



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 11:20 AM UTC

PDB ID : 7WPA / pdb_00007wpa
EMDB ID : EMD-32680
Title : SARS-CoV-2 Omicron Variant SPIKE trimer complexed with ACE2
Authors : Yin, W.; Xu, Y.; Xu, P.; Cao, X.; Wu, C.; Gu, C.; He, X.; Wang, X.; Huang, S.; Yuan, Q.; Wu, K.; Hu, W.; Huang, Z.; Liu, J.; Wang, Z.; Jia, F.; Xia, K.; Liu, P.; Wang, X.; Song, B.; Zheng, J.; Jiang, H.; Cheng, X.; Jiang, Y.; Deng, S.; Xu, E.
Deposited on : 2022-01-23
Resolution : 2.77 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

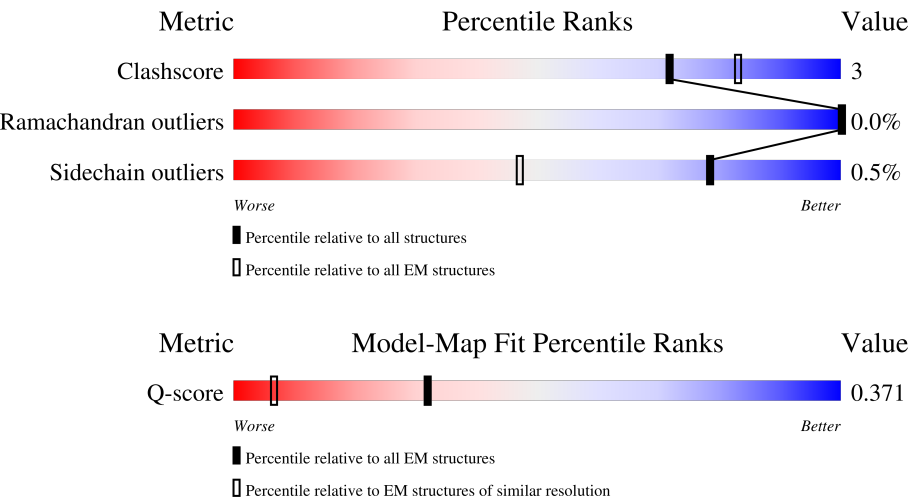
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.77 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	10695 (2.27 - 3.27)

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	1205	6% 80% 8% 12%
1	B	1205	5% 80% 8% 12%
1	C	1205	10% 80% 8% 12%
2	D	805	43% 64% 10% 26%

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Mol	Chain	Length	Quality of chain
3	E	2	 50% 50%
3	F	2	 100%
3	G	2	 50% 50%
3	H	2	 50% 50%
3	I	2	 50% 100%
3	J	2	 50% 100%
3	K	2	 50% 50%
3	L	2	 50% 100%
3	M	2	 50% 50%
3	N	2	 50% 50%
3	O	2	 50% 50%
3	P	2	 100%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 30444 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
1	A	1061	Total	C	N	O	S	0	0
			8329	5333	1385	1573	38		
1	B	1061	Total	C	N	O	S	0	0
			8329	5333	1385	1573	38		
1	C	1057	Total	C	N	O	S	0	0
			8297	5312	1381	1567	37		

There are 135 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	VAL	ALA	variant	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	93	ILE	THR	variant	UNP P0DTC2
A	140	ASP	GLY	variant	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	206	ILE	-	insertion	UNP P0DTC2
A	207	VAL	-	insertion	UNP P0DTC2
A	208	ARG	ASN	variant	UNP P0DTC2
A	209	GLU	LEU	variant	UNP P0DTC2
A	210	PRO	VAL	variant	UNP P0DTC2
A	211	GLU	ARG	variant	UNP P0DTC2
A	336	ASP	GLY	variant	UNP P0DTC2
A	368	LEU	SER	variant	UNP P0DTC2
A	370	PRO	SER	variant	UNP P0DTC2
A	372	PHE	SER	variant	UNP P0DTC2
A	414	ASN	LYS	variant	UNP P0DTC2
A	437	LYS	ASN	variant	UNP P0DTC2
A	443	SER	GLY	variant	UNP P0DTC2
A	474	ASN	SER	variant	UNP P0DTC2
A	475	LYS	THR	variant	UNP P0DTC2
A	481	ALA	GLU	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	490	ARG	GLN	variant	UNP P0DTC2
A	493	SER	GLY	variant	UNP P0DTC2
A	495	ARG	GLN	variant	UNP P0DTC2
A	498	TYR	ASN	variant	UNP P0DTC2
A	502	HIS	TYR	variant	UNP P0DTC2
A	544	LYS	THR	variant	UNP P0DTC2
A	611	GLY	ASP	variant	UNP P0DTC2
A	652	TYR	HIS	variant	UNP P0DTC2
A	676	LYS	ASN	variant	UNP P0DTC2
A	678	HIS	PRO	variant	UNP P0DTC2
A	679	GLY	ARG	engineered mutation	UNP P0DTC2
A	680	SER	ARG	engineered mutation	UNP P0DTC2
A	682	SER	ARG	engineered mutation	UNP P0DTC2
A	761	LYS	ASN	variant	UNP P0DTC2
A	793	TYR	ASP	variant	UNP P0DTC2
A	853	LYS	ASN	variant	UNP P0DTC2
A	951	HIS	GLN	variant	UNP P0DTC2
A	966	LYS	ASN	variant	UNP P0DTC2
A	978	PHE	LEU	variant	UNP P0DTC2
A	983	PRO	LYS	engineered mutation	UNP P0DTC2
A	984	PRO	VAL	engineered mutation	UNP P0DTC2
B	67	VAL	ALA	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	93	ILE	THR	variant	UNP P0DTC2
B	140	ASP	GLY	variant	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	206	ILE	-	insertion	UNP P0DTC2
B	207	VAL	-	insertion	UNP P0DTC2
B	208	ARG	ASN	variant	UNP P0DTC2
B	209	GLU	LEU	variant	UNP P0DTC2
B	210	PRO	VAL	variant	UNP P0DTC2
B	211	GLU	ARG	variant	UNP P0DTC2
B	336	ASP	GLY	variant	UNP P0DTC2
B	368	LEU	SER	variant	UNP P0DTC2
B	370	PRO	SER	variant	UNP P0DTC2
B	372	PHE	SER	variant	UNP P0DTC2
B	414	ASN	LYS	variant	UNP P0DTC2
B	437	LYS	ASN	variant	UNP P0DTC2
B	443	SER	GLY	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	474	ASN	SER	variant	UNP P0DTC2
B	475	LYS	THR	variant	UNP P0DTC2
B	481	ALA	GLU	variant	UNP P0DTC2
B	490	ARG	GLN	variant	UNP P0DTC2
B	493	SER	GLY	variant	UNP P0DTC2
B	495	ARG	GLN	variant	UNP P0DTC2
B	498	TYR	ASN	variant	UNP P0DTC2
B	502	HIS	TYR	variant	UNP P0DTC2
B	544	LYS	THR	variant	UNP P0DTC2
B	611	GLY	ASP	variant	UNP P0DTC2
B	652	TYR	HIS	variant	UNP P0DTC2
B	676	LYS	ASN	variant	UNP P0DTC2
B	678	HIS	PRO	variant	UNP P0DTC2
B	679	GLY	ARG	engineered mutation	UNP P0DTC2
B	680	SER	ARG	engineered mutation	UNP P0DTC2
B	682	SER	ARG	engineered mutation	UNP P0DTC2
B	761	LYS	ASN	variant	UNP P0DTC2
B	793	TYR	ASP	variant	UNP P0DTC2
B	853	LYS	ASN	variant	UNP P0DTC2
B	951	HIS	GLN	variant	UNP P0DTC2
B	966	LYS	ASN	variant	UNP P0DTC2
B	978	PHE	LEU	variant	UNP P0DTC2
B	983	PRO	LYS	engineered mutation	UNP P0DTC2
B	984	PRO	VAL	engineered mutation	UNP P0DTC2
C	67	VAL	ALA	variant	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	93	ILE	THR	variant	UNP P0DTC2
C	140	ASP	GLY	variant	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	206	ILE	-	insertion	UNP P0DTC2
C	207	VAL	-	insertion	UNP P0DTC2
C	208	ARG	ASN	variant	UNP P0DTC2
C	209	GLU	LEU	variant	UNP P0DTC2
C	210	PRO	VAL	variant	UNP P0DTC2
C	211	GLU	ARG	variant	UNP P0DTC2
C	336	ASP	GLY	variant	UNP P0DTC2
C	368	LEU	SER	variant	UNP P0DTC2
C	370	PRO	SER	variant	UNP P0DTC2
C	372	PHE	SER	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	414	ASN	LYS	variant	UNP P0DTC2
C	437	LYS	ASN	variant	UNP P0DTC2
C	443	SER	GLY	variant	UNP P0DTC2
C	474	ASN	SER	variant	UNP P0DTC2
C	475	LYS	THR	variant	UNP P0DTC2
C	481	ALA	GLU	variant	UNP P0DTC2
C	490	ARG	GLN	variant	UNP P0DTC2
C	493	SER	GLY	variant	UNP P0DTC2
C	495	ARG	GLN	variant	UNP P0DTC2
C	498	TYR	ASN	variant	UNP P0DTC2
C	502	HIS	TYR	variant	UNP P0DTC2
C	544	LYS	THR	variant	UNP P0DTC2
C	611	GLY	ASP	variant	UNP P0DTC2
C	652	TYR	HIS	variant	UNP P0DTC2
C	676	LYS	ASN	variant	UNP P0DTC2
C	678	HIS	PRO	variant	UNP P0DTC2
C	679	GLY	ARG	engineered mutation	UNP P0DTC2
C	680	SER	ARG	engineered mutation	UNP P0DTC2
C	682	SER	ARG	engineered mutation	UNP P0DTC2
C	761	LYS	ASN	variant	UNP P0DTC2
C	793	TYR	ASP	variant	UNP P0DTC2
C	853	LYS	ASN	variant	UNP P0DTC2
C	951	HIS	GLN	variant	UNP P0DTC2
C	966	LYS	ASN	variant	UNP P0DTC2
C	978	PHE	LEU	variant	UNP P0DTC2
C	983	PRO	LYS	engineered mutation	UNP P0DTC2
C	984	PRO	VAL	engineered mutation	UNP P0DTC2

- Molecule 2 is a protein called Angiotensin-converting enzyme 2.

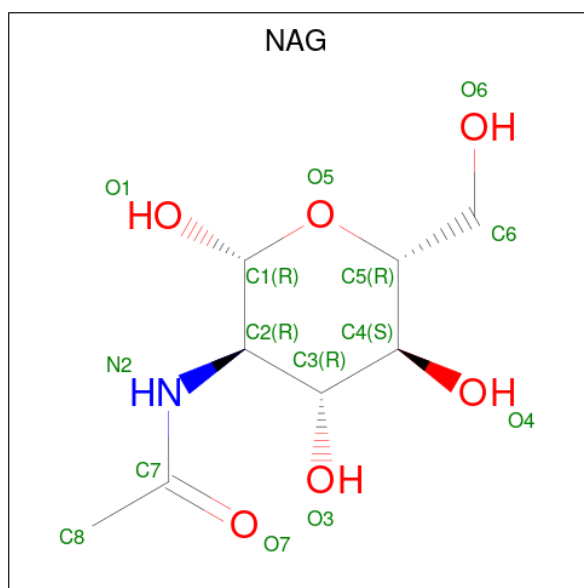
Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	595	Total	C	N	O	S	0	0
			4857	3108	804	916	29		

- Molecule 3 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
3	E	2	Total	C	N	O	0	0
			28	16	2	10		
3	F	2	Total	C	N	O	0	0
			28	16	2	10		
3	G	2	Total	C	N	O	0	0
			28	16	2	10		
3	H	2	Total	C	N	O	0	0
			28	16	2	10		
3	I	2	Total	C	N	O	0	0
			28	16	2	10		
3	J	2	Total	C	N	O	0	0
			28	16	2	10		
3	K	2	Total	C	N	O	0	0
			28	16	2	10		
3	L	2	Total	C	N	O	0	0
			28	16	2	10		
3	M	2	Total	C	N	O	0	0
			28	16	2	10		
3	N	2	Total	C	N	O	0	0
			28	16	2	10		
3	O	2	Total	C	N	O	0	0
			28	16	2	10		
3	P	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest")

by depositor).

Mol	Chain	Residues	Atoms		AltConf
5	D	1	Total	Cl	0
			1	1	

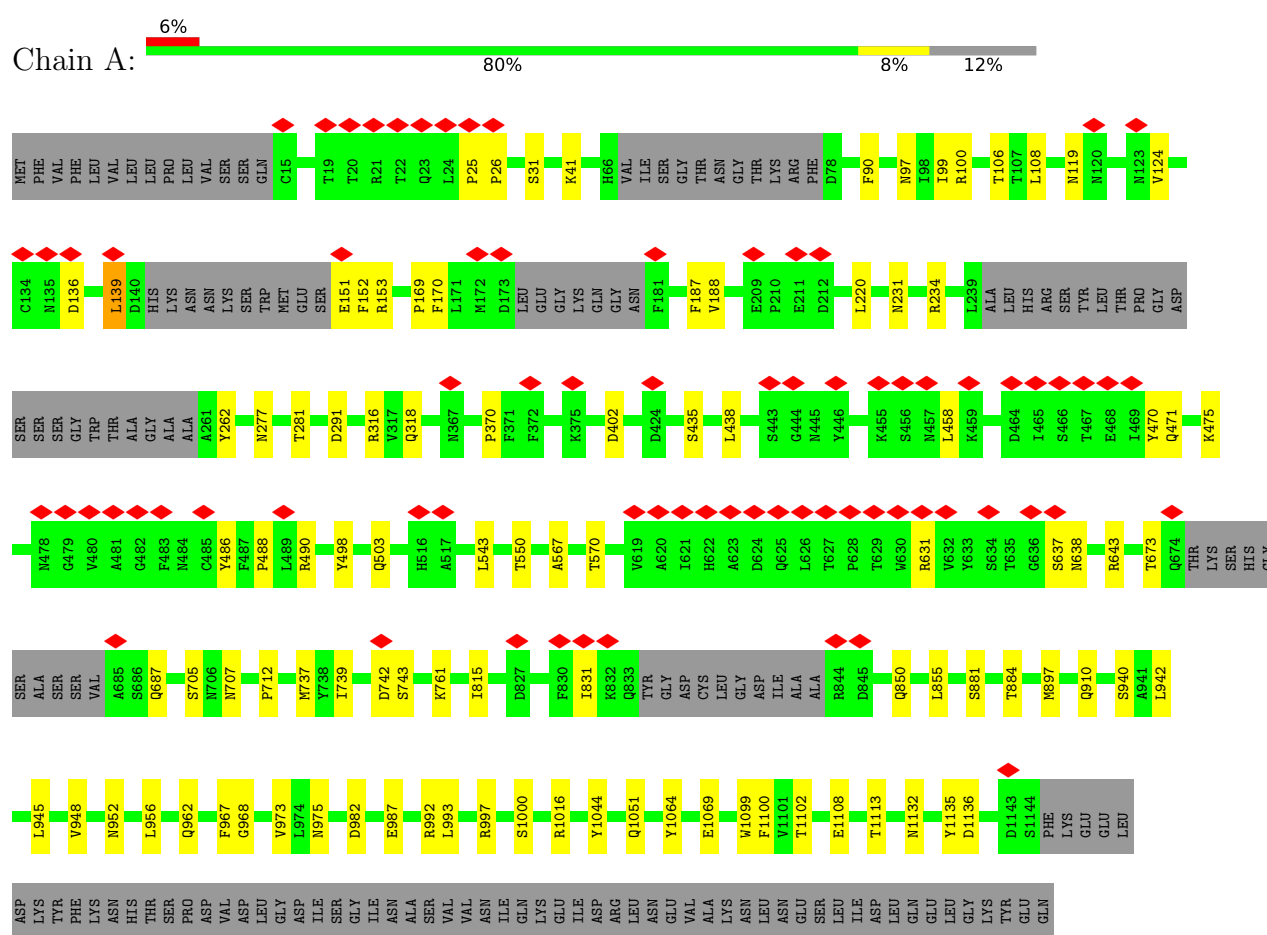
- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
6	D	1	Total	Zn	0
			1	1	

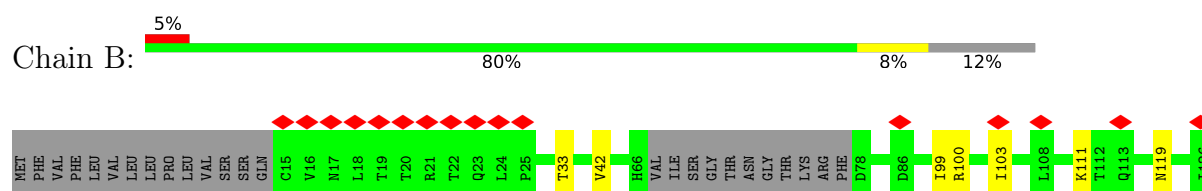
3 Residue-property plots

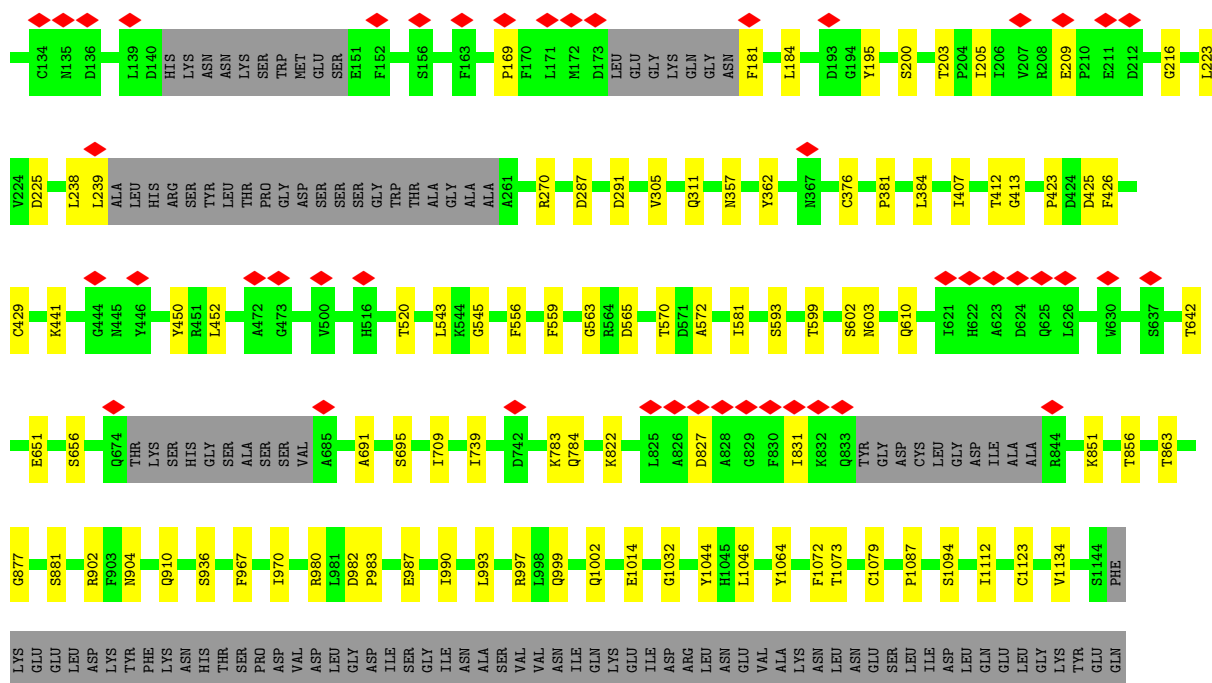
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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

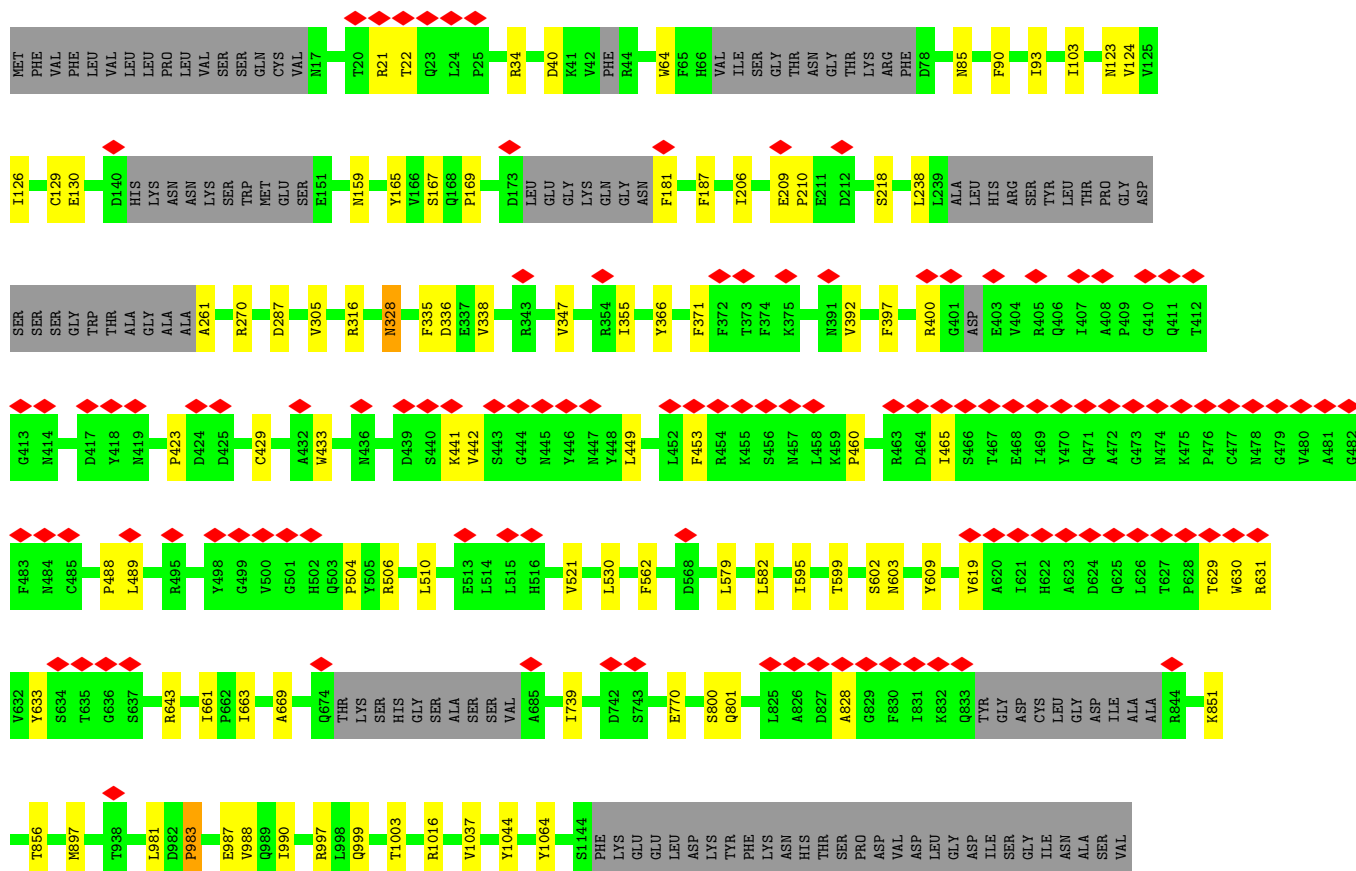
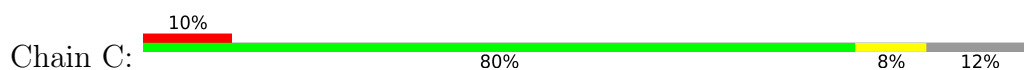


• Molecule 1: Spike glycoprotein



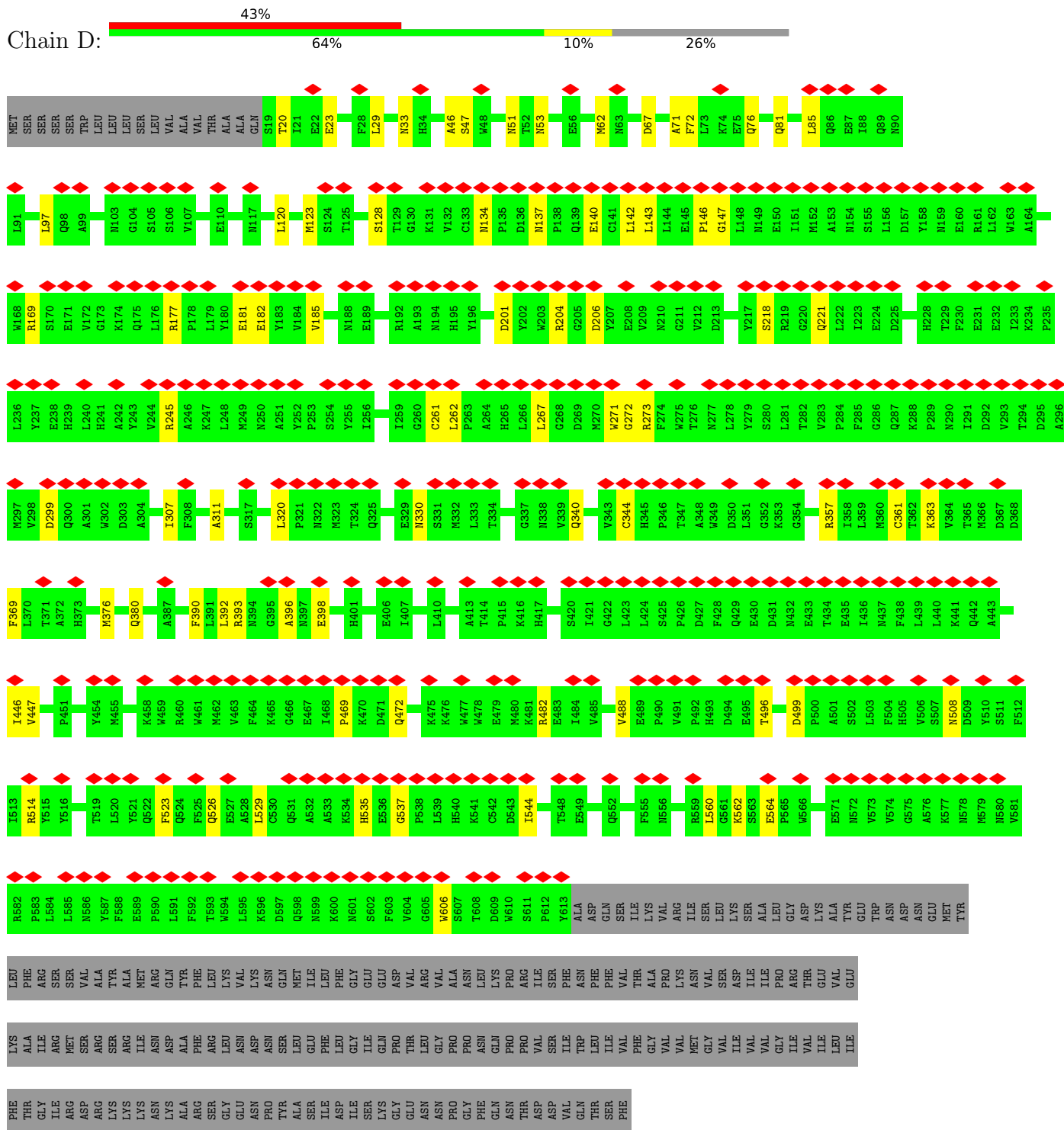


• Molecule 1: Spike glycoprotein



VAL ASN SER SER SER SER TRP LEU LEU ASP ARG ASN VAL VAL ALA ASN ASN ASN GLU SER SER LEU LEU ASP ASP LEU GLN GLU LEU GLY TYR TYR GLU GLN

• Molecule 2: Angiotensin-converting enzyme 2



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  50% 50%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100% 100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  50% 50% 50%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  50% 50%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  50% 100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  50% 100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	242374	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 ($e^-/\text{\AA}^2$)	1.389	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	5000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.752	Depositor
Minimum map value	-0.022	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.05	Depositor
Map size ($\text{\AA}$)	395.52, 395.52, 395.52	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.824, 0.824, 0.824	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.20	0/8524	0.51	6/11599 (0.1%)
1	B	0.19	0/8524	0.47	1/11599 (0.0%)
1	C	0.19	0/8489	0.48	3/11548 (0.0%)
2	D	0.19	0/4994	0.50	4/6785 (0.1%)
All	All	0.19	0/30531	0.49	14/41531 (0.0%)

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

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	983	PRO	N-CA-C	9.92	122.80	110.70
1	A	458	LEU	CA-C-N	7.45	143.58	122.21
1	A	458	LEU	C-N-CA	7.45	143.58	122.21
1	C	619	VAL	N-CA-C	-6.80	104.70	111.77
1	B	983	PRO	N-CA-C	6.79	118.98	110.70
1	A	31	SER	CA-C-N	5.96	129.08	120.79
1	A	31	SER	C-N-CA	5.96	129.08	120.79
2	D	128	SER	CA-C-N	5.94	132.89	121.54
2	D	128	SER	C-N-CA	5.94	132.89	121.54
1	C	206	ILE	N-CA-C	-5.87	107.54	113.47
1	A	567	ALA	CA-C-N	5.58	130.33	122.46
1	A	567	ALA	C-N-CA	5.58	130.33	122.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	508	ASN	CA-C-N	5.28	129.90	122.46
2	D	508	ASN	C-N-CA	5.28	129.90	122.46

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	370	PRO	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8329	0	8160	56	0
1	B	8329	0	8166	60	0
1	C	8297	0	8139	55	0
2	D	4857	0	4637	45	0
3	E	28	0	25	0	0
3	F	28	0	25	0	0
3	G	28	0	25	0	0
3	H	28	0	25	1	0
3	I	28	0	25	0	0
3	J	28	0	25	0	0
3	K	28	0	25	0	0
3	L	28	0	25	0	0
3	M	28	0	25	1	0
3	N	28	0	25	0	0
3	O	28	0	25	0	0
3	P	28	0	25	0	0
4	A	84	0	78	1	0
4	B	84	0	78	1	0
4	C	84	0	78	1	0
4	D	42	0	39	0	0
5	D	1	0	0	0	0
6	D	1	0	0	0	0
All	All	30444	0	29675	203	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 3.

All (203) 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:1044:TYR:HB2	1:B:1064:TYR:HB3	1.83	0.60
2:D:201:ASP:HA	2:D:204:ARG:HB2	1.84	0.60
1:C:21:ARG:HG3	1:C:22:THR:HG23	1.85	0.59
2:D:446:ILE:HG13	2:D:447:VAL:HG23	1.84	0.59
2:D:523:PHE:HA	2:D:526:GLN:HB2	1.83	0.58
1:C:1044:TYR:HB2	1:C:1064:TYR:HB3	1.85	0.58
2:D:344:CYS:HB3	2:D:361:CYS:HB3	1.87	0.57
1:C:347:VAL:HG12	1:C:397:PHE:HB2	1.86	0.57
1:A:1044:TYR:HB2	1:A:1064:TYR:HB3	1.88	0.55
1:B:99:ILE:HA	1:B:239:LEU:HA	1.87	0.55
1:B:987:GLU:HA	1:B:990:ILE:HG12	1.88	0.55
2:D:469:PRO:HD2	2:D:472:GLN:HE21	1.72	0.55
1:A:973:VAL:HG12	1:A:975:ASN:H	1.72	0.55
1:B:1002:GLN:NE2	1:C:1003:THR:OG1	2.38	0.54
2:D:120:LEU:HA	2:D:123:MET:HG2	1.89	0.54
1:B:100:ARG:O	1:B:119:ASN:ND2	2.41	0.54
1:B:181:PHE:HB2	1:B:209:GLU:HG2	1.89	0.54
2:D:311:ALA:HB1	2:D:376:MET:HG3	1.90	0.54
1:A:100:ARG:HB3	1:A:139:LEU:HD23	1.90	0.53
1:A:318:GLN:HE21	1:A:631:ARG:HH12	1.53	0.53
2:D:72:PHE:O	2:D:76:GLN:HB2	2.08	0.53
1:A:90:PHE:HB3	1:A:187:PHE:HB2	1.90	0.53
1:C:602:SER:OG	1:C:603:ASN:N	2.42	0.53
1:A:1113:THR:HG22	1:A:1135:TYR:HB3	1.90	0.53
1:A:712:PRO:HA	1:A:1069:GLU:HA	1.91	0.53
1:C:124:VAL:HB	1:C:169:PRO:HA	1.90	0.53
1:C:129:CYS:SG	1:C:130:GLU:N	2.81	0.53
1:B:863:THR:HG22	1:C:643:ARG:HH12	1.74	0.53
1:C:770:GLU:OE1	1:C:1016:ARG:NH1	2.42	0.52
1:A:815:ILE:HB	1:A:1051:GLN:HE21	1.74	0.52
1:B:904:ASN:HD21	1:B:910:GLN:HE21	1.57	0.52
1:B:1073:THR:HB	1:B:1094:SER:HB3	1.91	0.52
1:C:739:ILE:O	1:C:997:ARG:NH2	2.42	0.52
1:C:800:SER:OG	1:C:801:GLN:NE2	2.43	0.52
2:D:29:LEU:O	2:D:33:ASN:ND2	2.42	0.52
1:A:470:TYR:H	1:A:488:PRO:HD3	1.73	0.52
2:D:53:ASN:OD1	2:D:340:GLN:NE2	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:355:ILE:HB	1:C:392:VAL:HB	1.90	0.51
2:D:177:ARG:NH1	2:D:181:GLU:OE2	2.43	0.51
2:D:261:CYS:SG	2:D:262:LEU:N	2.80	0.51
1:B:200:SER:HB3	1:B:223:LEU:HD22	1.92	0.51
1:A:673:THR:HG22	1:A:687:GLN:HG2	1.93	0.51
1:A:855:LEU:HD13	1:A:956:LEU:HD22	1.93	0.51
1:B:357:ASN:H	1:B:520:THR:HB	1.76	0.51
2:D:20:THR:HG23	2:D:23:GLU:H	1.76	0.50
1:B:783:LYS:HG3	1:B:784:GLN:HG3	1.92	0.50
1:C:209:GLU:HG2	1:C:210:PRO:HD3	1.94	0.50
1:A:471:GLN:NE2	1:A:475:LYS:O	2.45	0.50
1:A:942:LEU:HD23	1:A:945:LEU:HD12	1.93	0.50
1:B:225:ASP:N	1:B:225:ASP:OD1	2.44	0.50
1:A:543:LEU:HD21	1:A:570:THR:HG21	1.93	0.50
1:B:169:PRO:HB2	1:B:223:LEU:HD23	1.93	0.50
1:B:423:PRO:HD2	1:B:426:PHE:HB2	1.94	0.49
1:B:827:ASP:OD1	1:B:827:ASP:N	2.44	0.49
1:A:316:ARG:O	1:A:631:ARG:NH1	2.45	0.49
1:A:962:GLN:NE2	1:A:1000:SER:OG	2.44	0.49
1:B:593:SER:OG	1:B:610:GLN:NE2	2.44	0.49
1:A:910:GLN:NE2	1:B:1087:PRO:O	2.45	0.49
1:C:123:ASN:ND2	1:C:167:SER:O	2.45	0.49
1:B:563:GLY:HA3	1:B:572:ALA:HB3	1.94	0.49
1:A:739:ILE:O	1:A:997:ARG:NH2	2.44	0.49
1:C:126:ILE:HB	1:C:165:TYR:HB3	1.93	0.49
1:C:130:GLU:OE2	1:C:159:ASN:ND2	2.46	0.49
1:C:328:ASN:OD1	1:C:328:ASN:N	2.44	0.49
1:A:1100:PHE:HZ	3:H:1:NAG:H62	1.78	0.49
1:B:1032:GLY:HA3	1:C:1037:VAL:HG21	1.94	0.49
1:A:136:ASP:OD1	1:A:136:ASP:N	2.46	0.48
2:D:535:HIS:HB3	2:D:537:GLY:H	1.77	0.48
1:C:181:PHE:HB2	1:C:209:GLU:HB3	1.94	0.48
2:D:143:LEU:HB2	2:D:146:PRO:HD2	1.95	0.48
1:A:486:TYR:HB3	1:A:490:ARG:HH12	1.77	0.48
1:A:119:ASN:HD21	1:A:170:PHE:HB2	1.78	0.48
1:A:968:GLY:O	1:A:992:ARG:NH2	2.47	0.48
1:A:761:LYS:NZ	1:B:311:GLN:O	2.45	0.48
1:A:435:SER:HB2	1:A:438:LEU:HB2	1.96	0.47
1:B:831:ILE:HB	4:C:1302:NAG:H82	1.96	0.47
2:D:134:ASN:ND2	2:D:137:ASN:OD1	2.46	0.47
1:A:910:GLN:HE22	1:B:1087:PRO:HG2	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1112:ILE:HG22	1:B:1134:VAL:HG13	1.97	0.47
2:D:47:SER:O	2:D:51:ASN:ND2	2.47	0.47
1:B:33:THR:OG1	1:B:216:GLY:O	2.33	0.47
1:C:90:PHE:O	1:C:187:PHE:N	2.47	0.47
1:C:336:ASP:OD1	1:C:336:ASP:N	2.47	0.47
1:C:453:PHE:H	1:C:488:PRO:HB3	1.80	0.47
2:D:46:ALA:HB1	2:D:62:MET:HA	1.95	0.47
2:D:177:ARG:HD3	2:D:496:THR:HA	1.97	0.47
1:A:291:ASP:OD1	1:A:291:ASP:N	2.47	0.47
1:B:291:ASP:N	1:B:291:ASP:OD1	2.47	0.47
2:D:142:LEU:HB2	2:D:147:GLY:HA3	1.96	0.47
1:B:184:LEU:HB2	1:B:205:ILE:HD11	1.97	0.47
2:D:218:SER:HB2	2:D:221:GLN:HB2	1.97	0.47
1:A:737:MET:HG3	1:A:742:ASP:HB2	1.97	0.46
2:D:140:GLU:HB3	2:D:142:LEU:HD22	1.97	0.46
2:D:396:ALA:HA	2:D:562:LYS:HD3	1.95	0.46
2:D:392:LEU:HG	2:D:562:LYS:HB3	1.96	0.46
1:B:999:GLN:NE2	1:C:999:GLN:OE1	2.48	0.46
1:A:948:VAL:O	1:A:952:ASN:ND2	2.49	0.46
1:A:881:SER:OG	1:A:884:THR:OG1	2.33	0.46
1:B:305:VAL:HG22	1:B:599:THR:HG23	1.97	0.46
1:C:270:ARG:NH2	1:C:287:ASP:OD2	2.49	0.46
1:C:103:ILE:HG23	1:C:238:LEU:HD11	1.97	0.46
1:C:366:TYR:HA	1:C:371:PHE:HZ	1.80	0.46
1:A:277:ASN:OD1	1:A:281:THR:N	2.47	0.45
1:B:412:THR:OG1	1:B:413:GLY:N	2.49	0.45
2:D:271:TRP:HE1	2:D:273:ARG:HH21	1.64	0.45
2:D:245:ARG:HA	2:D:262:LEU:HD21	1.98	0.45
1:C:897:MET:HE2	1:C:897:MET:HB3	1.87	0.45
2:D:182:GLU:HA	2:D:185:VAL:HG22	1.98	0.45
1:A:831:ILE:HB	4:B:1302:NAG:H83	1.98	0.45
1:B:42:VAL:HG22	1:C:562:PHE:HB2	1.98	0.45
1:B:656:SER:HB3	1:B:695:SER:HB2	1.99	0.45
1:B:184:LEU:HB3	1:B:203:THR:HB	1.98	0.45
1:C:629:THR:OG1	1:C:630:TRP:N	2.50	0.45
1:A:106:THR:OG1	1:A:231:ASN:O	2.34	0.45
1:A:1136:ASP:N	1:A:1136:ASP:OD1	2.47	0.45
1:B:195:TYR:HE2	1:B:225:ASP:HB2	1.82	0.45
1:C:335:PHE:HA	1:C:338:VAL:HG12	1.98	0.45
2:D:529:LEU:HD13	2:D:544:ILE:HD11	1.98	0.45
1:A:108:LEU:HD12	1:A:234:ARG:HH11	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:29:LEU:HD21	2:D:97:LEU:HD12	1.98	0.45
1:A:498:TYR:O	1:A:503:GLN:NE2	2.50	0.44
1:C:530:LEU:HD21	1:C:582:LEU:HD11	1.98	0.44
1:C:609:TYR:OH	1:C:631:ARG:NH1	2.50	0.44
1:B:270:ARG:NH2	1:B:287:ASP:OD2	2.50	0.44
1:C:316:ARG:HB2	1:C:630:TRP:HB2	1.98	0.44
1:C:579:LEU:HB2	3:M:2:NAG:H81	1.99	0.44
2:D:67:ASP:O	2:D:71:ALA:N	2.49	0.44
1:C:93:ILE:HG22	1:C:181:PHE:HB3	1.99	0.44
1:C:40:ASP:OD1	1:C:40:ASP:N	2.49	0.44
1:C:631:ARG:NH2	1:C:633:TYR:O	2.49	0.44
1:B:602:SER:OG	1:B:603:ASN:N	2.48	0.44
1:A:850:GLN:HB2	1:A:855:LEU:HD12	2.00	0.44
1:C:64:TRP:HE1	1:C:261:ALA:HB1	1.83	0.44
1:C:305:VAL:HG22	1:C:599:THR:HG23	1.99	0.44
1:C:400:ARG:HA	1:C:504:PRO:HA	2.00	0.44
1:C:423:PRO:HB3	1:C:460:PRO:HB3	2.00	0.44
2:D:482:ARG:HD2	2:D:488:VAL:HA	2.00	0.44
2:D:482:ARG:O	2:D:606:TRP:NE1	2.48	0.44
1:B:739:ILE:O	1:B:997:ARG:NH2	2.51	0.43
1:A:97:ASN:OD1	1:A:97:ASN:N	2.51	0.43
1:B:851:LYS:HG3	1:B:856:THR:HG22	2.01	0.43
1:A:1102:THR:HG23	1:A:1108:GLU:H	1.82	0.43
1:B:877:GLY:O	1:B:881:SER:HB2	2.19	0.43
1:C:34:ARG:NH1	1:C:218:SER:OG	2.52	0.43
2:D:560:LEU:HD12	2:D:564:GLU:HB3	1.99	0.43
1:C:429:CYS:HB2	1:C:510:LEU:HB2	2.00	0.43
2:D:307:ILE:HG23	2:D:369:PHE:HD1	1.82	0.43
1:A:188:VAL:HG23	1:A:220:LEU:HD12	1.99	0.43
1:A:1016:ARG:HH22	1:B:1014:GLU:HG2	1.84	0.43
1:B:425:ASP:OD1	1:B:425:ASP:N	2.52	0.43
1:C:987:GLU:HA	1:C:990:ILE:HG22	2.00	0.43
1:A:124:VAL:HG23	1:A:169:PRO:HA	2.00	0.43
1:A:637:SER:OG	1:A:638:ASN:N	2.52	0.43
2:D:320:LEU:HD13	2:D:380:GLN:HG2	1.99	0.43
1:A:25:PRO:HA	1:A:26:PRO:HD3	1.85	0.42
1:A:739:ILE:HA	1:A:997:ARG:HD3	2.01	0.42
1:B:970:ILE:HD12	1:B:980:ARG:HD3	2.01	0.42
1:C:85:ASN:OD1	1:C:85:ASN:N	2.52	0.42
1:A:402:ASP:N	1:A:402:ASP:OD1	2.52	0.42
1:B:822:LYS:HE3	1:B:936:SER:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:72:PHE:O	2:D:76:GLN:CB	2.67	0.42
1:B:376:CYS:HA	1:B:429:CYS:HA	2.01	0.42
1:B:565:ASP:OD1	1:B:565:ASP:N	2.53	0.42
2:D:390:PHE:HA	2:D:393:ARG:HB2	2.02	0.42
2:D:398:GLU:HB3	2:D:514:ARG:HB3	2.01	0.42
1:A:705:SER:HA	4:A:1303:NAG:H82	2.01	0.42
1:B:450:TYR:HE2	1:B:452:LEU:HD13	1.85	0.42
1:B:556:PHE:HB2	1:B:581:ILE:HD13	2.02	0.42
1:B:967:PHE:HD2	1:B:993:LEU:HD23	1.85	0.42
2:D:85:LEU:HD11	2:D:97:LEU:HD13	2.01	0.42
1:B:103:ILE:HG23	1:B:238:LEU:HD11	2.01	0.42
1:A:41:LYS:HG2	1:B:559:PHE:HD2	1.85	0.42
1:B:441:LYS:HD2	1:B:441:LYS:HA	1.82	0.42
1:B:543:LEU:HD21	1:B:570:THR:HG21	2.01	0.42
1:B:709:ILE:HG22	1:B:1072:PHE:HB2	2.01	0.41
2:D:267:LEU:HD13	2:D:272:GLY:HA3	2.02	0.41
2:D:169:ARG:NH1	2:D:499:ASP:OD1	2.45	0.41
1:A:707:ASN:N	1:A:707:ASN:OD1	2.52	0.41
1:B:651:GLU:O	1:B:691:ALA:N	2.48	0.41
2:D:206:ASP:OD1	2:D:206:ASP:N	2.47	0.41
1:A:643:ARG:HH11	1:C:828:ALA:HB3	1.86	0.41
1:C:595:ILE:HG23	1:C:661:ILE:HG21	2.02	0.41
1:A:940:SER:O	1:A:940:SER:OG	2.33	0.41
1:C:663:ILE:HD11	1:C:669:ALA:HB2	2.02	0.41
1:C:851:LYS:HG3	1:C:856:THR:HG22	2.02	0.41
1:A:1099:TRP:HB2	1:A:1132:ASN:ND2	2.36	0.41
1:B:1079:CYS:HB2	1:B:1123:CYS:HB2	1.68	0.41
1:C:441:LYS:HB3	1:C:442:VAL:H	1.75	0.41
2:D:363:LYS:HD2	2:D:363:LYS:HA	1.86	0.41
1:A:99:ILE:HD13	1:A:262:TYR:HB3	2.02	0.41
1:C:433:TRP:HB2	1:C:506:ARG:HB3	2.01	0.41
1:B:362:TYR:HD2	1:B:384:LEU:HB3	1.85	0.41
1:C:449:LEU:HB3	1:C:489:LEU:HD12	2.03	0.41
2:D:299:ASP:N	2:D:299:ASP:OD1	2.53	0.41
1:C:441:LYS:HD2	1:C:441:LYS:HA	1.85	0.41
1:A:897:MET:HE2	1:A:897:MET:HB3	1.91	0.40
1:B:902:ARG:NH2	1:B:1046:LEU:O	2.48	0.40
1:C:90:PHE:HB3	1:C:187:PHE:HB2	2.03	0.40
1:A:743:SER:HB3	1:B:545:GLY:HA2	2.02	0.40
1:A:967:PHE:HD2	1:A:993:LEU:HD23	1.87	0.40
1:B:111:LYS:HG3	1:C:465:ILE:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:381:PRO:HA	1:B:384:LEU:HD12	2.03	0.40
2:D:330:ASN:HB2	2:D:357:ARG:HE	1.86	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	1047/1205 (87%)	975 (93%)	72 (7%)	0	100	100
1	B	1047/1205 (87%)	1005 (96%)	42 (4%)	0	100	100
1	C	1039/1205 (86%)	991 (95%)	47 (4%)	1 (0%)	48	75
2	D	593/805 (74%)	551 (93%)	42 (7%)	0	100	100
All	All	3726/4420 (84%)	3522 (94%)	203 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	983	PRO

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	931/1053 (88%)	924 (99%)	7 (1%)	73	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	931/1053 (88%)	928 (100%)	3 (0%)	86	94
1	C	927/1053 (88%)	923 (100%)	4 (0%)	84	93
2	D	526/711 (74%)	525 (100%)	1 (0%)	87	96
All	All	3315/3870 (86%)	3300 (100%)	15 (0%)	78	92

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	139	LEU
1	A	151	GLU
1	A	152	PHE
1	A	153	ARG
1	A	550	THR
1	A	982	ASP
1	A	987	GLU
1	B	407	ILE
1	B	642	THR
1	B	982	ASP
1	C	328	ASN
1	C	521	VAL
1	C	981	LEU
1	C	988	VAL
2	D	81	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (93) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	ASN
1	A	311	GLN
1	A	318	GLN
1	A	357	ASN
1	A	391	ASN
1	A	406	GLN
1	A	419	ASN
1	A	447	ASN
1	A	471	GLN
1	A	474	ASN
1	A	503	GLN
1	A	533	ASN
1	A	604	GLN

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Mol	Chain	Res	Type
1	A	641	GLN
1	A	687	GLN
1	A	898	GLN
1	A	910	GLN
1	A	952	ASN
1	A	954	GLN
1	A	957	ASN
1	A	962	GLN
1	A	1002	GLN
1	A	1007	GLN
1	A	1033	GLN
1	A	1051	GLN
1	A	1068	GLN
1	A	1080	HIS
1	A	1103	GLN
1	A	1105	ASN
1	A	1139	GLN
1	B	23	GLN
1	B	66	HIS
1	B	97	ASN
1	B	160	ASN
1	B	215	GLN
1	B	236	GLN
1	B	268	GLN
1	B	340	ASN
1	B	357	ASN
1	B	411	GLN
1	B	471	GLN
1	B	478	ASN
1	B	537	ASN
1	B	610	GLN
1	B	672	GLN
1	B	748	ASN
1	B	801	GLN
1	B	850	GLN
1	B	898	GLN
1	B	910	GLN
1	B	946	GLN
1	B	1002	GLN
1	B	1008	GLN
1	B	1020	ASN
1	B	1033	GLN

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Mol	Chain	Res	Type
1	B	1068	GLN
1	B	1110	GLN
1	B	1132	ASN
1	C	66	HIS
1	C	97	ASN
1	C	191	ASN
1	C	236	GLN
1	C	340	ASN
1	C	411	GLN
1	C	447	ASN
1	C	503	GLN
1	C	537	ASN
1	C	539	ASN
1	C	577	GLN
1	C	641	GLN
1	C	687	GLN
1	C	774	ASN
1	C	801	GLN
1	C	821	ASN
1	C	898	GLN
1	C	917	GLN
1	C	932	GLN
1	C	1007	GLN
1	C	1020	ASN
1	C	1061	HIS
2	D	33	ASN
2	D	42	GLN
2	D	58	ASN
2	D	76	GLN
2	D	86	GLN
2	D	102	GLN
2	D	134	ASN
2	D	239	HIS
2	D	437	ASN
2	D	472	GLN
2	D	522	GLN
2	D	524	GLN
2	D	578	ASN

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	$\# Z > 2$
3	NAG	E	1	3,1	14,14,15	0.35	0	17,19,21	0.59	0
3	NAG	E	2	3	14,14,15	0.45	0	17,19,21	0.59	1 (5%)
3	NAG	F	1	3,1	14,14,15	0.37	0	17,19,21	0.54	0
3	NAG	F	2	3	14,14,15	0.48	0	17,19,21	0.51	0
3	NAG	G	1	3,1	14,14,15	0.39	0	17,19,21	0.60	1 (5%)
3	NAG	G	2	3	14,14,15	0.38	0	17,19,21	0.49	0
3	NAG	H	1	3,1	14,14,15	0.35	0	17,19,21	0.56	0
3	NAG	H	2	3	14,14,15	0.39	0	17,19,21	0.47	0
3	NAG	I	1	3,1	14,14,15	0.42	0	17,19,21	1.43	3 (17%)
3	NAG	I	2	3	14,14,15	0.41	0	17,19,21	0.77	1 (5%)
3	NAG	J	1	3,1	14,14,15	0.36	0	17,19,21	0.58	0
3	NAG	J	2	3	14,14,15	0.38	0	17,19,21	0.48	0
3	NAG	K	1	3,1	14,14,15	0.44	0	17,19,21	0.64	1 (5%)
3	NAG	K	2	3	14,14,15	0.40	0	17,19,21	0.48	0
3	NAG	L	1	3,1	14,14,15	0.36	0	17,19,21	0.56	0
3	NAG	L	2	3	14,14,15	0.49	0	17,19,21	0.51	0
3	NAG	M	1	3,1	14,14,15	0.39	0	17,19,21	0.54	0
3	NAG	M	2	3	14,14,15	0.42	0	17,19,21	0.50	0
3	NAG	N	1	3,1	14,14,15	0.33	0	17,19,21	0.60	1 (5%)
3	NAG	N	2	3	14,14,15	0.39	0	17,19,21	0.46	0
3	NAG	O	1	3,1	14,14,15	0.41	0	17,19,21	0.62	1 (5%)
3	NAG	O	2	3	14,14,15	0.40	0	17,19,21	0.47	0
3	NAG	P	1	3,1	14,14,15	0.34	0	17,19,21	0.56	0
3	NAG	P	2	3	14,14,15	0.43	0	17,19,21	0.51	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	E	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	NAG	G	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	NAG	I	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1
3	NAG	J	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
3	NAG	K	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	K	2	3	-	0/6/23/26	0/1/1/1
3	NAG	L	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
3	NAG	M	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	2/6/23/26	0/1/1/1
3	NAG	N	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	N	2	3	-	2/6/23/26	0/1/1/1
3	NAG	O	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	O	2	3	-	0/6/23/26	0/1/1/1
3	NAG	P	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	P	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	1	NAG	C1-O5-C5	3.42	116.77	112.19
3	I	1	NAG	C2-N2-C7	3.31	127.33	122.90
3	I	2	NAG	C1-O5-C5	2.84	115.99	112.19
3	K	1	NAG	C1-O5-C5	2.17	115.10	112.19
3	O	1	NAG	C1-O5-C5	2.09	114.98	112.19
3	I	1	NAG	O4-C4-C5	2.05	114.38	109.32
3	E	2	NAG	C1-O5-C5	2.04	114.92	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1	NAG	C1-O5-C5	2.03	114.91	112.19
3	N	1	NAG	C1-O5-C5	2.01	114.88	112.19

There are no chirality outliers.

All (21) torsion outliers are listed below:

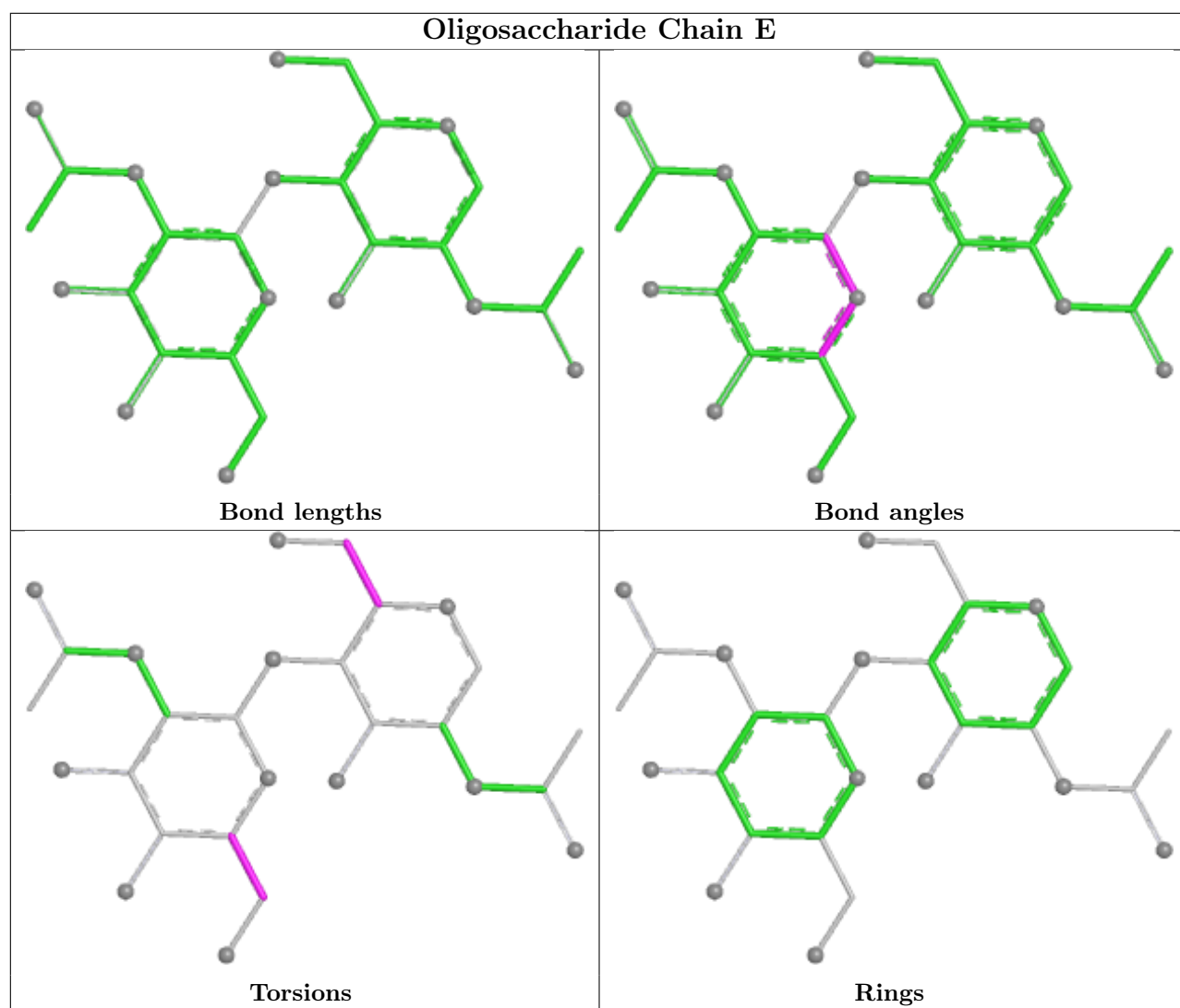
Mol	Chain	Res	Type	Atoms
3	I	1	NAG	C4-C5-C6-O6
3	M	2	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
3	M	2	NAG	C4-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
3	I	2	NAG	O5-C5-C6-O6
3	I	2	NAG	C4-C5-C6-O6
3	N	2	NAG	C4-C5-C6-O6
3	L	1	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	L	1	NAG	O5-C5-C6-O6
3	N	2	NAG	O5-C5-C6-O6
3	L	2	NAG	C4-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	L	2	NAG	O5-C5-C6-O6
3	I	1	NAG	C1-C2-N2-C7
3	I	1	NAG	C3-C2-N2-C7
3	E	2	NAG	O5-C5-C6-O6

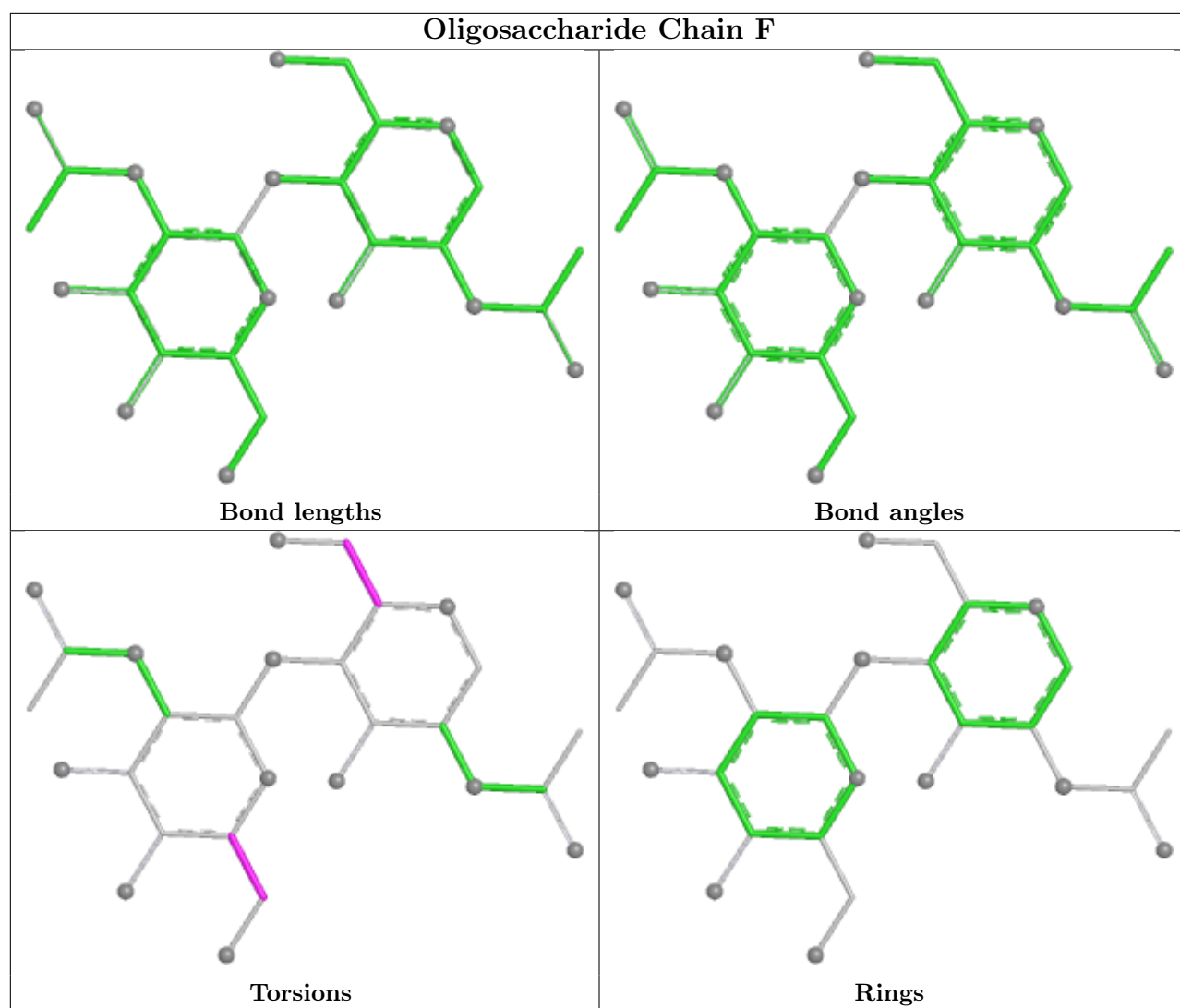
There are no ring outliers.

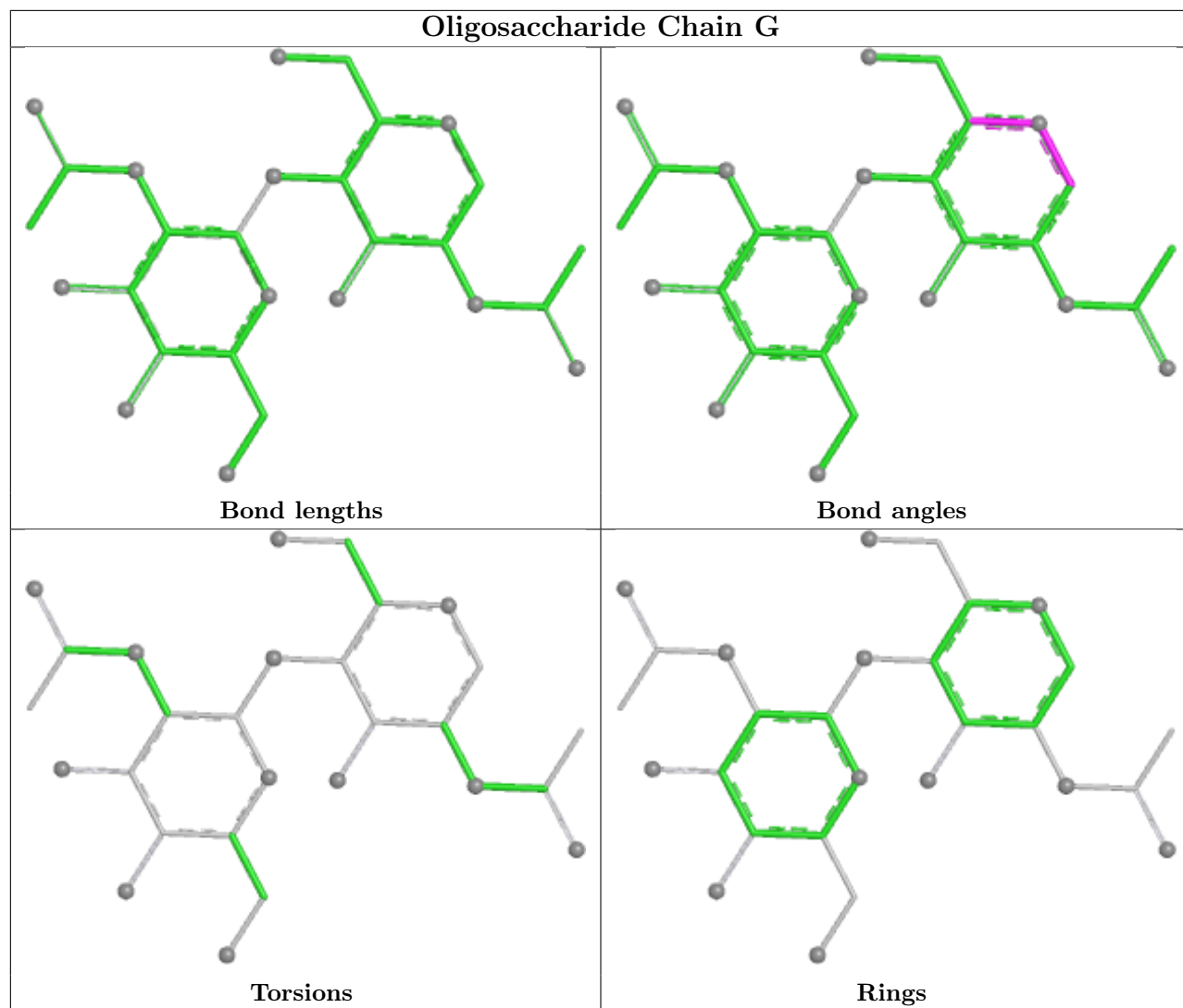
2 monomers are involved in 2 short contacts:

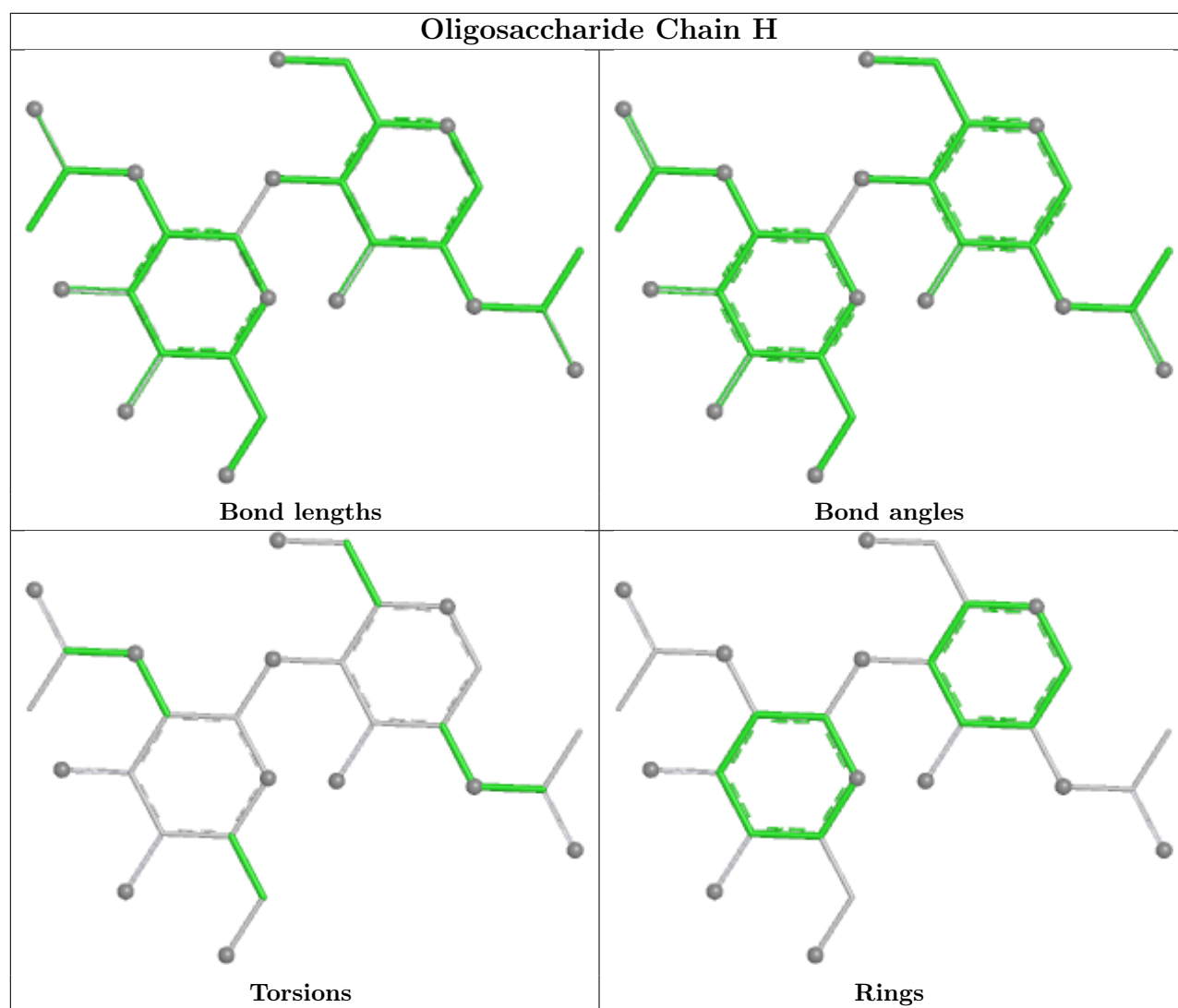
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	2	NAG	1	0
3	H	1	NAG	1	0

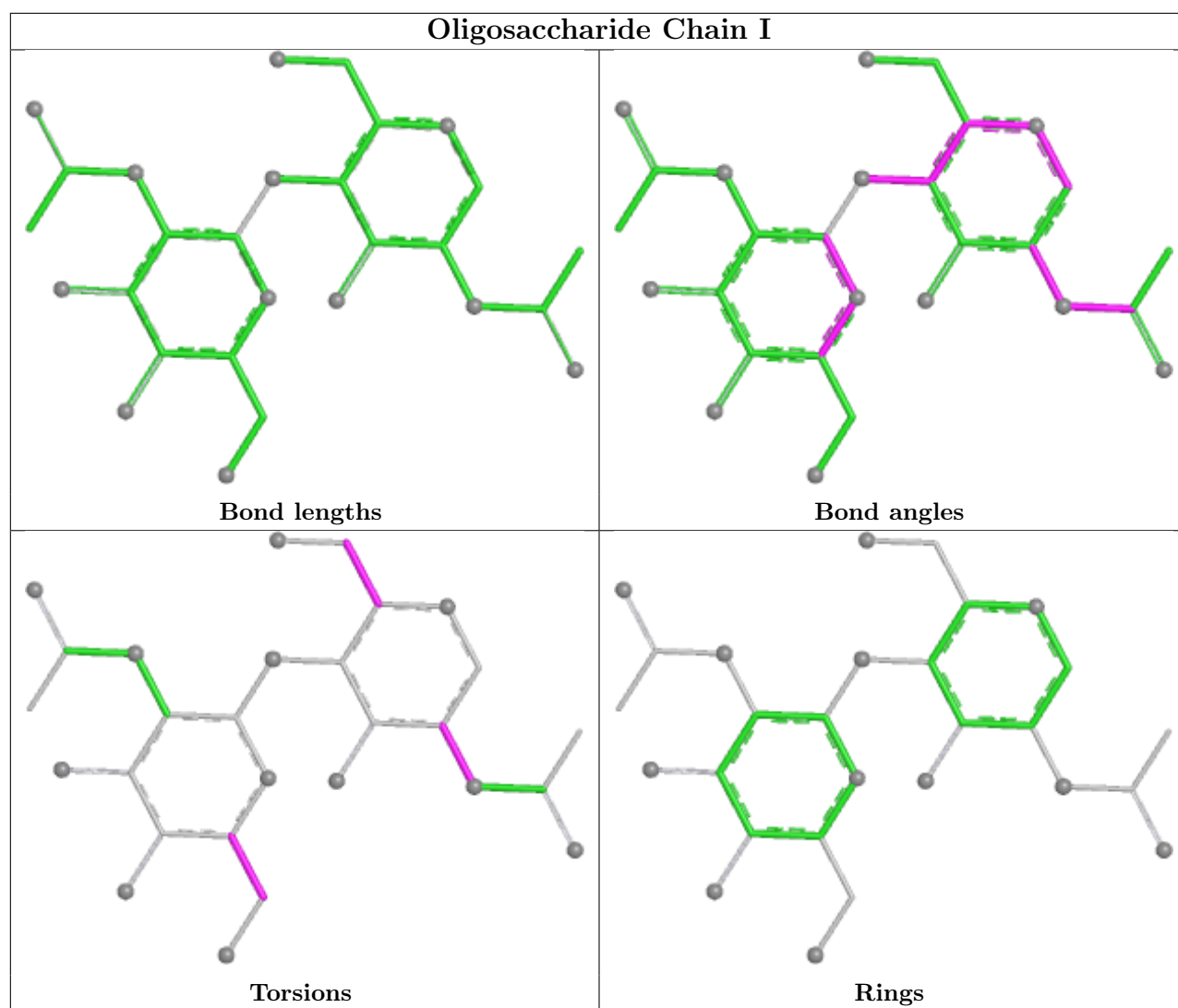
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

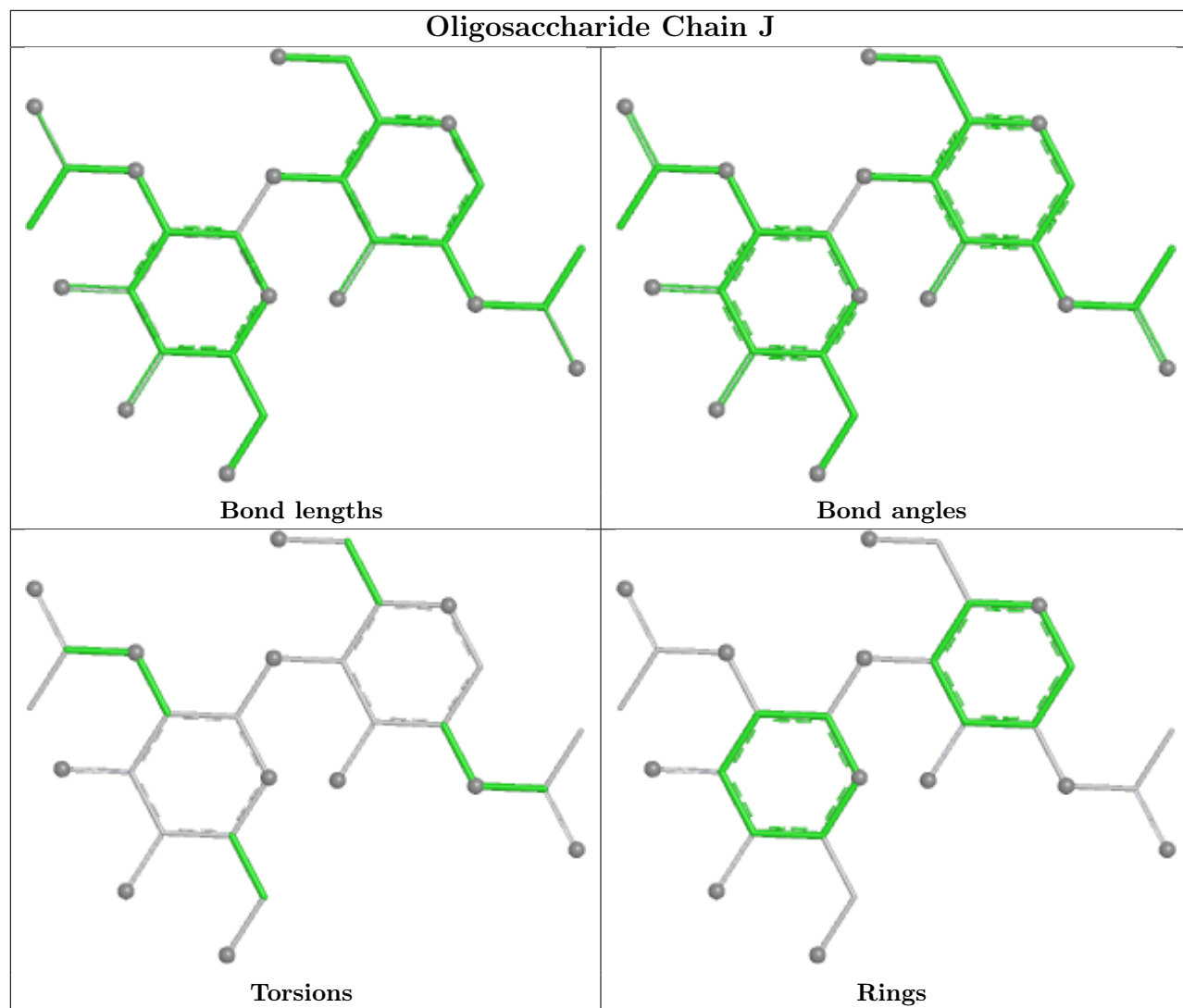


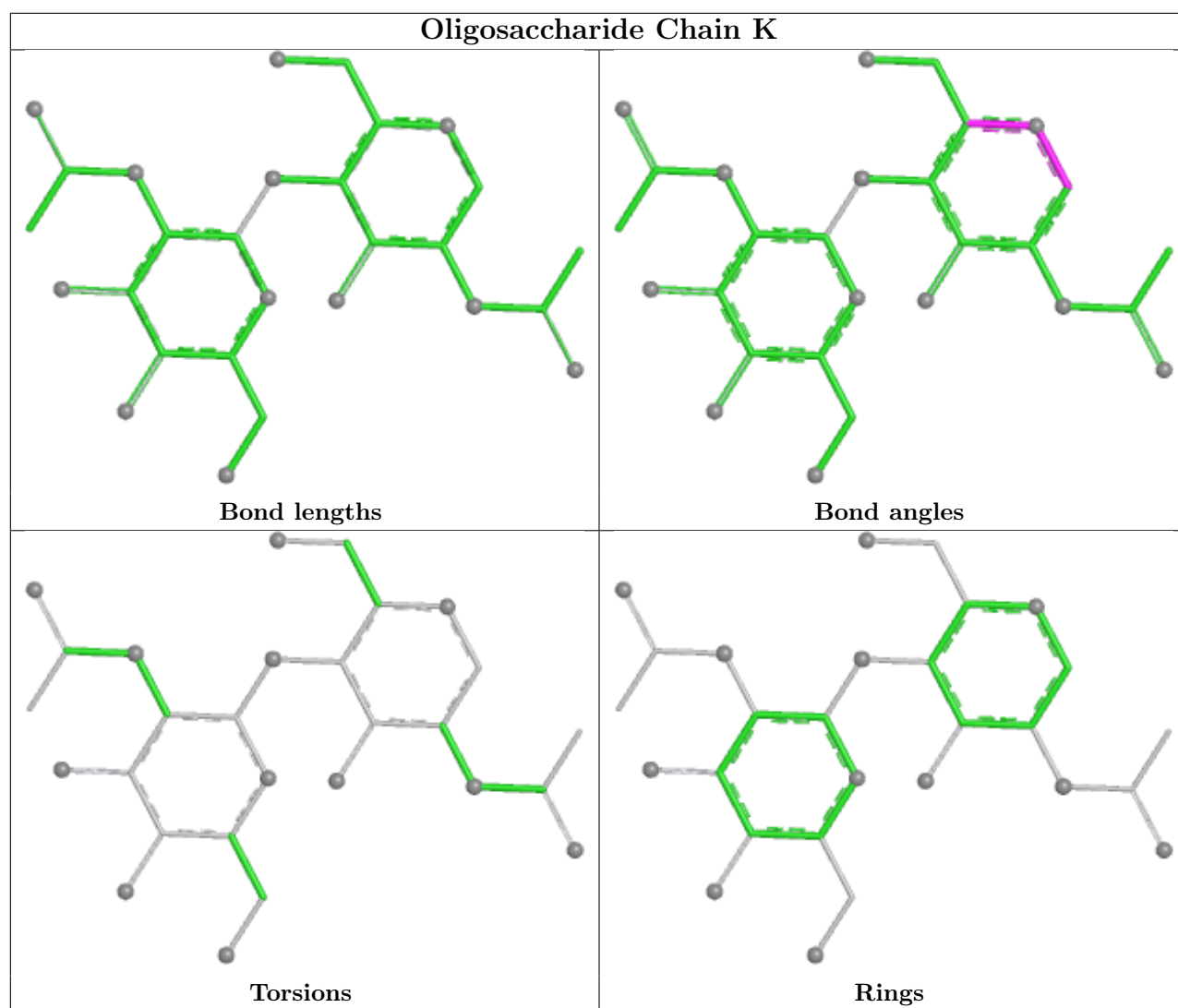


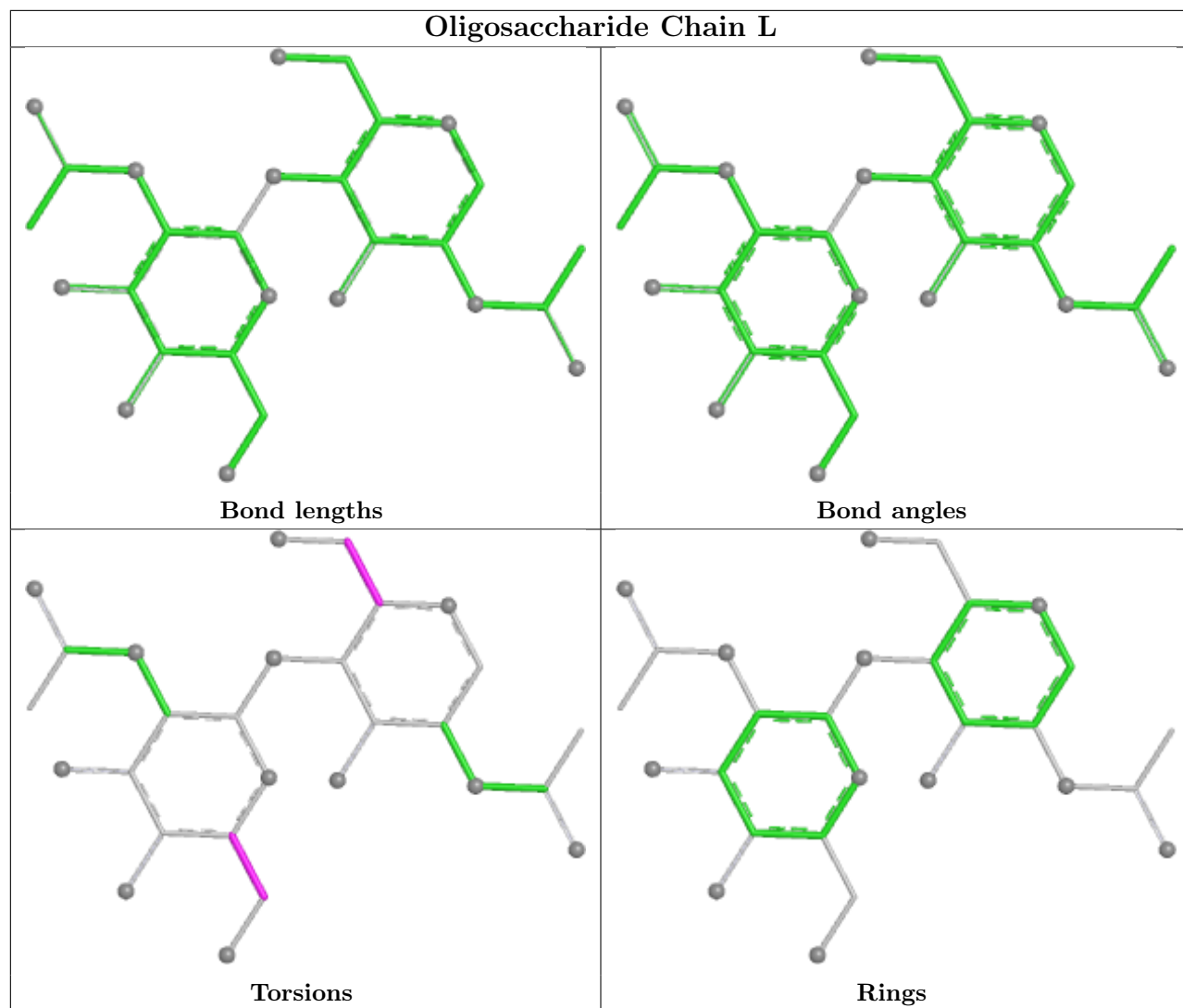


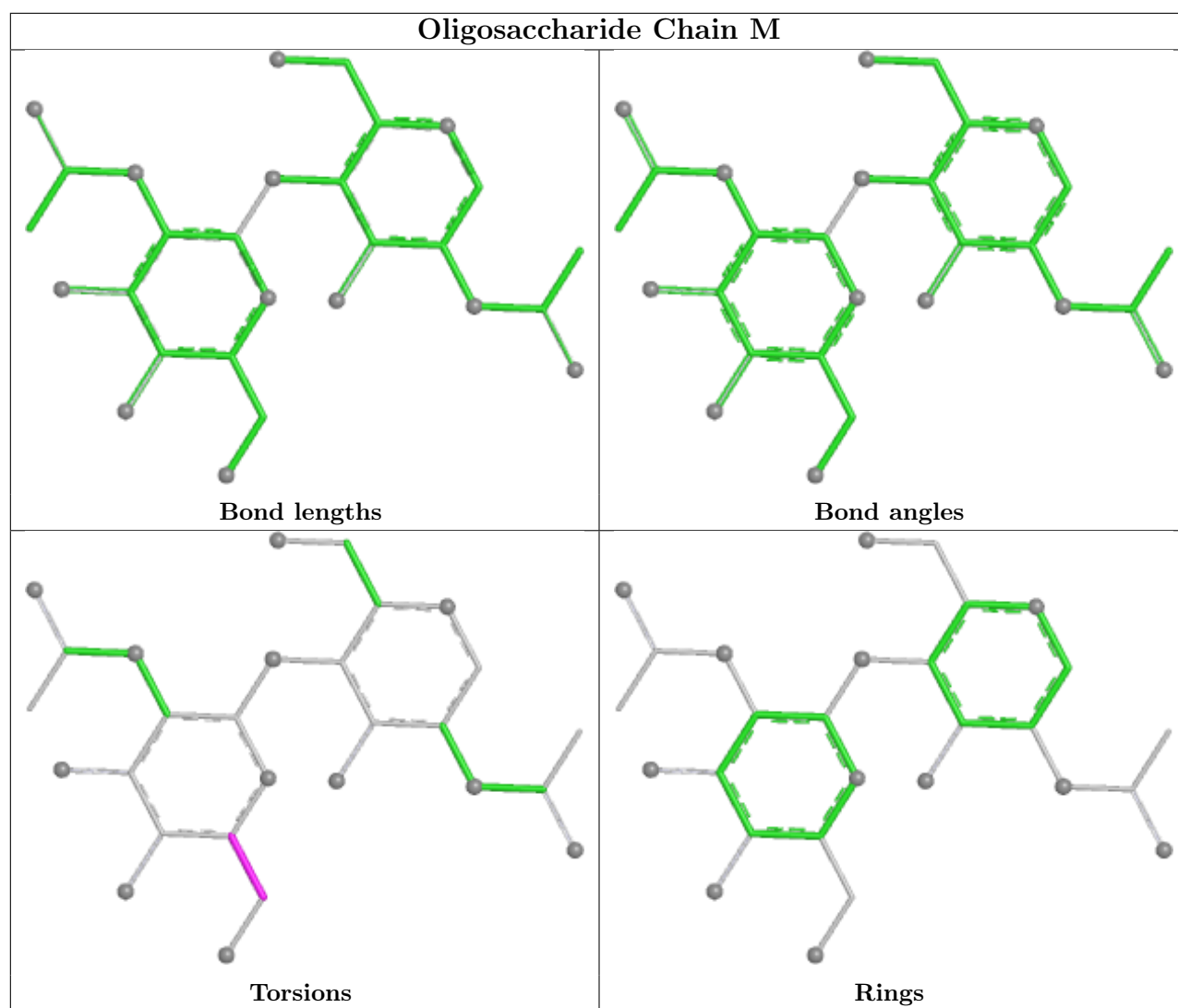


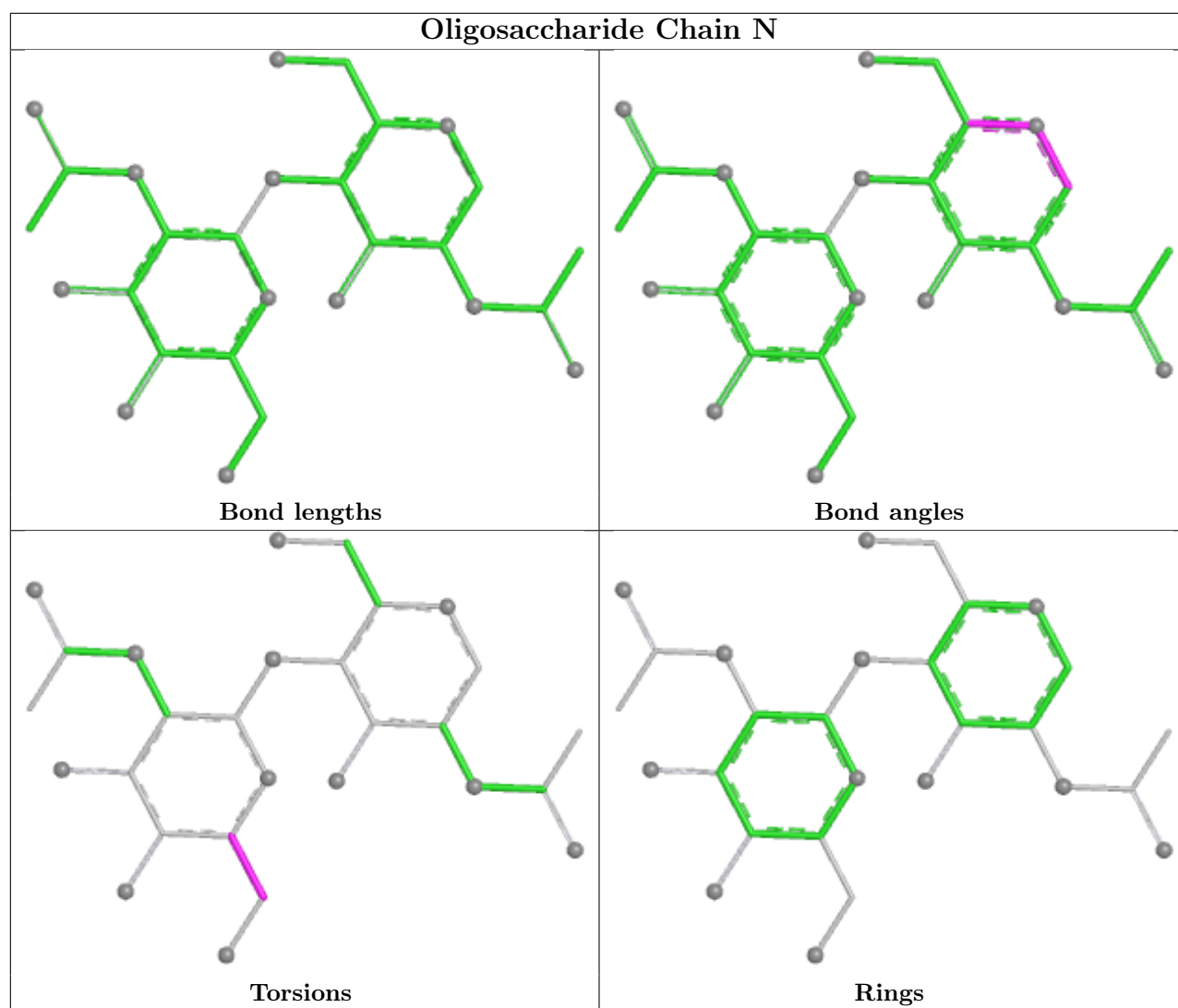


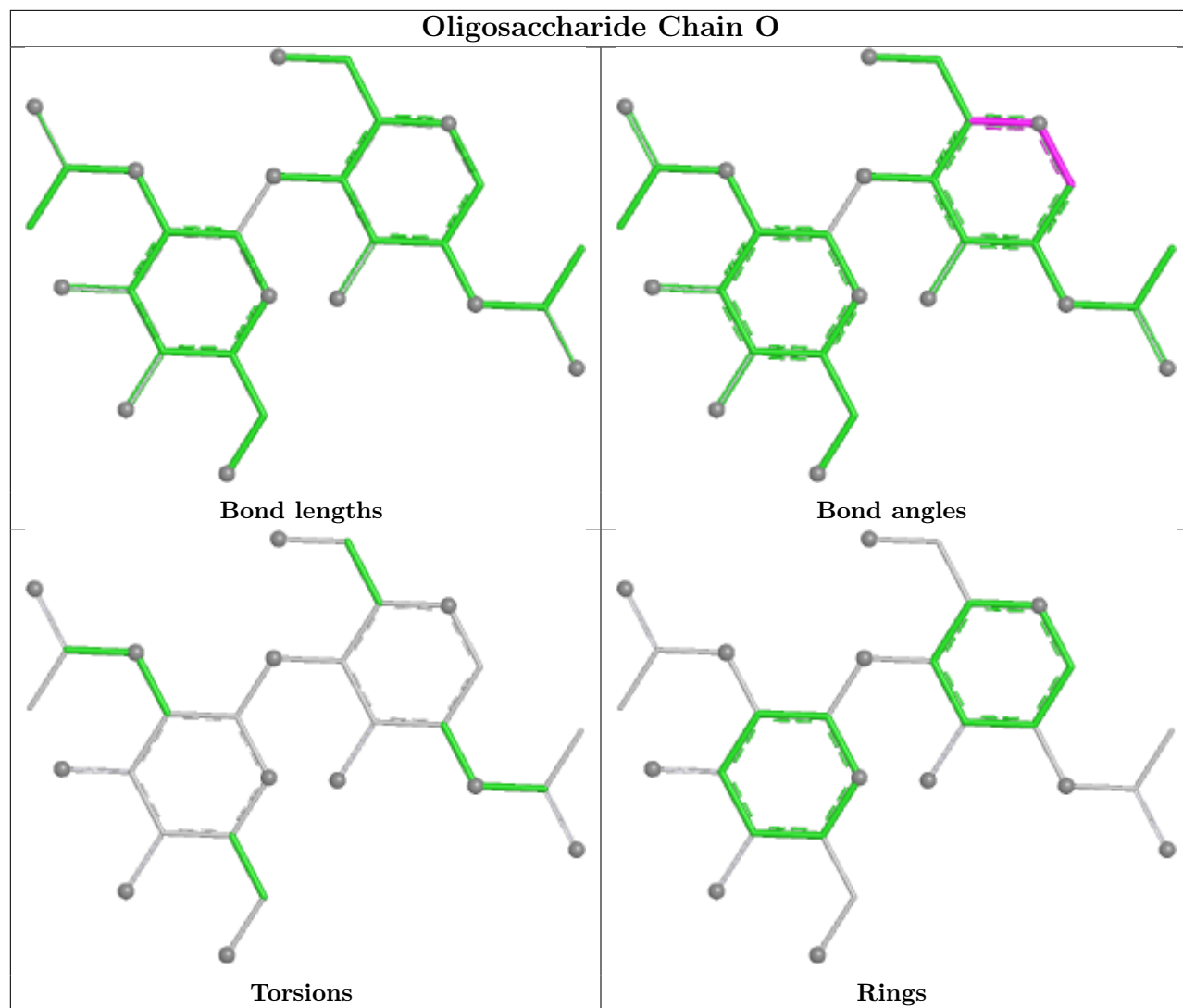


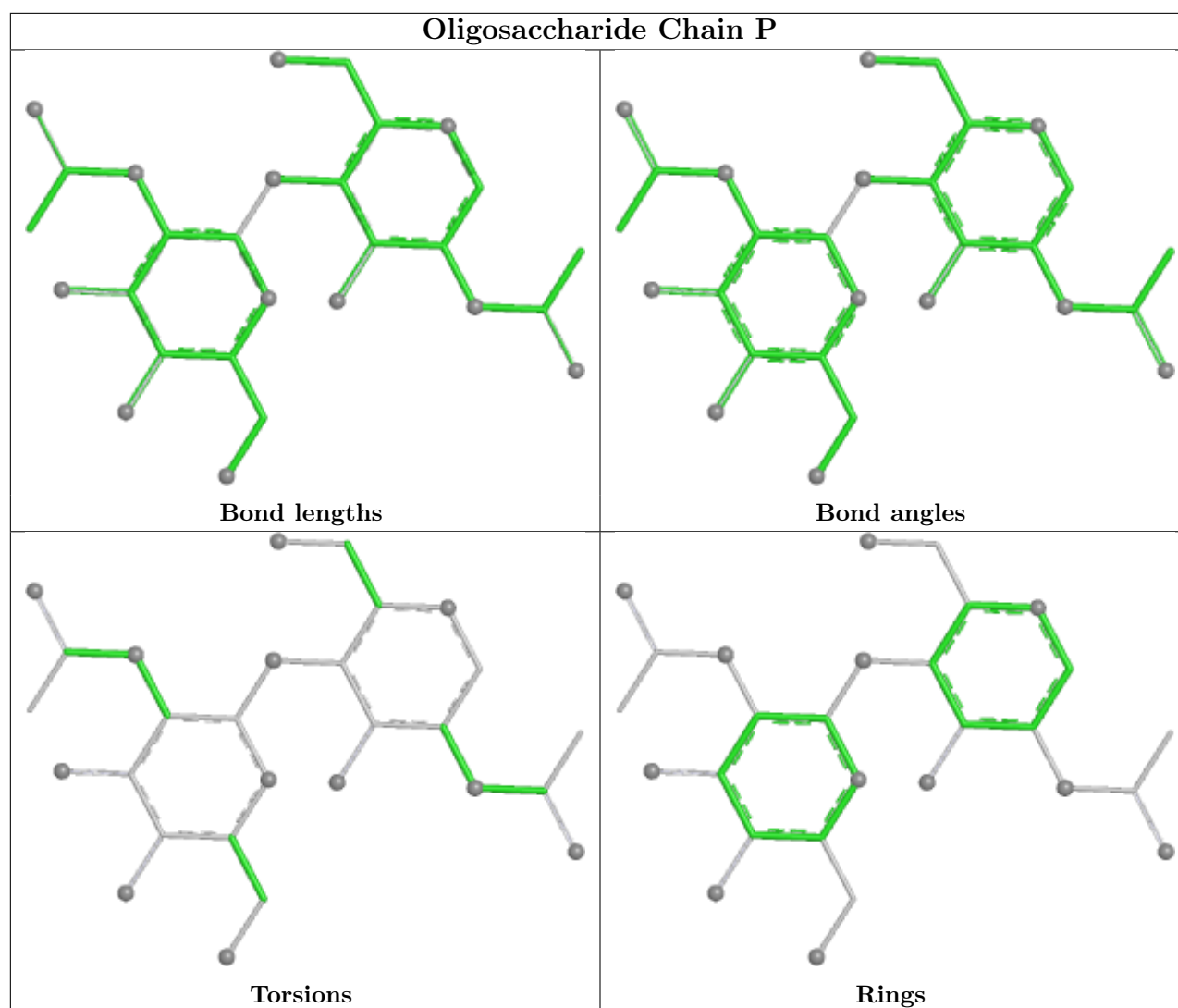












5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 2 are monoatomic - leaving 21 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	A	1302	1	14,14,15	0.40	0	17,19,21	0.56	0
4	NAG	B	1303	1	14,14,15	0.40	0	17,19,21	0.53	0
4	NAG	B	1306	1	14,14,15	0.37	0	17,19,21	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	1304	1	14,14,15	0.43	0	17,19,21	0.57	0
4	NAG	D	904	2	14,14,15	0.39	0	17,19,21	0.53	0
4	NAG	A	1301	1	14,14,15	0.38	0	17,19,21	0.50	0
4	NAG	D	903	2	14,14,15	0.38	0	17,19,21	0.51	0
4	NAG	C	1304	1	14,14,15	0.40	0	17,19,21	0.53	0
4	NAG	A	1303	1	14,14,15	0.41	0	17,19,21	0.55	0
4	NAG	C	1306	1	14,14,15	0.39	0	17,19,21	0.52	0
4	NAG	D	905	2	14,14,15	0.38	0	17,19,21	0.59	1 (5%)
4	NAG	C	1302	1	14,14,15	0.39	0	17,19,21	0.55	0
4	NAG	C	1305	1	14,14,15	0.50	0	17,19,21	0.62	1 (5%)
4	NAG	B	1304	1	14,14,15	0.39	0	17,19,21	0.56	0
4	NAG	C	1301	1	14,14,15	0.39	0	17,19,21	0.49	0
4	NAG	A	1306	1	14,14,15	0.38	0	17,19,21	0.52	0
4	NAG	B	1302	1	14,14,15	0.56	0	17,19,21	1.05	1 (5%)
4	NAG	B	1305	1	14,14,15	0.38	0	17,19,21	0.49	0
4	NAG	A	1305	1	14,14,15	0.39	0	17,19,21	0.50	0
4	NAG	B	1301	1	14,14,15	0.37	0	17,19,21	0.59	0
4	NAG	C	1303	1	14,14,15	0.39	0	17,19,21	0.54	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1302	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1303	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1306	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1304	1	-	1/6/23/26	0/1/1/1
4	NAG	D	904	2	-	0/6/23/26	0/1/1/1
4	NAG	A	1301	1	-	0/6/23/26	0/1/1/1
4	NAG	D	903	2	-	2/6/23/26	0/1/1/1
4	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1303	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1306	1	-	0/6/23/26	0/1/1/1
4	NAG	D	905	2	-	2/6/23/26	0/1/1/1
4	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1305	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1304	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1301	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1302	1	-	4/6/23/26	0/1/1/1
4	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1305	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1303	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1302	NAG	C2-N2-C7	3.20	127.18	122.90
4	C	1305	NAG	C1-O5-C5	2.14	115.05	112.19
4	D	905	NAG	C1-O5-C5	2.02	114.90	112.19

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1302	NAG	O5-C5-C6-O6
4	A	1305	NAG	O5-C5-C6-O6
4	D	905	NAG	O5-C5-C6-O6
4	C	1301	NAG	C4-C5-C6-O6
4	A	1305	NAG	C4-C5-C6-O6
4	D	905	NAG	C4-C5-C6-O6
4	A	1302	NAG	C4-C5-C6-O6
4	D	903	NAG	O5-C5-C6-O6
4	B	1301	NAG	C8-C7-N2-C2
4	B	1301	NAG	O7-C7-N2-C2
4	D	903	NAG	C4-C5-C6-O6
4	C	1301	NAG	O5-C5-C6-O6
4	B	1305	NAG	C4-C5-C6-O6
4	C	1304	NAG	O5-C5-C6-O6
4	C	1304	NAG	C4-C5-C6-O6
4	A	1306	NAG	O5-C5-C6-O6
4	A	1306	NAG	C4-C5-C6-O6
4	B	1302	NAG	C4-C5-C6-O6
4	B	1305	NAG	O5-C5-C6-O6
4	C	1305	NAG	O5-C5-C6-O6
4	C	1305	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	B	1302	NAG	O5-C5-C6-O6
4	B	1306	NAG	C4-C5-C6-O6
4	C	1302	NAG	C4-C5-C6-O6
4	B	1306	NAG	O5-C5-C6-O6
4	C	1302	NAG	O5-C5-C6-O6
4	A	1304	NAG	C1-C2-N2-C7
4	B	1302	NAG	C1-C2-N2-C7
4	B	1302	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1303	NAG	1	0
4	C	1302	NAG	1	0
4	B	1302	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

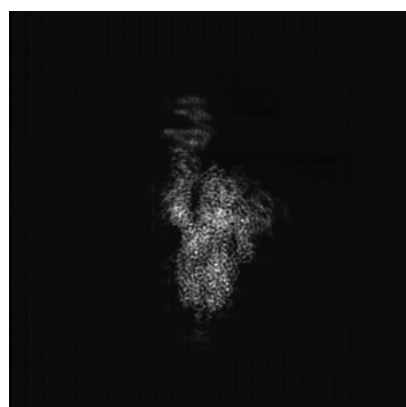
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32680. These allow visual inspection of the internal detail of the map and identification of artifacts.

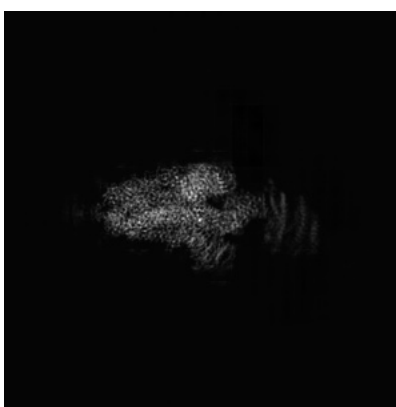
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

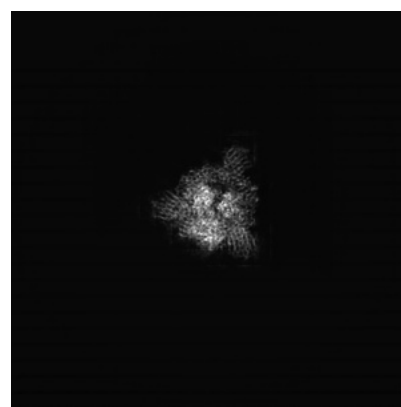
6.1.1 Primary map



X



Y



Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 240



Y Index: 240



Z Index: 240

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: 239



Y Index: 246

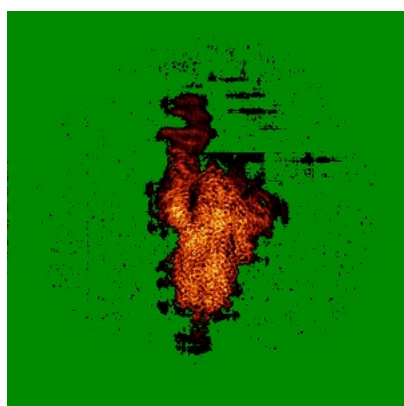


Z Index: 227

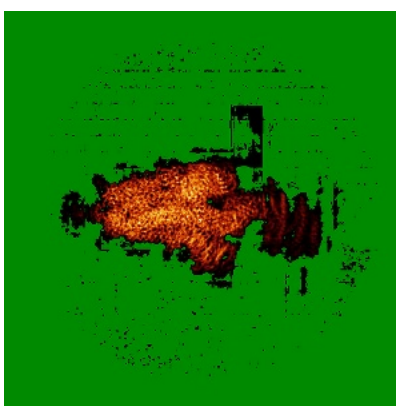
The images above show the largest variance slices of the map in three orthogonal directions.

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

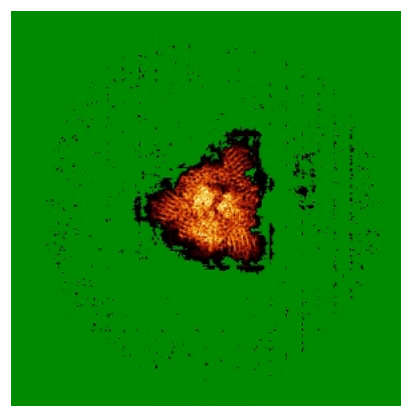
6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



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

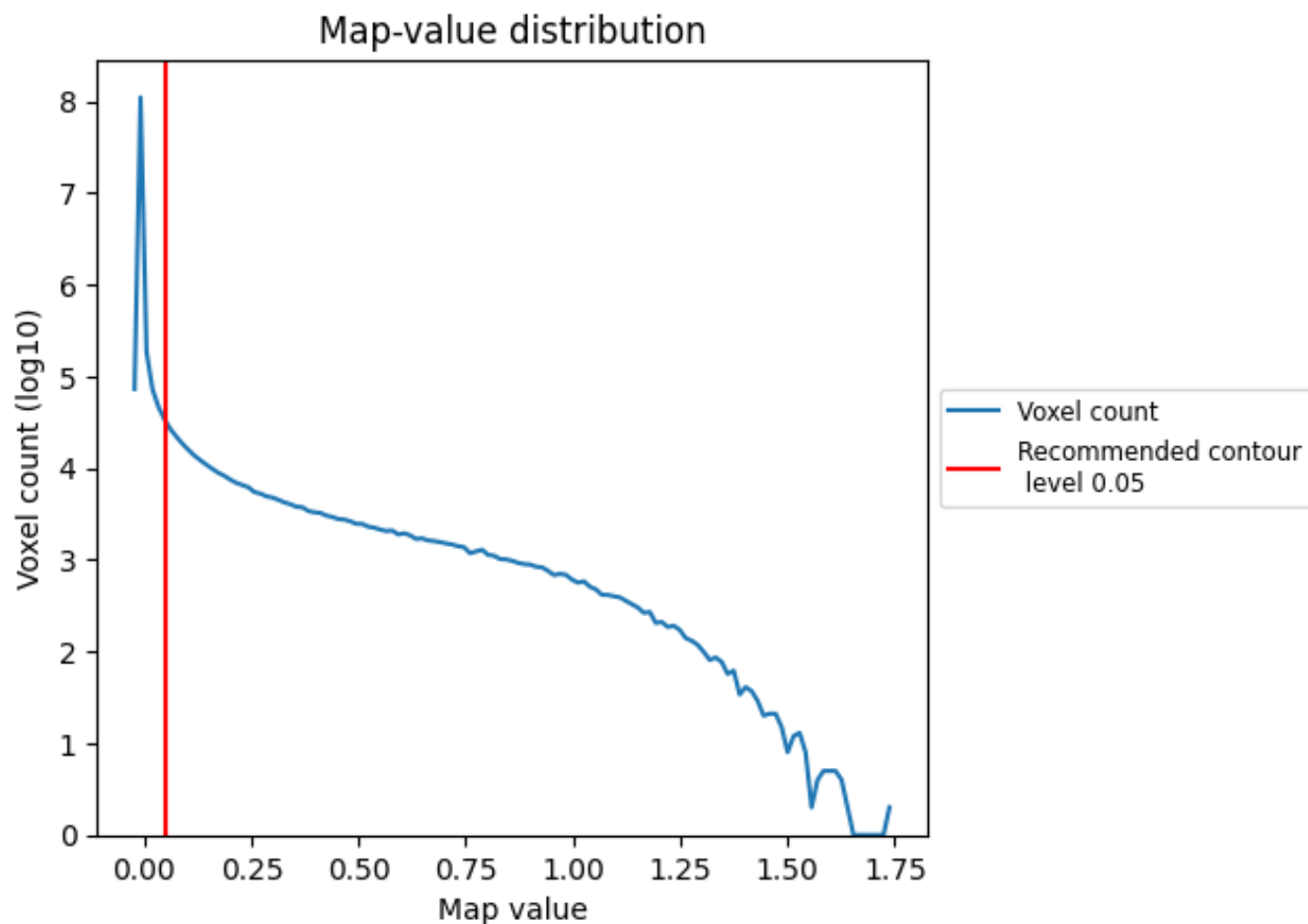
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

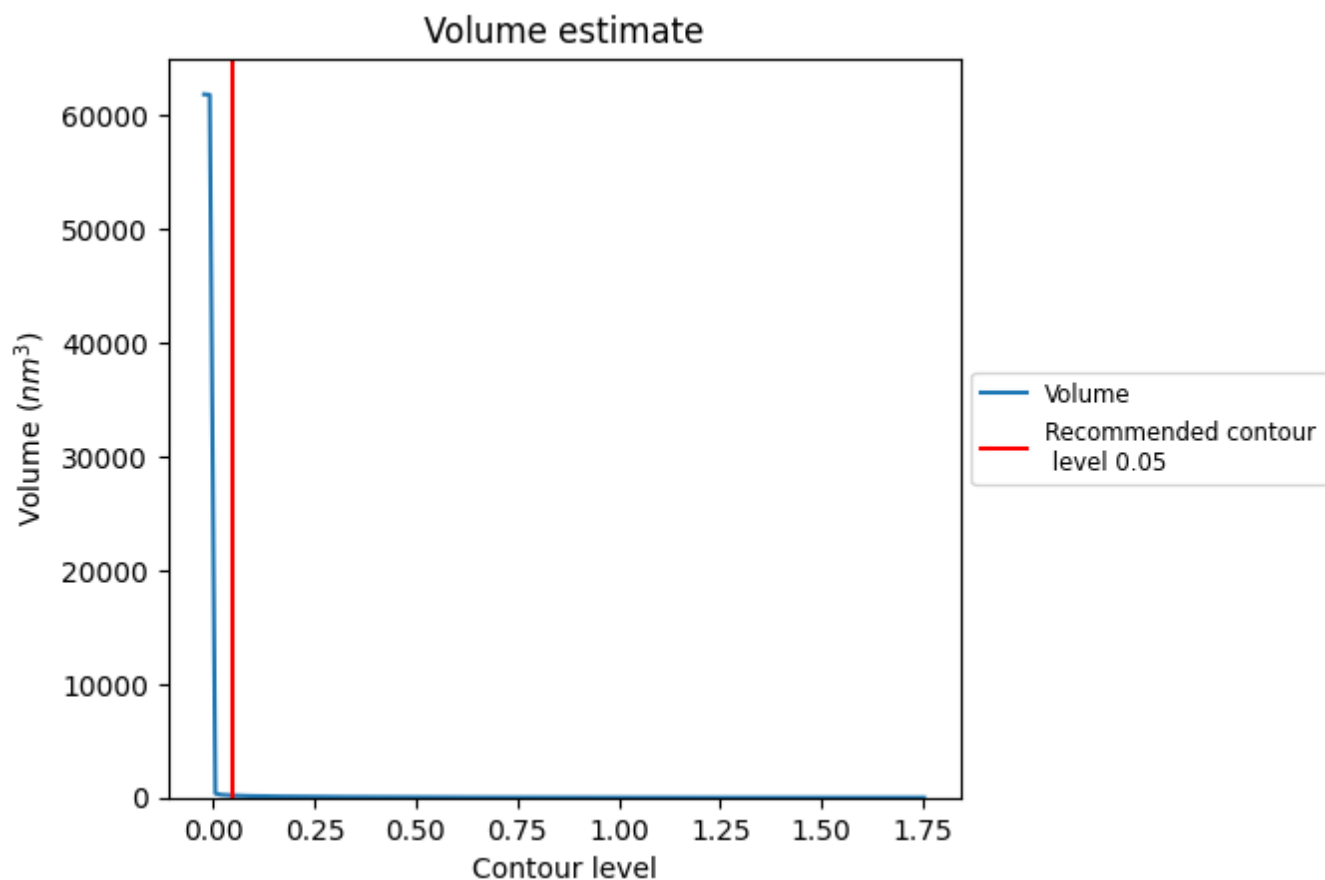
This section contains the results of statistical analysis of the map.

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.

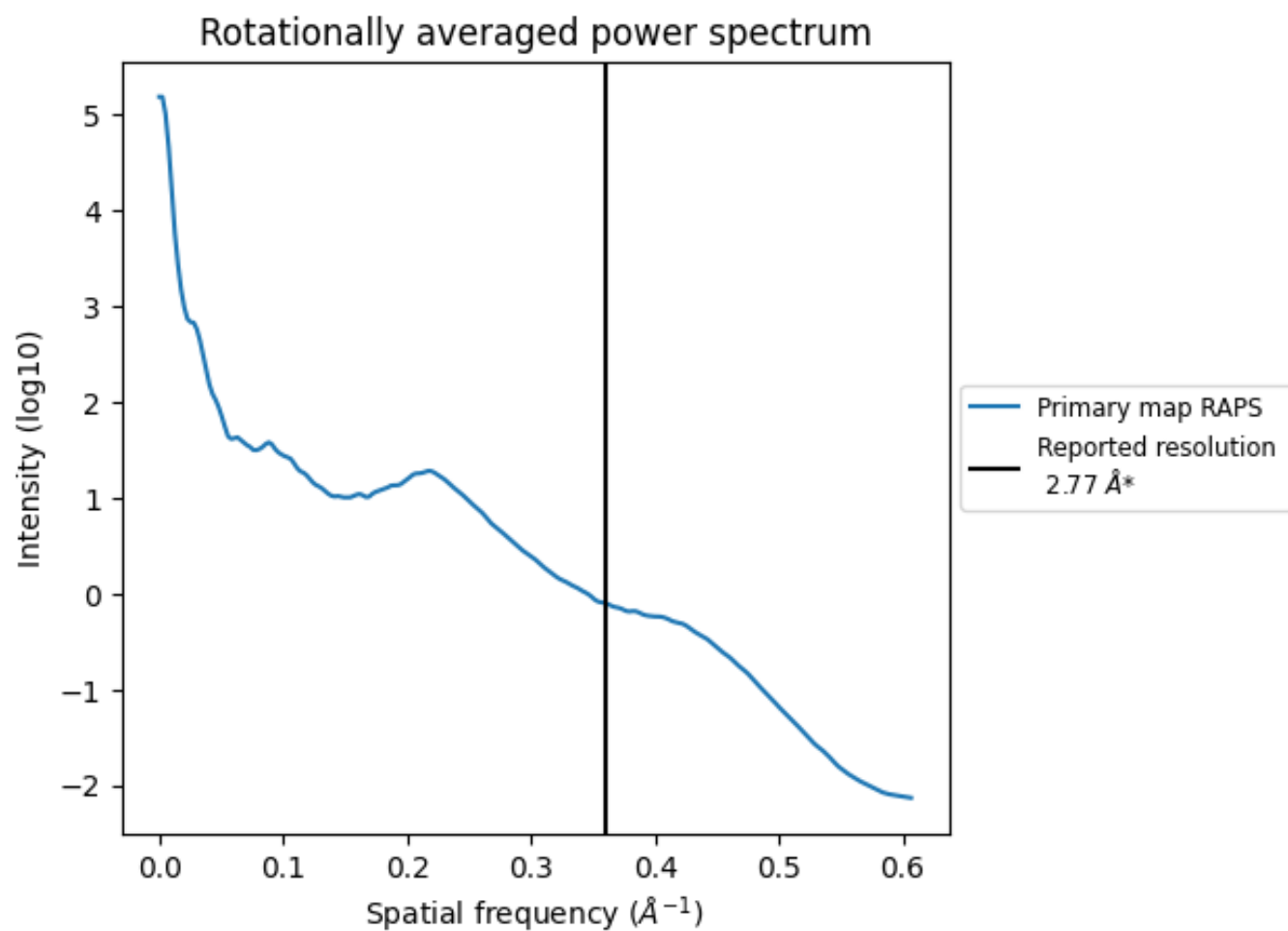
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 182 nm³; this corresponds to an approximate mass of 165 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 0.361 Å⁻¹

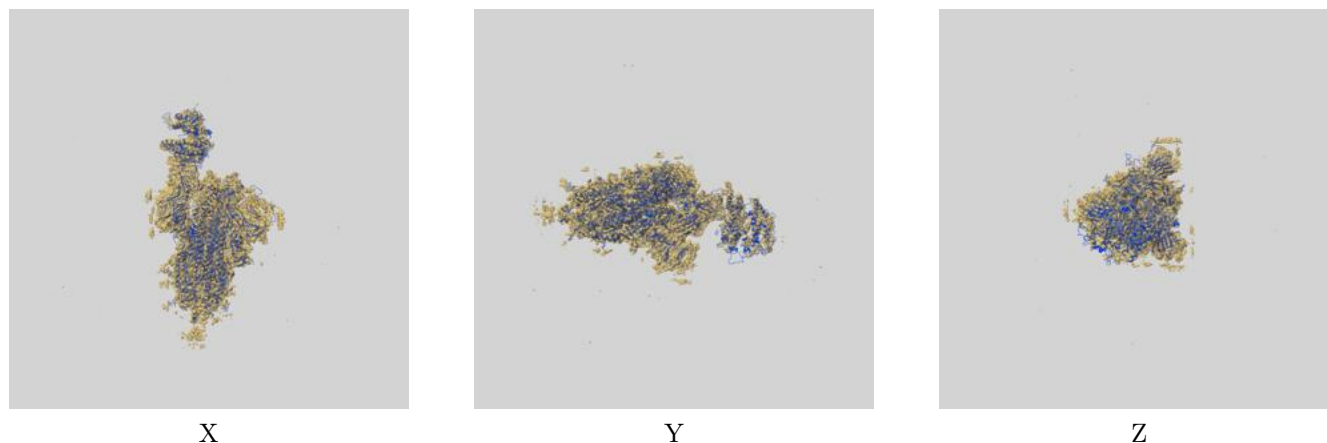
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

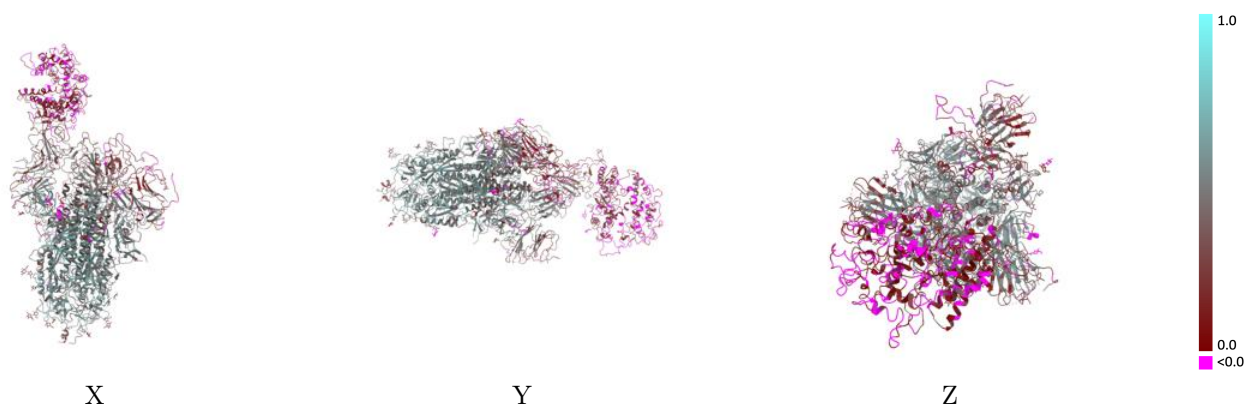
This section contains information regarding the fit between EMDB map EMD-32680 and PDB model 7WPA. Per-residue inclusion information can be found in section [3](#) on page [11](#).

9.1 Map-model overlay [i](#)



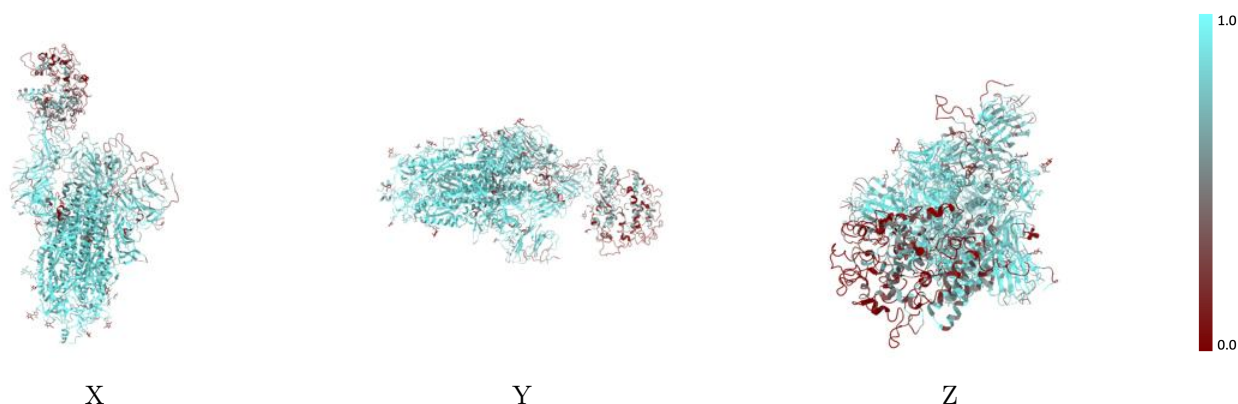
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



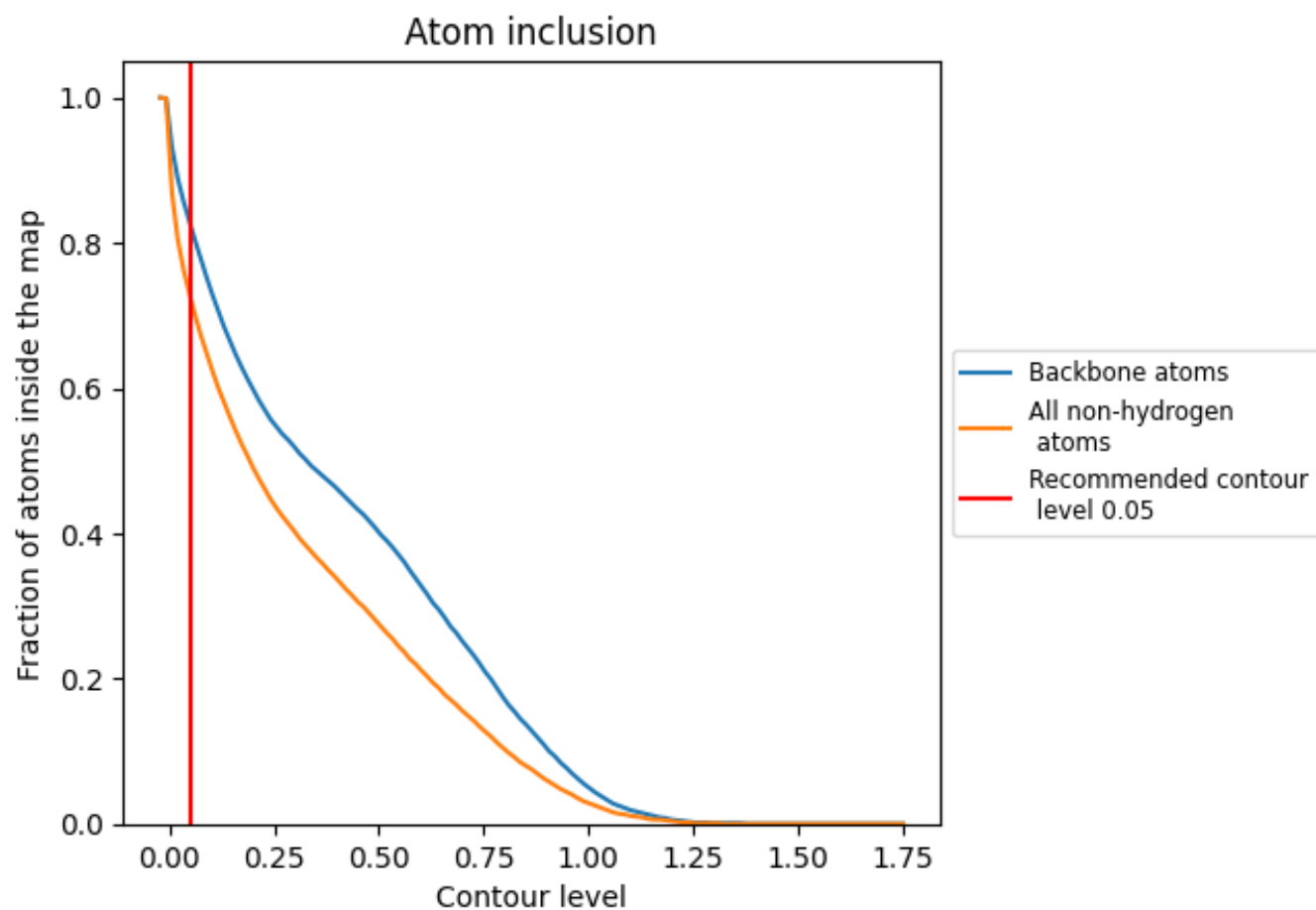
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.05).

9.4 Atom inclusion ⓘ



At the recommended contour level, 82% of all backbone atoms, 72% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7240	 0.3710
A	 0.8070	 0.4310
B	 0.8190	 0.4400
C	 0.7750	 0.4180
D	 0.3470	 0.0780
E	 0.7860	 0.3490
F	 0.1790	 0.1230
G	 0.2860	 0.2570
H	 0.6790	 0.3970
I	 0.2500	 0.0770
J	 0.3570	 0.3530
K	 0.5000	 0.2550
L	 0.3930	 0.2930
M	 0.2140	 -0.0530
N	 0.6790	 0.2490
O	 0.3930	 0.2960
P	 0.3930	 0.2760

