:HD13	1:B:308:VAL:HG23	1.97	0.46
1:B:618:THR:HG23	1:B:619:GLU:HG3	1.96	0.46
1:B:1141:LEU:HG	1:B:1145:LEU:HD23	1.97	0.46
1:C:1052:PHE:HB2	1:C:1063:LEU:HB2	1.97	0.46
1:B:406:GLU:HB3	1:B:418:ILE:HG21	1.97	0.46
1:B:152:TRP:HD1	1:B:154:GLU:HG3	1.80	0.46
1:B:578:ASP:OD1	1:B:578:ASP:N	2.46	0.46
1:A:788:ILE:HG23	1:A:876:ALA:HB2	1.98	0.46
1:B:97:LYS:HA	1:B:97:LYS:HD2	1.77	0.46
1:A:204:TYR:HB3	1:A:223:LEU:HB3	1.96	0.46
1:A:544:ASN:HD21	1:A:579:PRO:HG3	1.81	0.46
1:C:317:ASN:HA	1:C:594:GLY:HA2	1.97	0.46
1:B:914:ASN:ND2	1:C:1123:SER:OG	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:ILE:HD11	1:B:78:ARG:HG3	1.98	0.45
1:C:129:LYS:HE3	1:C:129:LYS:HB3	1.73	0.45
1:A:708:SER:HG	1:A:711:SER:HG	1.63	0.45
1:C:736:VAL:HG12	1:C:858:LEU:HA	1.98	0.45
1:C:825:LYS:HA	1:C:825:LYS:HD3	1.76	0.45
1:B:112:SER:HB3	1:B:134:GLN:HB3	1.99	0.45
1:A:99:ASN:N	1:A:99:ASN:OD1	2.49	0.45
1:C:71:SER:HB2	1:C:244:LEU:HD23	1.99	0.45
1:C:552:LEU:HD12	1:C:585:LEU:HB3	1.99	0.45
1:B:866:THR:HG22	1:B:869:MET:HG3	1.97	0.45
1:B:986:PRO:HA	1:B:989:ALA:HB3	1.98	0.45
1:A:715:PRO:HD3	1:C:894:LEU:HD13	1.98	0.45
1:C:93:ALA:HB3	1:C:266:TYR:HB2	1.99	0.45
1:C:618:THR:HG23	1:C:619:GLU:HG3	1.98	0.45
1:B:84:LEU:HD12	1:B:238:PHE:HD2	1.81	0.45
1:B:451:TYR:HB2	1:B:495:TYR:H	1.81	0.45
1:A:273:ARG:HA	1:A:273:ARG:HD3	1.68	0.45
1:A:452:LEU:HD12	1:A:492:LEU:HB3	1.99	0.45
1:A:605:SER:OG	1:A:606:ASN:N	2.50	0.45
1:B:319:ARG:NH1	1:B:591:SER:O	2.50	0.45
1:B:176:LEU:HB3	1:B:190:ARG:HH21	1.82	0.45
1:A:1052:PHE:HB2	1:A:1063:LEU:HB2	1.99	0.45
1:C:532:ASN:N	1:C:532:ASN:OD1	2.47	0.45
1:A:345:THR:O	1:A:509:ARG:NH2	2.47	0.45
1:C:137:ASN:N	1:C:159:VAL:O	2.48	0.45
1:B:547:THR:HG1	1:A:978:ASN:HB3	1.74	0.44
1:C:816:SER:OG	1:C:819:GLU:OE1	2.33	0.44
1:B:42:VAL:O	1:C:563:GLN:NE2	2.51	0.44
1:B:124:THR:O	1:B:124:THR:OG1	2.35	0.44
1:B:1047:TYR:HB2	1:B:1067:TYR:HB3	1.98	0.44
1:C:310:LYS:HB2	1:C:310:LYS:HE3	1.79	0.44
1:B:792:PRO:HG2	1:C:707:TYR:HB3	1.99	0.44
1:C:101:ILE:HD12	1:C:242:LEU:HB2	1.98	0.44
1:B:133:PHE:HE2	1:B:160:TYR:HB2	1.82	0.44
1:C:454:ARG:O	1:C:456:PHE:N	2.50	0.44
1:C:986:PRO:HA	1:C:989:ALA:HB3	1.99	0.44
1:B:321:GLN:N	1:B:628:GLN:O	2.50	0.44
1:A:231:ILE:HD12	1:A:233:ILE:H	1.82	0.44
1:A:406:GLU:HA	1:A:409:GLN:HB2	1.98	0.44
1:B:129:LYS:HB3	1:B:129:LYS:HE3	1.77	0.44
1:B:787:GLN:OE1	1:C:703:ASN:ND2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1006:THR:OG1	1:C:1005:GLN:OE1	2.30	0.44
1:B:99:ASN:OD1	1:B:99:ASN:N	2.50	0.44
1:B:323:THR:HG23	1:B:324:GLU:HG2	1.99	0.44
1:A:655:TYR:OH	1:A:696:THR:OG1	2.32	0.44
1:A:736:VAL:HG11	1:A:1004:LEU:HD11	1.99	0.44
1:A:1053:PRO:O	1:A:1054:GLN:NE2	2.47	0.44
1:C:710:ASN:N	1:C:710:ASN:OD1	2.49	0.44
1:A:398:ASP:HB2	1:A:512:VAL:HG22	1.99	0.44
1:C:358:ILE:HB	1:C:395:VAL:HB	2.00	0.44
1:B:707:TYR:HB2	1:A:883:THR:HG23	1.99	0.43
1:A:96:GLU:OE1	1:A:101:ILE:N	2.47	0.43
1:C:296:LEU:O	1:C:299:THR:OG1	2.29	0.43
1:A:299:THR:O	1:A:303:LEU:HD12	2.18	0.43
1:C:157:PHE:HB3	1:C:158:ARG:H	1.64	0.43
1:C:308:VAL:HG22	1:C:602:THR:HG23	2.00	0.43
1:C:97:LYS:HA	1:C:97:LYS:HD2	1.76	0.43
1:A:127:VAL:HG22	1:A:171:VAL:HG22	2.00	0.43
1:A:567:ARG:HG2	1:A:573:THR:HG22	1.99	0.43
1:C:502:GLY:O	1:C:504:GLY:N	2.41	0.43
1:A:210:ILE:HG22	1:A:212:ILE:HG23	2.01	0.43
1:C:244:LEU:HD12	1:C:244:LEU:HA	1.87	0.43
1:B:244:LEU:HD12	1:B:244:LEU:HA	1.89	0.43
1:B:645:THR:HG22	1:B:647:ALA:H	1.81	0.43
1:B:884:SER:OG	1:B:887:THR:OG1	2.33	0.43
1:A:864:LEU:HA	1:A:864:LEU:HD23	1.82	0.43
1:A:133:PHE:HD1	1:A:161:SER:HB3	1.83	0.43
1:A:173:GLN:HG2	1:A:174:PRO:HD2	2.00	0.43
1:B:662:CYS:HB2	1:B:671:CYS:HB3	1.71	0.43
1:A:415:THR:HA	1:A:419:ALA:HB3	2.01	0.43
1:B:438:SER:HB2	1:B:507:PRO:HG2	2.01	0.42
1:A:105:ILE:HG12	1:A:241:LEU:HD11	2.00	0.42
1:A:369:TYR:HB3	1:A:375:PHE:HZ	1.84	0.42
1:A:555:SER:OG	1:A:584:ILE:O	2.32	0.42
1:C:1082:CYS:HB2	1:C:1126:CYS:HB2	1.87	0.42
1:B:83:VAL:HG21	1:B:237:ARG:HH21	1.83	0.42
1:A:310:LYS:HE2	1:A:310:LYS:HB3	1.92	0.42
1:C:621:PRO:HB2	1:C:622:VAL:H	1.59	0.42
1:C:916:LEU:HD12	1:C:923:ILE:HD13	2.01	0.42
1:B:709:ASN:OD1	1:B:709:ASN:N	2.51	0.42
1:C:197:ILE:HB	1:C:202:LYS:HZ1	1.84	0.42
1:B:989:ALA:O	1:B:993:ILE:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:913:GLN:H	1:A:913:GLN:HG2	1.51	0.42
1:C:454:ARG:NH2	1:C:467:ASP:O	2.52	0.42
1:C:566:GLY:N	1:C:575:ALA:O	2.40	0.42
1:C:776:LYS:O	1:C:780:GLU:HG2	2.20	0.42
1:B:39:PRO:HG3	1:B:55:PHE:HZ	1.85	0.42
1:B:105:ILE:HB	1:B:239:GLN:HB3	2.02	0.42
1:B:715:PRO:HA	1:B:1072:GLU:HA	2.01	0.42
1:A:67:VAL:HG22	1:A:80:ASP:HB3	2.00	0.42
1:A:550:GLY:HA2	1:A:589:PRO:HA	2.00	0.42
1:A:1082:CYS:HB2	1:A:1126:CYS:HB2	1.69	0.42
1:B:122:ASN:HA	1:B:157:PHE:HZ	1.84	0.42
1:B:351:TYR:HB2	1:B:454:ARG:HG2	2.02	0.42
1:B:625:HIS:ND1	1:B:628:GLN:OE1	2.52	0.42
1:A:621:PRO:O	1:A:623:ALA:N	2.53	0.42
1:A:1093:GLY:O	1:C:904:TYR:OH	2.38	0.42
1:C:130:VAL:HG11	1:C:231:ILE:HG21	2.02	0.42
1:C:337:PRO:HD2	1:C:358:ILE:HG23	2.02	0.42
1:C:438:SER:HB2	1:C:507:PRO:HG2	2.02	0.42
1:C:640:SER:HB3	1:C:652:GLY:HA2	2.01	0.42
1:B:61:ASN:OD1	1:B:61:ASN:N	2.53	0.42
1:B:117:LEU:HA	1:B:130:VAL:HG22	2.00	0.42
1:A:41:LYS:HE3	1:A:41:LYS:HB3	1.84	0.41
1:C:153:MET:O	1:C:154:GLU:HG3	2.19	0.41
1:A:640:SER:OG	1:A:654:GLU:OE2	2.39	0.41
1:A:1107:ARG:HH22	1:C:896:ILE:HD12	1.86	0.41
1:C:34:ARG:NH1	1:C:217:PRO:O	2.49	0.41
1:C:202:LYS:HE2	1:C:202:LYS:HB2	1.72	0.41
1:B:533:LEU:HD23	1:B:533:LEU:HA	1.89	0.41
1:B:1094:VAL:HG23	1:A:900:MET:HE1	2.03	0.41
1:A:722:VAL:HG22	1:A:1065:VAL:HG22	2.02	0.41
1:C:776:LYS:NZ	1:C:780:GLU:OE1	2.54	0.41
1:B:130:VAL:HG11	1:B:231:ILE:HD13	2.03	0.41
1:A:566:GLY:HA3	1:A:575:ALA:HB3	2.03	0.41
1:A:1004:LEU:HD12	1:A:1004:LEU:HA	1.90	0.41
1:C:715:PRO:HA	1:C:1072:GLU:HA	2.01	0.41
1:B:43:PHE:HE1	1:B:283:GLY:HA3	1.85	0.41
1:C:497:PHE:HZ	1:C:507:PRO:HD3	1.86	0.41
1:A:327:VAL:HG13	1:A:542:ASN:HB3	2.03	0.41
1:C:74:ASN:OD1	1:C:74:ASN:N	2.49	0.41
1:B:152:TRP:HD1	1:B:154:GLU:CG	2.34	0.41
1:B:454:ARG:NH1	1:B:469:SER:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:744:GLY:H	1:B:977:LEU:HD22	1.86	0.41
1:C:339:ASP:HA	1:C:371:PHE:HE2	1.84	0.41
1:B:130:VAL:HG21	1:B:231:ILE:HD13	2.03	0.41
1:B:365:TYR:HA	1:B:368:LEU:HD23	2.03	0.41
1:B:408:ARG:HH12	1:A:377:PHE:HB2	1.85	0.41
1:B:431:GLY:HA2	1:B:515:PHE:CE2	2.56	0.41
1:A:412:PRO:HD3	1:A:425:LEU:HD23	2.03	0.41
1:C:406:GLU:HA	1:C:409:GLN:HB2	2.02	0.41
1:B:152:TRP:CD1	1:B:154:GLU:OE2	2.74	0.40
1:A:318:PHE:HA	1:A:631:PRO:HG2	2.03	0.40
1:A:561:PRO:O	1:A:577:ARG:NH1	2.50	0.40
1:B:894:LEU:HB3	1:C:713:ALA:HB3	2.02	0.40
1:A:406:GLU:HG3	1:A:418:ILE:HG13	2.02	0.40
1:C:84:LEU:HA	1:C:85:PRO:HD3	1.91	0.40
1:B:1072:GLU:N	1:B:1072:GLU:OE2	2.53	0.40
1:A:58:PHE:HB2	1:A:293:LEU:HD22	2.03	0.40
1:A:557:LYS:HD3	1:A:557:LYS:HA	1.81	0.40
1:C:201:PHE:HB3	1:C:229:LEU:HB2	2.02	0.40
1:C:435:ALA:HB2	1:C:510:VAL:HG22	2.03	0.40
1:B:869:MET:HB3	1:C:699:LEU:HD21	2.02	0.40
1:A:323:THR:HG23	1:A:324:GLU:HG2	2.02	0.40
1:A:472:ILE:HA	1:A:472:ILE:HD12	1.86	0.40
1:C:127:VAL:HG12	1:C:171:VAL:HG22	2.03	0.40
1:B:356:LYS:HB3	1:B:356:LYS:HE3	1.94	0.40
1:B:640:SER:HB3	1:B:652:GLY:HA2	2.03	0.40
1:B:715:PRO:HD3	1:A:894:LEU:HD13	2.02	0.40
1:C:936:ASP:OD1	1:C:936:ASP:N	2.54	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	1079/1267 (85%)	997 (92%)	79 (7%)	3 (0%)	41	71
1	B	1079/1267 (85%)	992 (92%)	83 (8%)	4 (0%)	34	66
1	C	1079/1267 (85%)	990 (92%)	85 (8%)	4 (0%)	34	66
All	All	3237/3801 (85%)	2979 (92%)	247 (8%)	11 (0%)	44	71

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	621	PRO
1	A	621	PRO
1	C	621	PRO
1	B	455	LEU
1	A	455	LEU
1	C	503	VAL
1	C	403	ARG
1	B	344	ALA
1	B	622	VAL
1	A	622	VAL
1	C	622	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	945/1109 (85%)	918 (97%)	27 (3%)	42	70
1	B	945/1109 (85%)	910 (96%)	35 (4%)	34	65
1	C	945/1109 (85%)	906 (96%)	39 (4%)	30	62
All	All	2835/3327 (85%)	2734 (96%)	101 (4%)	38	66

All (101) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	56	LEU
1	B	87	ASN

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Mol	Chain	Res	Type
1	B	92	PHE
1	B	111	ASP
1	B	129	LYS
1	B	186	PHE
1	B	189	LEU
1	B	198	ASP
1	B	241	LEU
1	B	258	TRP
1	B	276	LEU
1	B	278	LYS
1	B	294	ASP
1	B	305	SER
1	B	319	ARG
1	B	375	PHE
1	B	390	LEU
1	B	440	LYS
1	B	442	ASP
1	B	469	SER
1	B	498	ARG
1	B	515	PHE
1	B	532	ASN
1	B	633	TRP
1	B	636	TYR
1	B	660	TYR
1	B	821	LEU
1	B	855	PHE
1	B	927	PHE
1	B	939	SER
1	B	940	SER
1	B	969	LYS
1	B	1041	ASP
1	B	1097	SER
1	B	1098	ASN
1	A	79	PHE
1	A	175	PHE
1	A	186	PHE
1	A	190	ARG
1	A	301	CYS
1	A	365	TYR
1	A	375	PHE
1	A	387	LEU
1	A	389	ASP

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Mol	Chain	Res	Type
1	A	409	GLN
1	A	423	TYR
1	A	440	LYS
1	A	441	LEU
1	A	456	PHE
1	A	515	PHE
1	A	517	LEU
1	A	611	LEU
1	A	637	SER
1	A	677	GLN
1	A	759	PHE
1	A	808	ASP
1	A	873	TYR
1	A	960	ASN
1	A	994	ASP
1	A	1050	MET
1	A	1101	HIS
1	A	1127	ASP
1	C	122	ASN
1	C	160	TYR
1	C	175	PHE
1	C	186	PHE
1	C	189	LEU
1	C	207	HIS
1	C	220	PHE
1	C	258	TRP
1	C	278	LYS
1	C	316	SER
1	C	325	SER
1	C	356	LYS
1	C	375	PHE
1	C	403	ARG
1	C	408	ARG
1	C	440	LYS
1	C	456	PHE
1	C	473	TYR
1	C	498	ARG
1	C	535	LYS
1	C	544	ASN
1	C	559	PHE
1	C	590	CYS
1	C	592	PHE

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Mol	Chain	Res	Type
1	C	605	SER
1	C	660	TYR
1	C	662	CYS
1	C	740	MET
1	C	765	ARG
1	C	779	GLN
1	C	790	LYS
1	C	869	MET
1	C	873	TYR
1	C	936	ASP
1	C	995	ARG
1	C	1019	ARG
1	C	1050	MET
1	C	1098	ASN
1	C	1101	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

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
2	NAG	D	1	1,2	14,14,15	0.22	0	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	D	2	2	14,14,15	0.26	0	17,19,21	0.44	0
2	NAG	E	1	1,2	14,14,15	0.23	0	17,19,21	0.50	0
2	NAG	E	2	2	14,14,15	0.24	0	17,19,21	0.48	0
2	NAG	F	1	1,2	14,14,15	0.38	0	17,19,21	0.45	0
2	NAG	F	2	2	14,14,15	0.29	0	17,19,21	0.42	0
2	NAG	G	1	1,2	14,14,15	0.21	0	17,19,21	0.50	0
2	NAG	G	2	2	14,14,15	0.24	0	17,19,21	0.46	0
2	NAG	H	1	1,2	14,14,15	0.25	0	17,19,21	0.48	0
2	NAG	H	2	2	14,14,15	0.28	0	17,19,21	0.43	0
2	NAG	I	1	1,2	14,14,15	0.24	0	17,19,21	0.43	0
2	NAG	I	2	2	14,14,15	0.32	0	17,19,21	0.49	0
2	NAG	J	1	1,2	14,14,15	0.40	0	17,19,21	0.46	0
2	NAG	J	2	2	14,14,15	0.27	0	17,19,21	0.42	0
2	NAG	K	1	1,2	14,14,15	0.20	0	17,19,21	0.45	0
2	NAG	K	2	2	14,14,15	0.29	0	17,19,21	0.49	0
2	NAG	L	1	1,2	14,14,15	0.25	0	17,19,21	0.46	0
2	NAG	L	2	2	14,14,15	0.27	0	17,19,21	0.44	0
2	NAG	M	1	1,2	14,14,15	0.22	0	17,19,21	0.45	0
2	NAG	M	2	2	14,14,15	0.23	0	17,19,21	0.51	0
2	NAG	N	1	1,2	14,14,15	0.42	0	17,19,21	0.43	0
2	NAG	N	2	2	14,14,15	0.29	0	17,19,21	0.42	0
2	NAG	O	1	1,2	14,14,15	0.20	0	17,19,21	0.47	0
2	NAG	O	2	2	14,14,15	0.28	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
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	NAG	J	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	NAG	K	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	K	2	2	-	0/6/23/26	0/1/1/1
2	NAG	L	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	L	2	2	-	0/6/23/26	0/1/1/1
2	NAG	M	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	M	2	2	-	2/6/23/26	0/1/1/1
2	NAG	N	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	N	2	2	-	2/6/23/26	0/1/1/1
2	NAG	O	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	O	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	N	2	NAG	O5-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
2	M	2	NAG	O5-C5-C6-O6
2	J	2	NAG	C4-C5-C6-O6
2	M	1	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
2	M	2	NAG	C4-C5-C6-O6
2	N	1	NAG	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	O	1	NAG	O5-C5-C6-O6
2	N	2	NAG	C4-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	N	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C8-C7-N2-C2
2	F	1	NAG	O7-C7-N2-C2
2	J	1	NAG	C8-C7-N2-C2
2	J	1	NAG	O7-C7-N2-C2
2	N	1	NAG	C8-C7-N2-C2
2	N	1	NAG	O7-C7-N2-C2

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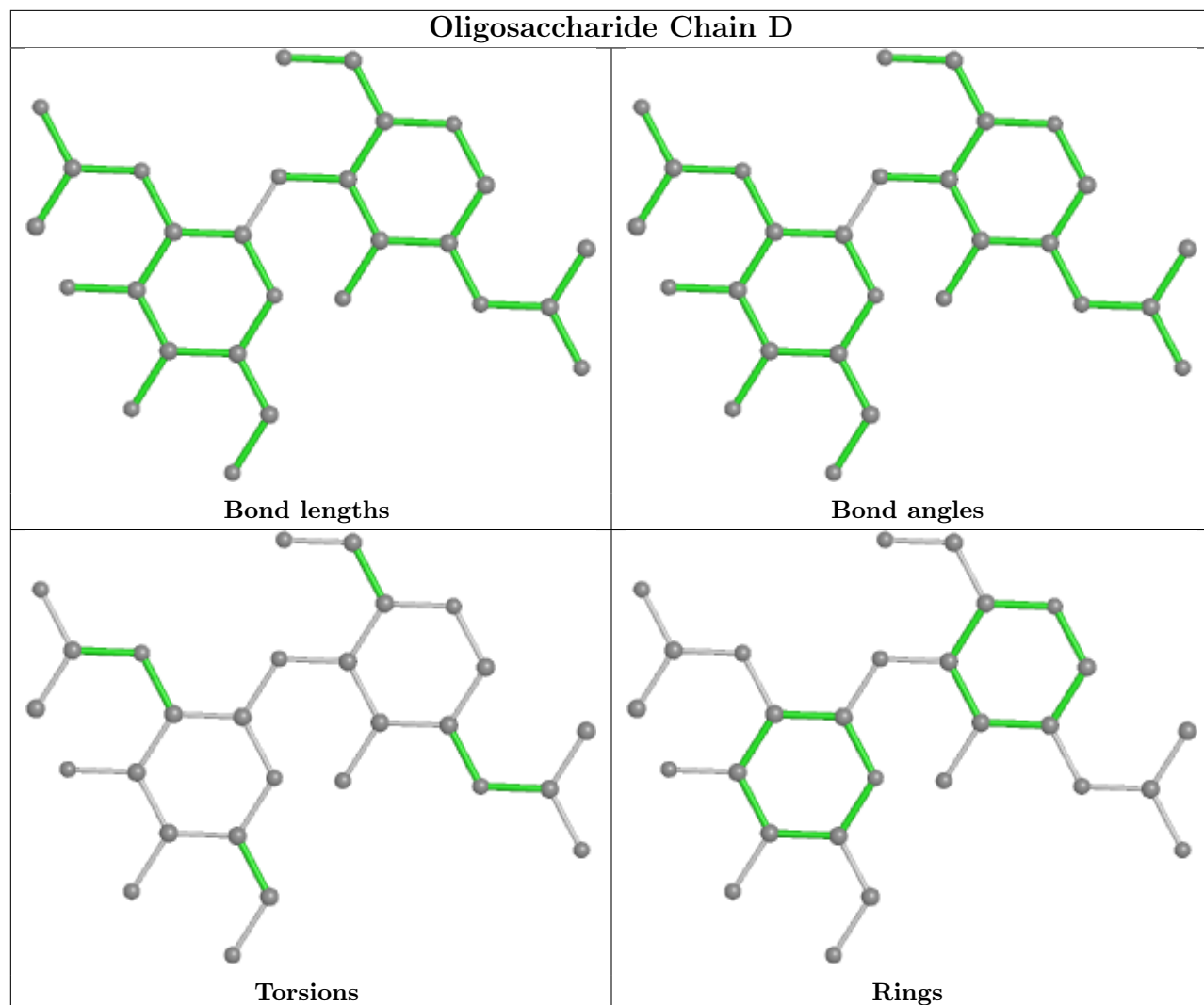
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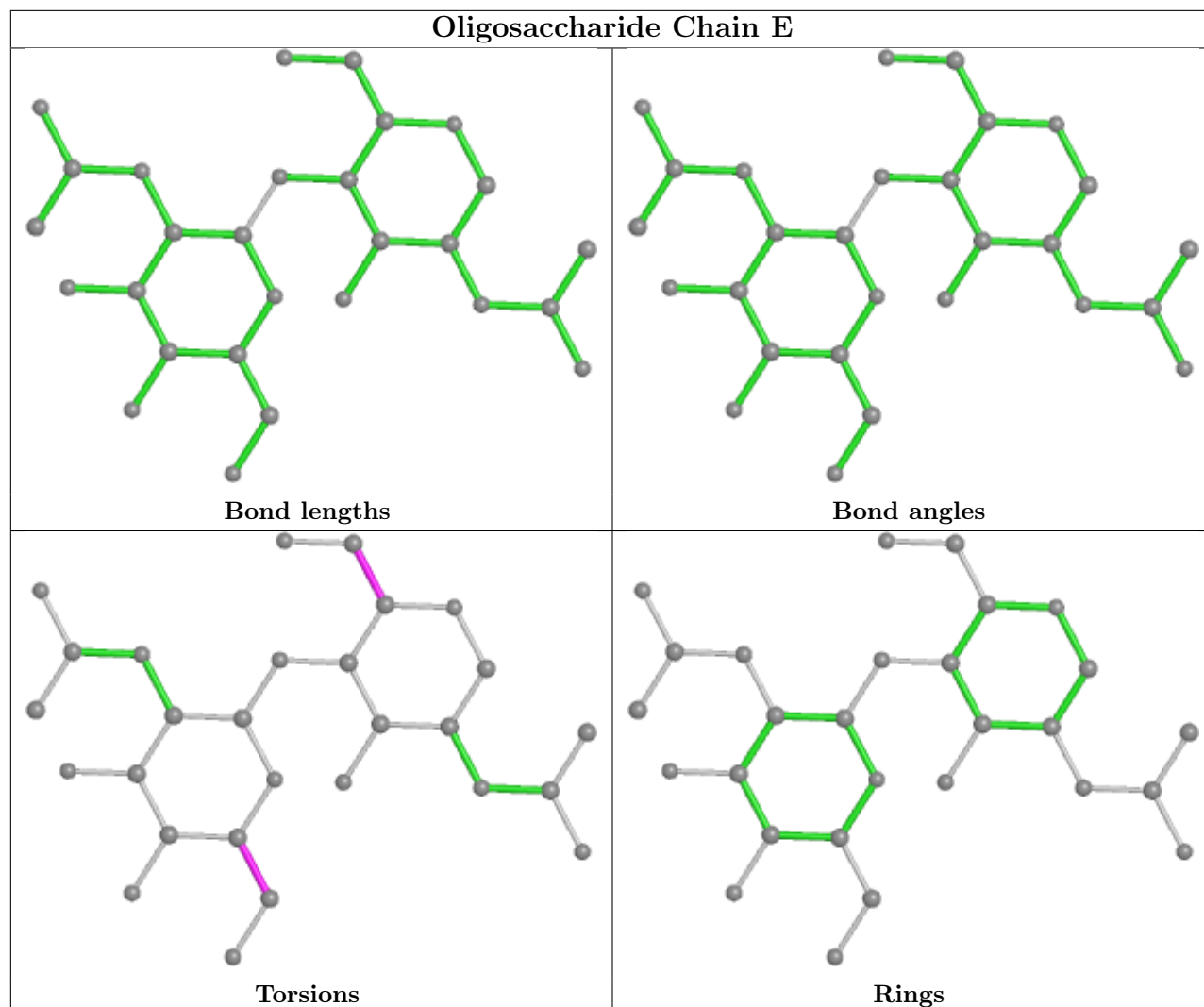
Mol	Chain	Res	Type	Atoms
2	M	1	NAG	C4-C5-C6-O6
2	O	1	NAG	C4-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6

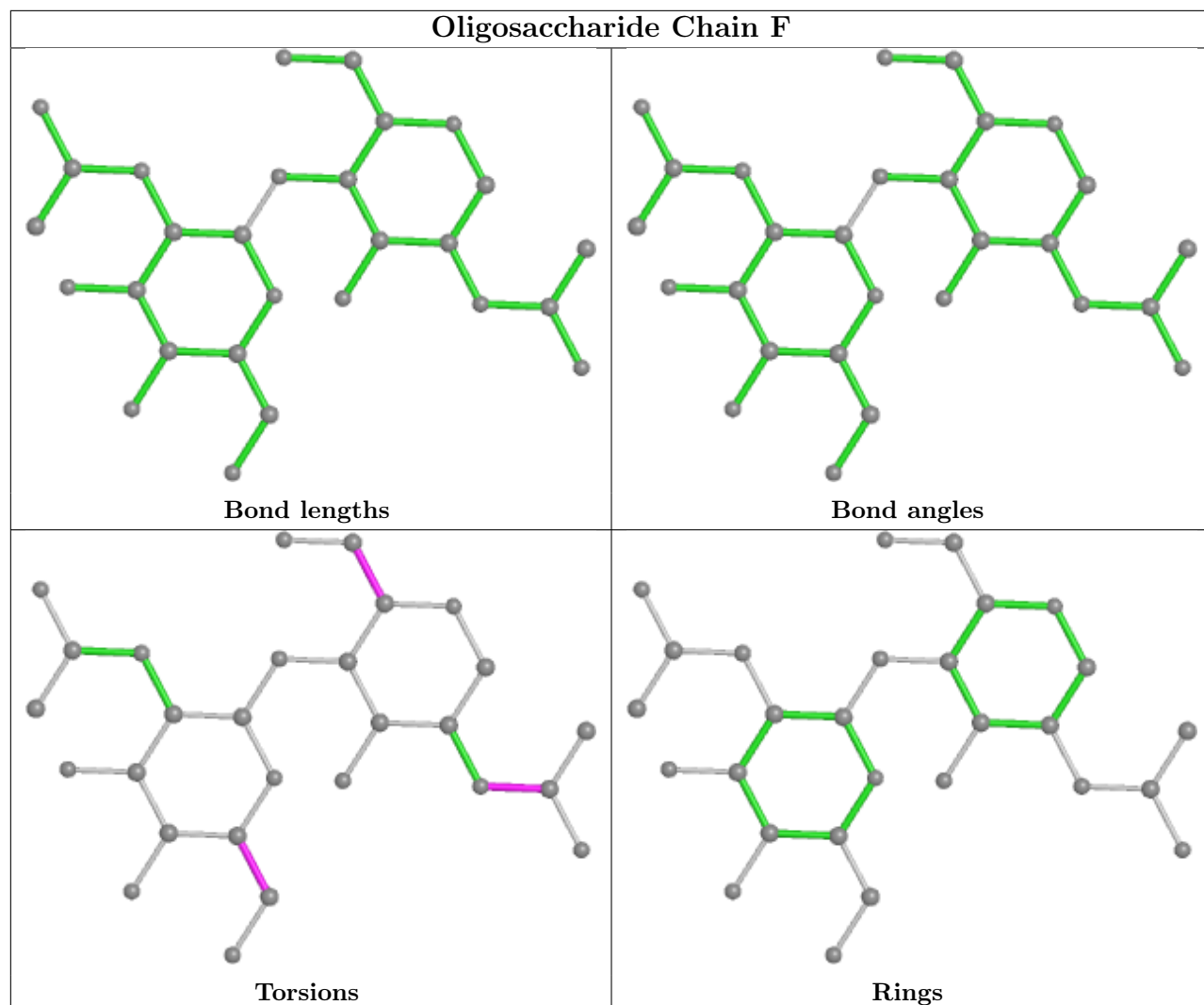
There are no ring outliers.

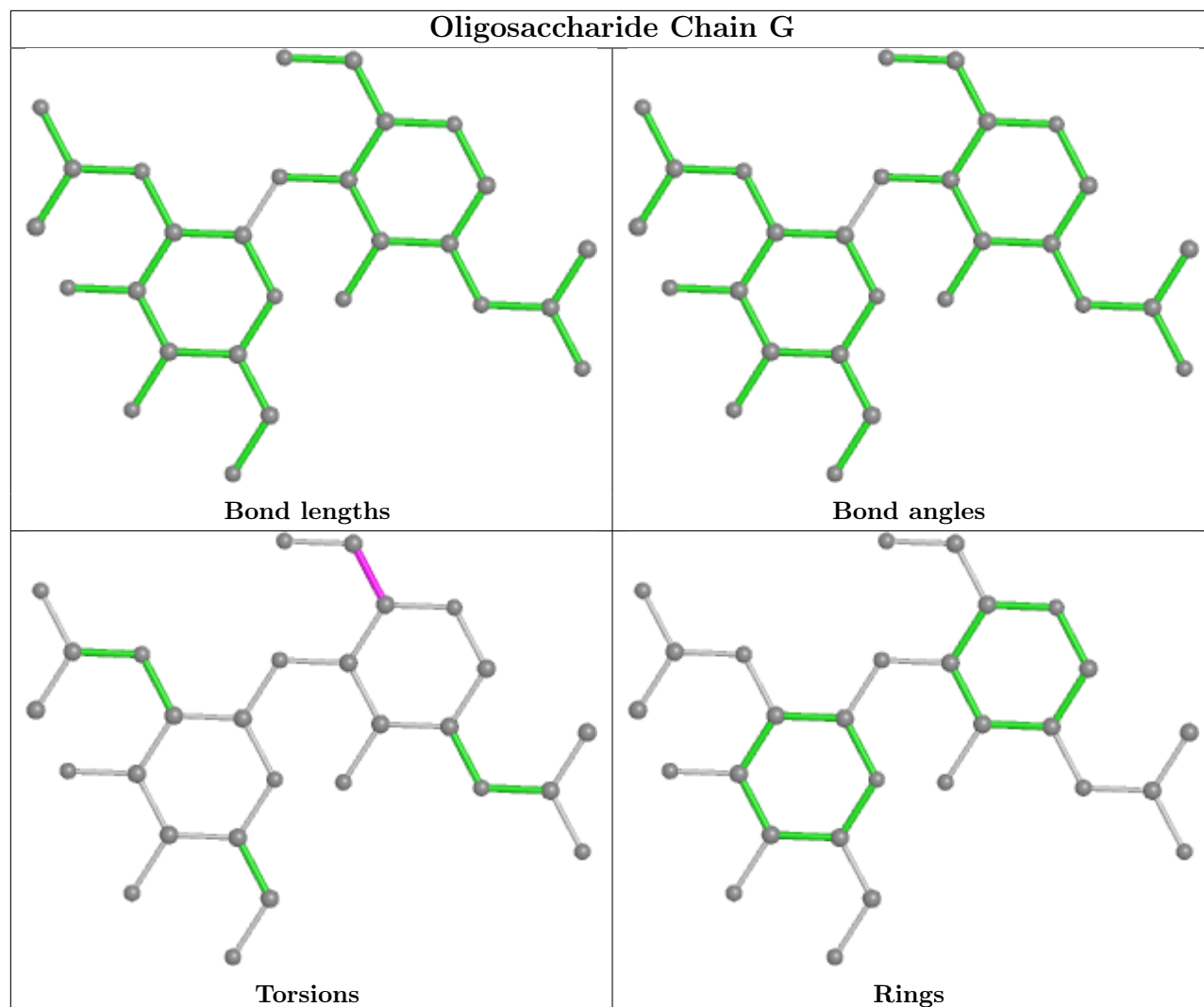
No monomer is involved in short contacts.

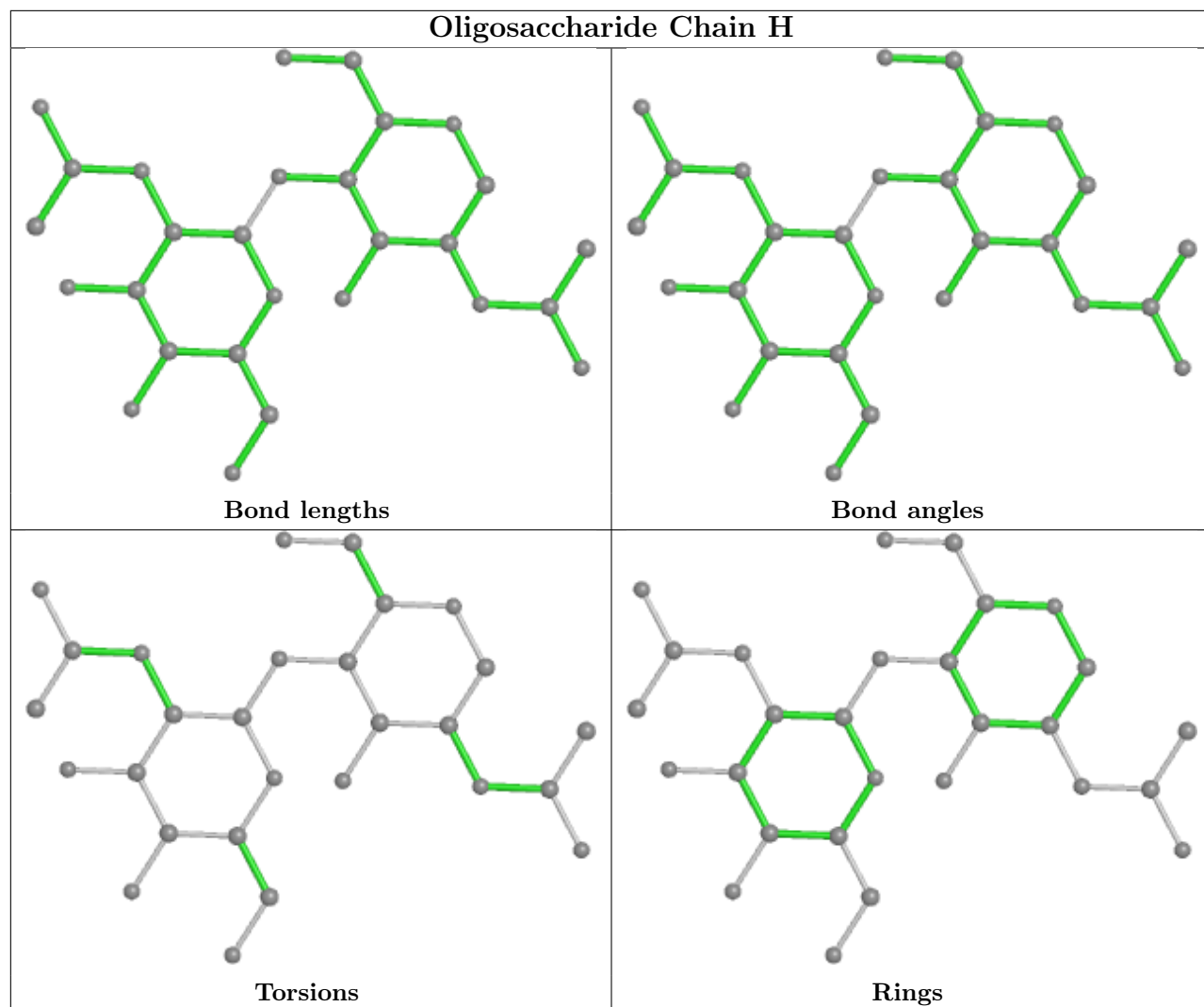
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

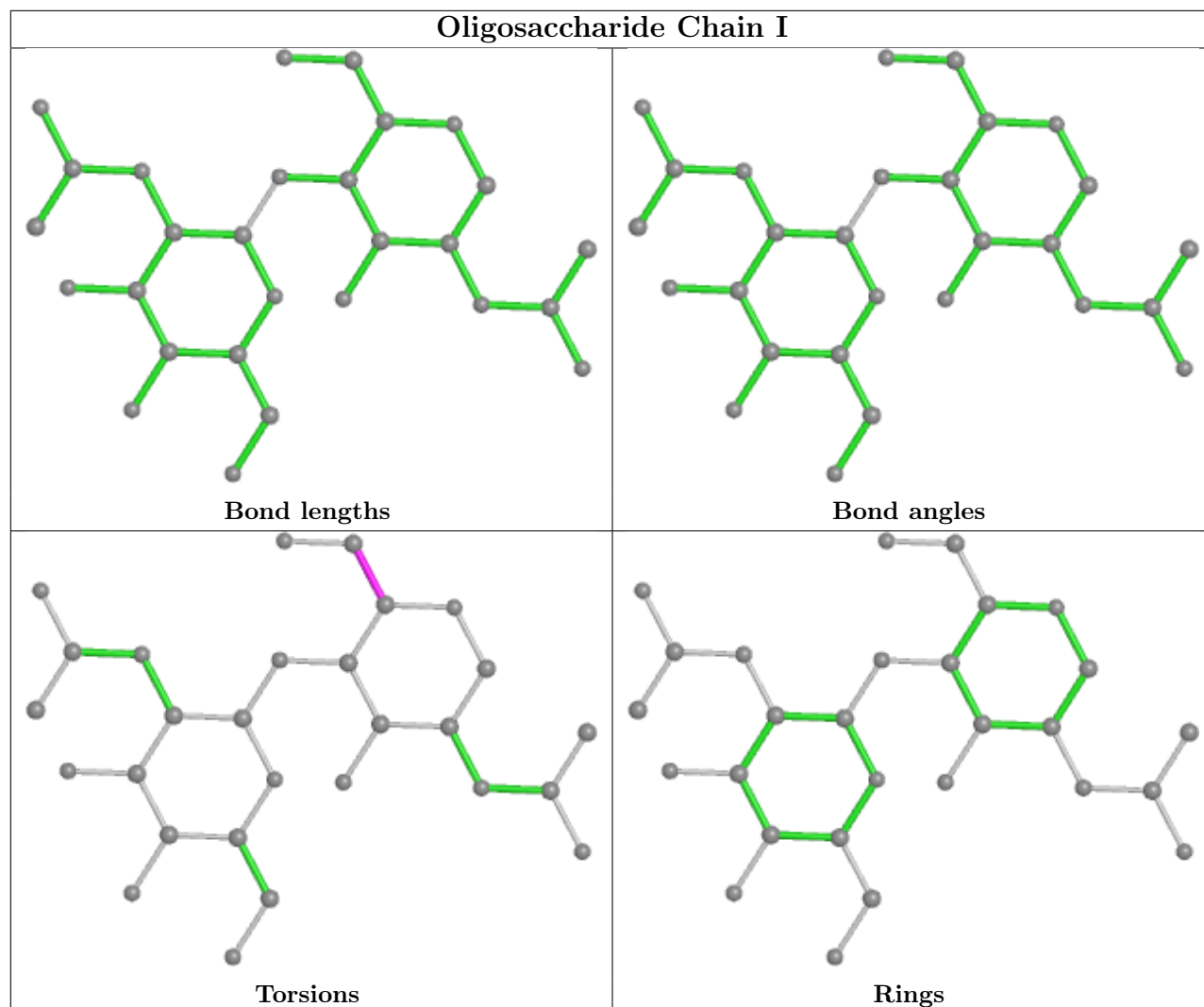


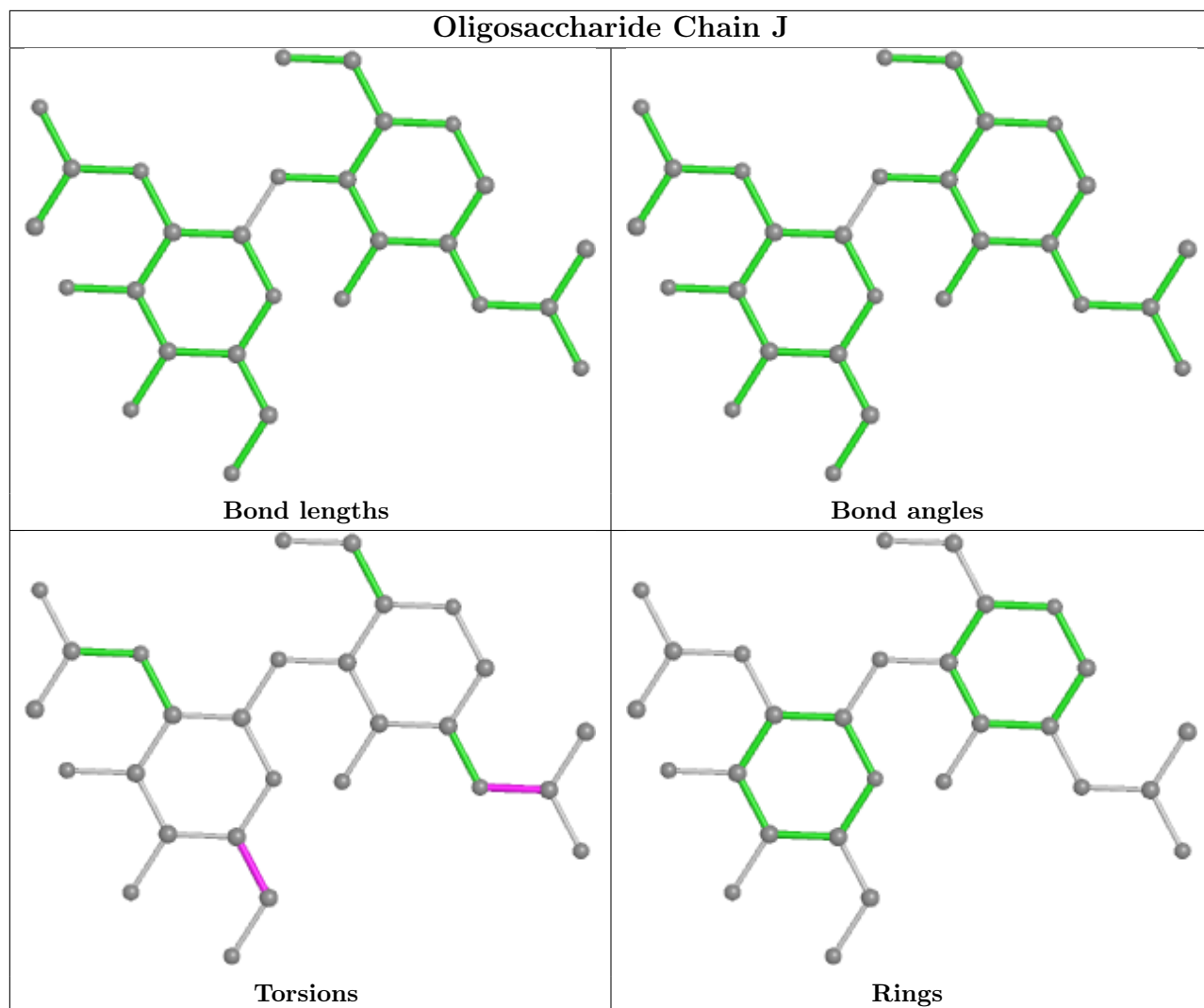


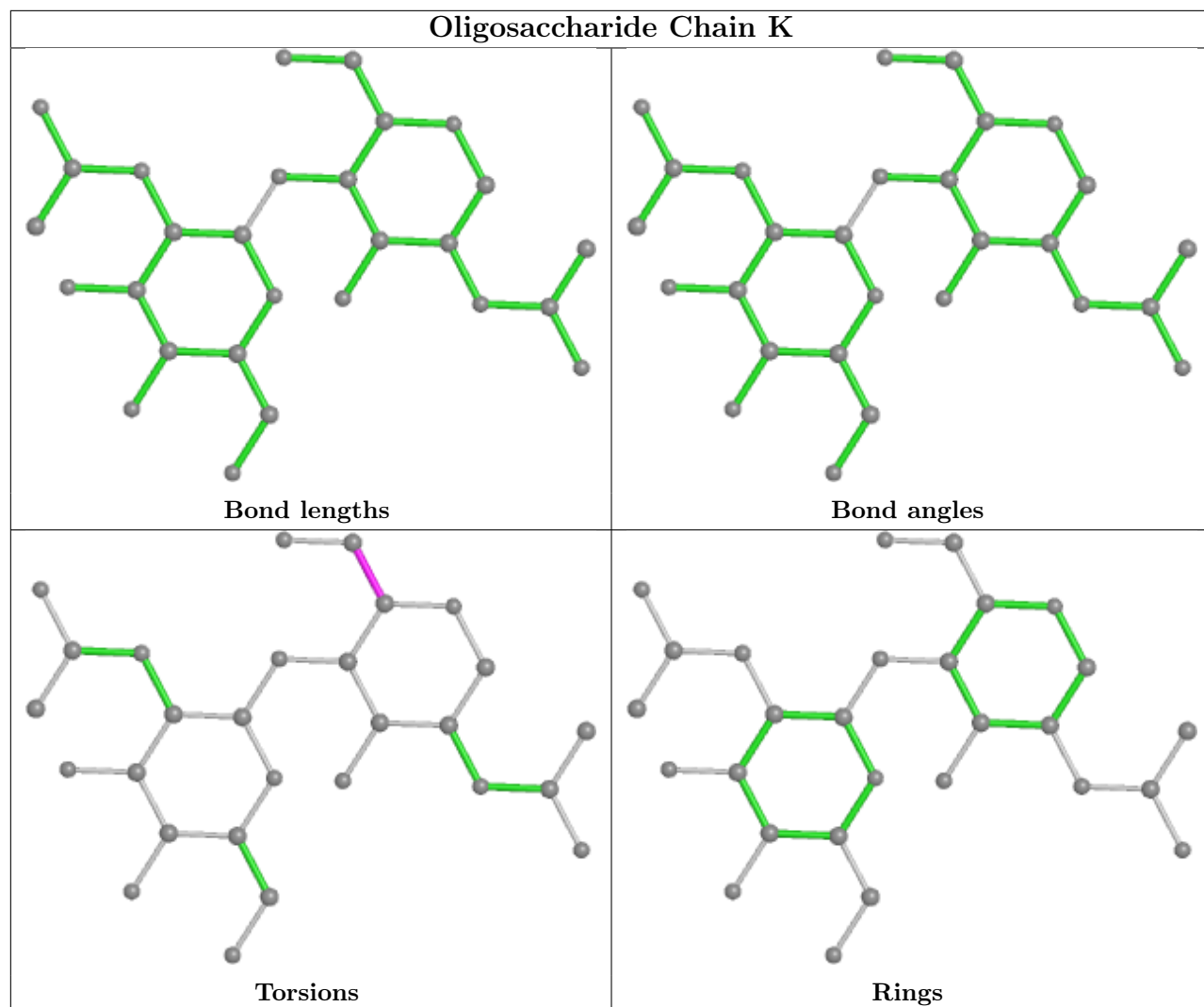


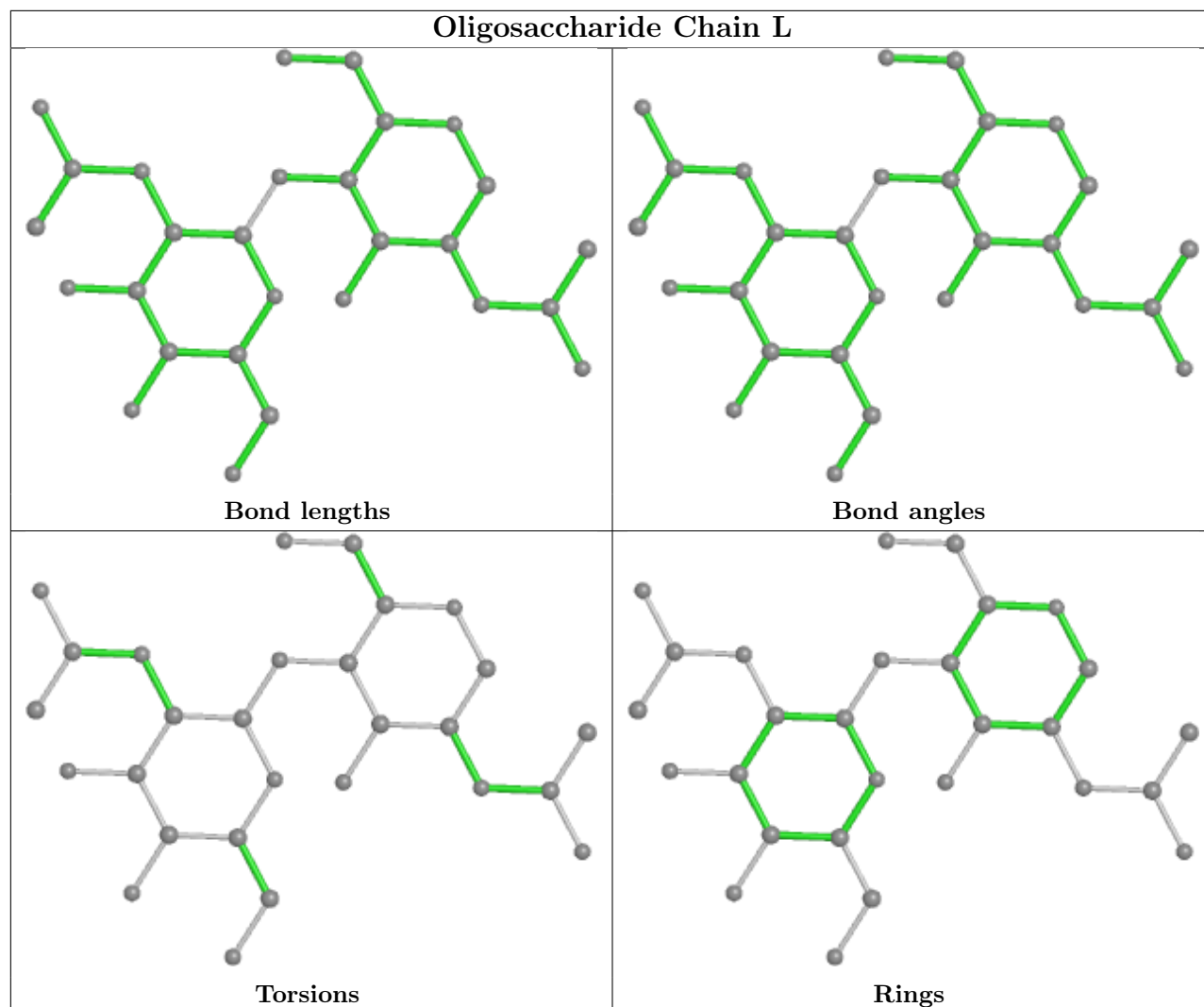


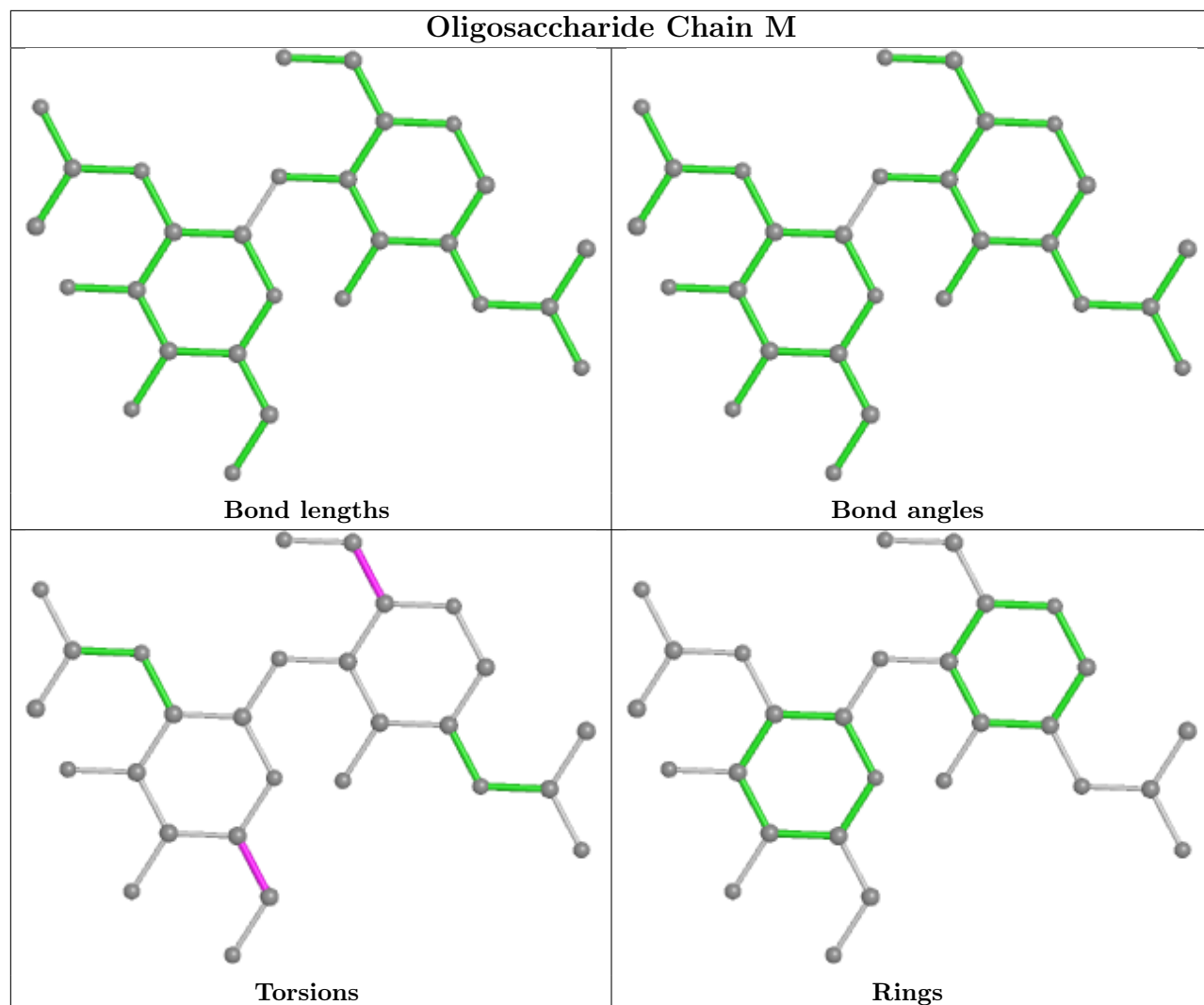


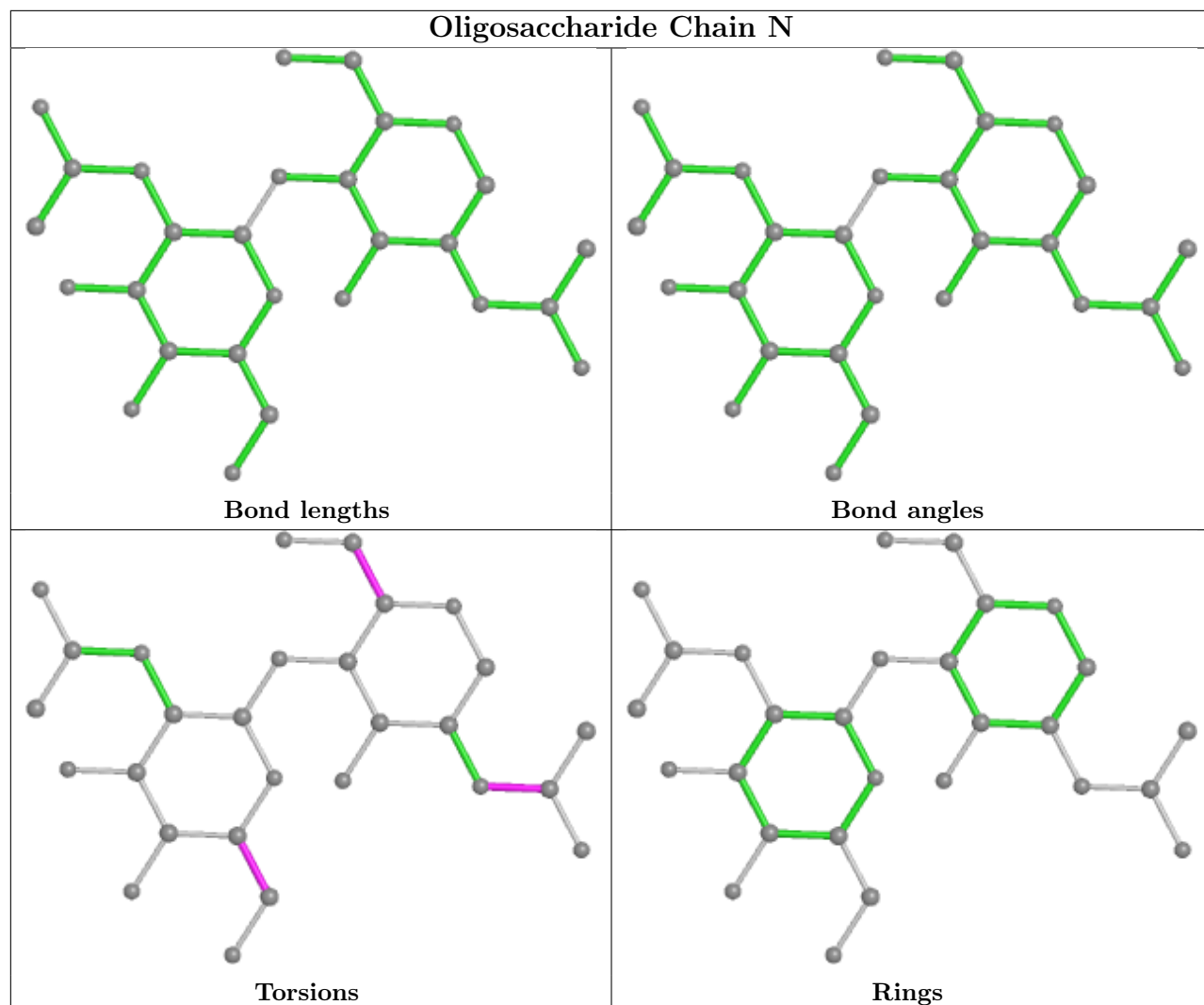


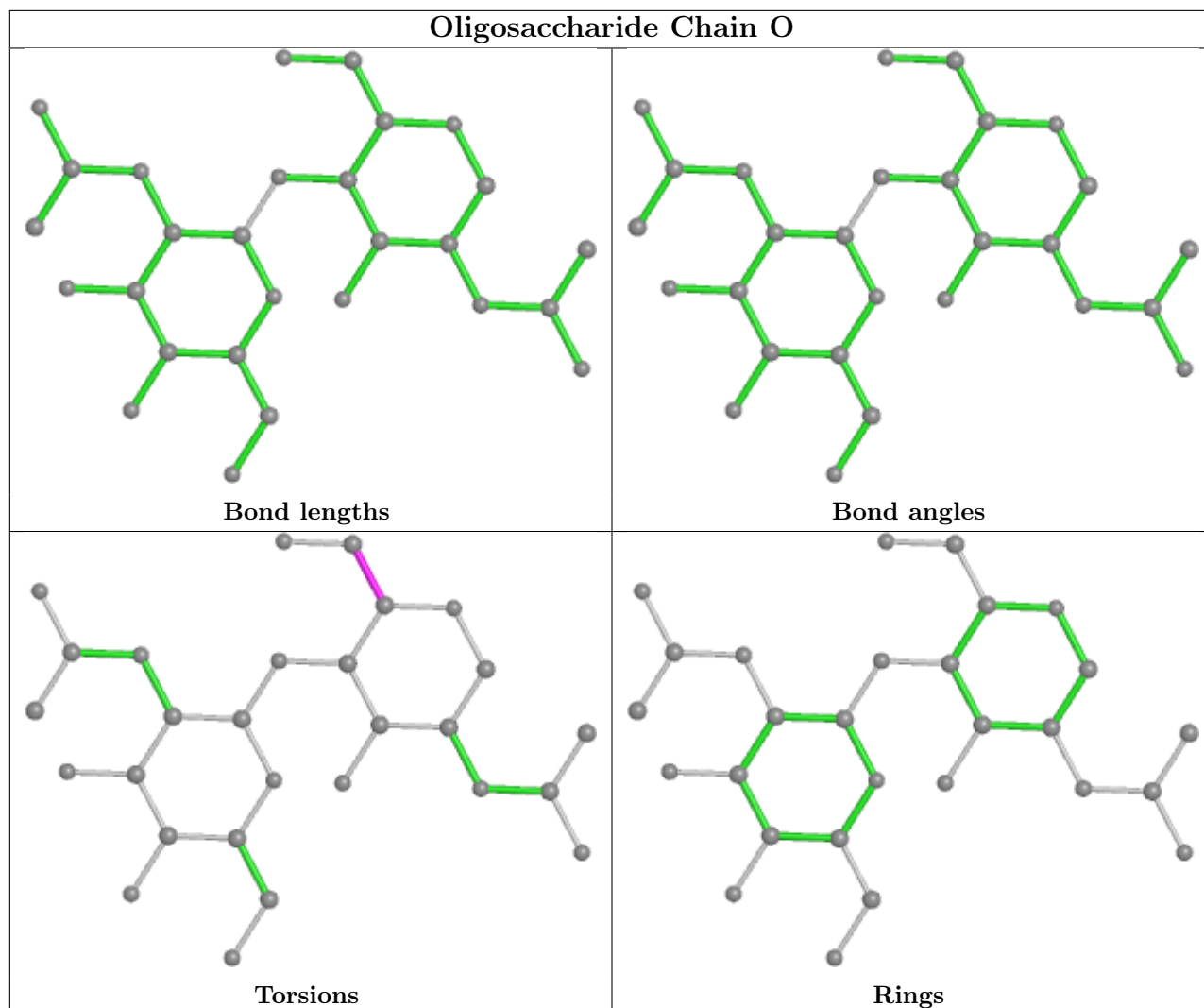












5.6 Ligand geometry [i](#)

23 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	B	1306	1	14,14,15	0.67	0	17,19,21	1.96	2 (11%)
3	NAG	C	1301	1	14,14,15	0.32	0	17,19,21	0.42	0
3	NAG	B	1303	1	14,14,15	0.39	0	17,19,21	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	1307	1	14,14,15	0.34	0	17,19,21	0.51	0
3	NAG	C	1306	1	14,14,15	0.72	1 (7%)	17,19,21	1.96	2 (11%)
3	NAG	B	1305	1	14,14,15	0.62	0	17,19,21	0.43	0
3	NAG	C	1304	1	14,14,15	0.30	0	17,19,21	0.46	0
3	NAG	C	1307	1	14,14,15	0.34	0	17,19,21	0.51	0
3	NAG	B	1302	1	14,14,15	1.26	2 (14%)	17,19,21	1.19	1 (5%)
3	NAG	C	1305	1	14,14,15	0.55	0	17,19,21	0.43	0
3	NAG	A	1304	1	14,14,15	0.27	0	17,19,21	0.43	0
3	NAG	B	1301	1	14,14,15	0.28	0	17,19,21	0.44	0
3	NAG	A	1307	1	14,14,15	0.29	0	17,19,21	0.51	0
3	NAG	A	1305	1	14,14,15	0.59	0	17,19,21	0.50	0
3	NAG	C	1302	1	14,14,15	1.31	2 (14%)	17,19,21	1.20	1 (5%)
3	NAG	A	1302	1	14,14,15	1.35	2 (14%)	17,19,21	1.20	1 (5%)
3	NAG	A	1303	1	14,14,15	0.39	0	17,19,21	0.36	0
3	NAG	A	1306	1	14,14,15	0.79	1 (7%)	17,19,21	1.97	2 (11%)
3	NAG	A	1301	1	14,14,15	0.30	0	17,19,21	0.47	0
3	NAG	B	1308	1	14,14,15	0.30	0	17,19,21	0.52	0
3	NAG	C	1303	1	14,14,15	0.38	0	17,19,21	0.39	0
3	NAG	C	1308	1	14,14,15	0.32	0	17,19,21	0.51	0
3	NAG	B	1304	1	14,14,15	0.26	0	17,19,21	0.42	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1306	1	-	5/6/23/26	0/1/1/1
3	NAG	C	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1303	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1307	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1306	1	-	3/6/23/26	0/1/1/1
3	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1307	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1302	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1305	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1304	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1301	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	1307	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1305	1	-	1/6/23/26	0/1/1/1
3	NAG	C	1302	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1302	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1303	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1306	1	-	3/6/23/26	0/1/1/1
3	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1308	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1308	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1304	1	-	2/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1302	NAG	O5-C1	4.19	1.50	1.43
3	C	1302	NAG	O5-C1	3.98	1.50	1.43
3	B	1302	NAG	O5-C1	3.81	1.49	1.43
3	A	1302	NAG	C1-C2	2.63	1.56	1.52
3	C	1302	NAG	C1-C2	2.62	1.56	1.52
3	B	1302	NAG	C1-C2	2.55	1.56	1.52
3	A	1306	NAG	C1-C2	2.42	1.55	1.52
3	C	1306	NAG	C1-C2	2.15	1.55	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1306	NAG	C2-N2-C7	6.86	132.68	122.90
3	A	1306	NAG	C2-N2-C7	6.84	132.64	122.90
3	C	1306	NAG	C2-N2-C7	6.84	132.64	122.90
3	C	1302	NAG	C1-O5-C5	4.66	118.50	112.19
3	A	1302	NAG	C1-O5-C5	4.65	118.50	112.19
3	B	1302	NAG	C1-O5-C5	4.64	118.48	112.19
3	B	1306	NAG	C1-C2-N2	3.12	115.82	110.49
3	C	1306	NAG	C1-C2-N2	3.11	115.80	110.49
3	A	1306	NAG	C1-C2-N2	3.02	115.65	110.49

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1304	NAG	C4-C5-C6-O6
3	B	1305	NAG	C4-C5-C6-O6
3	A	1304	NAG	O5-C5-C6-O6
3	C	1304	NAG	C4-C5-C6-O6
3	C	1305	NAG	O5-C5-C6-O6
3	A	1307	NAG	O5-C5-C6-O6
3	B	1305	NAG	O5-C5-C6-O6
3	B	1307	NAG	O5-C5-C6-O6
3	C	1304	NAG	O5-C5-C6-O6
3	B	1307	NAG	C4-C5-C6-O6
3	C	1303	NAG	C4-C5-C6-O6
3	C	1305	NAG	C4-C5-C6-O6
3	B	1306	NAG	C8-C7-N2-C2
3	B	1306	NAG	O7-C7-N2-C2
3	B	1308	NAG	C8-C7-N2-C2
3	B	1308	NAG	O7-C7-N2-C2
3	A	1306	NAG	C8-C7-N2-C2
3	A	1306	NAG	O7-C7-N2-C2
3	C	1306	NAG	C8-C7-N2-C2
3	C	1306	NAG	O7-C7-N2-C2
3	B	1306	NAG	O5-C5-C6-O6
3	C	1301	NAG	O5-C5-C6-O6
3	C	1308	NAG	O5-C5-C6-O6
3	B	1304	NAG	O5-C5-C6-O6
3	A	1307	NAG	C4-C5-C6-O6
3	A	1302	NAG	O5-C5-C6-O6
3	C	1303	NAG	O5-C5-C6-O6
3	A	1301	NAG	O5-C5-C6-O6
3	A	1302	NAG	C4-C5-C6-O6
3	B	1306	NAG	C4-C5-C6-O6
3	C	1308	NAG	C4-C5-C6-O6
3	B	1301	NAG	O5-C5-C6-O6
3	A	1305	NAG	O5-C5-C6-O6
3	C	1301	NAG	C4-C5-C6-O6
3	A	1306	NAG	C3-C2-N2-C7
3	C	1306	NAG	C3-C2-N2-C7
3	B	1304	NAG	C4-C5-C6-O6
3	A	1301	NAG	C4-C5-C6-O6
3	B	1306	NAG	C3-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

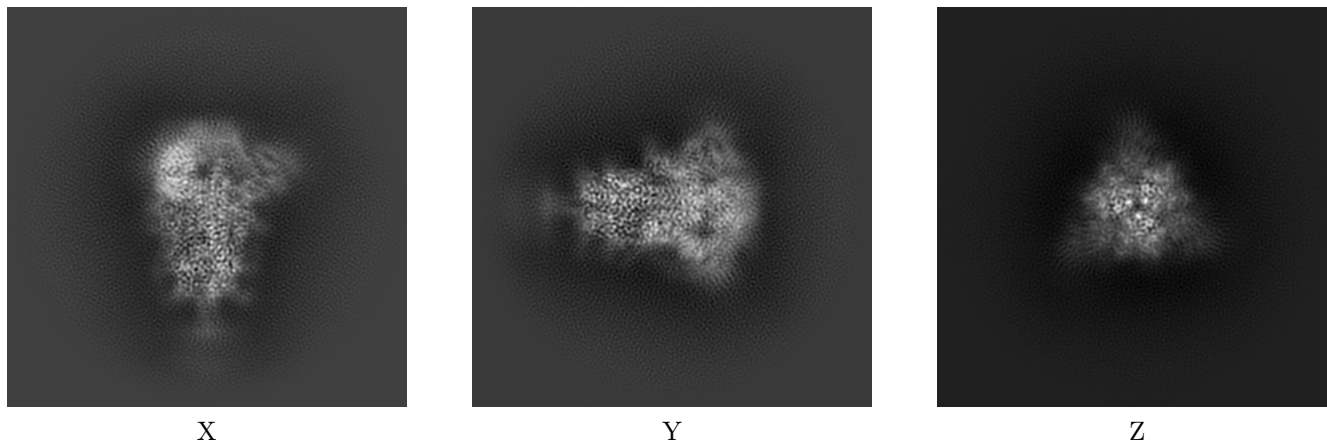
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33212. These allow visual inspection of the internal detail of the map and identification of artifacts.

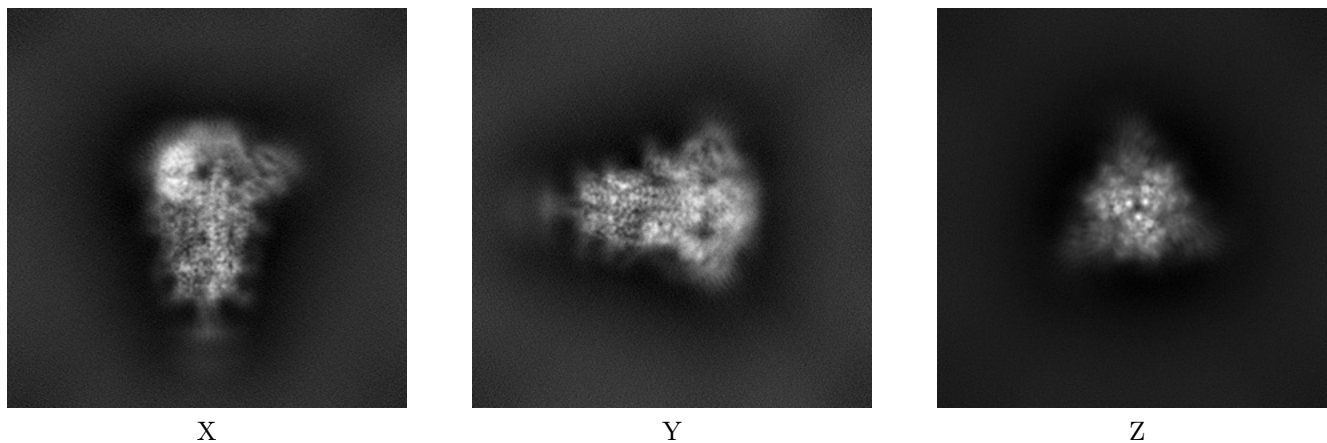
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



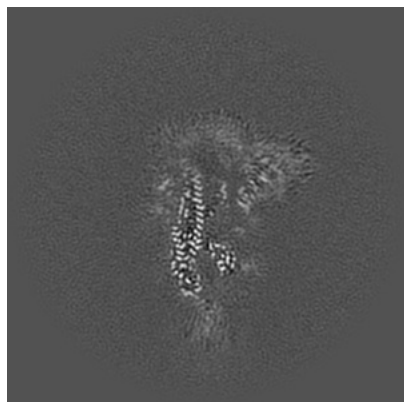
6.1.2 Raw map



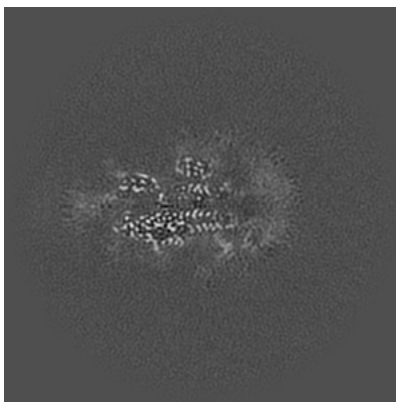
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

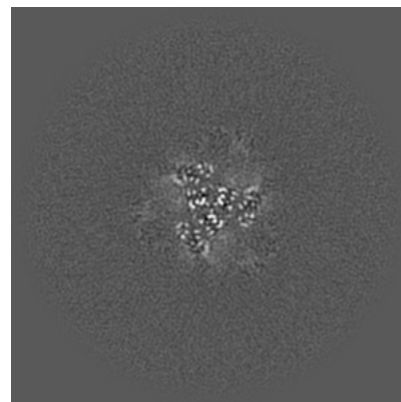
6.2.1 Primary map



X Index: 160

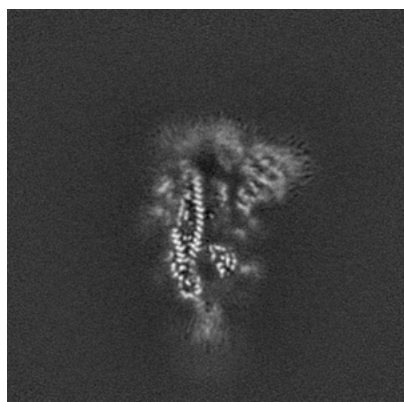


Y Index: 160

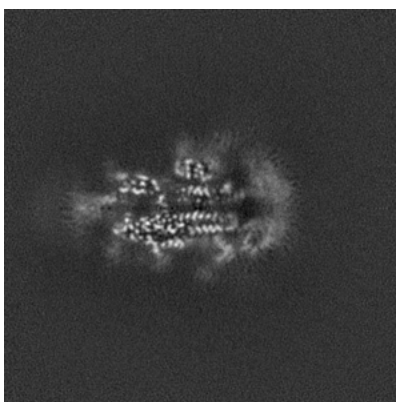


Z Index: 160

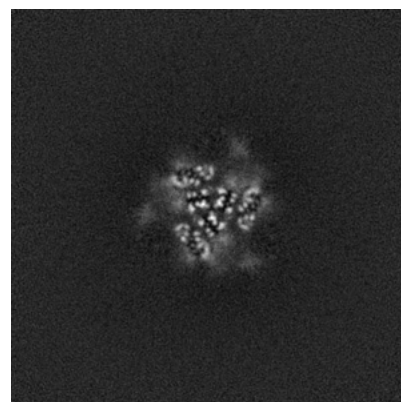
6.2.2 Raw map



X Index: 160



Y Index: 160

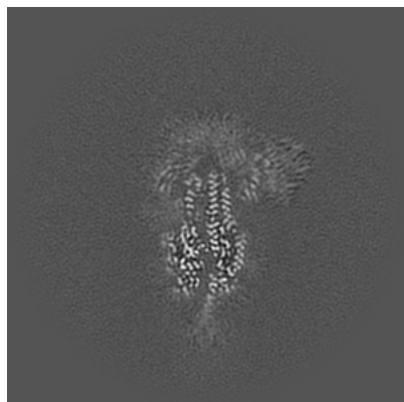


Z Index: 160

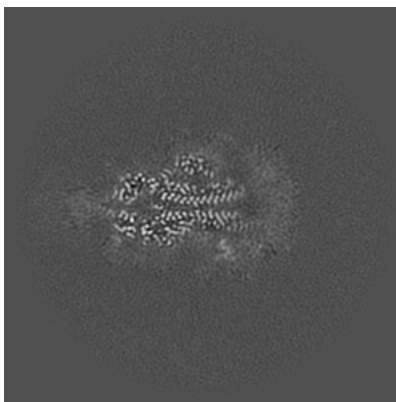
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

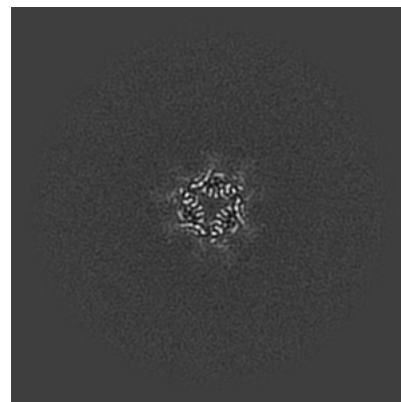
6.3.1 Primary map



X Index: 167

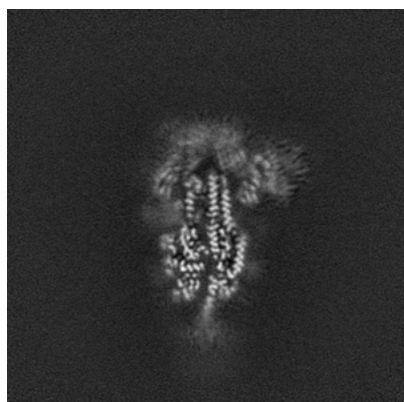


Y Index: 164

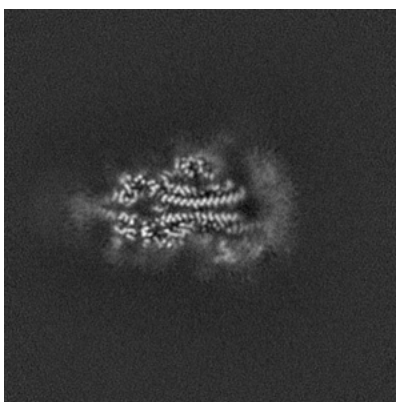


Z Index: 111

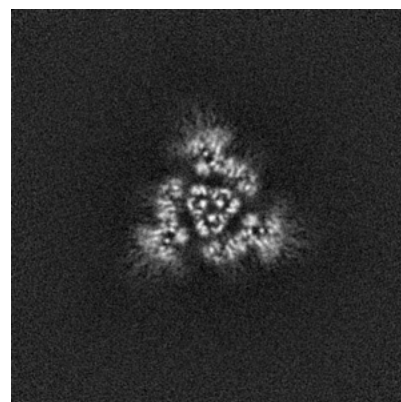
6.3.2 Raw map



X Index: 167



Y Index: 164

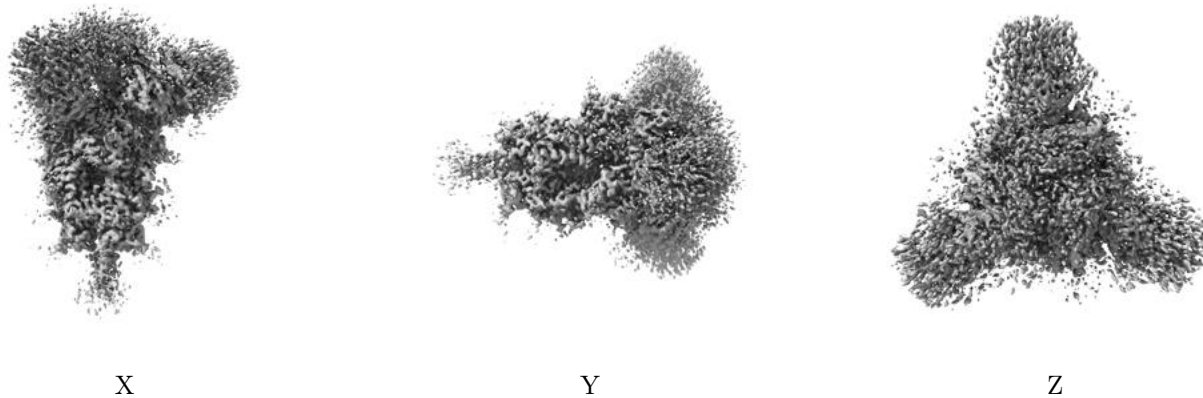


Z Index: 179

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

6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.4. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 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.

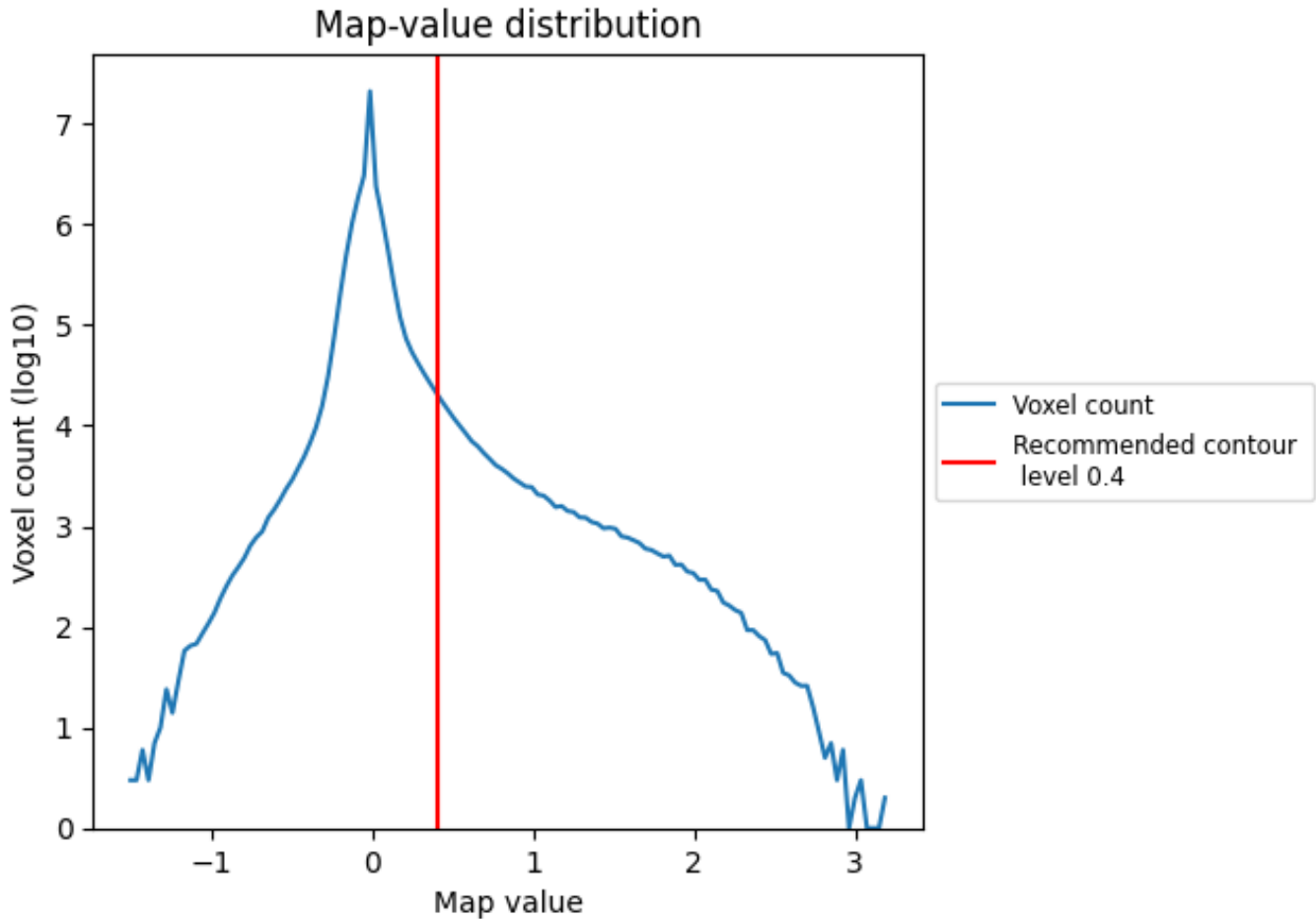
6.5 Mask visualisation [i](#)

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

7 Map analysis [i](#)

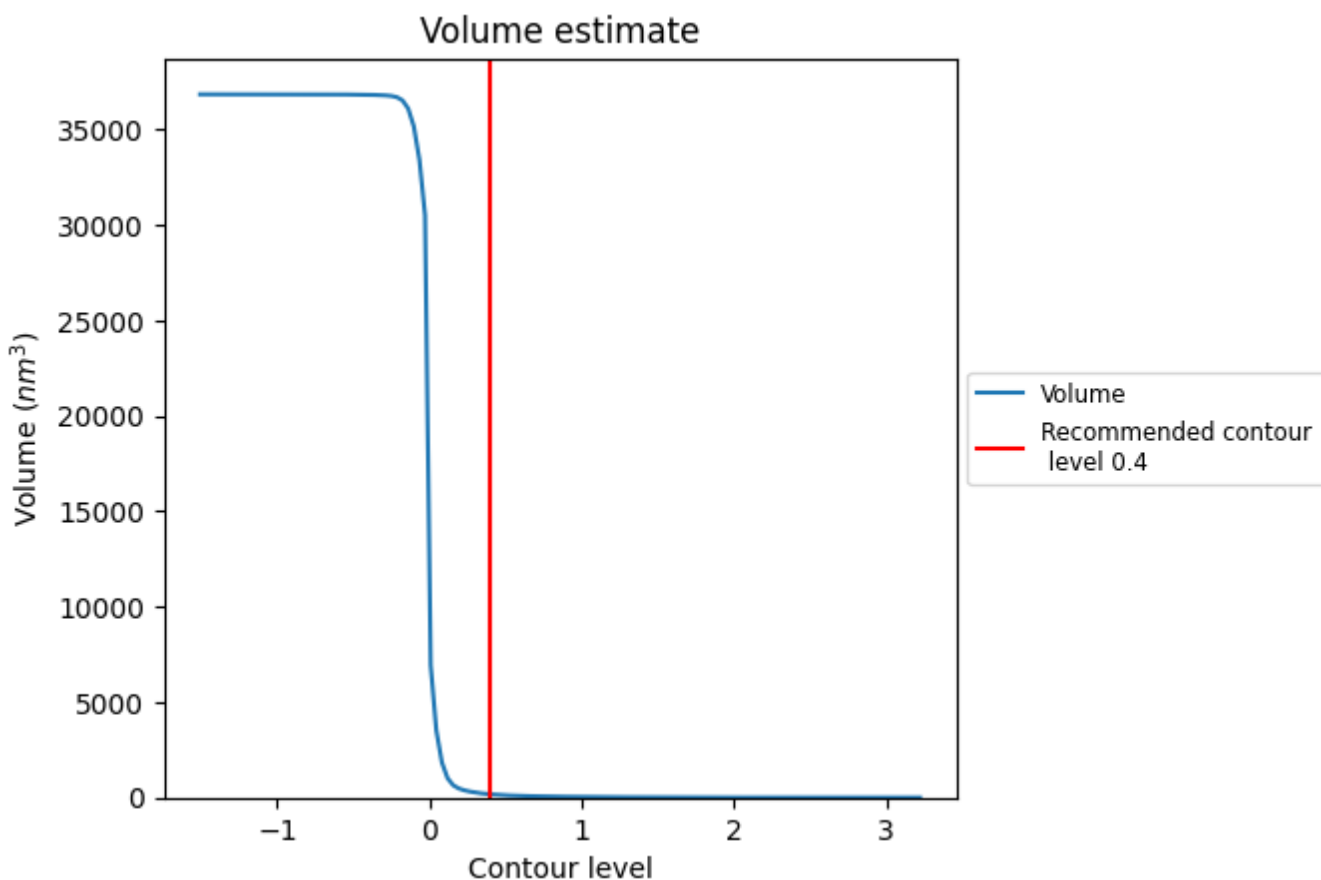
This section contains the results of statistical analysis of the map.

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.

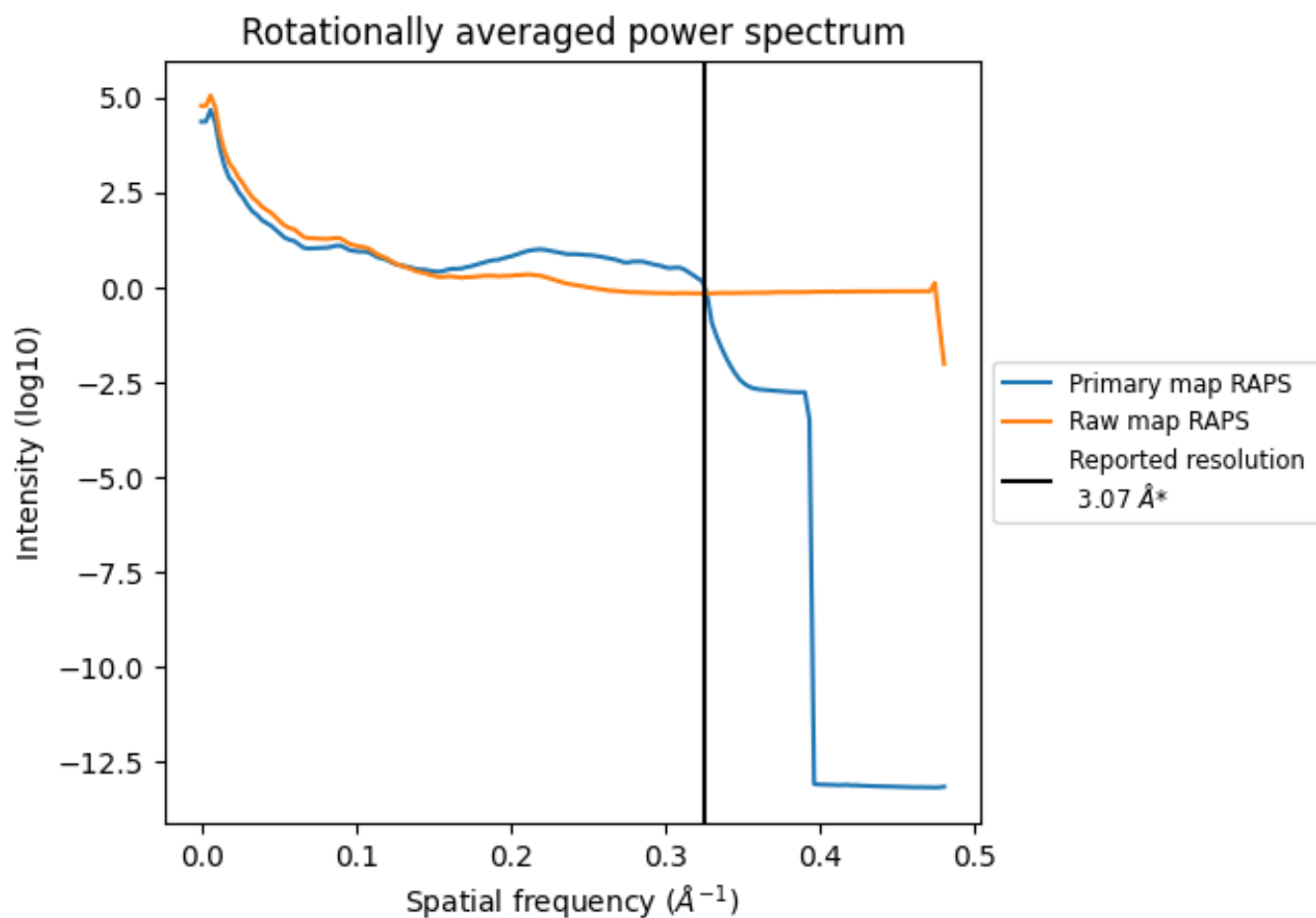
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 174 nm^3 ; this corresponds to an approximate mass of 157 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 i

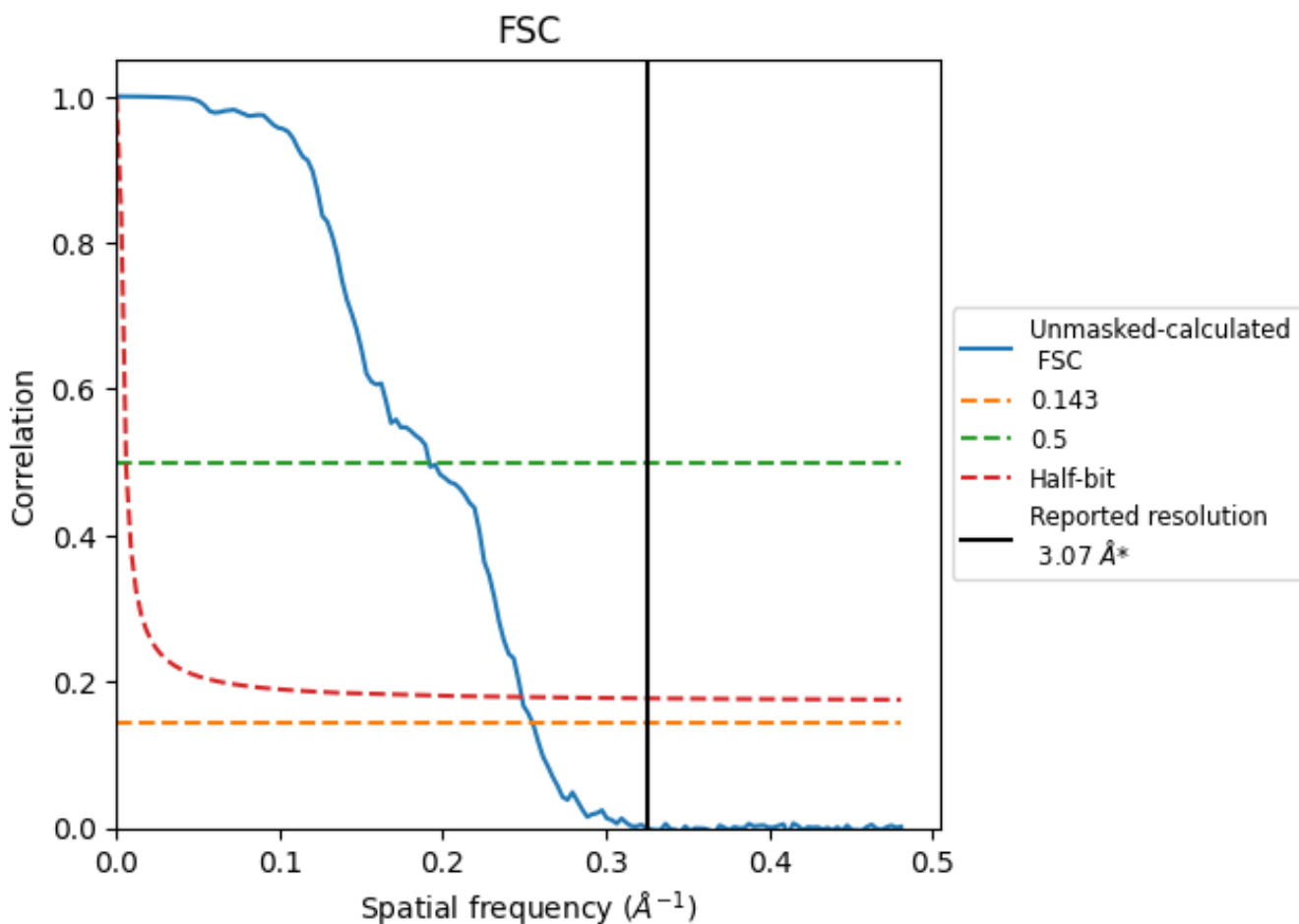


*Reported resolution corresponds to spatial frequency of 0.326 Å⁻¹

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 0.326 \AA^{-1}

8.2 Resolution estimates [i](#)

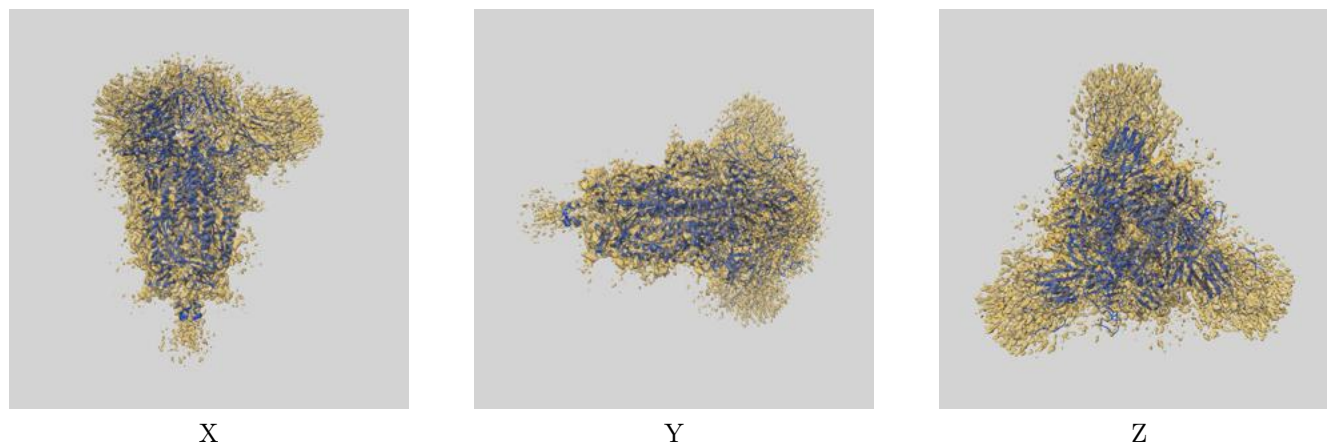
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.07	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.92	5.22	4.03

*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 3.92 differs from the reported value 3.07 by more than 10 %

9 Map-model fit [i](#)

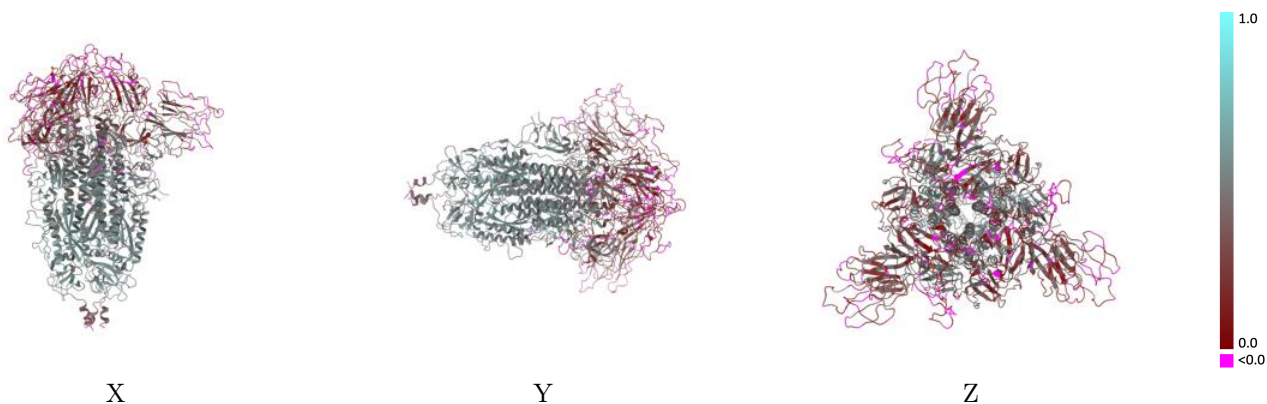
This section contains information regarding the fit between EMDB map EMD-33212 and PDB model 7XIY. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay [i](#)



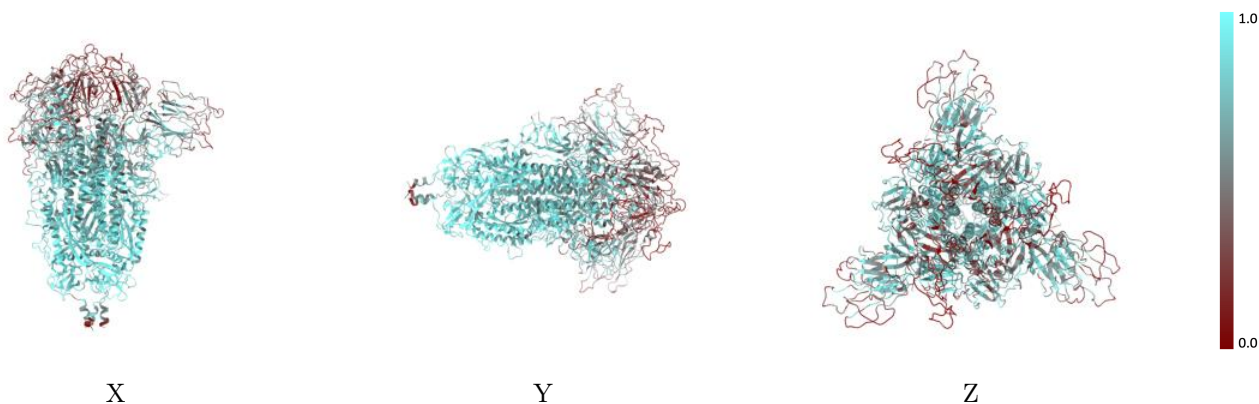
The images above show the 3D surface view of the map at the recommended contour level 0.4 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



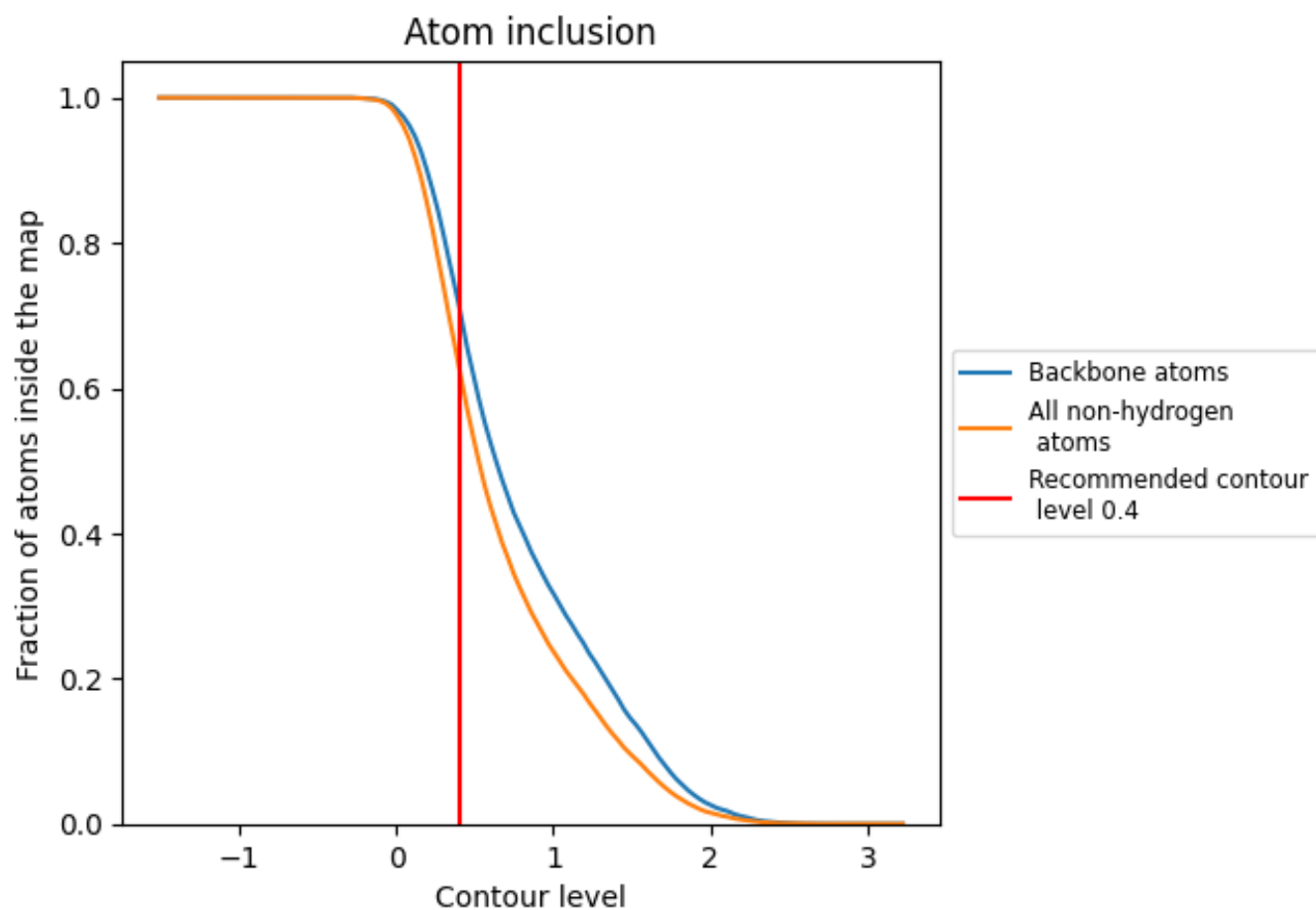
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.4).

































9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 63% 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 (0.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6300	 0.3590
A	 0.6273	 0.3570
B	 0.6315	 0.3590
C	 0.6304	 0.3610
D	 0.7143	 0.4480
E	 0.6786	 0.4470
F	 0.6071	 0.2520
G	 0.6786	 0.3390
H	 0.7500	 0.4420
I	 0.6071	 0.3830
J	 0.6071	 0.2600
K	 0.6429	 0.3330
L	 0.7143	 0.4230
M	 0.5714	 0.3510
N	 0.6071	 0.2340
O	 0.6786	 0.3570

