



wwPDB EM Validation Summary Report ⓘ

Mar 28, 2026 – 03:05 PM UTC

PDB ID : 7UGQ / pdb_00007ugq
EMDB ID : EMD-26496
Title : Cryo-EM structure of BG24 Fabs with an inferred germline CDRL1 and 10-1074 Fabs in complex with HIV-1 Env 6405-SOSIP.664
Authors : Dam, K.A.; Bjorkman, P.J.
Deposited on : 2022-03-25
Resolution : 3.40 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

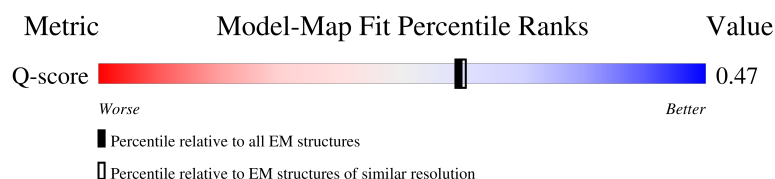
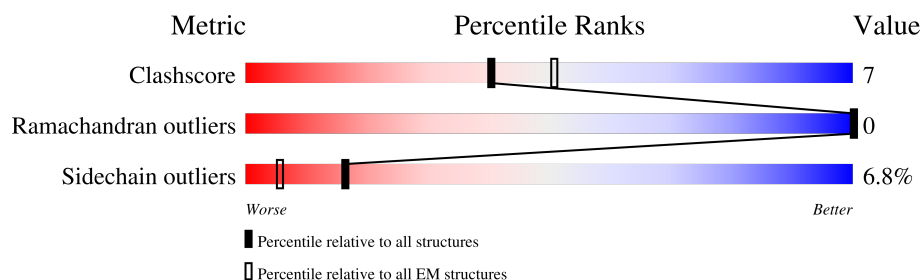
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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.



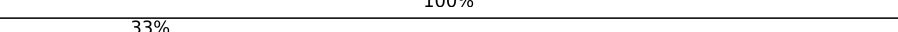
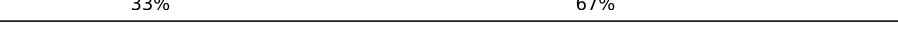
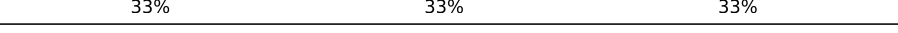
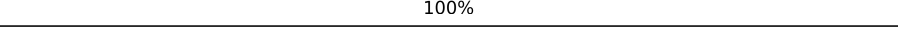
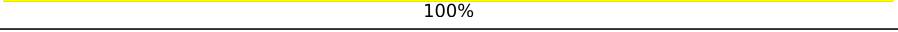
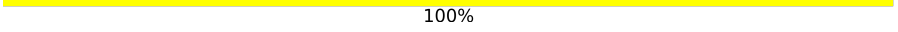
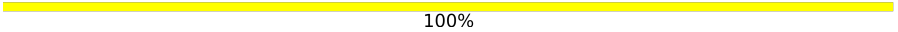
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14717 (2.90 - 3.90)

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	447	
1	B	447	
1	C	447	
2	D	129	

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Mol	Chain	Length	Quality of chain
2	E	129	 71% 26% .
2	F	129	 84% 14% .
3	G	125	 85% 14% .
3	H	125	 82% 16% .
3	I	125	 85% 15% .
4	J	105	 86% 13% .
4	K	105	 88% 11% .
4	L	105	 87% 13% .
5	M	133	 20% 78% 20% .
5	N	133	 19% 73% 26% .
5	O	133	 17% 79% 20% .
6	P	107	 8% 87% 12% .
6	Q	107	 7% 76% 21% .
6	R	107	 9% 83% 13% .
7	0	3	 33% 67%
7	S	3	 100%
7	V	3	 33% 33% 67%
7	c	3	 33% 33% 33%
7	e	3	 100%
7	h	3	 100%
7	o	3	 33% 67%
7	q	3	 100%
7	t	3	 33% 67%
8	T	2	 100%
8	U	2	 100%

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Mol	Chain	Length	Quality of chain
8	X	2	100%
8	Y	2	100%
8	Z	2	50% 50% 50%
8	a	2	50% 50%
8	b	2	100%
8	f	2	100%
8	g	2	100%
8	j	2	100%
8	k	2	100%
8	l	2	50% 50% 50%
8	m	2	50% 50%
8	n	2	100%
8	r	2	100%
8	s	2	100%
8	v	2	100%
8	w	2	100%
8	x	2	50% 50% 50%
8	y	2	100%
8	z	2	100%
9	W	5	100%
9	i	5	100%
9	u	5	100%
10	1	9	11% 89%
10	d	9	11% 89%
11	p	8	12% 62% 25%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	NAG	0	1	X	-	-	-
7	NAG	c	1	X	-	X	-
7	NAG	o	1	X	-	-	-

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 26113 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Envelope glycoprotein gp120.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	447	Total	C	N	O	S	0	0
			3498	2191	616	662	29		
1	B	447	Total	C	N	O	S	0	0
			3498	2191	616	662	29		
1	C	447	Total	C	N	O	S	0	0
			3498	2191	616	662	29		

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	ASN	ASP	conflict	UNP D7S1H2
A	?	-	GLU	deletion	UNP D7S1H2
A	?	-	THR	deletion	UNP D7S1H2
A	?	-	GLU	deletion	UNP D7S1H2
A	?	-	ARG	deletion	UNP D7S1H2
A	?	-	GLY	deletion	UNP D7S1H2
A	?	-	ALA	deletion	UNP D7S1H2
A	?	-	SER	deletion	UNP D7S1H2
A	?	-	ASN	deletion	UNP D7S1H2
A	?	-	SER	deletion	UNP D7S1H2
A	?	-	THR	deletion	UNP D7S1H2
A	?	-	GLY	deletion	UNP D7S1H2
A	?	-	THR	deletion	UNP D7S1H2
A	?	-	ASN	deletion	UNP D7S1H2
A	?	-	ASN	deletion	UNP D7S1H2
A	?	-	GLY	deletion	UNP D7S1H2
A	?	-	THR	deletion	UNP D7S1H2
A	?	-	ILE	deletion	UNP D7S1H2
A	?	-	SER	deletion	UNP D7S1H2
A	?	-	SER	deletion	UNP D7S1H2
A	?	-	THR	deletion	UNP D7S1H2
A	?	-	GLY	deletion	UNP D7S1H2
A	?	-	PRO	deletion	UNP D7S1H2
A	?	-	GLY	deletion	UNP D7S1H2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLN	deletion	UNP D7S1H2
A	?	-	ALA	deletion	UNP D7S1H2
A	?	-	ILE	deletion	UNP D7S1H2
A	?	-	TRP	deletion	UNP D7S1H2
A	496	VAL	ILE	conflict	UNP D7S1H2
A	500	ARG	LYS	conflict	UNP D7S1H2
A	501	CYS	ALA	conflict	UNP D7S1H2
A	502	LYS	ARG	conflict	UNP D7S1H2
B	33	ASN	ASP	conflict	UNP D7S1H2
B	?	-	GLU	deletion	UNP D7S1H2
B	?	-	THR	deletion	UNP D7S1H2
B	?	-	GLU	deletion	UNP D7S1H2
B	?	-	ARG	deletion	UNP D7S1H2
B	?	-	GLY	deletion	UNP D7S1H2
B	?	-	ALA	deletion	UNP D7S1H2
B	?	-	SER	deletion	UNP D7S1H2
B	?	-	ASN	deletion	UNP D7S1H2
B	?	-	SER	deletion	UNP D7S1H2
B	?	-	THR	deletion	UNP D7S1H2
B	?	-	GLY	deletion	UNP D7S1H2
B	?	-	THR	deletion	UNP D7S1H2
B	?	-	ASN	deletion	UNP D7S1H2
B	?	-	ASN	deletion	UNP D7S1H2
B	?	-	GLY	deletion	UNP D7S1H2
B	?	-	THR	deletion	UNP D7S1H2
B	?	-	ILE	deletion	UNP D7S1H2
B	?	-	SER	deletion	UNP D7S1H2
B	?	-	SER	deletion	UNP D7S1H2
B	?	-	THR	deletion	UNP D7S1H2
B	?	-	GLY	deletion	UNP D7S1H2
B	?	-	PRO	deletion	UNP D7S1H2
B	?	-	GLY	deletion	UNP D7S1H2
B	?	-	GLN	deletion	UNP D7S1H2
B	?	-	ALA	deletion	UNP D7S1H2
B	?	-	ILE	deletion	UNP D7S1H2
B	?	-	TRP	deletion	UNP D7S1H2
B	496	VAL	ILE	conflict	UNP D7S1H2
B	500	ARG	LYS	conflict	UNP D7S1H2
B	501	CYS	ALA	conflict	UNP D7S1H2
B	502	LYS	ARG	conflict	UNP D7S1H2
C	33	ASN	ASP	conflict	UNP D7S1H2
C	?	-	GLU	deletion	UNP D7S1H2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	THR	deletion	UNP D7S1H2
C	?	-	GLU	deletion	UNP D7S1H2
C	?	-	ARG	deletion	UNP D7S1H2
C	?	-	GLY	deletion	UNP D7S1H2
C	?	-	ALA	deletion	UNP D7S1H2
C	?	-	SER	deletion	UNP D7S1H2
C	?	-	ASN	deletion	UNP D7S1H2
C	?	-	SER	deletion	UNP D7S1H2
C	?	-	THR	deletion	UNP D7S1H2
C	?	-	GLY	deletion	UNP D7S1H2
C	?	-	THR	deletion	UNP D7S1H2
C	?	-	ASN	deletion	UNP D7S1H2
C	?	-	ASN	deletion	UNP D7S1H2
C	?	-	GLY	deletion	UNP D7S1H2
C	?	-	THR	deletion	UNP D7S1H2
C	?	-	ILE	deletion	UNP D7S1H2
C	?	-	SER	deletion	UNP D7S1H2
C	?	-	SER	deletion	UNP D7S1H2
C	?	-	THR	deletion	UNP D7S1H2
C	?	-	GLY	deletion	UNP D7S1H2
C	?	-	PRO	deletion	UNP D7S1H2
C	?	-	GLY	deletion	UNP D7S1H2
C	?	-	GLN	deletion	UNP D7S1H2
C	?	-	ALA	deletion	UNP D7S1H2
C	?	-	ILE	deletion	UNP D7S1H2
C	?	-	TRP	deletion	UNP D7S1H2
C	496	VAL	ILE	conflict	UNP D7S1H2
C	500	ARG	LYS	conflict	UNP D7S1H2
C	501	CYS	ALA	conflict	UNP D7S1H2
C	502	LYS	ARG	conflict	UNP D7S1H2

- Molecule 2 is a protein called Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	129	Total	C	N	O	S	0	0
			1030	655	176	193	6		
2	E	129	Total	C	N	O	S	0	0
			1030	655	176	193	6		
2	F	129	Total	C	N	O	S	0	0
			1030	655	176	193	6		

- Molecule 3 is a protein called BG24 with an inferred germline CDRL1 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	125	Total	C	N	O	S	0	0
			966	605	172	183	6		
3	H	125	Total	C	N	O	S	0	0
			966	605	172	183	6		
3	I	125	Total	C	N	O	S	0	0
			966	605	172	183	6		

- Molecule 4 is a protein called BG24 with an inferred germline CDRL1 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	J	105	Total	C	N	O	S	0	0
			758	468	135	153	2		
4	K	105	Total	C	N	O	S	0	0
			758	468	135	153	2		
4	L	105	Total	C	N	O	S	0	0
			758	468	135	153	2		

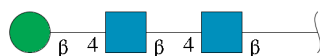
- Molecule 5 is a protein called 10-1074 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	M	133	Total	C	N	O	S	0	0
			1041	657	175	205	4		
5	N	133	Total	C	N	O	S	0	0
			1041	657	175	205	4		
5	O	133	Total	C	N	O	S	0	0
			1041	657	175	205	4		

- Molecule 6 is a protein called 10-1074 Fab light chain.

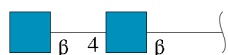
Mol	Chain	Residues	Atoms					AltConf	Trace
6	P	107	Total	C	N	O	S	0	0
			824	515	152	154	3		
6	Q	107	Total	C	N	O	S	0	0
			824	515	152	154	3		
6	R	107	Total	C	N	O	S	0	0
			824	515	152	154	3		

- Molecule 7 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
7	S	3	Total	C	N	O	0	0
			39	22	2	15		
7	V	3	Total	C	N	O	0	0
			39	22	2	15		
7	c	3	Total	C	N	O	0	0
			39	22	2	15		
7	e	3	Total	C	N	O	0	0
			39	22	2	15		
7	h	3	Total	C	N	O	0	0
			39	22	2	15		
7	o	3	Total	C	N	O	0	0
			39	22	2	15		
7	q	3	Total	C	N	O	0	0
			39	22	2	15		
7	t	3	Total	C	N	O	0	0
			39	22	2	15		
7	0	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 8 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



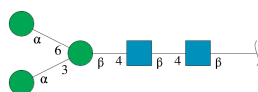
Mol	Chain	Residues	Atoms				AltConf	Trace
8	T	2	Total	C	N	O	0	0
			28	16	2	10		
8	U	2	Total	C	N	O	0	0
			28	16	2	10		
8	X	2	Total	C	N	O	0	0
			28	16	2	10		
8	Y	2	Total	C	N	O	0	0
			28	16	2	10		
8	Z	2	Total	C	N	O	0	0
			28	16	2	10		
8	a	2	Total	C	N	O	0	0
			28	16	2	10		
8	b	2	Total	C	N	O	0	0
			28	16	2	10		
8	f	2	Total	C	N	O	0	0
			28	16	2	10		

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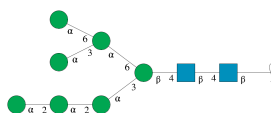
Mol	Chain	Residues	Atoms				AltConf	Trace
8	g	2	Total	C	N	O	0	0
			28	16	2	10		
8	j	2	Total	C	N	O	0	0
			28	16	2	10		
8	k	2	Total	C	N	O	0	0
			28	16	2	10		
8	l	2	Total	C	N	O	0	0
			28	16	2	10		
8	m	2	Total	C	N	O	0	0
			28	16	2	10		
8	n	2	Total	C	N	O	0	0
			28	16	2	10		
8	r	2	Total	C	N	O	0	0
			28	16	2	10		
8	s	2	Total	C	N	O	0	0
			28	16	2	10		
8	v	2	Total	C	N	O	0	0
			28	16	2	10		
8	w	2	Total	C	N	O	0	0
			28	16	2	10		
8	x	2	Total	C	N	O	0	0
			28	16	2	10		
8	y	2	Total	C	N	O	0	0
			28	16	2	10		
8	z	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



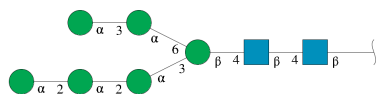
Mol	Chain	Residues	Atoms				AltConf	Trace
9	W	5	Total	C	N	O	0	0
			61	34	2	25		
9	i	5	Total	C	N	O	0	0
			61	34	2	25		
9	u	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 10 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
10	d	9	Total	C	N	O	0	0
			105	58	2	45		
10	1	9	Total	C	N	O	0	0
			105	58	2	45		

- Molecule 11 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
11	p	8	Total	C	N	O	0	0
			94	52	2	40		

- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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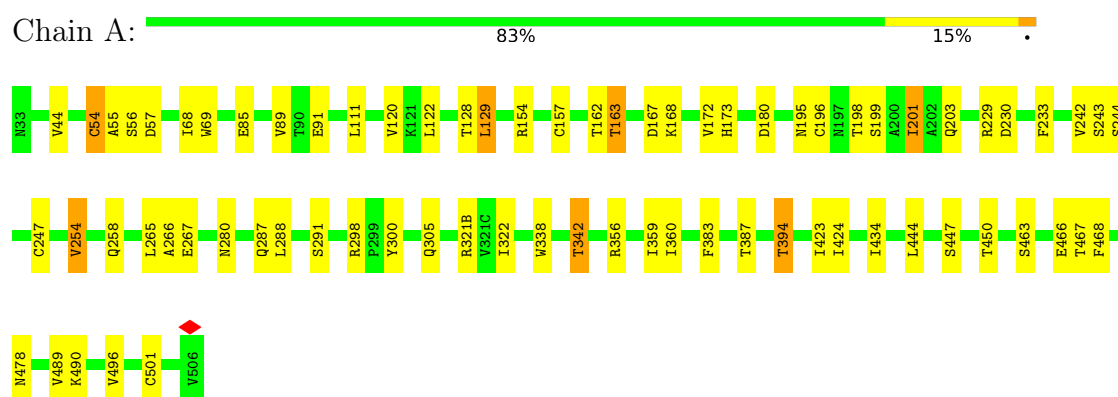
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Mol	Chain	Residues	Atoms				AltConf
12	C	1	Total 14	C 8	N 1	O 5	0
12	C	1	Total 14	C 8	N 1	O 5	0
12	C	1	Total 14	C 8	N 1	O 5	0
12	C	1	Total 14	C 8	N 1	O 5	0
12	C	1	Total 14	C 8	N 1	O 5	0
12	C	1	Total 14	C 8	N 1	O 5	0
12	C	1	Total 14	C 8	N 1	O 5	0
12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

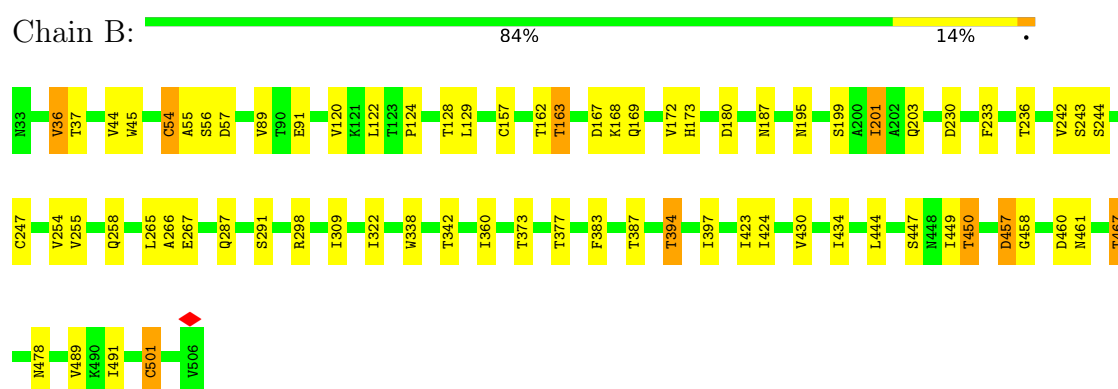
3 Residue-property plots

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.

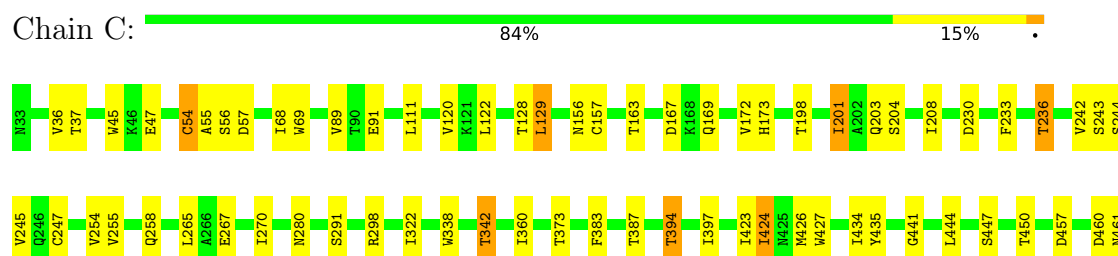
• Molecule 1: Envelope glycoprotein gp120



• Molecule 1: Envelope glycoprotein gp120

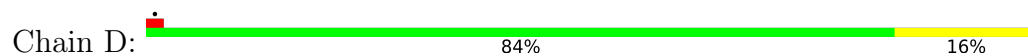


• Molecule 1: Envelope glycoprotein gp120

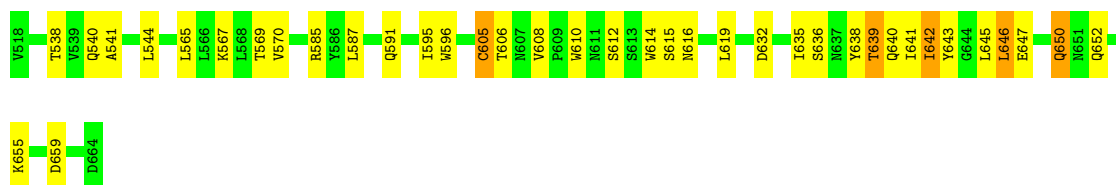




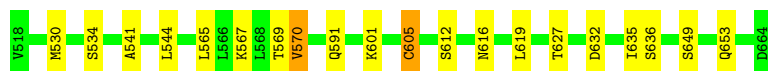
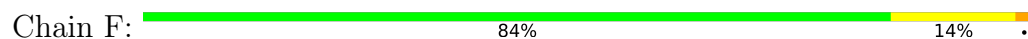
- Molecule 2: Envelope glycoprotein gp41



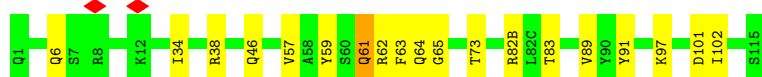
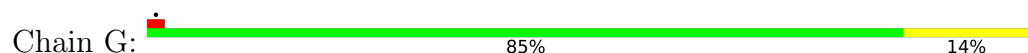
- Molecule 2: Envelope glycoprotein gp41



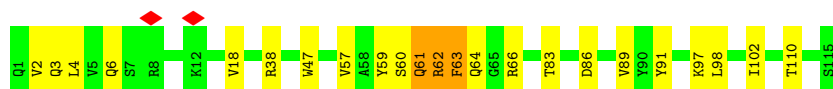
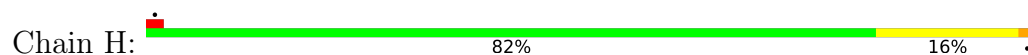
- Molecule 2: Envelope glycoprotein gp41



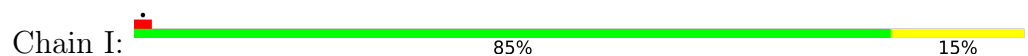
- Molecule 3: BG24 with an inferred germline CDRL1 Fab heavy chain



- Molecule 3: BG24 with an inferred germline CDRL1 Fab heavy chain

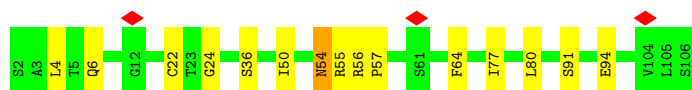


- Molecule 3: BG24 with an inferred germline CDRL1 Fab heavy chain



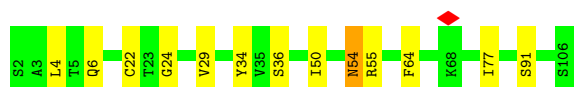
- Molecule 4: BG24 with an inferred germline CDRL1 Fab light chain

Chain J:  86% 13%



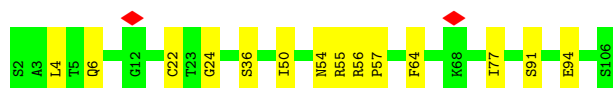
- Molecule 4: BG24 with an inferred germline CDRL1 Fab light chain

Chain K:  88% 11%



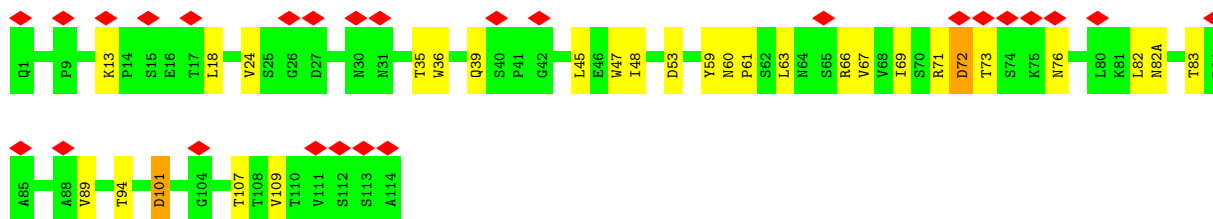
- Molecule 4: BG24 with an inferred germline CDRL1 Fab light chain

Chain L:  87% 13%



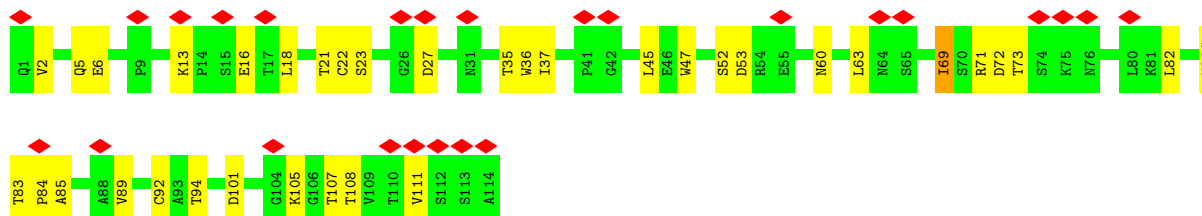
- Molecule 5: 10-1074 Fab heavy chain

Chain M:  20% 78% 20%



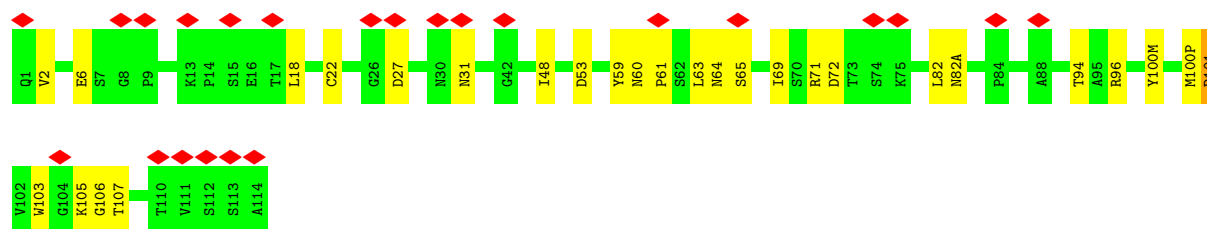
- Molecule 5: 10-1074 Fab heavy chain

Chain N:  19% 73% 26%

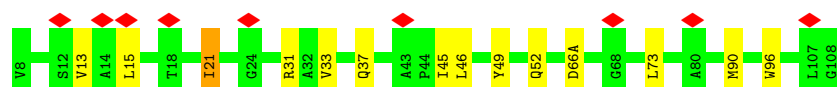
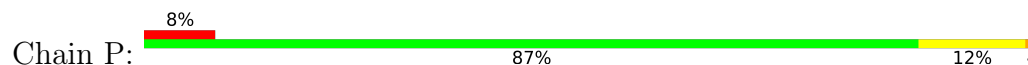


- Molecule 5: 10-1074 Fab heavy chain

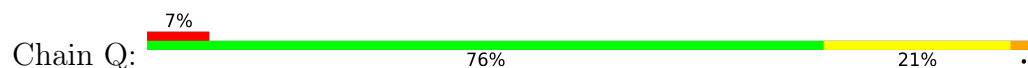
Chain O:  17% 79% 20%



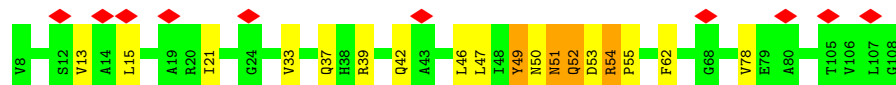
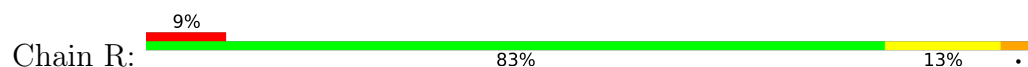
- Molecule 6: 10-1074 Fab light chain



- Molecule 6: 10-1074 Fab light chain



- Molecule 6: 10-1074 Fab light chain



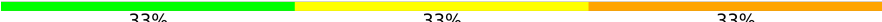
- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

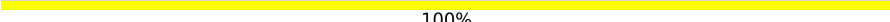


- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain c:  33% 33% 33%

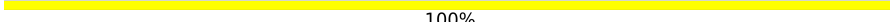
NAG1
NAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain e:  100%

NAG1
NAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain h:  100%

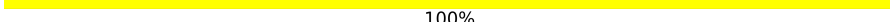
NAG1
NAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain o:  33% 67%

NAG1
NAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain q:  100%

NAG1
NAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain t:  33% 67%

NAG1
NAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 0:  33% 67%

NAG1
NAG2
BMA3

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain X:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:  50% 50%

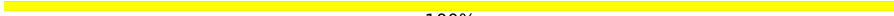

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a:  50% 50%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain f:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain g:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain j:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain k:  100%

NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain l:  50% 50%


NAG1
NAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain m:  50% 50%



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain n:  100%



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain r:  100%



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain s:  100%



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain v:  100%



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain w:  100%

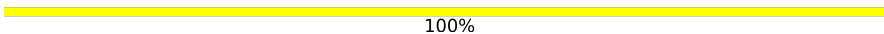


- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain x:  50% 50%

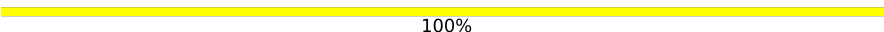


- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain y:  100%

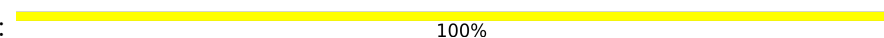
MAG1
MAG2

- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain z:  100%

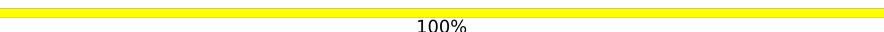
MAG1
MAG2

- Molecule 9: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:  100%

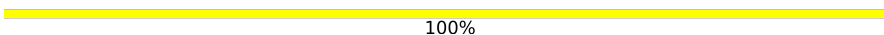
MAG1
MAG2
BMA3
MAN4
MAN5

- Molecule 9: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain i:  100%

MAG1
MAG2
BMA3
MAN4
MAN5

- Molecule 9: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain u:  100%

MAG1
MAG2
BMA3
MAN4
MAN5

- Molecule 10: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain d:  11% 89%

MAG1
MAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9

- Molecule 10: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 1:  11% 89%

MAN1
MAN2
MAN3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9

- Molecule 11: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain p:  12% 62% 25%

MAN1
MAN2
MAN3
MAN4
MAN5
MAN6
MAN7
MAN8

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	170897	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.119	Depositor
Minimum map value	-0.070	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0115	Depositor
Map size ($\text{\AA}$)	312.84, 312.84, 312.84	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.869, 0.869, 0.869	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, MAN, BMA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.26	0/3572	0.58	0/4859
1	B	0.26	0/3572	0.58	0/4859
1	C	0.26	0/3572	0.58	2/4859 (0.0%)
2	D	0.21	0/1048	0.55	0/1421
2	E	0.40	0/1048	0.65	1/1421 (0.1%)
2	F	0.22	0/1048	0.53	0/1421
3	G	0.25	0/990	0.52	0/1346
3	H	0.26	0/990	0.54	0/1346
3	I	0.23	0/990	0.49	0/1346
4	J	0.21	0/773	0.56	0/1047
4	K	0.21	0/773	0.58	0/1047
4	L	0.21	0/773	0.59	0/1047
5	M	0.21	0/1066	0.55	0/1451
5	N	0.20	0/1066	0.49	0/1451
5	O	0.22	0/1066	0.51	0/1451
6	P	0.17	0/845	0.47	0/1148
6	Q	0.21	0/845	0.51	0/1148
6	R	0.21	0/845	0.59	2/1148 (0.2%)
All	All	0.25	0/24882	0.56	5/33816 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	R	54	ARG	CA-C-N	7.73	127.66	119.85
6	R	54	ARG	C-N-CA	7.73	127.66	119.85
1	C	460	ASP	CA-C-N	5.53	132.11	121.54
1	C	460	ASP	C-N-CA	5.53	132.11	121.54
2	E	642	ILE	CB-CA-C	-5.39	105.08	111.97

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3498	0	3405	35	0
1	B	3498	0	3405	35	0
1	C	3498	0	3407	31	0
2	D	1030	0	1020	10	0
2	E	1030	0	1020	39	0
2	F	1030	0	1020	11	0
3	G	966	0	927	12	0
3	H	966	0	927	23	0
3	I	966	0	927	7	0
4	J	758	0	706	6	0
4	K	758	0	706	6	0
4	L	758	0	706	6	0
5	M	1041	0	1005	43	0
5	N	1041	0	1005	17	0
5	O	1041	0	1005	33	0
6	P	824	0	788	6	0
6	Q	824	0	788	33	0
6	R	824	0	788	21	0
7	0	39	0	34	6	0
7	S	39	0	34	0	0
7	V	39	0	34	0	0
7	c	39	0	34	7	0
7	e	39	0	34	0	0
7	h	39	0	34	0	0
7	o	39	0	32	5	0
7	q	39	0	34	0	0
7	t	39	0	34	0	0
8	T	28	0	25	0	0
8	U	28	0	25	0	0
8	X	28	0	25	0	0
8	Y	28	0	25	0	0
8	Z	28	0	25	0	0
8	a	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	b	28	0	25	0	0
8	f	28	0	25	0	0
8	g	28	0	25	0	0
8	j	28	0	25	0	0
8	k	28	0	25	1	0
8	l	28	0	25	0	0
8	m	28	0	25	0	0
8	n	28	0	25	0	0
8	r	28	0	25	0	0
8	s	28	0	25	0	0
8	v	28	0	25	0	0
8	w	28	0	25	1	0
8	x	28	0	25	0	0
8	y	28	0	25	1	0
8	z	28	0	25	0	0
9	W	61	0	52	0	0
9	i	61	0	52	0	0
9	u	61	0	52	0	0
10	l	105	0	88	0	0
10	d	105	0	88	0	0
11	p	94	0	79	6	0
12	A	98	0	91	0	0
12	B	98	0	91	1	0
12	C	98	0	91	0	0
12	D	14	0	13	0	0
12	E	14	0	13	0	0
12	F	14	0	13	0	0
All	All	26113	0	25107	360	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 7.

The worst 5 of 360 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)
2:E:596:TRP:CD1	2:E:647:GLU:HB2	1.63	1.33
5:M:59:TYR:CE1	5:M:69:ILE:HG22	1.71	1.25
2:E:596:TRP:NE1	2:E:647:GLU:HB2	1.48	1.24
5:O:59:TYR:HD2	5:O:65:SER:CB	1.58	1.14
5:M:63:LEU:HD21	5:M:67:VAL:CG1	1.79	1.12

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	439/447 (98%)	405 (92%)	34 (8%)	0	100	100
1	B	439/447 (98%)	404 (92%)	35 (8%)	0	100	100
1	C	439/447 (98%)	407 (93%)	32 (7%)	0	100	100
2	D	125/129 (97%)	116 (93%)	9 (7%)	0	100	100
2	E	125/129 (97%)	118 (94%)	7 (6%)	0	100	100
2	F	125/129 (97%)	115 (92%)	10 (8%)	0	100	100
3	G	123/125 (98%)	113 (92%)	10 (8%)	0	100	100
3	H	123/125 (98%)	112 (91%)	11 (9%)	0	100	100
3	I	123/125 (98%)	114 (93%)	9 (7%)	0	100	100
4	J	103/105 (98%)	91 (88%)	12 (12%)	0	100	100
4	K	103/105 (98%)	91 (88%)	12 (12%)	0	100	100
4	L	103/105 (98%)	91 (88%)	12 (12%)	0	100	100
5	M	131/133 (98%)	129 (98%)	2 (2%)	0	100	100
5	N	131/133 (98%)	127 (97%)	4 (3%)	0	100	100
5	O	131/133 (98%)	126 (96%)	5 (4%)	0	100	100
6	P	105/107 (98%)	99 (94%)	6 (6%)	0	100	100
6	Q	105/107 (98%)	100 (95%)	5 (5%)	0	100	100
6	R	105/107 (98%)	94 (90%)	11 (10%)	0	100	100
All	All	3078/3138 (98%)	2852 (93%)	226 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	397/397 (100%)	370 (93%)	27 (7%)	14	40
1	B	397/397 (100%)	367 (92%)	30 (8%)	12	38
1	C	397/397 (100%)	367 (92%)	30 (8%)	12	38
2	D	112/112 (100%)	105 (94%)	7 (6%)	16	42
2	E	112/112 (100%)	101 (90%)	11 (10%)	7	27
2	F	112/112 (100%)	106 (95%)	6 (5%)	20	47
3	G	102/102 (100%)	97 (95%)	5 (5%)	22	49
3	H	102/102 (100%)	94 (92%)	8 (8%)	11	37
3	I	102/102 (100%)	94 (92%)	8 (8%)	11	37
4	J	79/87 (91%)	74 (94%)	5 (6%)	16	42
4	K	79/87 (91%)	75 (95%)	4 (5%)	21	48
4	L	79/87 (91%)	76 (96%)	3 (4%)	29	54
5	M	116/116 (100%)	109 (94%)	7 (6%)	17	43
5	N	116/116 (100%)	108 (93%)	8 (7%)	14	40
5	O	116/116 (100%)	110 (95%)	6 (5%)	21	48
6	P	85/85 (100%)	81 (95%)	4 (5%)	23	50
6	Q	85/85 (100%)	80 (94%)	5 (6%)	18	44
6	R	85/85 (100%)	78 (92%)	7 (8%)	10	35
All	All	2673/2697 (99%)	2492 (93%)	181 (7%)	16	40

5 of 181 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	G	89	VAL
4	L	77	ILE
3	H	63	PHE
3	I	89	VAL
5	M	107	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 60 such sidechains are listed below:

Mol	Chain	Res	Type
3	G	43	GLN
6	Q	51	ASN
3	H	52	ASN
6	P	26	GLN
6	R	52	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

110 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	$\# Z > 2$
7	NAG	0	1	7,1	14,14,15	2.24	1 (7%)	17,19,21	1.93	2 (11%)
7	NAG	0	2	7	14,14,15	0.43	0	17,19,21	1.40	4 (23%)
7	BMA	0	3	7	11,11,12	0.23	0	15,15,17	0.82	0
10	NAG	1	1	10,1	14,14,15	0.27	0	17,19,21	0.76	0
10	NAG	1	2	10	14,14,15	0.28	0	17,19,21	0.83	1 (5%)
10	BMA	1	3	10	11,11,12	1.34	3 (27%)	15,15,17	1.78	3 (20%)
10	MAN	1	4	10	11,11,12	0.85	0	15,15,17	1.37	1 (6%)
10	MAN	1	5	10	11,11,12	0.84	0	15,15,17	1.28	1 (6%)
10	MAN	1	6	10	11,11,12	0.96	0	15,15,17	1.16	2 (13%)
10	MAN	1	7	10	11,11,12	1.59	3 (27%)	15,15,17	2.21	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	MAN	1	8	10	11,11,12	1.03	1 (9%)	15,15,17	1.28	2 (13%)
10	MAN	1	9	10	11,11,12	1.08	0	15,15,17	1.15	2 (13%)
7	NAG	S	1	7,1	14,14,15	0.57	0	17,19,21	0.83	1 (5%)
7	NAG	S	2	7	14,14,15	1.15	1 (7%)	17,19,21	2.77	3 (17%)
7	BMA	S	3	7	11,11,12	1.03	0	15,15,17	1.02	1 (6%)
8	NAG	T	1	8,1	14,14,15	1.01	1 (7%)	17,19,21	2.74	3 (17%)
8	NAG	T	2	8	14,14,15	0.62	0	17,19,21	0.64	1 (5%)
8	NAG	U	1	8,1	14,14,15	1.03	1 (7%)	17,19,21	2.78	4 (23%)
8	NAG	U	2	8	14,14,15	0.51	0	17,19,21	0.63	1 (5%)
7	NAG	V	1	7,1	14,14,15	0.41	0	17,19,21	0.77	0
7	NAG	V	2	7	14,14,15	0.78	1 (7%)	17,19,21	1.42	3 (17%)
7	BMA	V	3	7	11,11,12	1.02	1 (9%)	15,15,17	1.14	1 (6%)
9	NAG	W	1	9,1	14,14,15	0.78	0	17,19,21	2.89	4 (23%)
9	NAG	W	2	9	14,14,15	0.46	0	17,19,21	0.64	1 (5%)
9	BMA	W	3	9	11,11,12	0.89	0	15,15,17	1.06	1 (6%)
9	MAN	W	4	9	11,11,12	1.27	2 (18%)	15,15,17	1.74	1 (6%)
9	MAN	W	5	9	11,11,12	1.13	0	15,15,17	1.02	2 (13%)
8	NAG	X	1	8,1	14,14,15	0.39	0	17,19,21	0.68	1 (5%)
8	NAG	X	2	8	14,14,15	1.11	1 (7%)	17,19,21	2.76	3 (17%)
8	NAG	Y	1	8,1	14,14,15	0.89	1 (7%)	17,19,21	1.29	3 (17%)
8	NAG	Y	2	8	14,14,15	0.53	0	17,19,21	0.61	1 (5%)
8	NAG	Z	1	8,1	14,14,15	0.33	0	17,19,21	0.64	0
8	NAG	Z	2	8	14,14,15	0.55	0	17,19,21	0.87	2 (11%)
8	NAG	a	1	8,1	14,14,15	0.47	0	17,19,21	0.62	0
8	NAG	a	2	8	14,14,15	0.63	0	17,19,21	0.61	1 (5%)
8	NAG	b	1	8,1	14,14,15	0.39	0	17,19,21	0.67	1 (5%)
8	NAG	b	2	8	14,14,15	0.55	0	17,19,21	0.91	2 (11%)
7	NAG	c	1	7,1	14,14,15	2.31	1 (7%)	17,19,21	2.33	5 (29%)
7	NAG	c	2	7	14,14,15	0.40	0	17,19,21	1.48	3 (17%)
7	BMA	c	3	7	11,11,12	0.23	0	15,15,17	0.84	0
10	NAG	d	1	10,1	14,14,15	0.28	0	17,19,21	0.78	0
10	NAG	d	2	10	14,14,15	0.27	0	17,19,21	0.84	1 (5%)
10	BMA	d	3	10	11,11,12	1.30	2 (18%)	15,15,17	1.83	4 (26%)
10	MAN	d	4	10	11,11,12	0.84	0	15,15,17	1.42	1 (6%)
10	MAN	d	5	10	11,11,12	0.86	0	15,15,17	1.26	1 (6%)
10	MAN	d	6	10	11,11,12	0.90	0	15,15,17	1.16	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	MAN	d	7	10	11,11,12	1.50	3 (27%)	15,15,17	2.10	2 (13%)
10	MAN	d	8	10	11,11,12	1.04	1 (9%)	15,15,17	1.30	2 (13%)
10	MAN	d	9	10	11,11,12	1.05	0	15,15,17	1.17	2 (13%)
7	NAG	e	1	7,1	14,14,15	0.58	0	17,19,21	0.82	1 (5%)
7	NAG	e	2	7	14,14,15	1.13	1 (7%)	17,19,21	2.76	3 (17%)
7	BMA	e	3	7	11,11,12	1.02	0	15,15,17	1.04	1 (6%)
8	NAG	f	1	8,1	14,14,15	1.01	1 (7%)	17,19,21	2.75	3 (17%)
8	NAG	f	2	8	14,14,15	0.61	0	17,19,21	0.63	1 (5%)
8	NAG	g	1	8,1	14,14,15	1.03	1 (7%)	17,19,21	2.78	4 (23%)
8	NAG	g	2	8	14,14,15	0.49	0	17,19,21	0.83	2 (11%)
7	NAG	h	1	7,1	14,14,15	0.33	0	17,19,21	1.08	1 (5%)
7	NAG	h	2	7	14,14,15	0.68	0	17,19,21	0.75	1 (5%)
7	BMA	h	3	7	11,11,12	1.01	1 (9%)	15,15,17	1.14	1 (6%)
9	NAG	i	1	9,1	14,14,15	0.83	1 (7%)	17,19,21	2.87	4 (23%)
9	NAG	i	2	9	14,14,15	0.48	0	17,19,21	0.65	1 (5%)
9	BMA	i	3	9	11,11,12	0.85	0	15,15,17	1.09	1 (6%)
9	MAN	i	4	9	11,11,12	1.18	1 (9%)	15,15,17	1.77	1 (6%)
9	MAN	i	5	9	11,11,12	1.02	0	15,15,17	1.09	2 (13%)
8	NAG	j	1	8,1	14,14,15	0.39	0	17,19,21	0.68	1 (5%)
8	NAG	j	2	8	14,14,15	1.11	1 (7%)	17,19,21	2.75	3 (17%)
8	NAG	k	1	8,1	14,14,15	0.99	1 (7%)	17,19,21	1.39	4 (23%)
8	NAG	k	2	8	14,14,15	0.60	0	17,19,21	0.56	0
8	NAG	l	1	8,1	14,14,15	0.33	0	17,19,21	0.62	0
8	NAG	l	2	8	14,14,15	0.54	0	17,19,21	0.87	2 (11%)
8	NAG	m	1	8,1	14,14,15	0.31	0	17,19,21	0.82	0
8	NAG	m	2	8	14,14,15	0.72	1 (7%)	17,19,21	0.53	0
8	NAG	n	1	8,1	14,14,15	0.42	0	17,19,21	0.69	1 (5%)
8	NAG	n	2	8	14,14,15	0.54	0	17,19,21	0.89	2 (11%)
7	NAG	o	1	7,1	14,14,15	2.30	1 (7%)	17,19,21	1.91	2 (11%)
7	NAG	o	2	7	14,14,15	0.45	0	17,19,21	1.75	4 (23%)
7	BMA	o	3	7	11,11,12	0.23	0	15,15,17	0.81	0
11	NAG	p	1	11,1	14,14,15	0.28	0	17,19,21	0.79	0
11	NAG	p	2	11	14,14,15	0.28	0	17,19,21	0.86	2 (11%)
11	BMA	p	3	11	11,11,12	1.40	3 (27%)	15,15,17	1.86	4 (26%)
11	MAN	p	4	11	11,11,12	0.89	0	15,15,17	1.32	1 (6%)
11	MAN	p	5	11	11,11,12	0.82	0	15,15,17	1.24	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	MAN	p	6	11	11,11,12	0.97	0	15,15,17	1.16	2 (13%)
11	MAN	p	7	11	11,11,12	1.35	3 (27%)	15,15,17	1.92	2 (13%)
11	MAN	p	8	11	11,11,12	1.04	1 (9%)	15,15,17	1.31	2 (13%)
7	NAG	q	1	7,1	14,14,15	0.60	0	17,19,21	0.82	1 (5%)
7	NAG	q	2	7	14,14,15	1.15	1 (7%)	17,19,21	2.77	3 (17%)
7	BMA	q	3	7	11,11,12	1.03	0	15,15,17	1.02	1 (6%)
8	NAG	r	1	8,1	14,14,15	1.02	1 (7%)	17,19,21	2.75	3 (17%)
8	NAG	r	2	8	14,14,15	0.61	0	17,19,21	0.65	1 (5%)
8	NAG	s	1	8,1	14,14,15	1.04	1 (7%)	17,19,21	2.78	4 (23%)
8	NAG	s	2	8	14,14,15	0.52	0	17,19,21	0.64	1 (5%)
7	NAG	t	1	7,1	14,14,15	0.39	0	17,19,21	0.88	1 (5%)
7	NAG	t	2	7	14,14,15	0.50	0	17,19,21	0.76	0
7	BMA	t	3	7	11,11,12	1.04	1 (9%)	15,15,17	1.11	1 (6%)
9	NAG	u	1	9,1	14,14,15	0.79	1 (7%)	17,19,21	2.89	4 (23%)
9	NAG	u	2	9	14,14,15	0.48	0	17,19,21	0.65	1 (5%)
9	BMA	u	3	9	11,11,12	0.98	0	15,15,17	0.93	1 (6%)
9	MAN	u	4	9	11,11,12	1.11	1 (9%)	15,15,17	1.59	2 (13%)
9	MAN	u	5	9	11,11,12	0.99	0	15,15,17	1.16	2 (13%)
8	NAG	v	1	8,1	14,14,15	0.40	0	17,19,21	0.66	1 (5%)
8	NAG	v	2	8	14,14,15	1.12	1 (7%)	17,19,21	2.76	3 (17%)
8	NAG	w	1	8,1	14,14,15	0.99	1 (7%)	17,19,21	1.38	4 (23%)
8	NAG	w	2	8	14,14,15	0.57	0	17,19,21	0.57	0
8	NAG	x	1	8,1	14,14,15	0.34	0	17,19,21	0.64	0
8	NAG	x	2	8	14,14,15	0.55	0	17,19,21	0.87	2 (11%)
8	NAG	y	1	8,1	14,14,15	0.47	0	17,19,21	0.61	0
8	NAG	y	2	8	14,14,15	0.63	0	17,19,21	0.60	1 (5%)
8	NAG	z	1	8,1	14,14,15	0.42	0	17,19,21	0.69	1 (5%)
8	NAG	z	2	8	14,14,15	0.53	0	17,19,21	0.90	1 (5%)

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
7	NAG	0	1	7,1	1/1/5/7	6/6/23/26	0/1/1/1
7	NAG	0	2	7	-	5/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	BMA	0	3	7	-	1/2/19/22	0/1/1/1
10	NAG	1	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	1	2	10	-	0/6/23/26	0/1/1/1
10	BMA	1	3	10	-	1/2/19/22	0/1/1/1
10	MAN	1	4	10	-	2/2/19/22	0/1/1/1
10	MAN	1	5	10	-	2/2/19/22	0/1/1/1
10	MAN	1	6	10	-	0/2/19/22	0/1/1/1
10	MAN	1	7	10	-	2/2/19/22	0/1/1/1
10	MAN	1	8	10	-	0/2/19/22	0/1/1/1
10	MAN	1	9	10	-	0/2/19/22	0/1/1/1
7	NAG	S	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	S	2	7	-	6/6/23/26	0/1/1/1
7	BMA	S	3	7	-	0/2/19/22	0/1/1/1
8	NAG	T	1	8,1	-	4/6/23/26	0/1/1/1
8	NAG	T	2	8	-	2/6/23/26	0/1/1/1
8	NAG	U	1	8,1	-	6/6/23/26	0/1/1/1
8	NAG	U	2	8	-	0/6/23/26	0/1/1/1
7	NAG	V	1	7,1	-	2/6/23/26	0/1/1/1
7	NAG	V	2	7	-	4/6/23/26	0/1/1/1
7	BMA	V	3	7	-	2/2/19/22	0/1/1/1
9	NAG	W	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	W	2	9	-	0/6/23/26	0/1/1/1
9	BMA	W	3	9	-	0/2/19/22	0/1/1/1
9	MAN	W	4	9	-	0/2/19/22	0/1/1/1
9	MAN	W	5	9	-	1/2/19/22	0/1/1/1
8	NAG	X	1	8,1	-	1/6/23/26	0/1/1/1
8	NAG	X	2	8	-	6/6/23/26	0/1/1/1
8	NAG	Y	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	Y	2	8	-	2/6/23/26	0/1/1/1
8	NAG	Z	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	Z	2	8	-	4/6/23/26	0/1/1/1
8	NAG	a	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	a	2	8	-	0/6/23/26	0/1/1/1
8	NAG	b	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	b	2	8	-	2/6/23/26	0/1/1/1
7	NAG	c	1	7,1	1/1/5/7	3/6/23/26	0/1/1/1
7	NAG	c	2	7	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	BMA	c	3	7	-	2/2/19/22	0/1/1/1
10	NAG	d	1	10,1	-	1/6/23/26	0/1/1/1
10	NAG	d	2	10	-	0/6/23/26	0/1/1/1
10	BMA	d	3	10	-	1/2/19/22	0/1/1/1
10	MAN	d	4	10	-	0/2/19/22	0/1/1/1
10	MAN	d	5	10	-	2/2/19/22	0/1/1/1
10	MAN	d	6	10	-	0/2/19/22	0/1/1/1
10	MAN	d	7	10	-	2/2/19/22	0/1/1/1
10	MAN	d	8	10	-	0/2/19/22	0/1/1/1
10	MAN	d	9	10	-	0/2/19/22	0/1/1/1
7	NAG	e	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	e	2	7	-	6/6/23/26	0/1/1/1
7	BMA	e	3	7	-	0/2/19/22	0/1/1/1
8	NAG	f	1	8,1	-	4/6/23/26	0/1/1/1
8	NAG	f	2	8	-	2/6/23/26	0/1/1/1
8	NAG	g	1	8,1	-	6/6/23/26	0/1/1/1
8	NAG	g	2	8	-	2/6/23/26	0/1/1/1
7	NAG	h	1	7,1	-	4/6/23/26	0/1/1/1
7	NAG	h	2	7	-	2/6/23/26	0/1/1/1
7	BMA	h	3	7	-	2/2/19/22	0/1/1/1
9	NAG	i	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	i	2	9	-	1/6/23/26	0/1/1/1
9	BMA	i	3	9	-	0/2/19/22	0/1/1/1
9	MAN	i	4	9	-	0/2/19/22	0/1/1/1
9	MAN	i	5	9	-	0/2/19/22	0/1/1/1
8	NAG	j	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	j	2	8	-	6/6/23/26	0/1/1/1
8	NAG	k	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	k	2	8	-	2/6/23/26	0/1/1/1
8	NAG	l	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	l	2	8	-	4/6/23/26	0/1/1/1
8	NAG	m	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	m	2	8	-	1/6/23/26	0/1/1/1
8	NAG	n	1	8,1	-	1/6/23/26	0/1/1/1
8	NAG	n	2	8	-	2/6/23/26	0/1/1/1
7	NAG	o	1	7,1	1/1/5/7	4/6/23/26	0/1/1/1
7	NAG	o	2	7	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	BMA	o	3	7	-	2/2/19/22	0/1/1/1
11	NAG	p	1	11,1	-	2/6/23/26	0/1/1/1
11	NAG	p	2	11	-	0/6/23/26	0/1/1/1
11	BMA	p	3	11	-	1/2/19/22	0/1/1/1
11	MAN	p	4	11	-	0/2/19/22	0/1/1/1
11	MAN	p	5	11	-	2/2/19/22	0/1/1/1
11	MAN	p	6	11	-	0/2/19/22	0/1/1/1
11	MAN	p	7	11	-	2/2/19/22	0/1/1/1
11	MAN	p	8	11	-	1/2/19/22	0/1/1/1
7	NAG	q	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	q	2	7	-	6/6/23/26	0/1/1/1
7	BMA	q	3	7	-	0/2/19/22	0/1/1/1
8	NAG	r	1	8,1	-	4/6/23/26	0/1/1/1
8	NAG	r	2	8	-	2/6/23/26	0/1/1/1
8	NAG	s	1	8,1	-	6/6/23/26	0/1/1/1
8	NAG	s	2	8	-	0/6/23/26	0/1/1/1
7	NAG	t	1	7,1	-	4/6/23/26	0/1/1/1
7	NAG	t	2	7	-	2/6/23/26	0/1/1/1
7	BMA	t	3	7	-	2/2/19/22	0/1/1/1
9	NAG	u	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	u	2	9	-	2/6/23/26	0/1/1/1
9	BMA	u	3	9	-	2/2/19/22	0/1/1/1
9	MAN	u	4	9	-	0/2/19/22	0/1/1/1
9	MAN	u	5	9	-	0/2/19/22	0/1/1/1
8	NAG	v	1	8,1	-	1/6/23/26	0/1/1/1
8	NAG	v	2	8	-	6/6/23/26	0/1/1/1
8	NAG	w	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	w	2	8	-	2/6/23/26	0/1/1/1
8	NAG	x	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	x	2	8	-	4/6/23/26	0/1/1/1
8	NAG	y	1	8,1	-	1/6/23/26	0/1/1/1
8	NAG	y	2	8	-	0/6/23/26	0/1/1/1
8	NAG	z	1	8,1	-	1/6/23/26	0/1/1/1
8	NAG	z	2	8	-	2/6/23/26	0/1/1/1

The worst 5 of 49 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	c	1	NAG	C1-C2	8.57	1.64	1.52
7	o	1	NAG	C1-C2	8.50	1.63	1.52
7	0	1	NAG	C1-C2	8.31	1.63	1.52
10	1	7	MAN	O5-C5	3.34	1.49	1.43
7	q	2	NAG	C1-C2	3.24	1.56	1.52

The worst 5 of 184 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	r	1	NAG	C2-N2-C7	9.67	135.86	122.90
8	f	1	NAG	C2-N2-C7	9.66	135.84	122.90
9	W	1	NAG	C2-N2-C7	9.66	135.84	122.90
9	u	1	NAG	C2-N2-C7	9.65	135.84	122.90
9	i	1	NAG	C2-N2-C7	9.65	135.83	122.90

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	c	1	NAG	C1
7	o	1	NAG	C1
7	0	1	NAG	C1

5 of 208 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	c	1	NAG	C8-C7-N2-C2
7	c	1	NAG	O7-C7-N2-C2
7	c	2	NAG	C8-C7-N2-C2
7	c	2	NAG	O7-C7-N2-C2
7	o	1	NAG	C8-C7-N2-C2

There are no ring outliers.

10 monomers are involved in 27 short contacts:

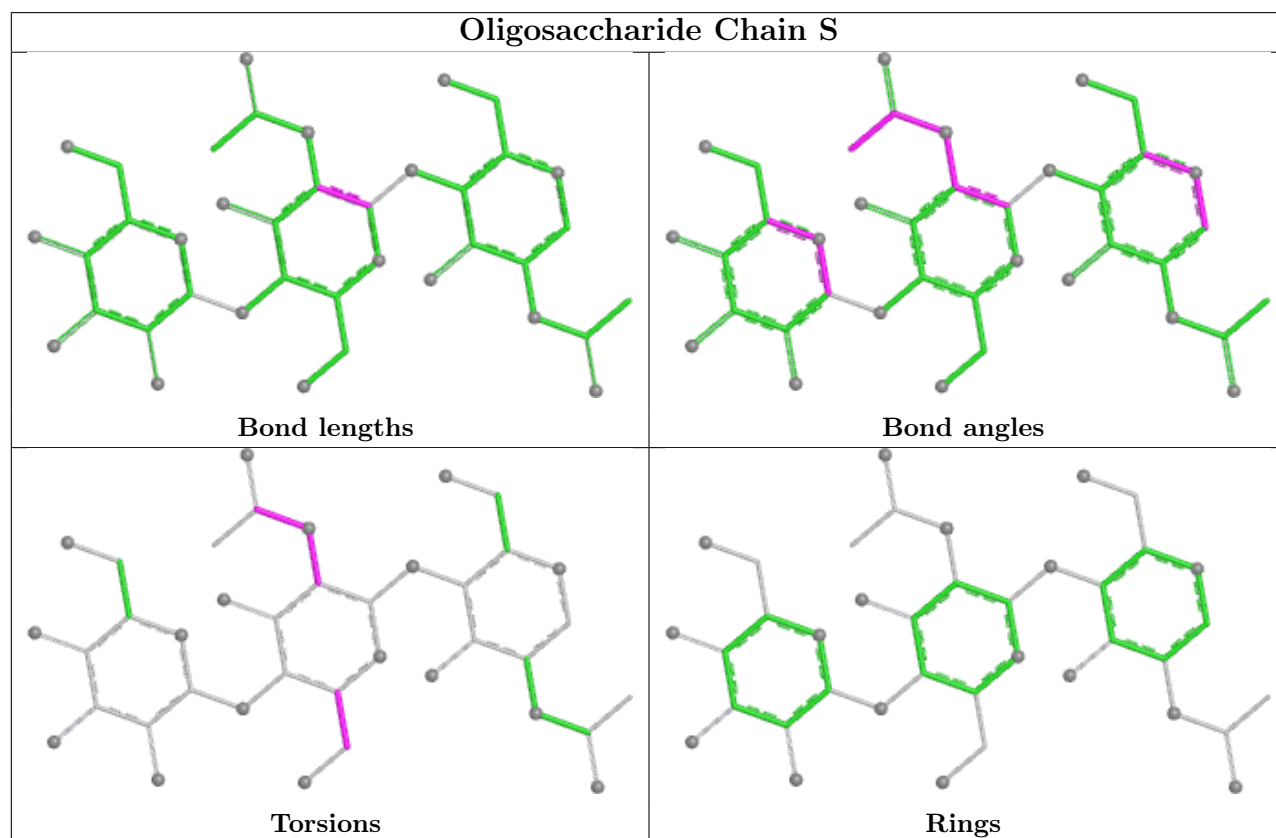
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	o	1	NAG	4	0
7	c	1	NAG	7	0
11	p	5	MAN	1	0
8	k	2	NAG	1	0
8	w	2	NAG	1	0
11	p	6	MAN	5	0
8	y	1	NAG	1	0

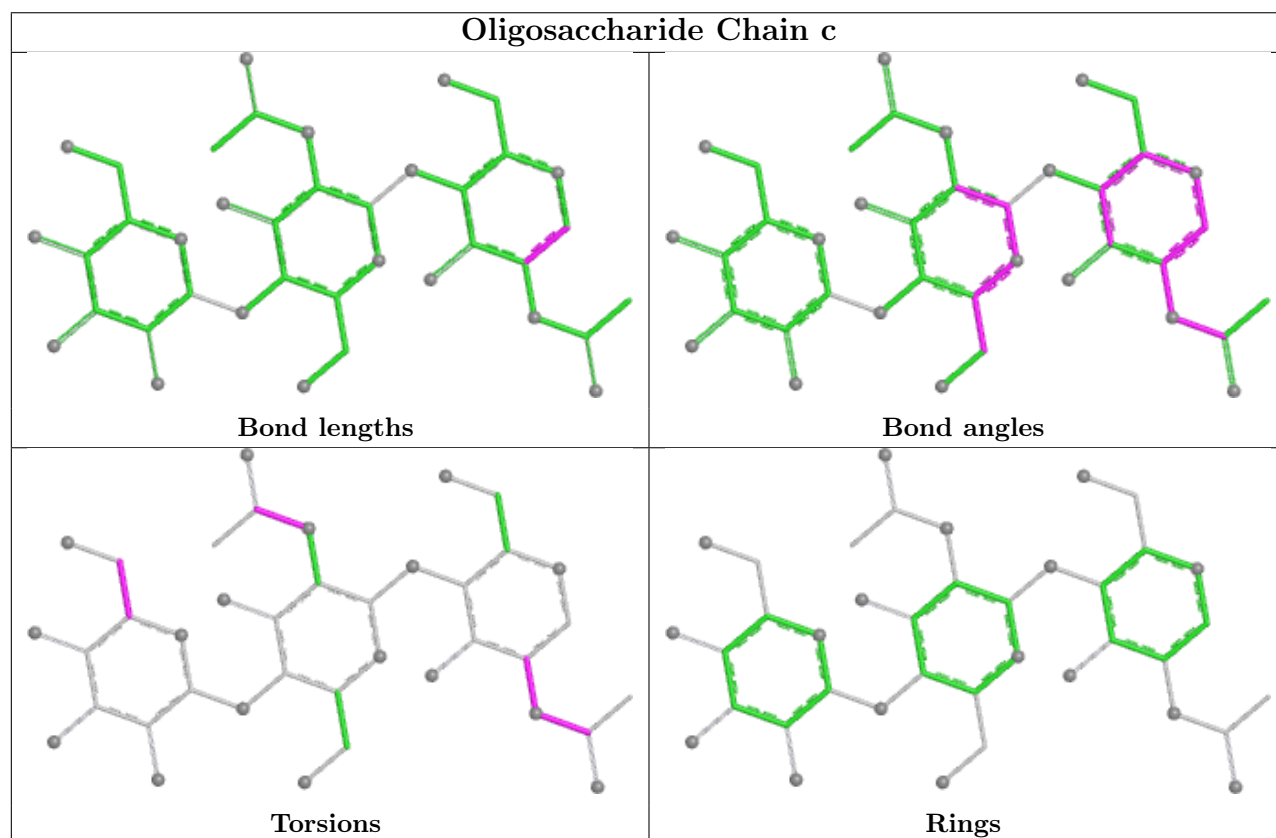
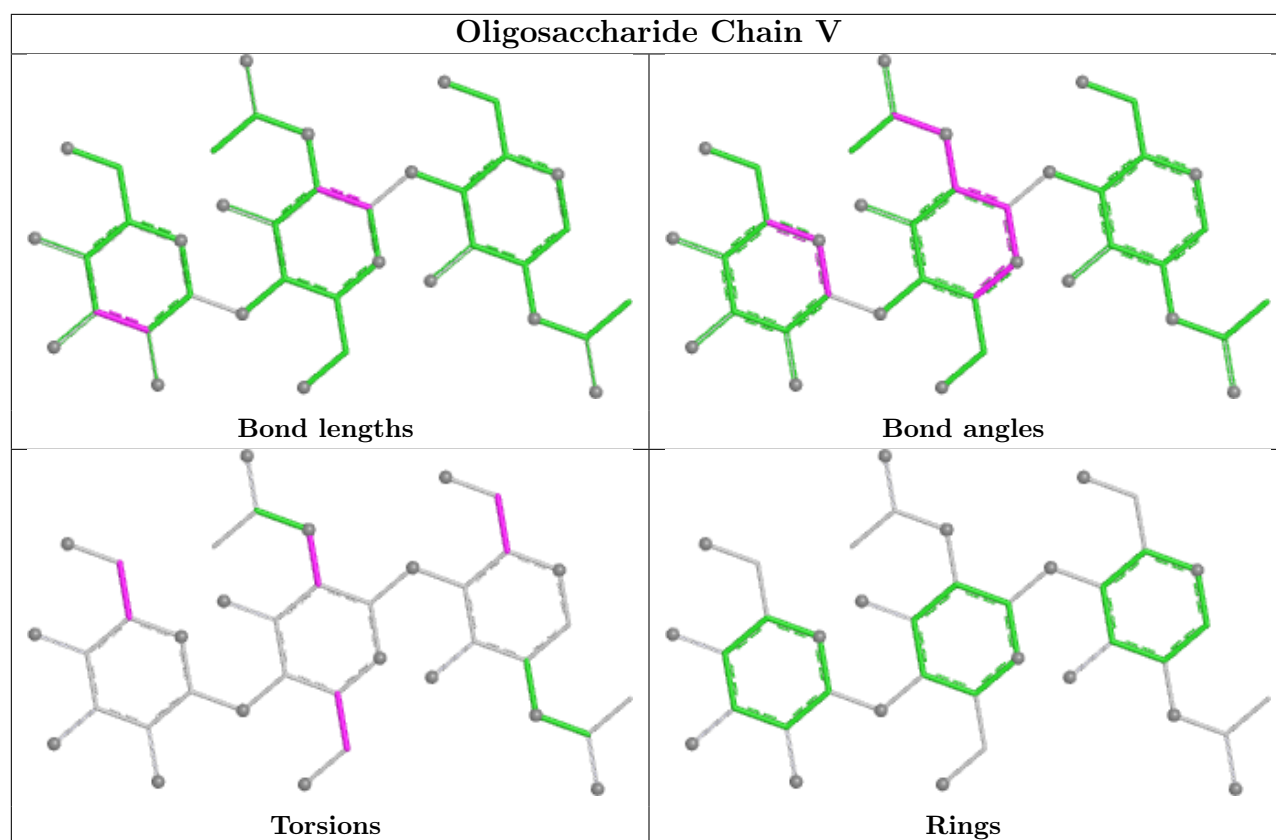
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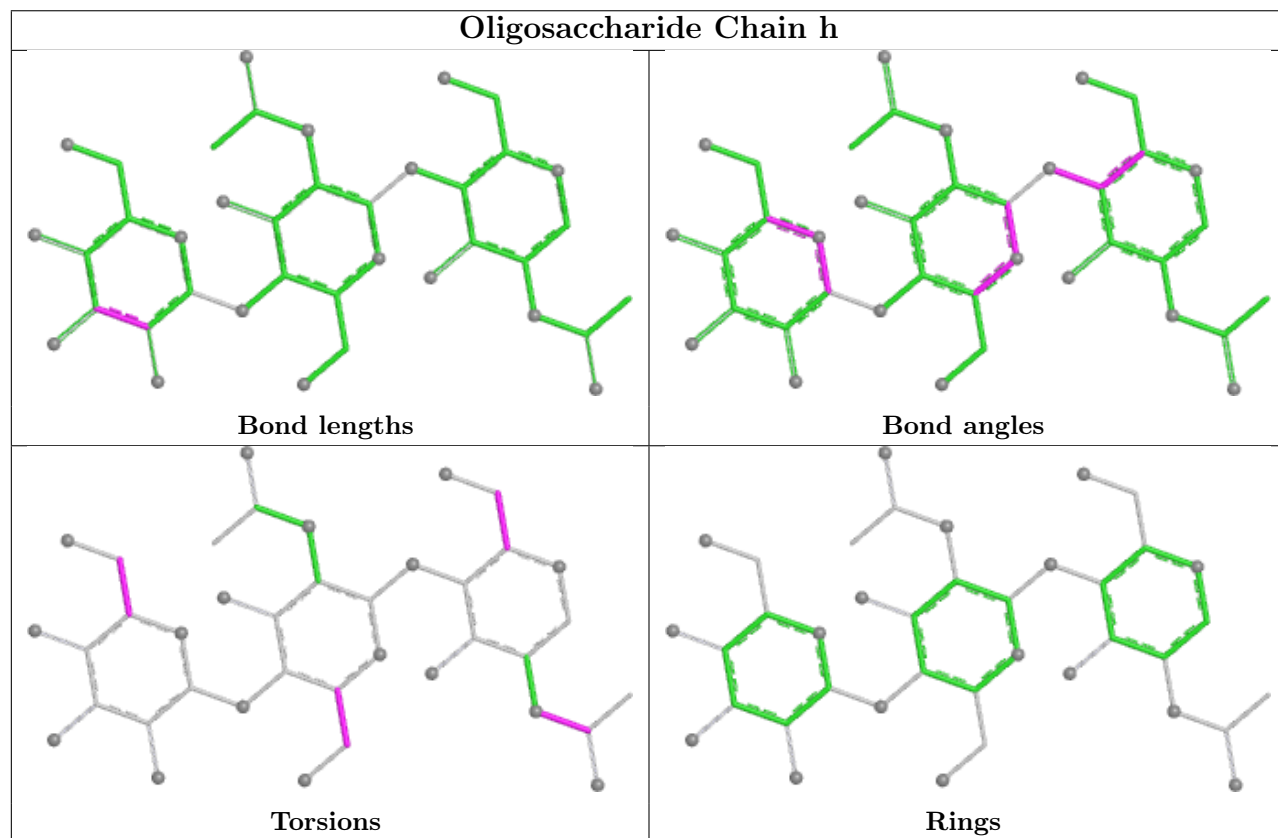
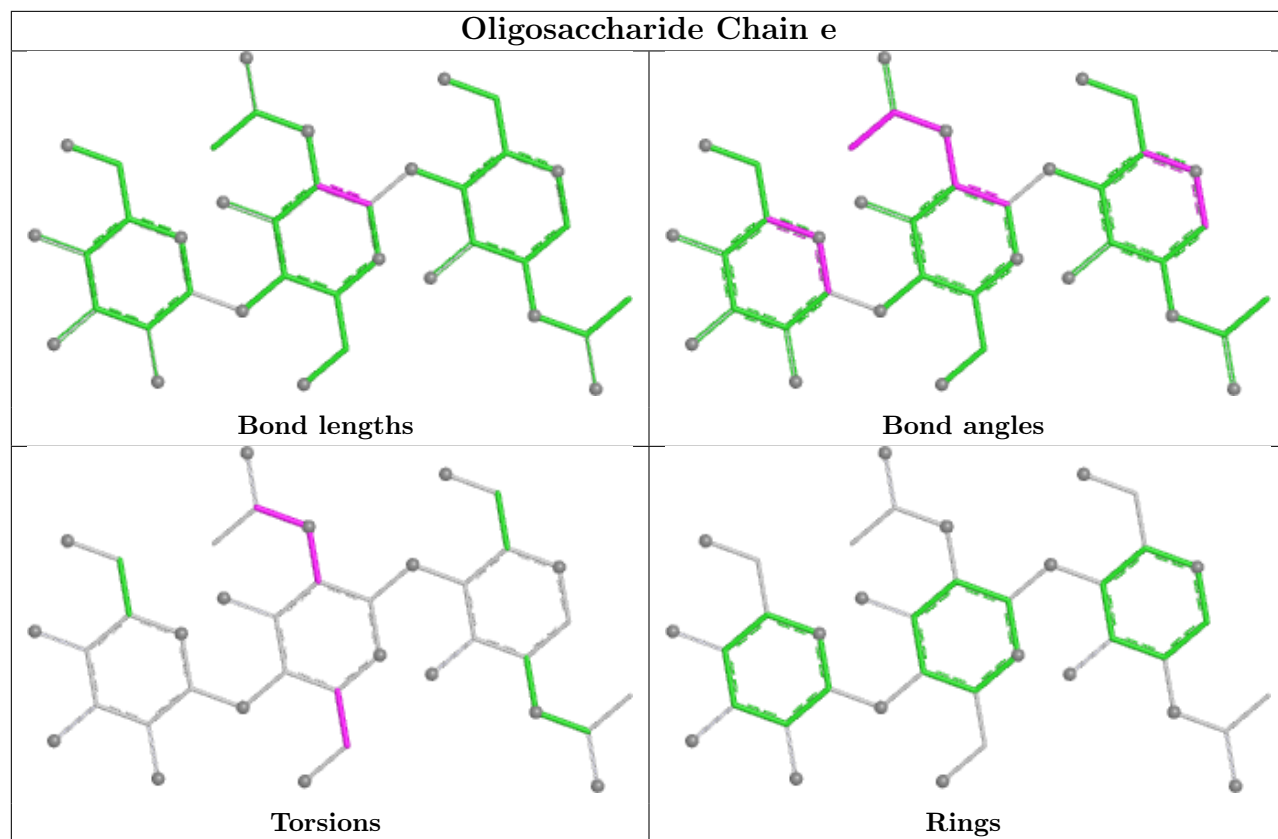
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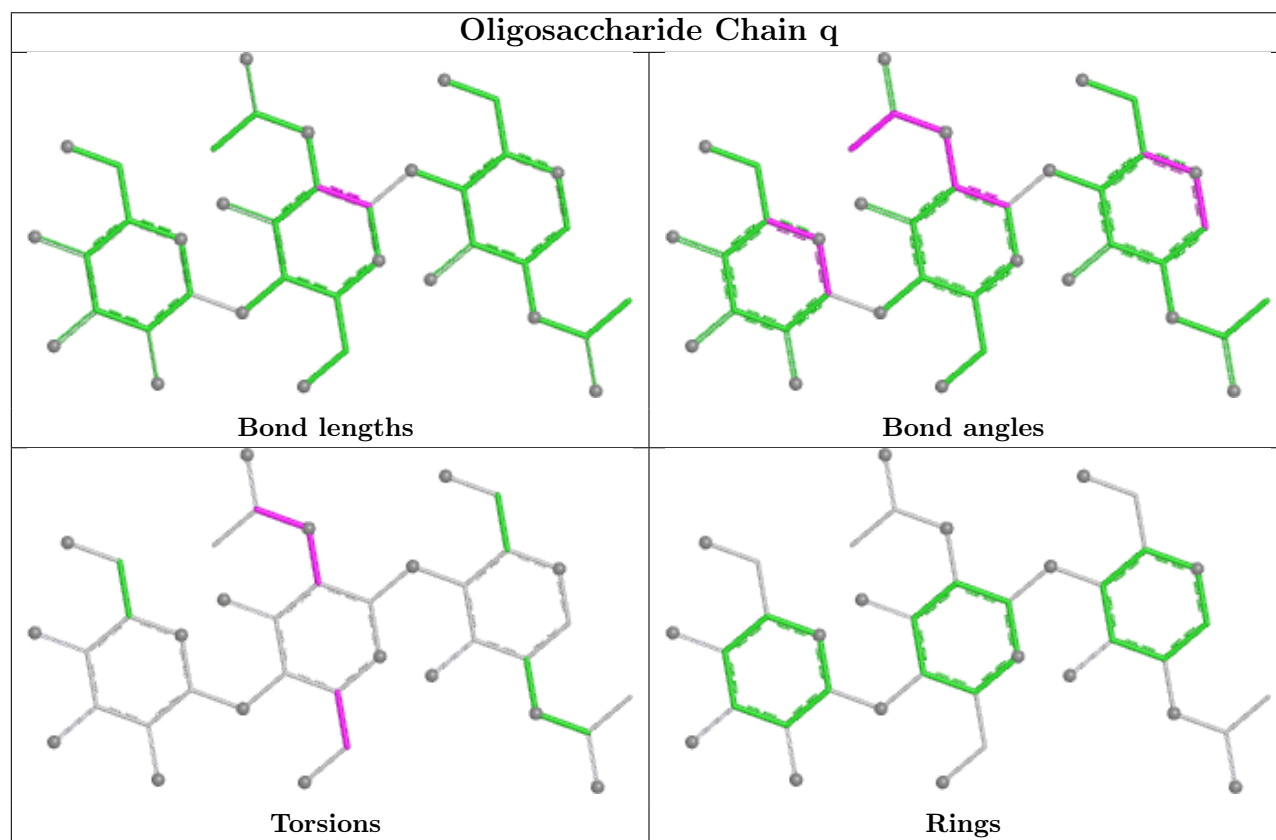
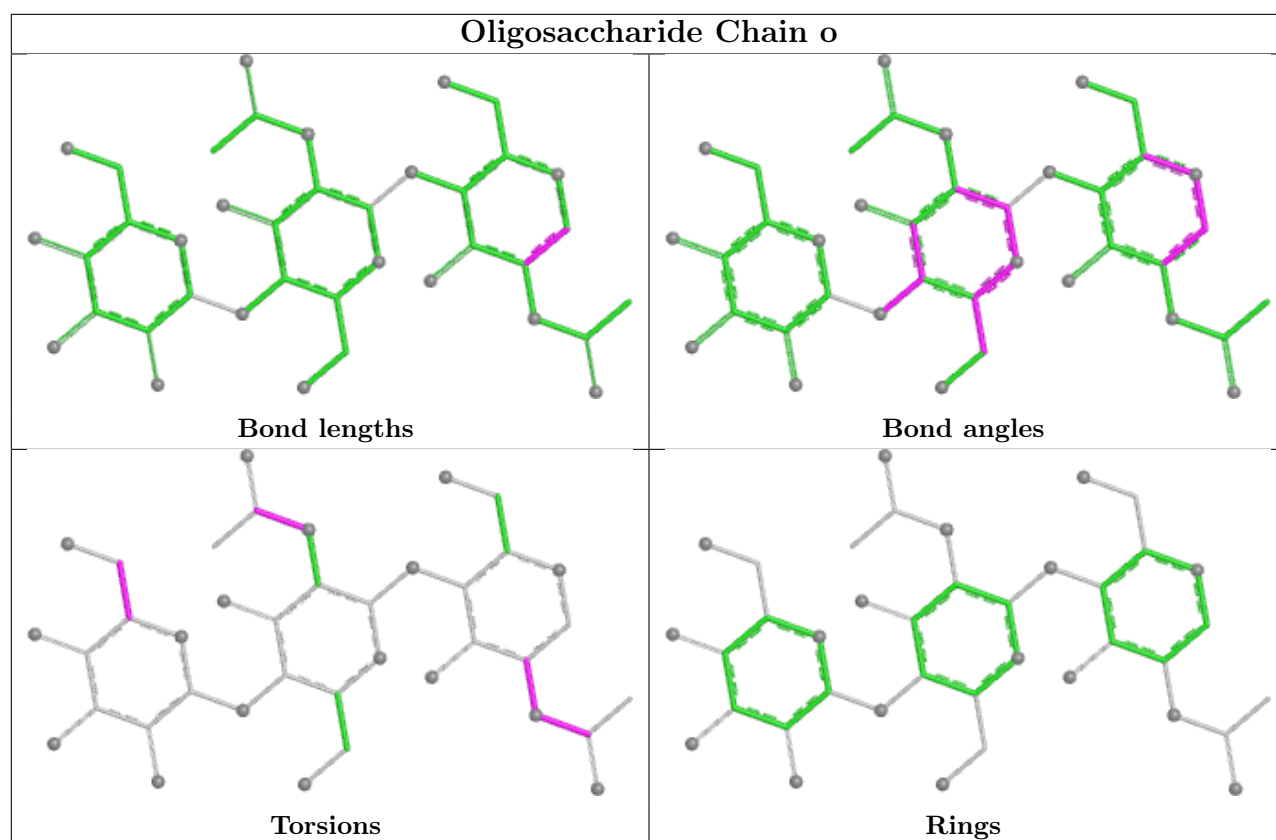
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	o	2	NAG	1	0
7	0	1	NAG	6	0
7	0	2	NAG	1	0

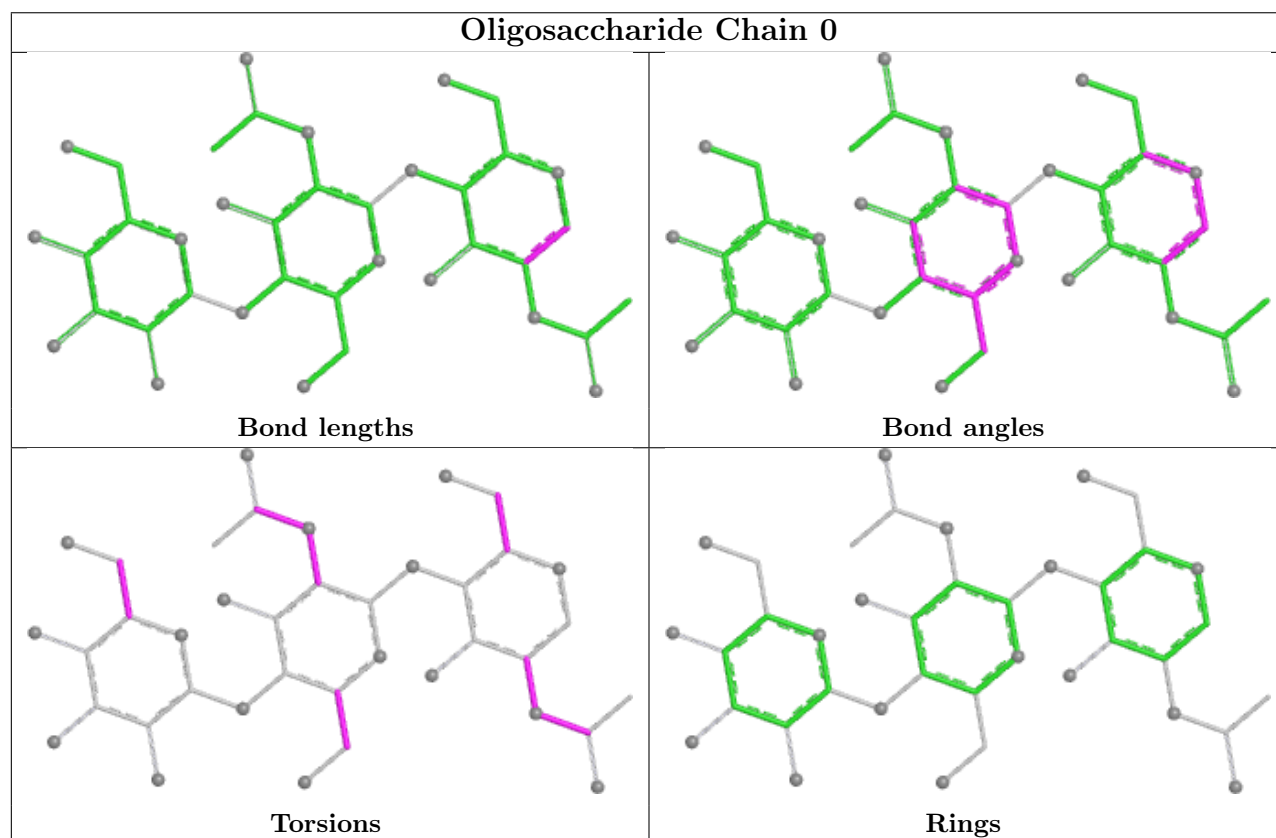
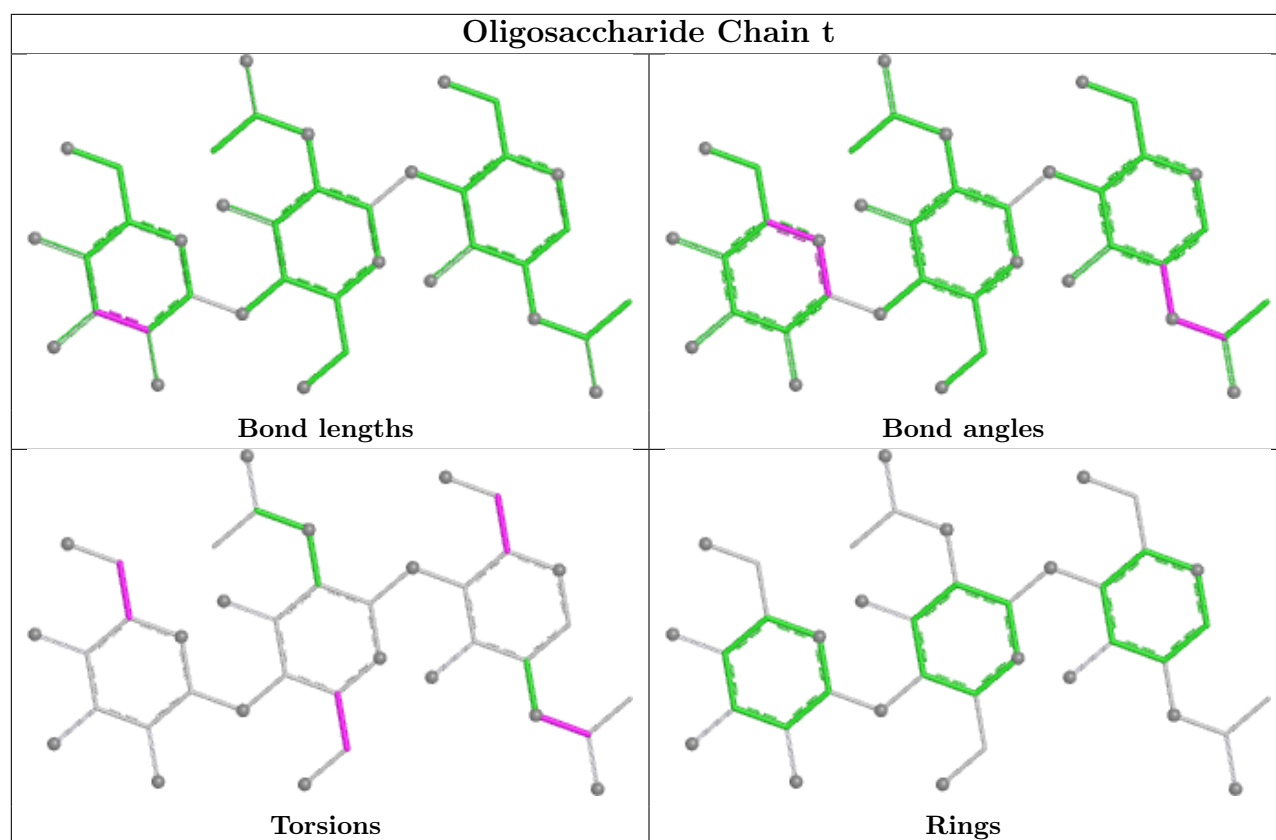
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

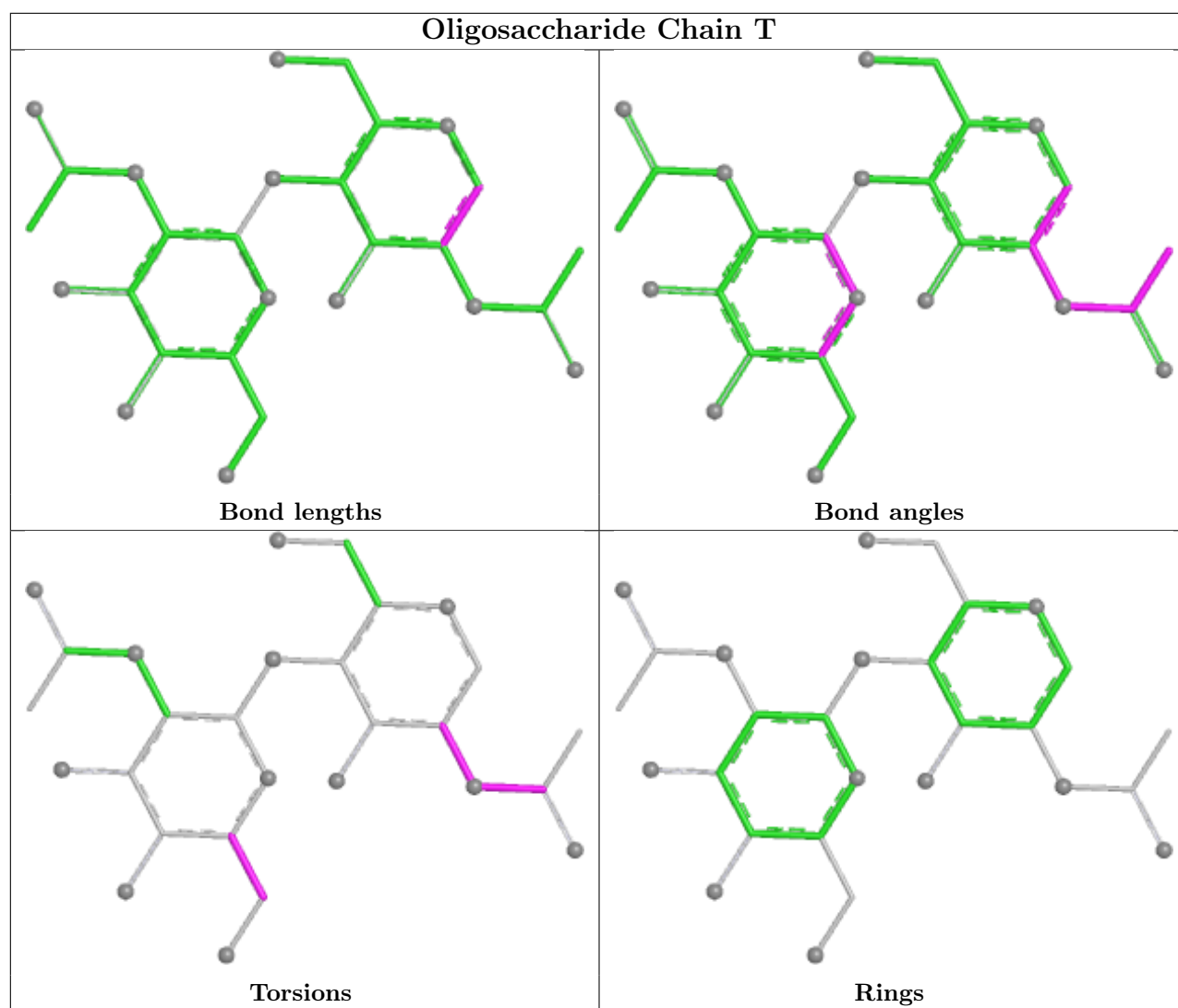


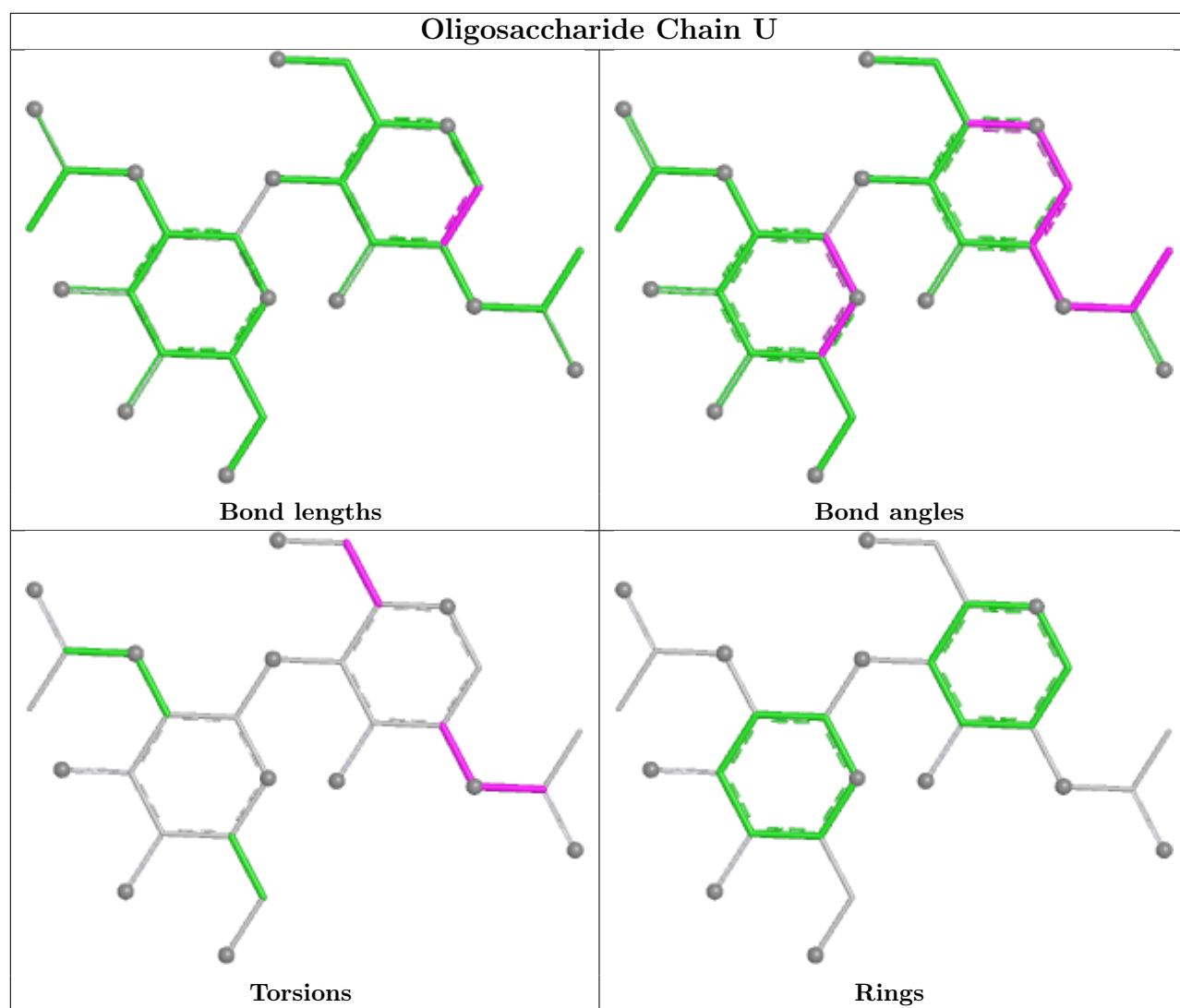


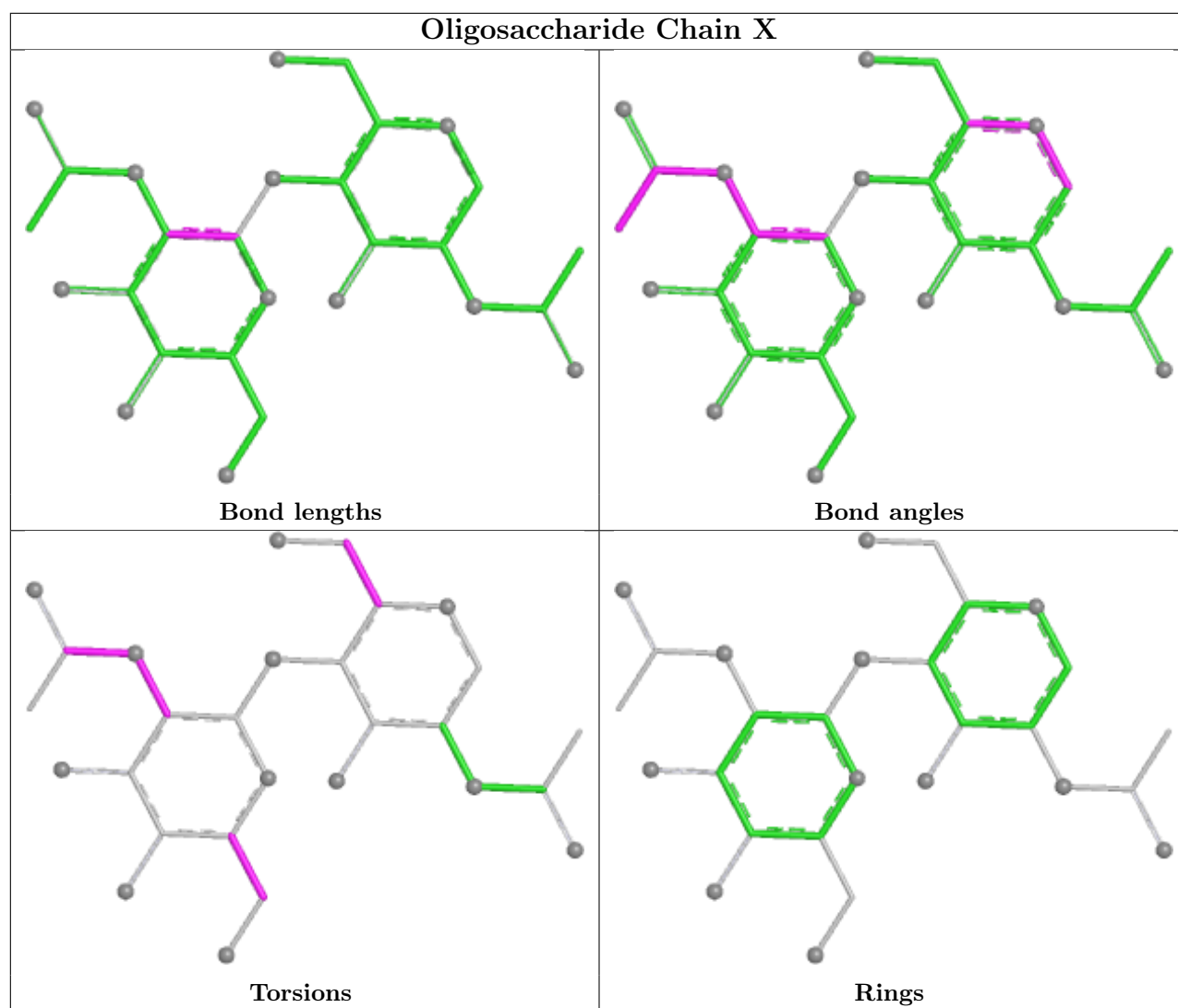


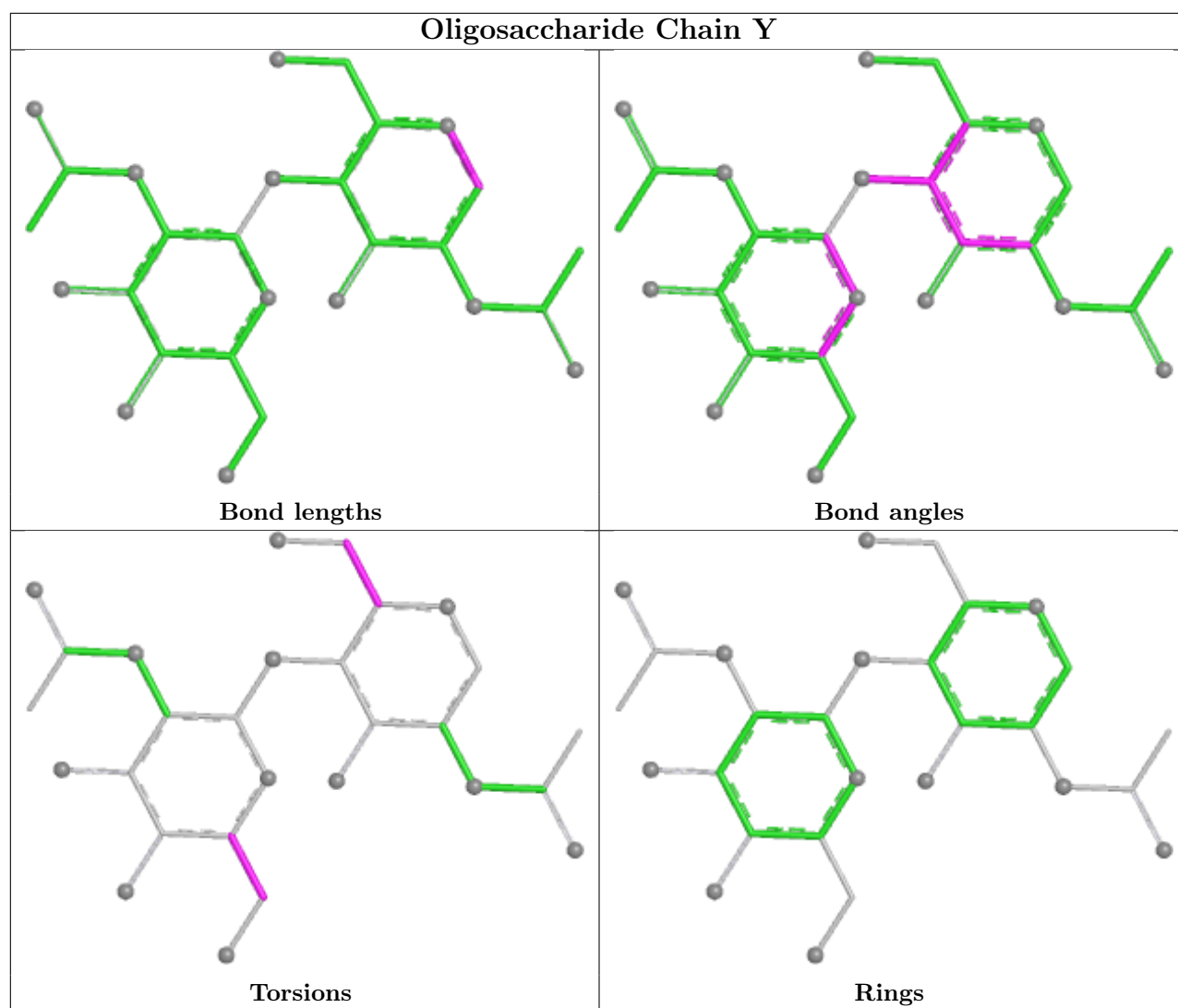


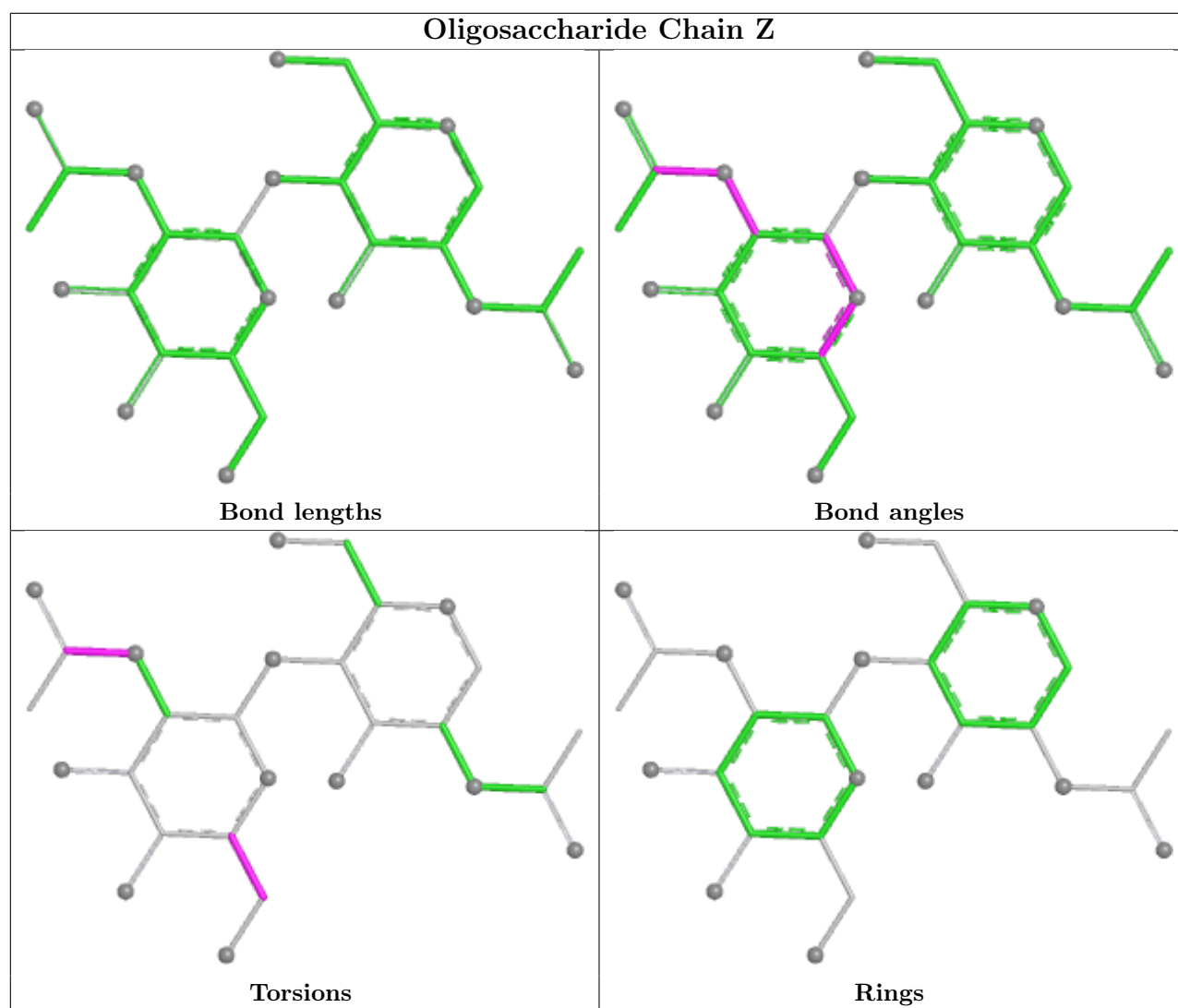


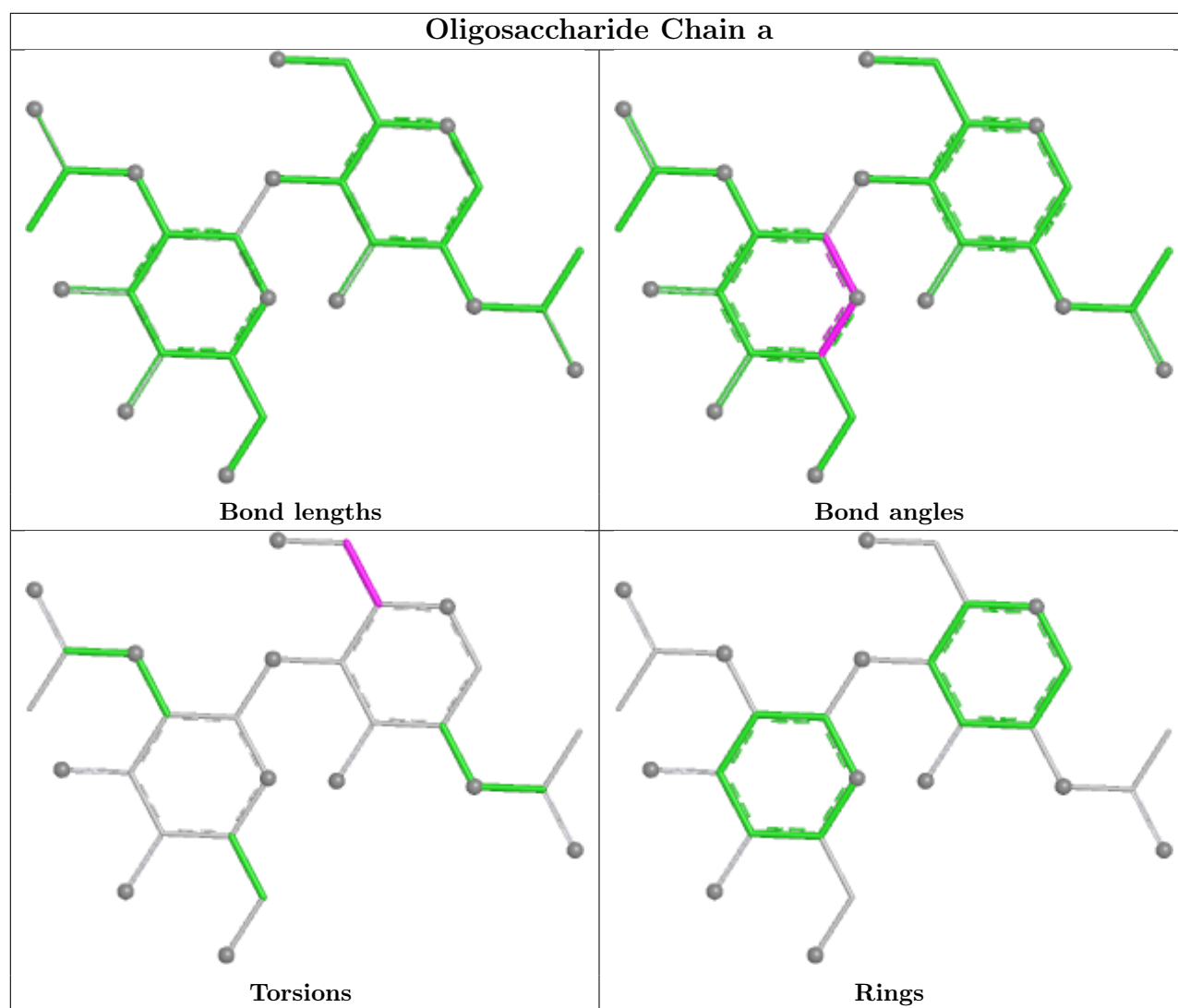


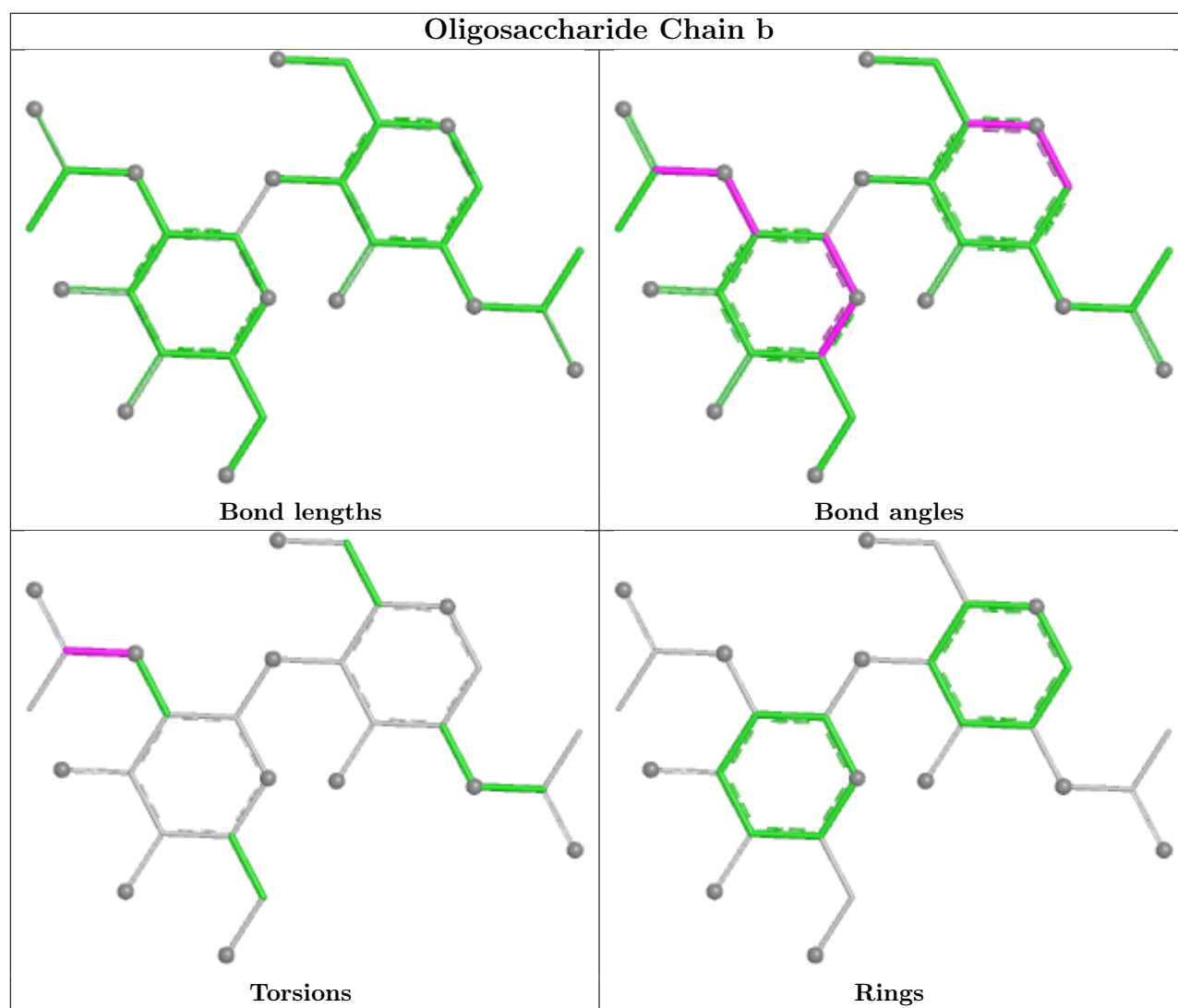


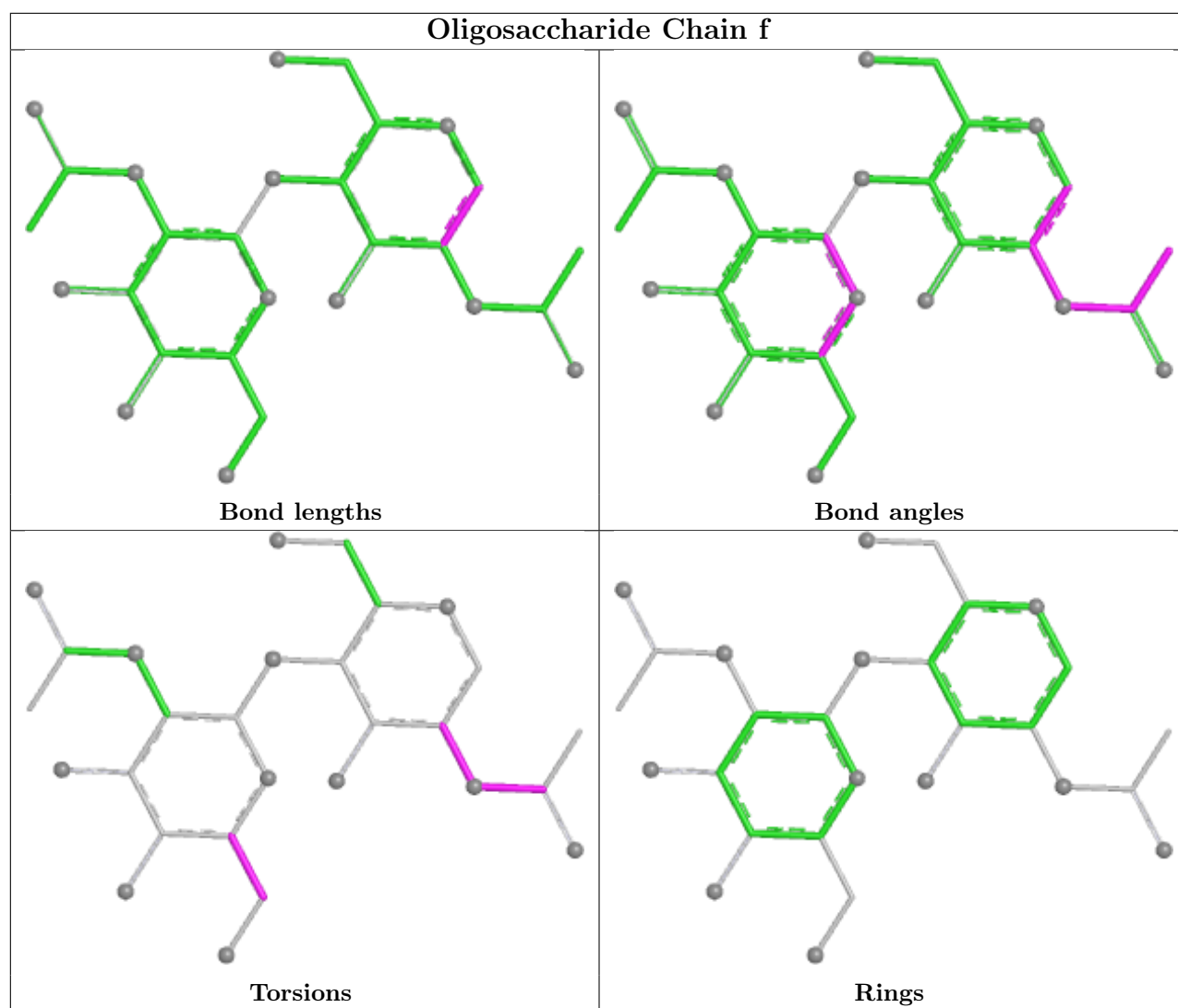


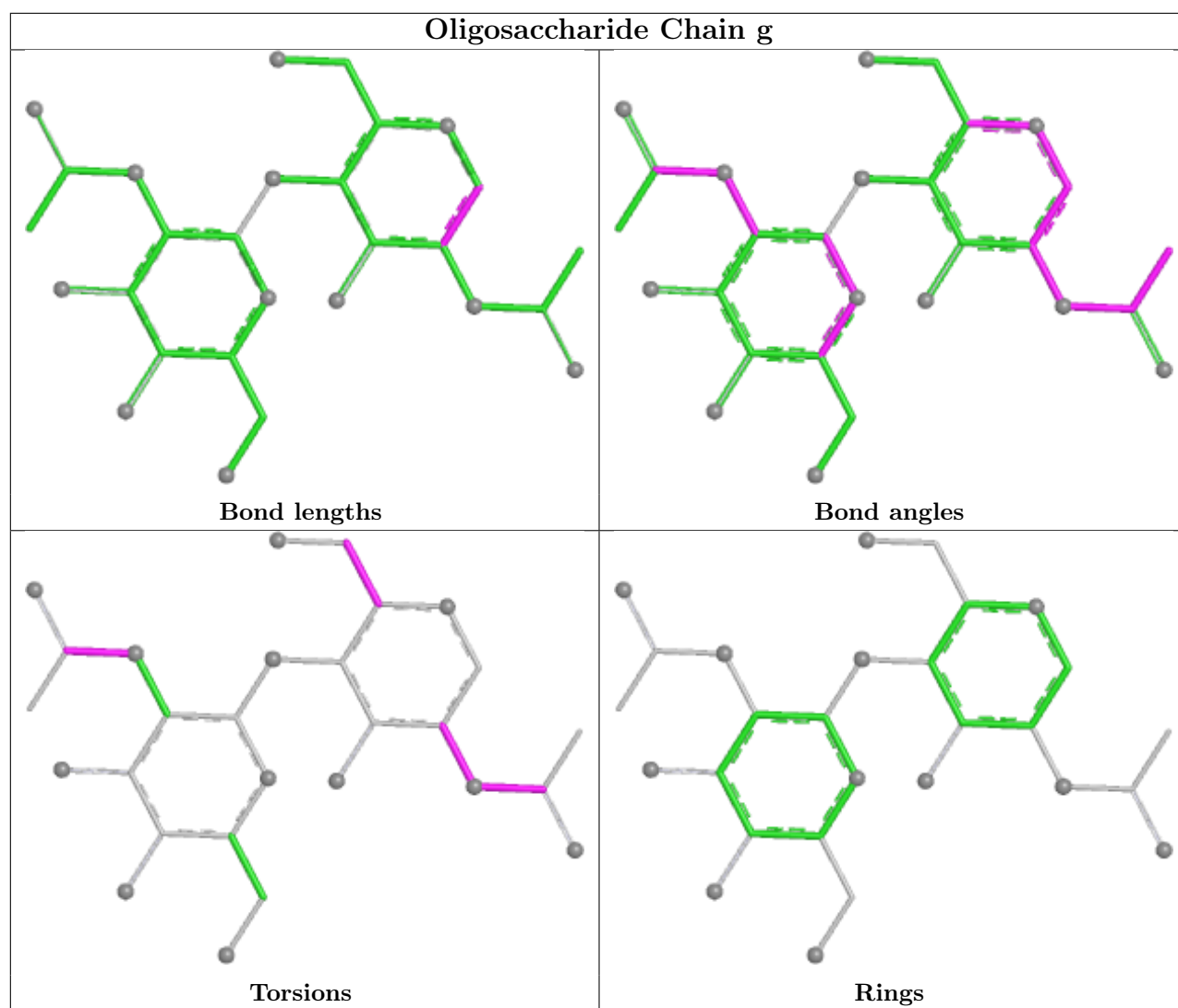


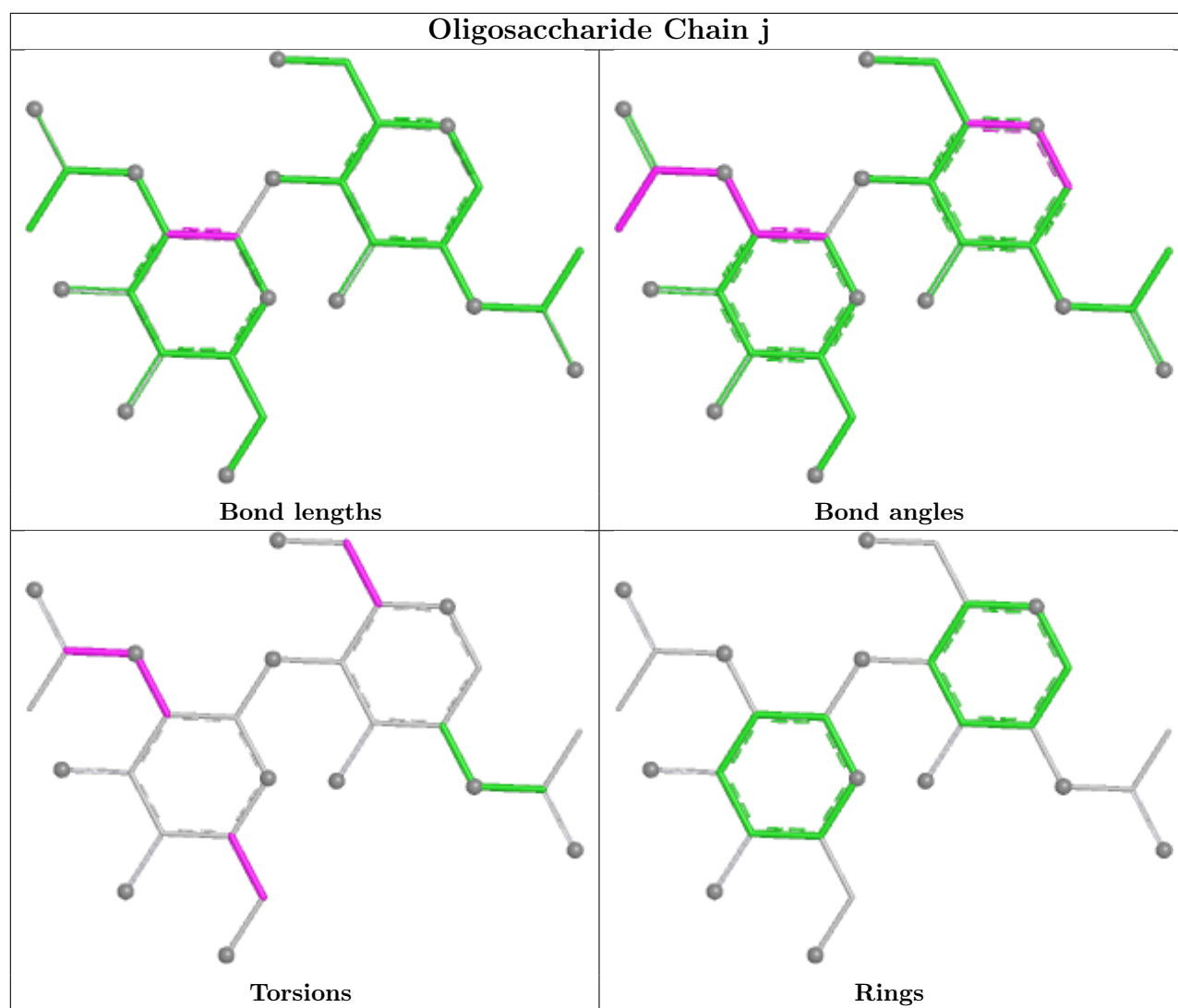


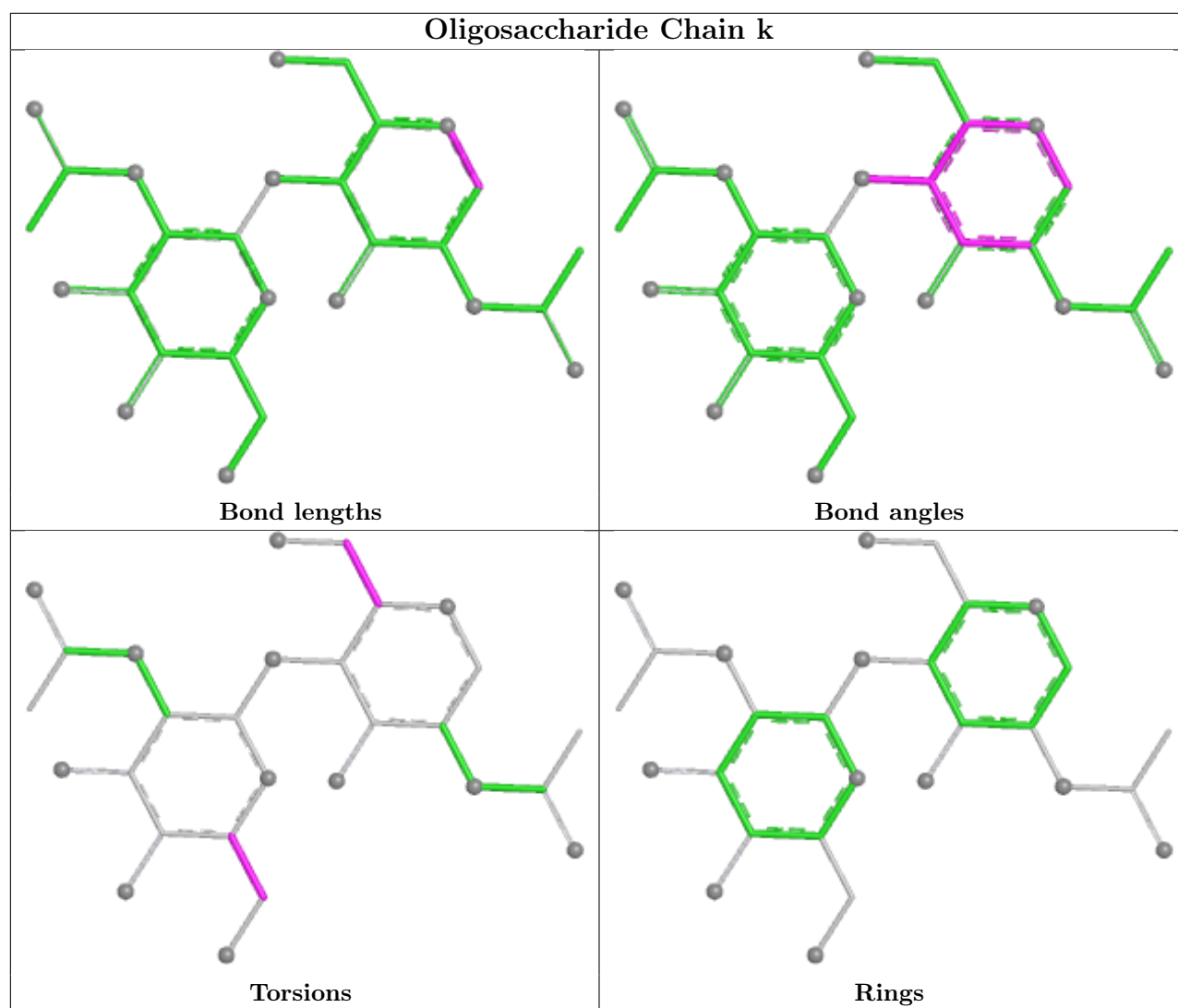


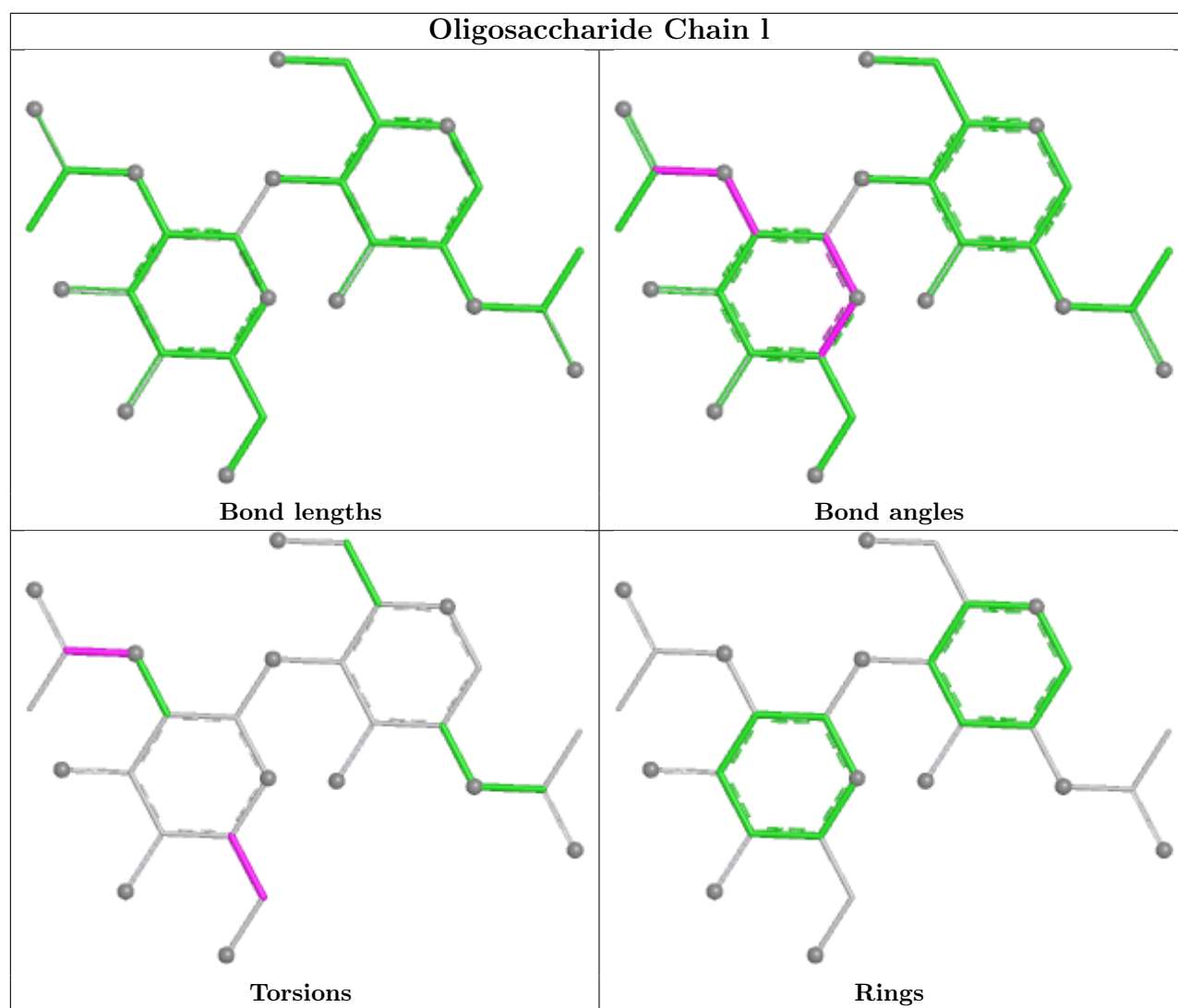


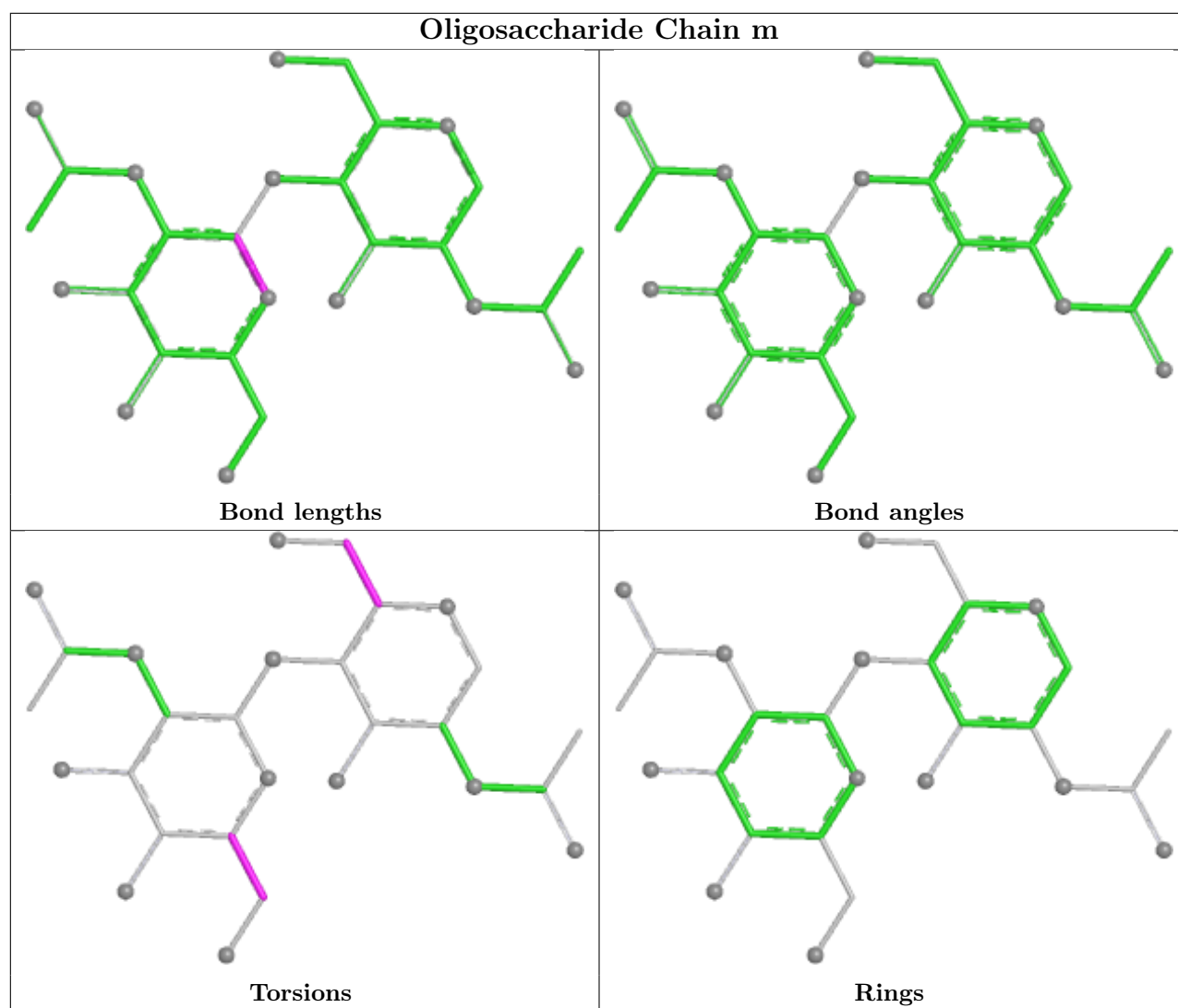


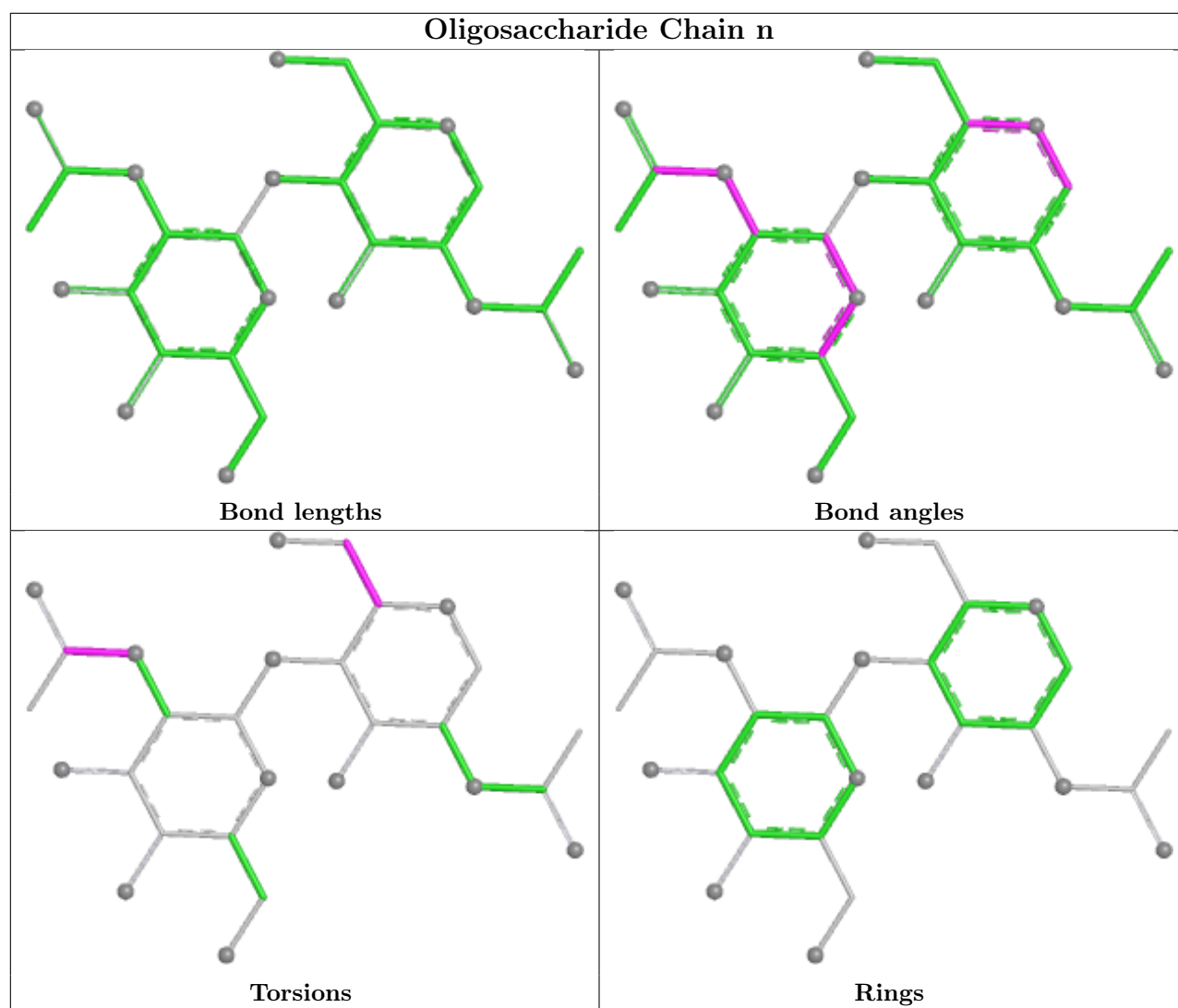


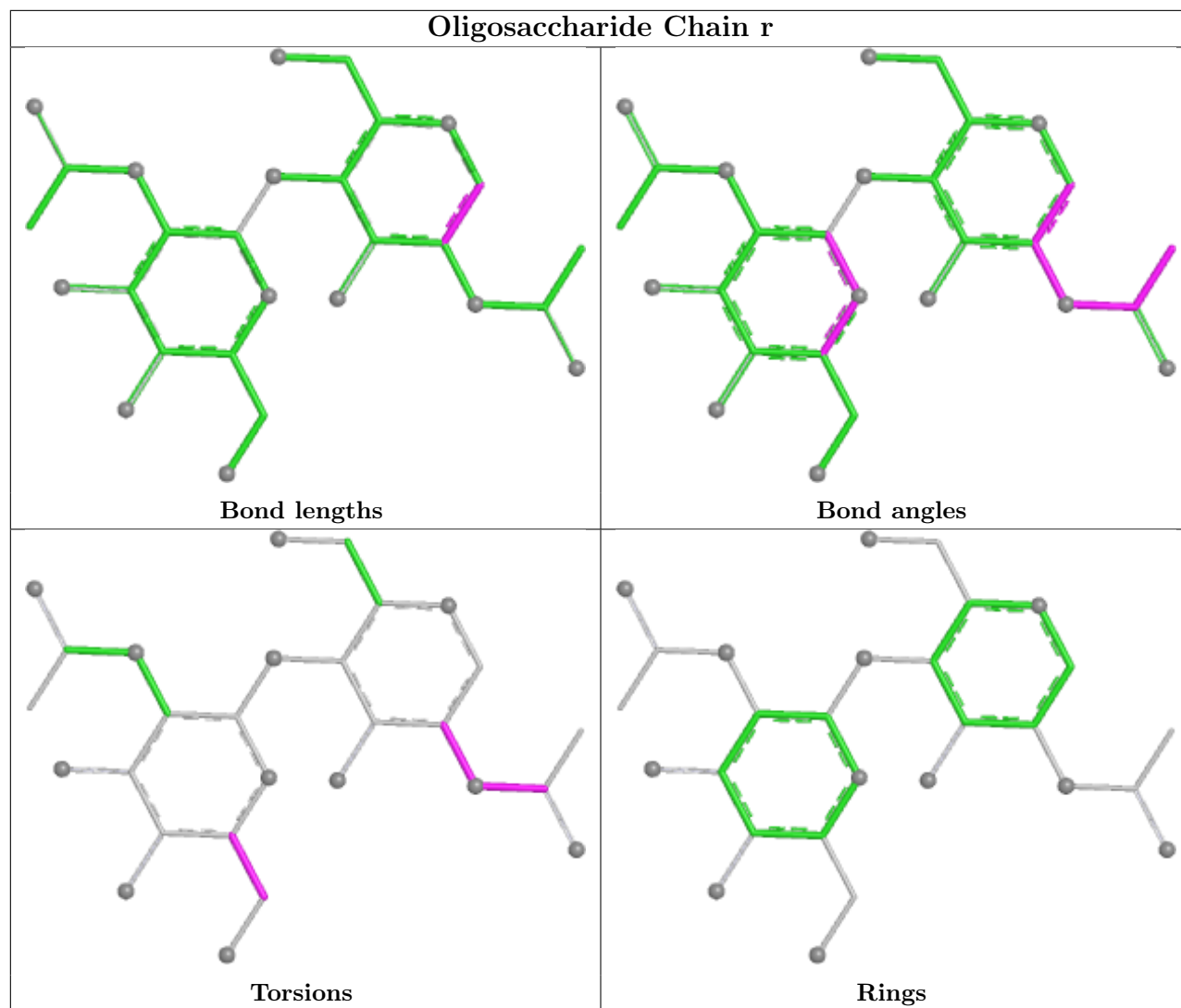


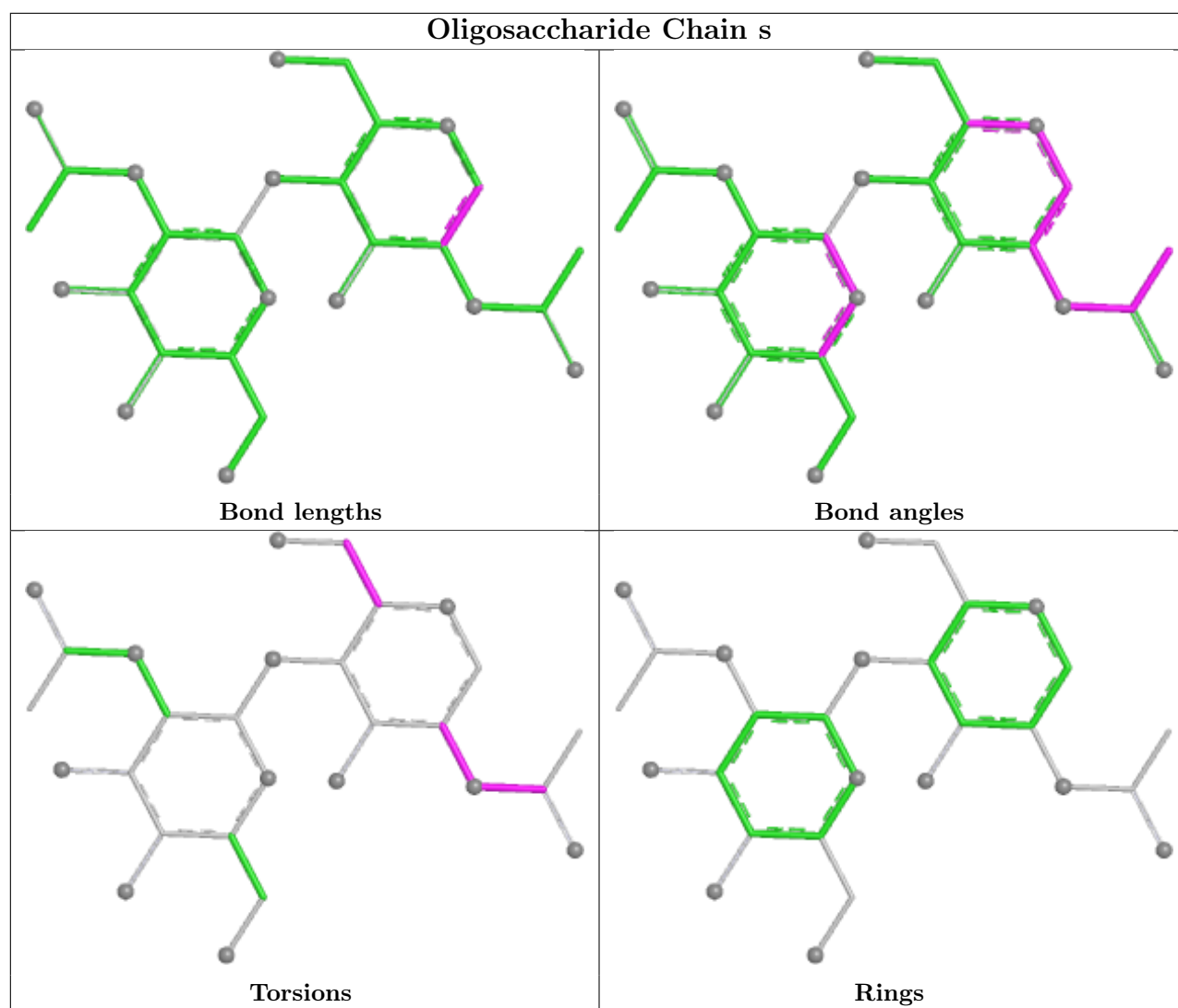


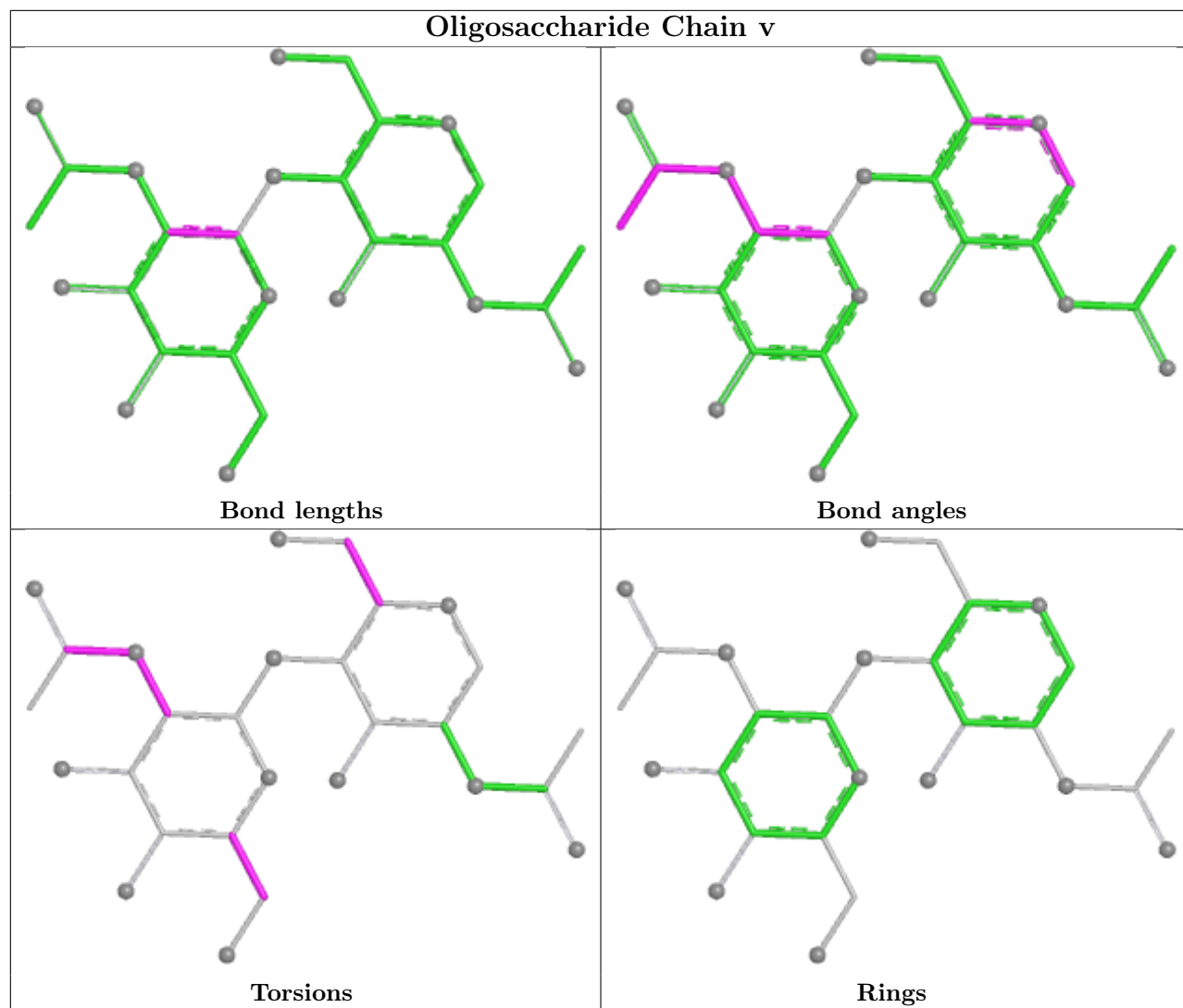


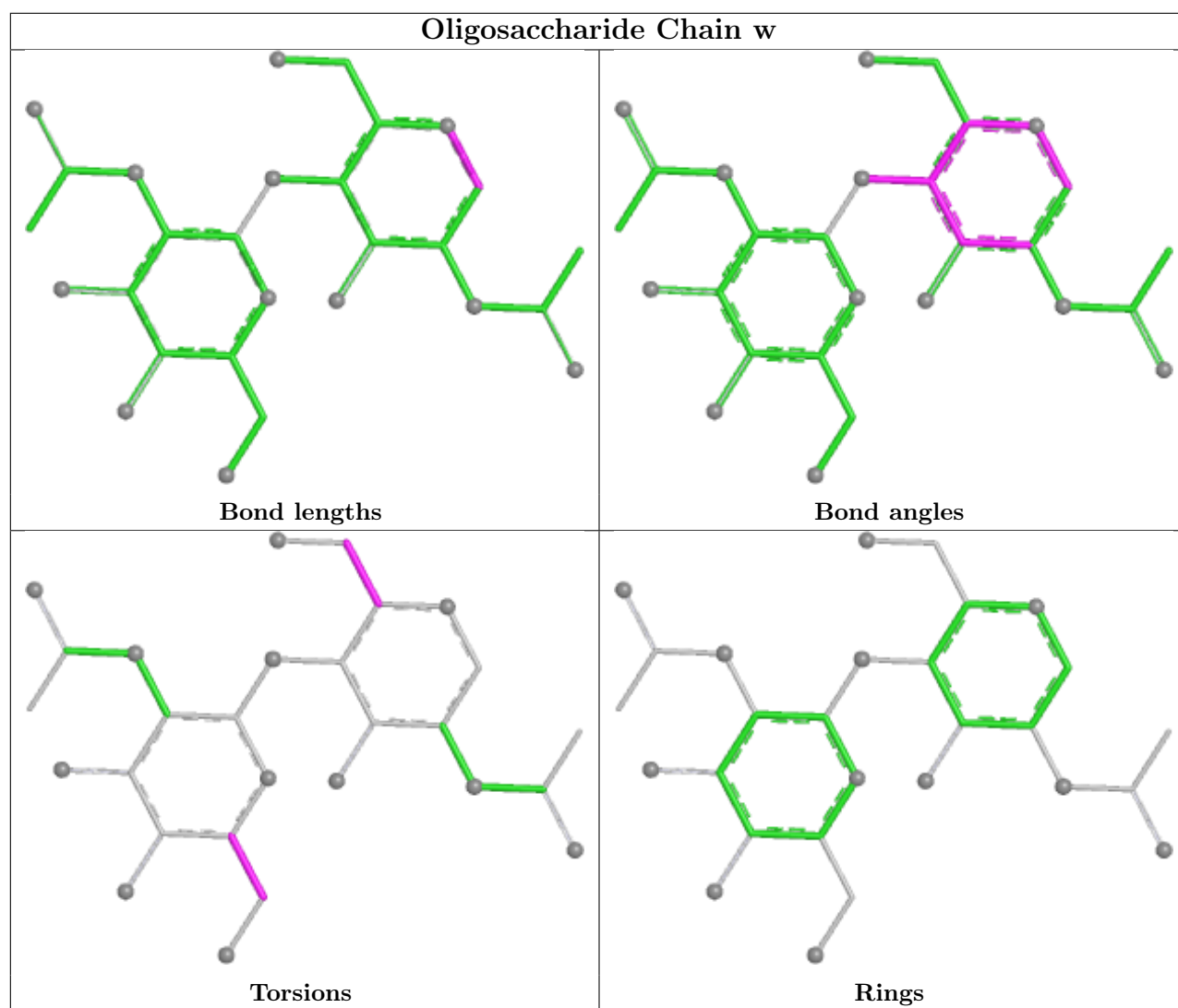


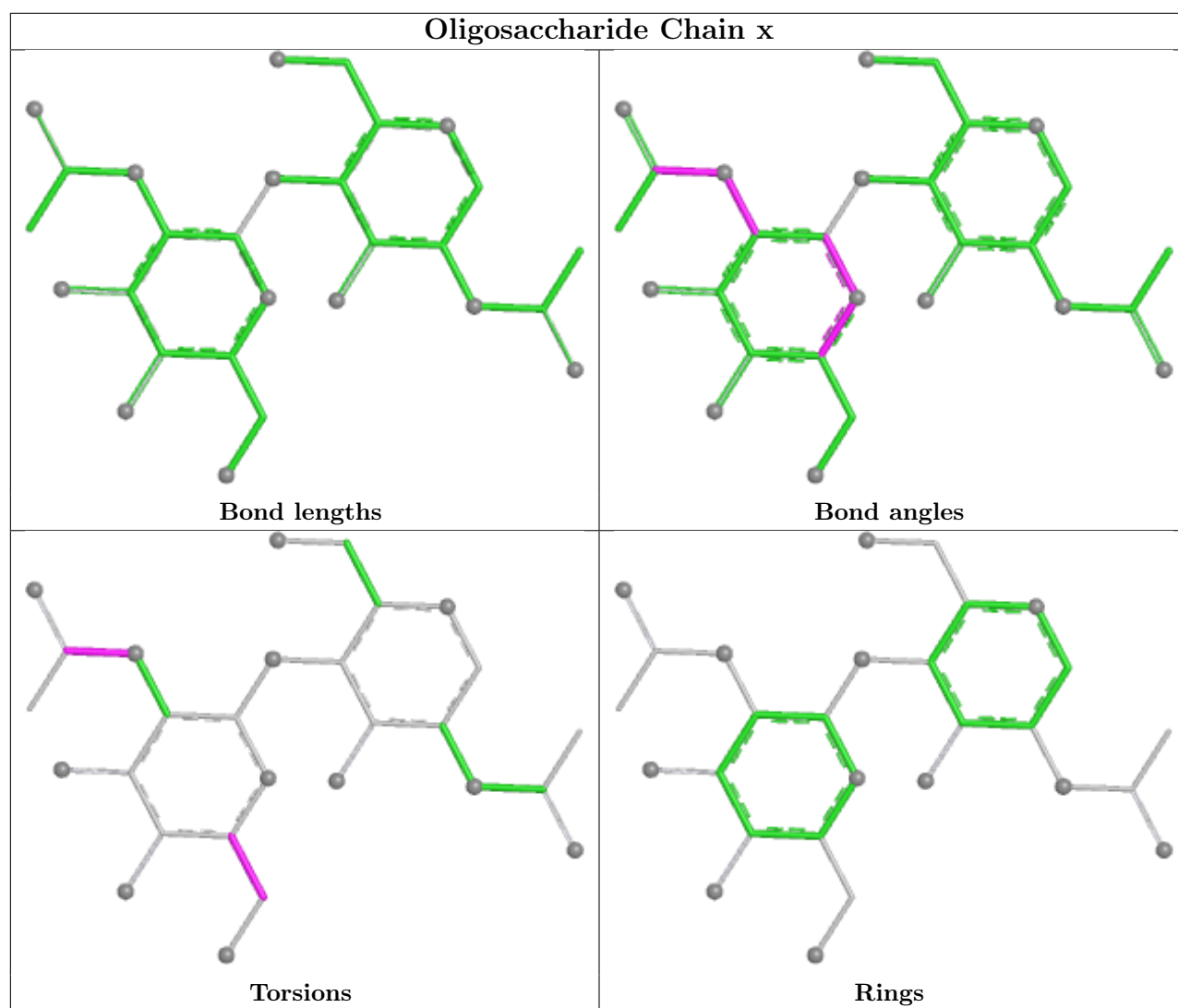


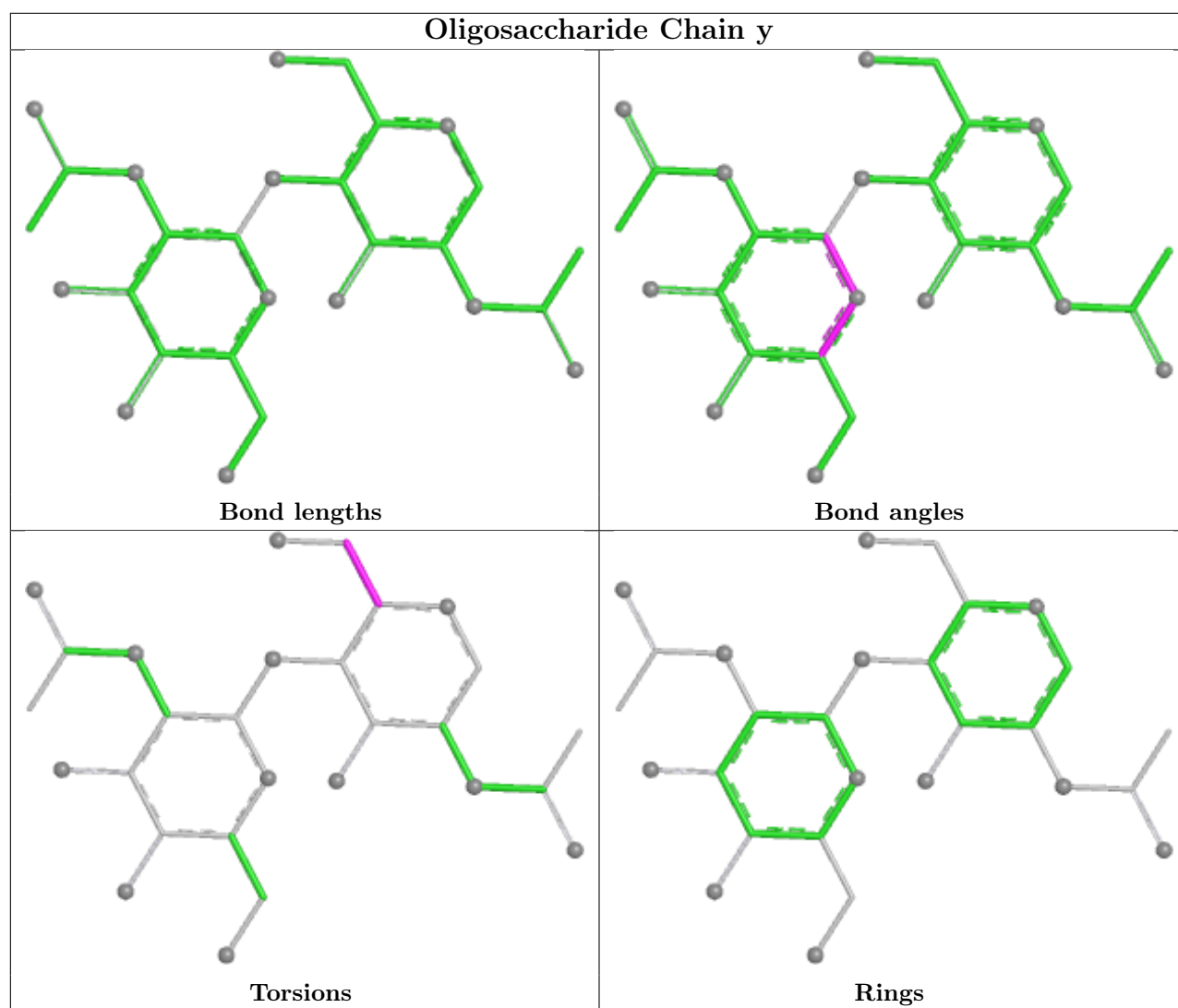


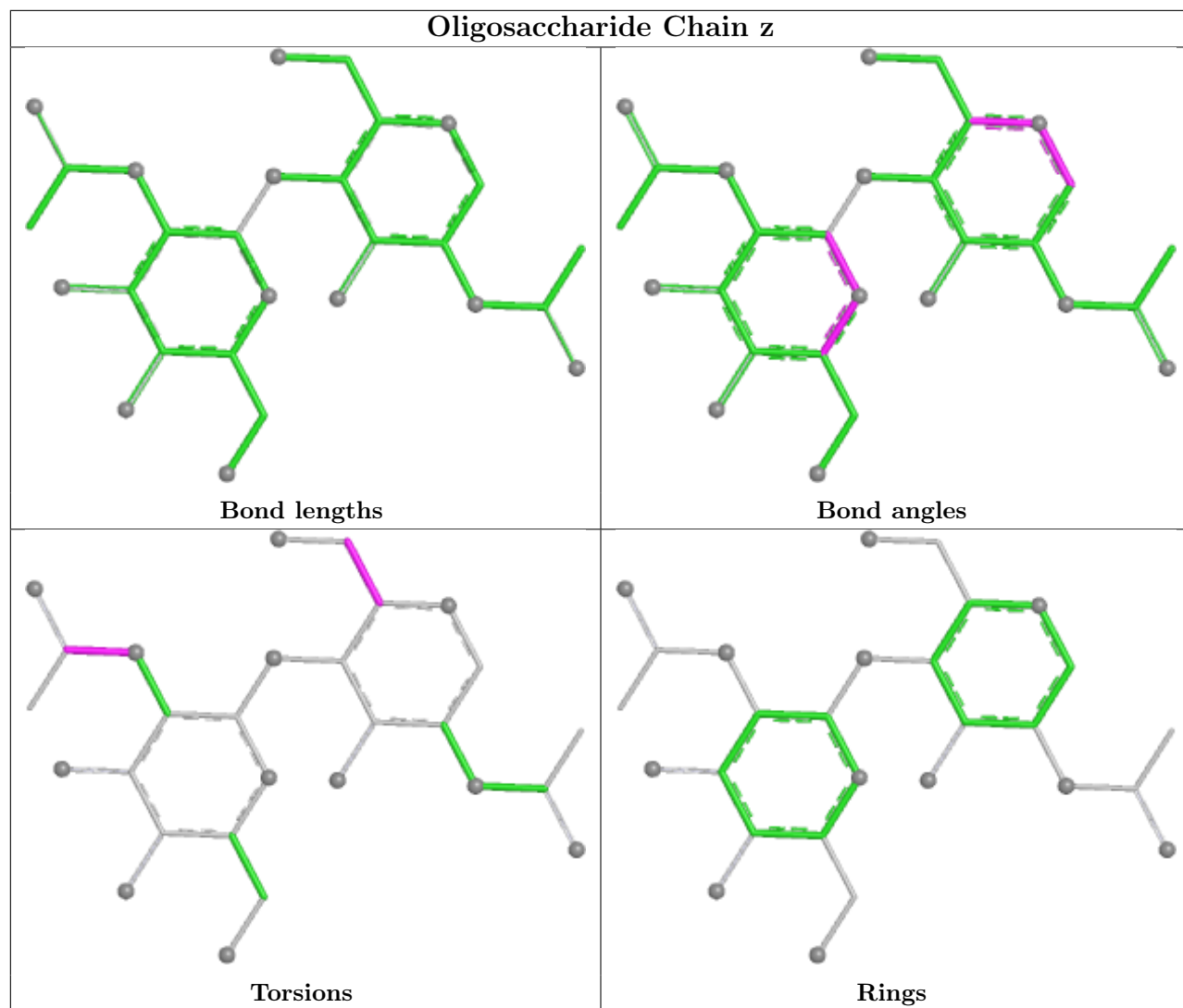


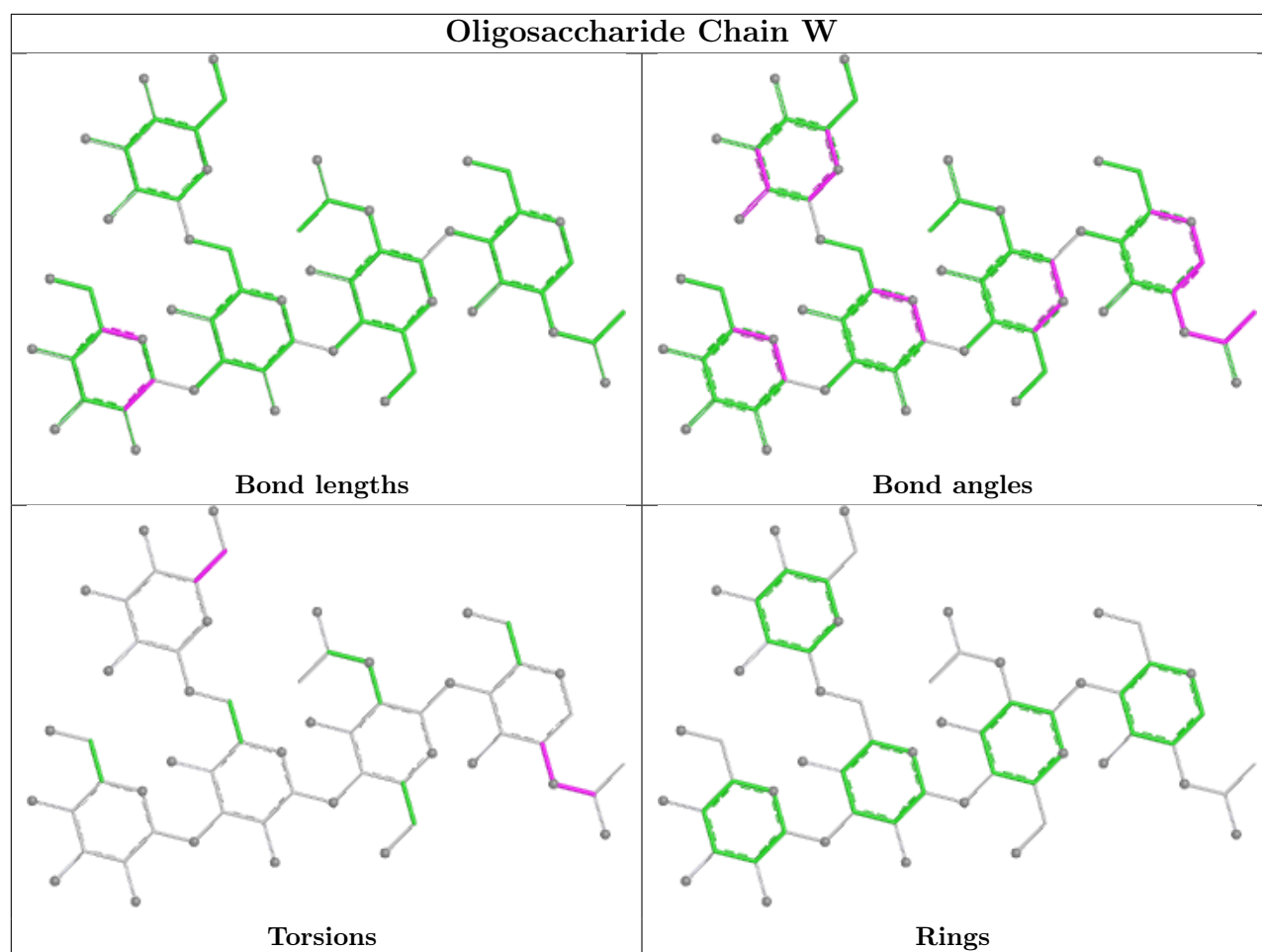


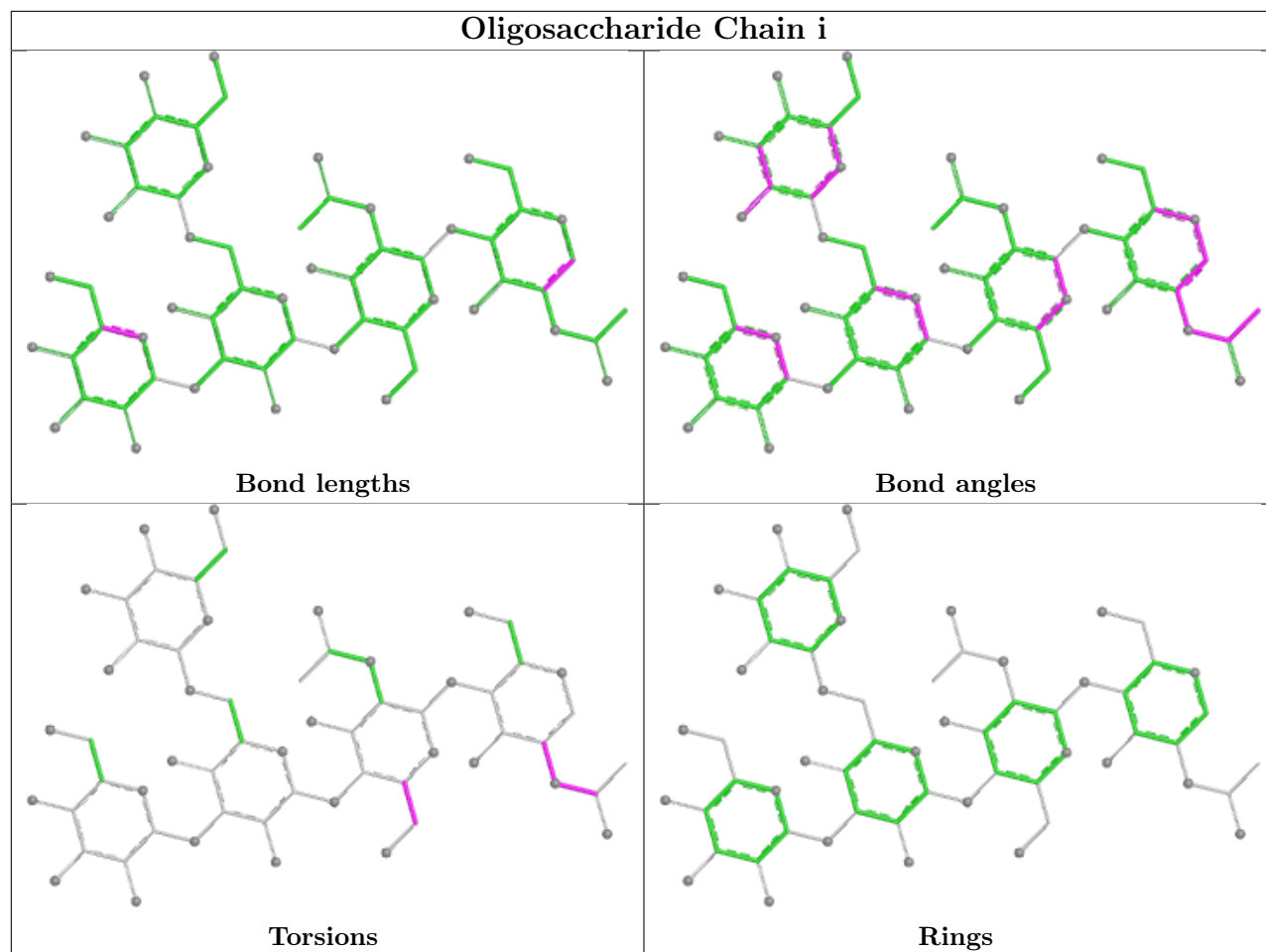


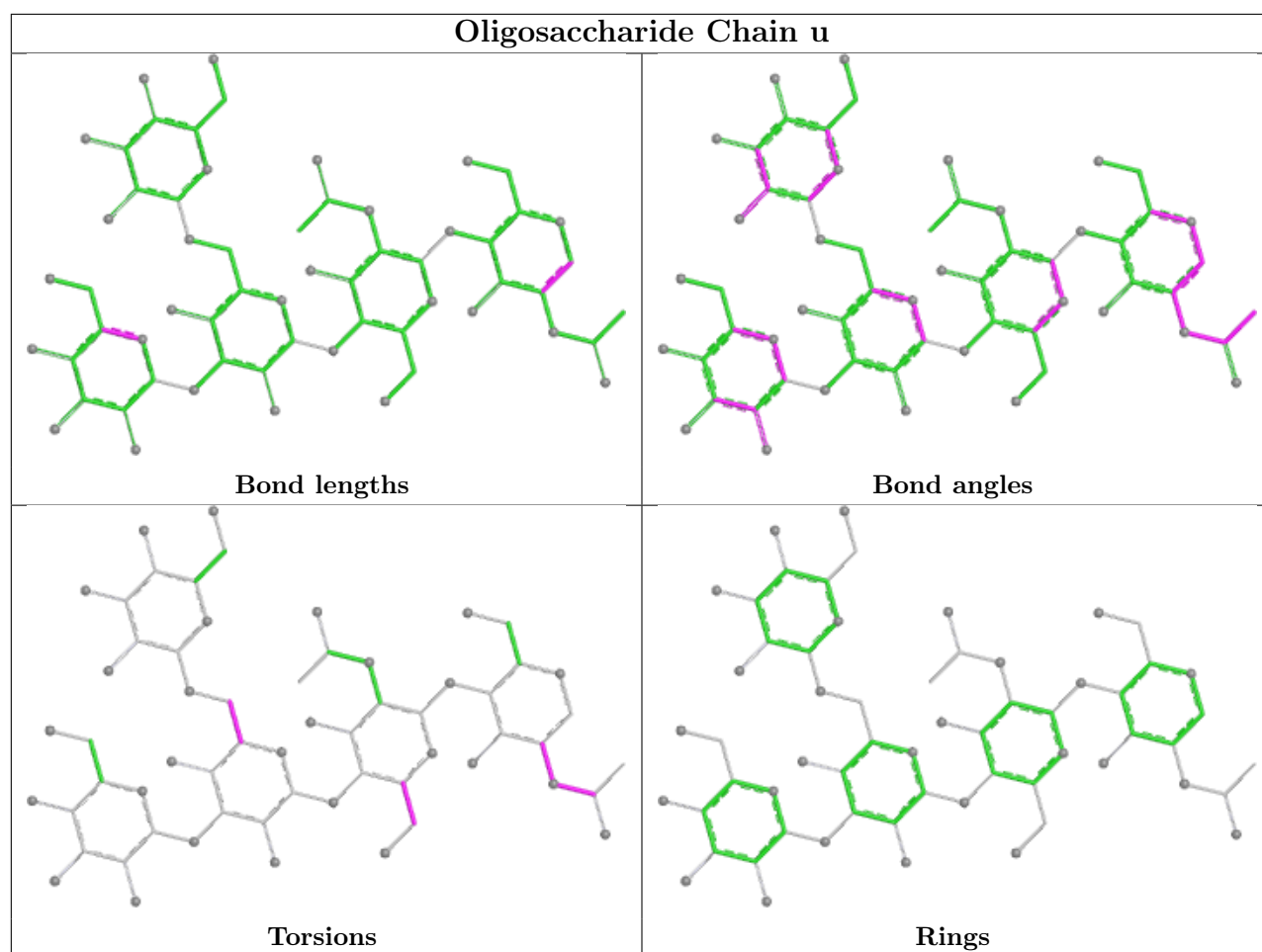


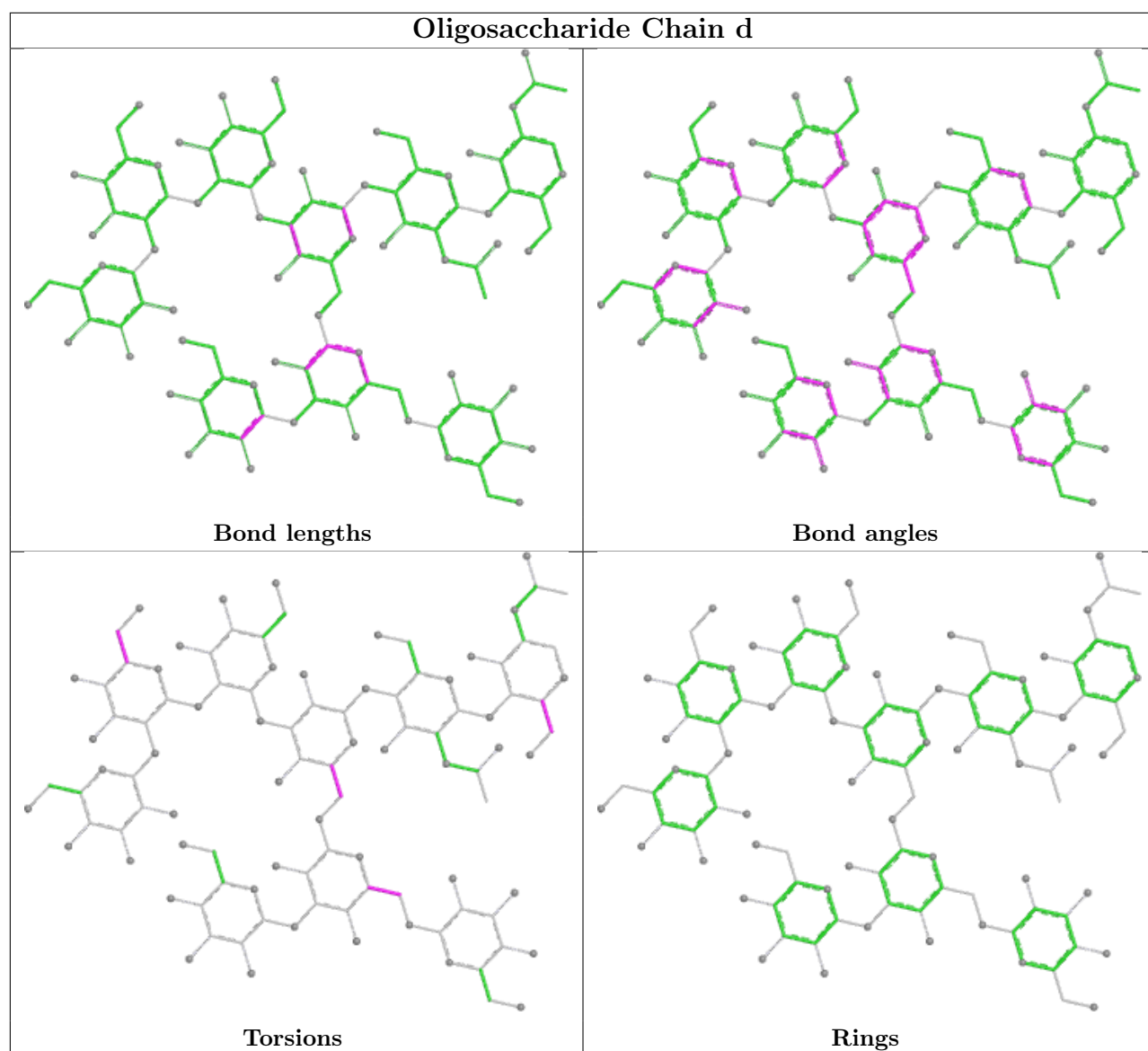


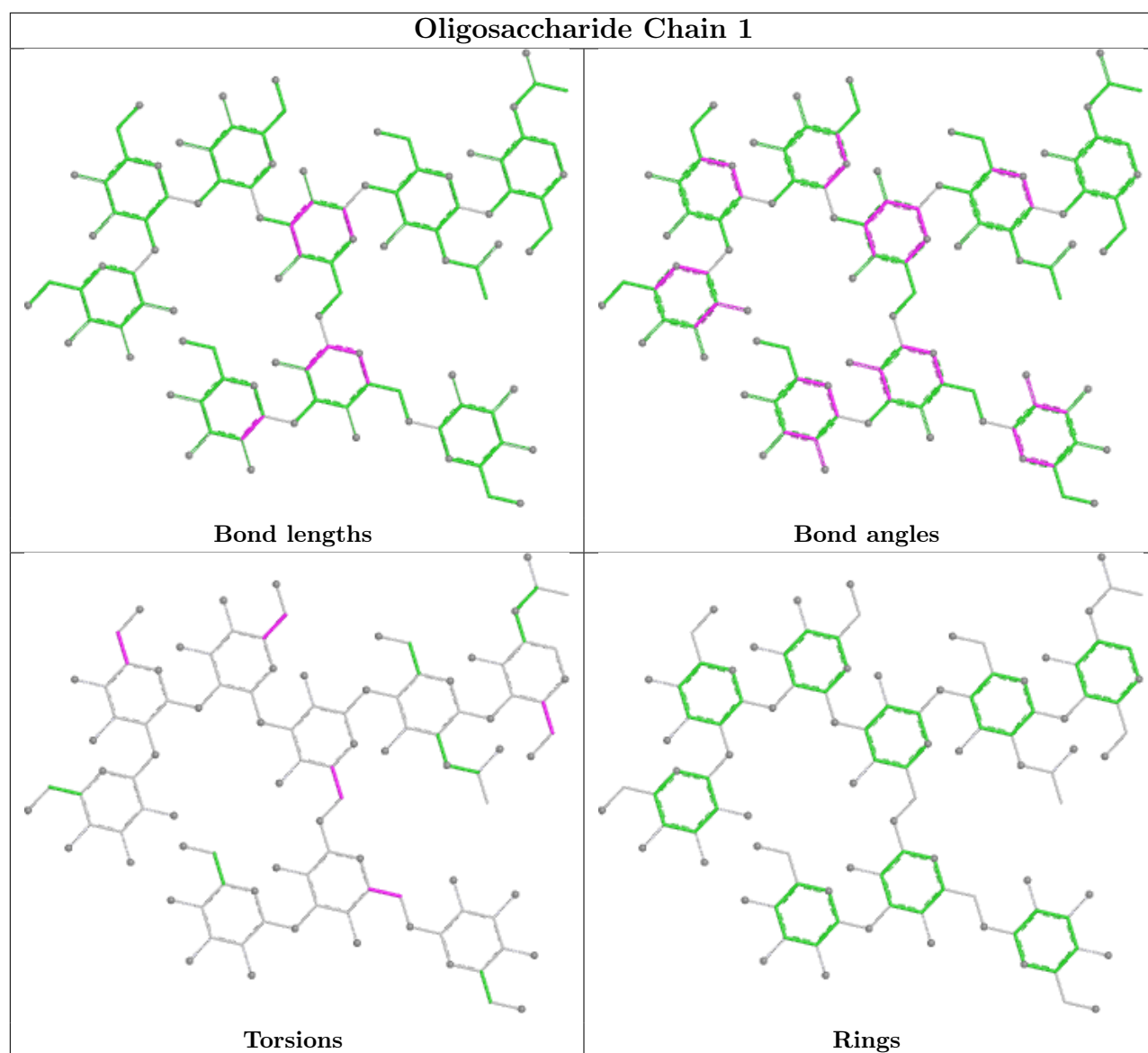


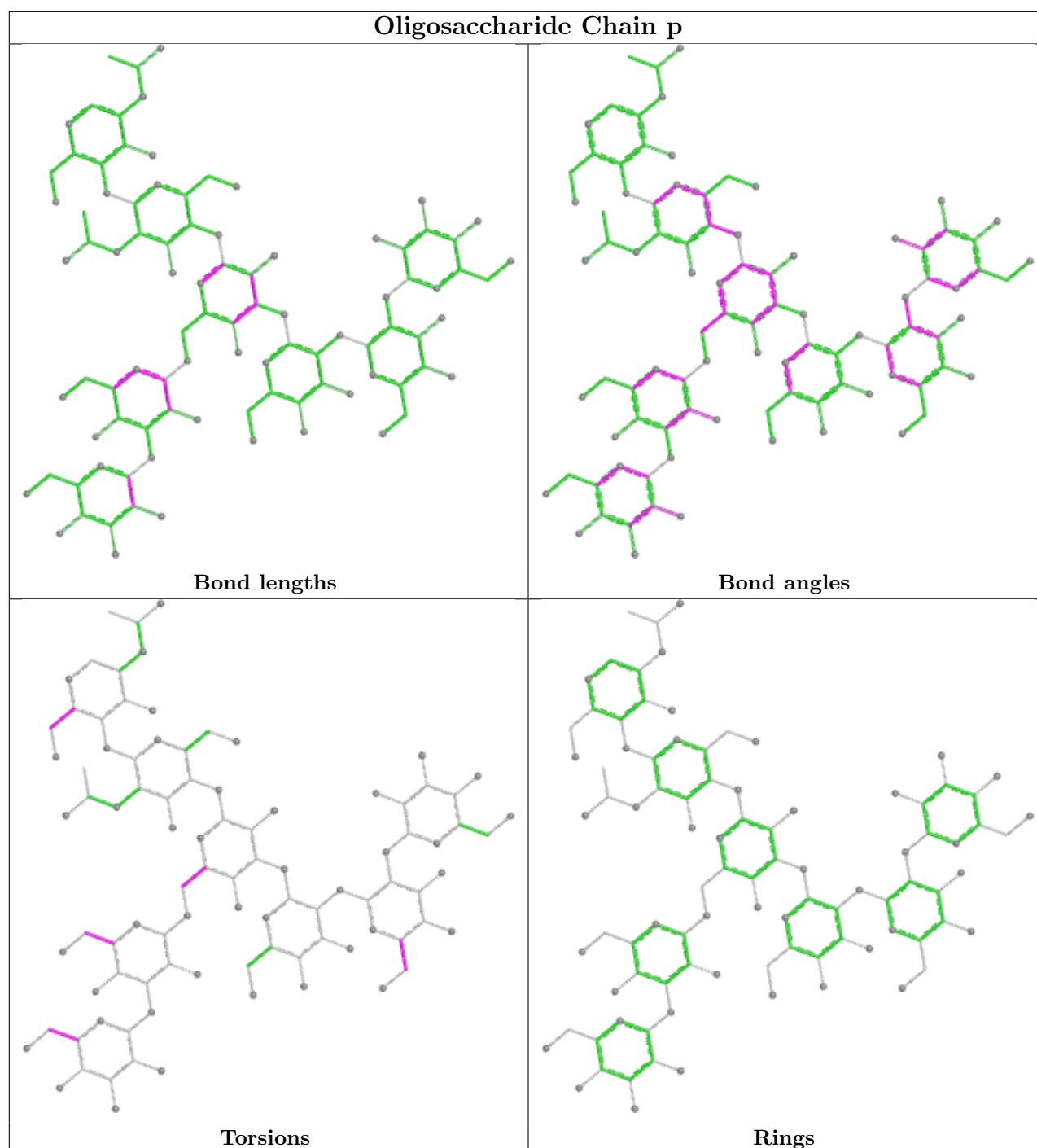












5.6 Ligand geometry [i](#)

24 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	$\# Z > 2$
12	NAG	C	605	1	14,14,15	0.57	0	17,19,21	0.73	1 (5%)
12	NAG	C	607	1	14,14,15	0.63	0	17,19,21	0.71	1 (5%)
12	NAG	F	701	2	14,14,15	0.88	1 (7%)	17,19,21	1.30	2 (11%)
12	NAG	A	604	1	14,14,15	0.72	0	17,19,21	1.36	3 (17%)
12	NAG	C	602	1	14,14,15	0.61	0	17,19,21	0.60	1 (5%)
12	NAG	B	607	1	14,14,15	0.62	0	17,19,21	0.71	1 (5%)
12	NAG	B	602	1	14,14,15	0.62	0	17,19,21	0.62	1 (5%)
12	NAG	A	603	1	14,14,15	0.60	0	17,19,21	0.66	1 (5%)
12	NAG	C	606	1	14,14,15	1.21	1 (7%)	17,19,21	2.75	3 (17%)
12	NAG	C	603	1	14,14,15	0.57	0	17,19,21	0.57	0
12	NAG	A	602	1	14,14,15	0.62	0	17,19,21	0.62	1 (5%)
12	NAG	B	601	1	14,14,15	0.76	1 (7%)	17,19,21	1.39	3 (17%)
12	NAG	B	603	1	14,14,15	0.56	0	17,19,21	0.56	0
12	NAG	D	701	2	14,14,15	0.86	1 (7%)	17,19,21	1.29	2 (11%)
12	NAG	A	601	1	14,14,15	0.77	1 (7%)	17,19,21	1.36	3 (17%)
12	NAG	B	605	1	14,14,15	0.58	0	17,19,21	0.72	1 (5%)
12	NAG	B	606	1	14,14,15	1.11	1 (7%)	17,19,21	2.77	4 (23%)
12	NAG	A	606	1	14,14,15	0.63	0	17,19,21	0.70	1 (5%)
12	NAG	A	605	1	14,14,15	0.59	0	17,19,21	0.73	1 (5%)
12	NAG	A	607	1	14,14,15	0.82	1 (7%)	17,19,21	0.68	1 (5%)
12	NAG	C	601	1	14,14,15	0.77	1 (7%)	17,19,21	1.40	3 (17%)
12	NAG	B	604	1	14,14,15	0.72	0	17,19,21	1.36	3 (17%)
12	NAG	C	604	1	14,14,15	0.72	0	17,19,21	1.36	3 (17%)
12	NAG	E	701	2	14,14,15	0.90	1 (7%)	17,19,21	1.29	2 (11%)

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
12	NAG	C	605	1	-	2/6/23/26	0/1/1/1
12	NAG	C	607	1	-	2/6/23/26	0/1/1/1
12	NAG	F	701	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	A	604	1	-	4/6/23/26	0/1/1/1
12	NAG	C	602	1	-	1/6/23/26	0/1/1/1
12	NAG	B	607	1	-	2/6/23/26	0/1/1/1
12	NAG	B	602	1	-	1/6/23/26	0/1/1/1
12	NAG	A	603	1	-	2/6/23/26	0/1/1/1
12	NAG	C	606	1	-	5/6/23/26	0/1/1/1
12	NAG	C	603	1	-	0/6/23/26	0/1/1/1
12	NAG	A	602	1	-	1/6/23/26	0/1/1/1
12	NAG	B	601	1	-	4/6/23/26	0/1/1/1
12	NAG	B	603	1	-	0/6/23/26	0/1/1/1
12	NAG	D	701	2	-	2/6/23/26	0/1/1/1
12	NAG	A	601	1	-	4/6/23/26	0/1/1/1
12	NAG	B	605	1	-	2/6/23/26	0/1/1/1
12	NAG	B	606	1	-	6/6/23/26	0/1/1/1
12	NAG	A	606	1	-	2/6/23/26	0/1/1/1
12	NAG	A	605	1	-	2/6/23/26	0/1/1/1
12	NAG	A	607	1	-	0/6/23/26	0/1/1/1
12	NAG	C	601	1	-	4/6/23/26	0/1/1/1
12	NAG	B	604	1	-	4/6/23/26	0/1/1/1
12	NAG	C	604	1	-	4/6/23/26	0/1/1/1
12	NAG	E	701	2	-	2/6/23/26	0/1/1/1

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	C	606	NAG	C1-C2	3.44	1.57	1.52
12	B	606	NAG	C1-C2	3.02	1.56	1.52
12	E	701	NAG	C1-C2	2.51	1.55	1.52
12	F	701	NAG	C1-C2	2.46	1.55	1.52
12	D	701	NAG	C1-C2	2.35	1.55	1.52

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	B	606	NAG	C2-N2-C7	9.60	135.77	122.90
12	C	606	NAG	C2-N2-C7	9.53	135.68	122.90
12	C	606	NAG	C1-C2-N2	4.91	118.18	110.43
12	B	606	NAG	C1-C2-N2	4.72	117.88	110.43
12	C	604	NAG	C2-N2-C7	4.25	128.60	122.90

There are no chirality outliers.

5 of 58 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	A	601	NAG	O5-C5-C6-O6
12	B	606	NAG	O5-C5-C6-O6
12	B	601	NAG	O5-C5-C6-O6
12	C	601	NAG	O5-C5-C6-O6
12	A	603	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	B	603	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	3
1	B	3
1	A	3
2	E	1
2	F	1
2	D	1

The worst 5 of 12 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	546:SER	C	565:LEU	N	25.08
1	F	546:SER	C	565:LEU	N	25.01
1	D	546:SER	C	565:LEU	N	24.75
1	C	137:GLY	C	152:THR	N	13.35
1	B	137:GLY	C	152:THR	N	13.34

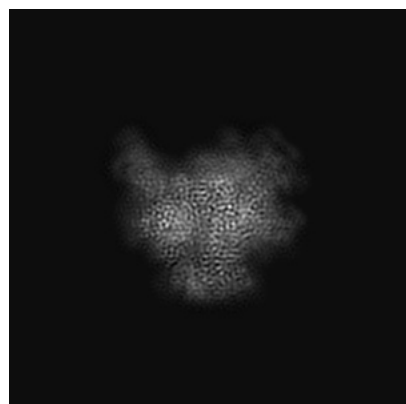
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26496. These allow visual inspection of the internal detail of the map and identification of artifacts.

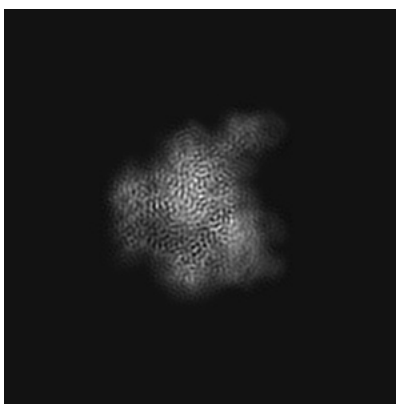
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

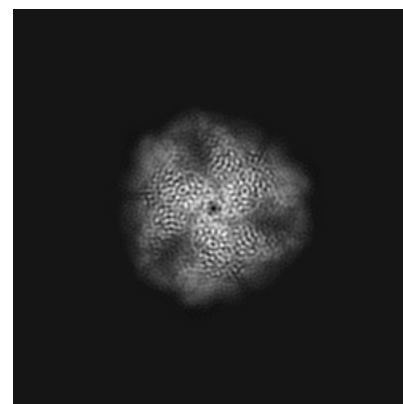
6.1.1 Primary map



X

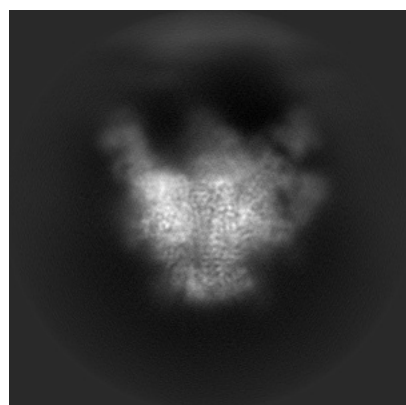


Y

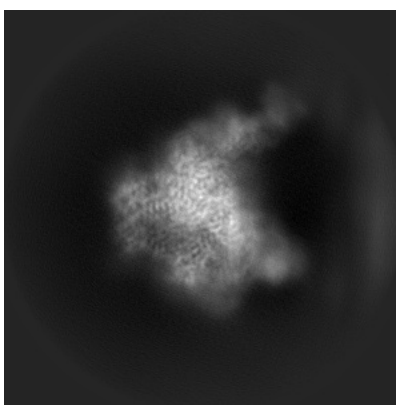


Z

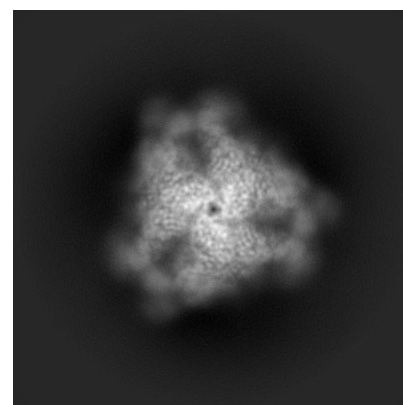
6.1.2 Raw map



X



Y

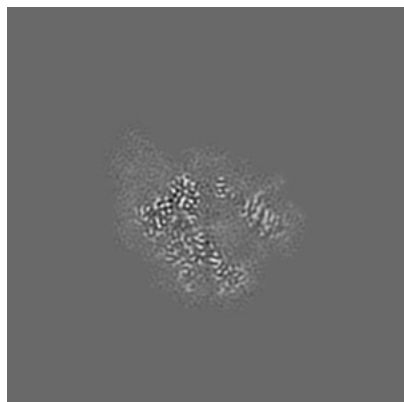


Z

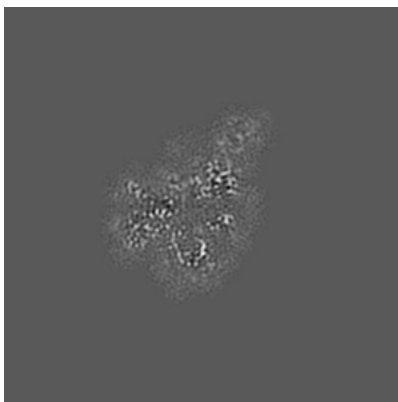
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

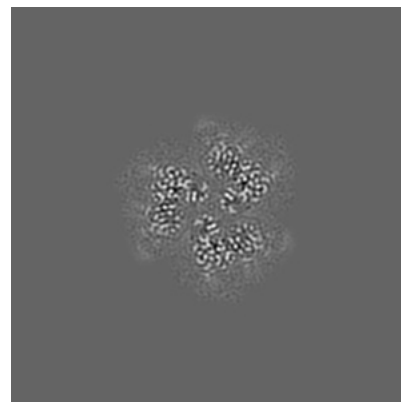
6.2.1 Primary map



X Index: 180

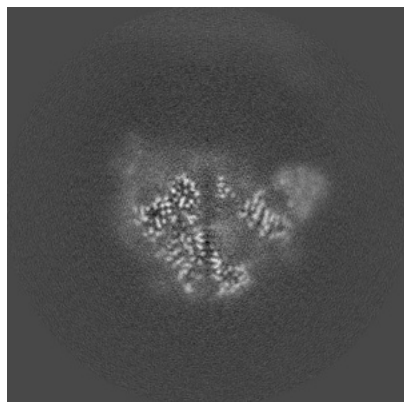


Y Index: 180

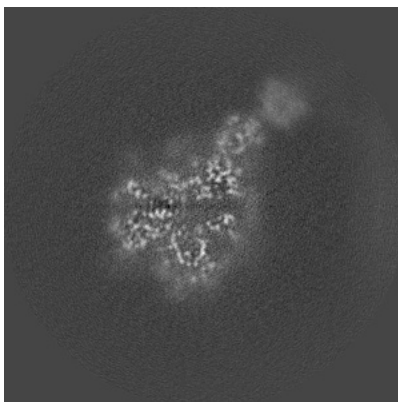


Z Index: 180

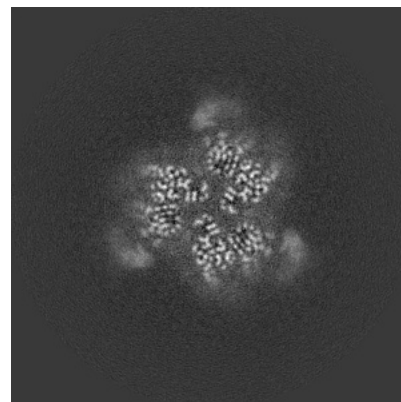
6.2.2 Raw map



X Index: 180



Y Index: 180

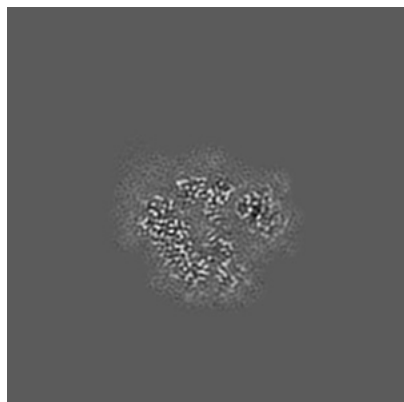


Z Index: 180

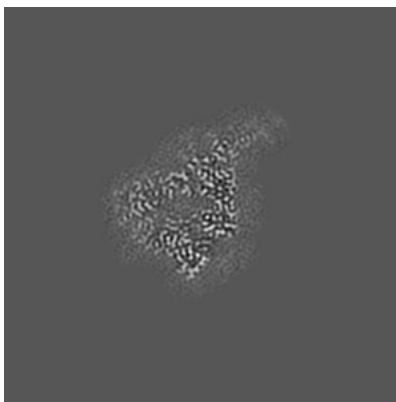
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

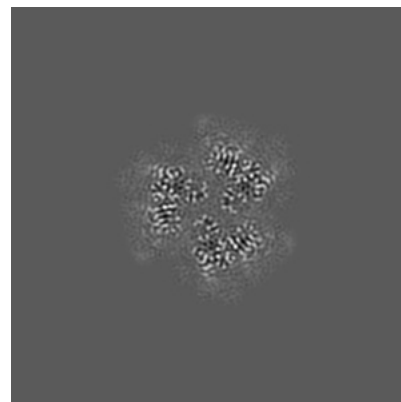
6.3.1 Primary map



X Index: 189

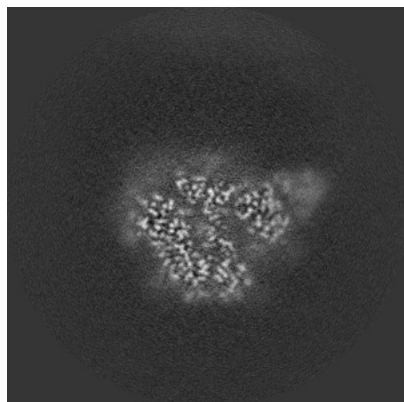


Y Index: 191

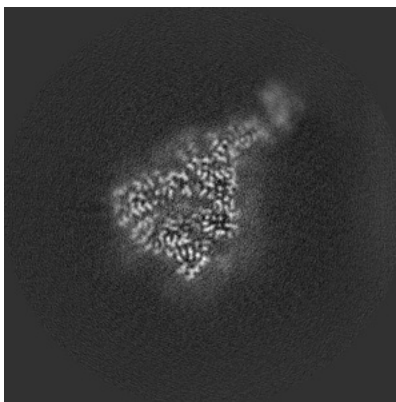


Z Index: 179

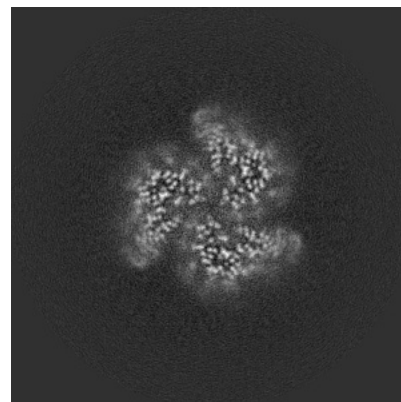
6.3.2 Raw map



X Index: 189



Y Index: 191

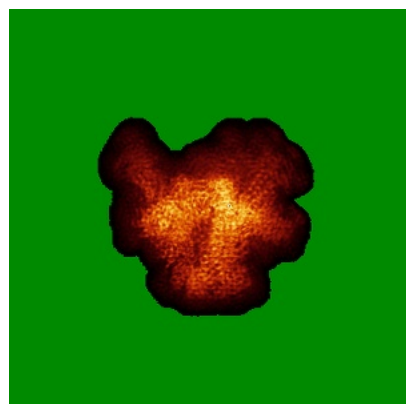


Z Index: 168

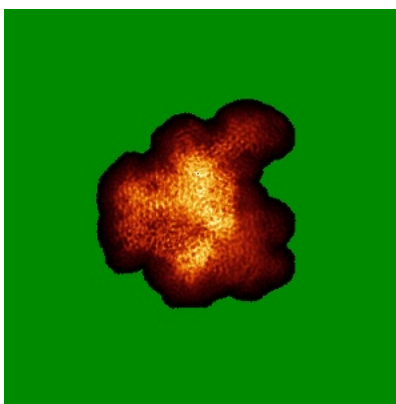
The images above show the largest variance slices of the map in three orthogonal directions.

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

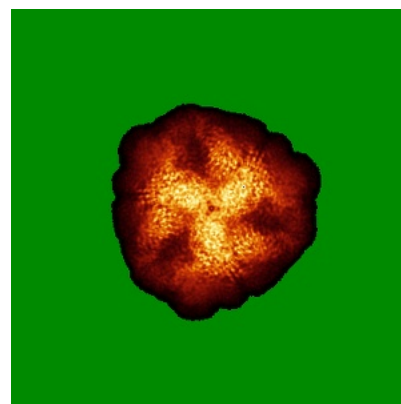
6.4.1 Primary map



X

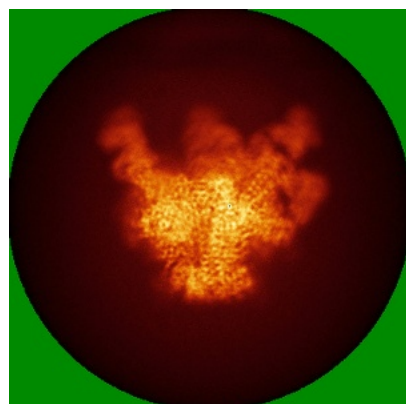


Y

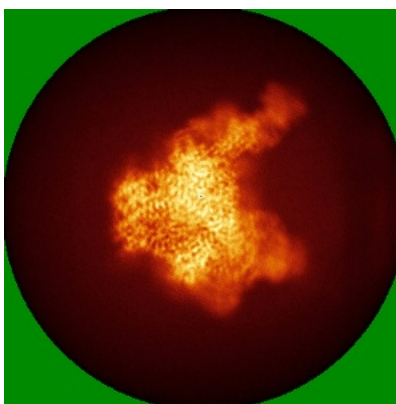


Z

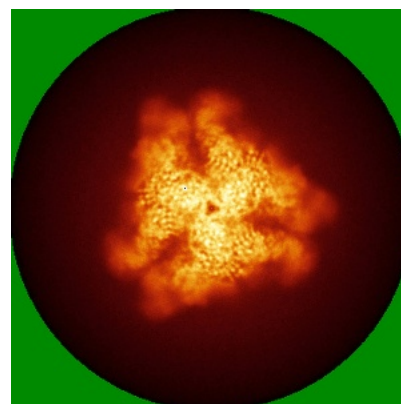
6.4.2 Raw map



X



Y



Z

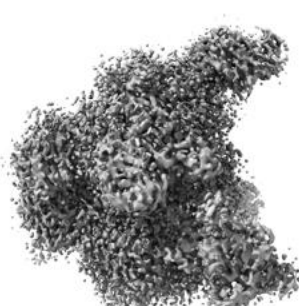
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

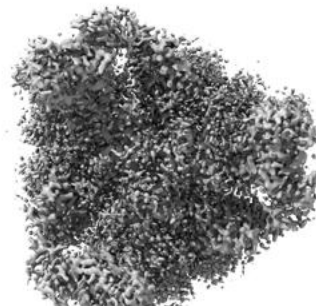
6.5.1 Primary map



X



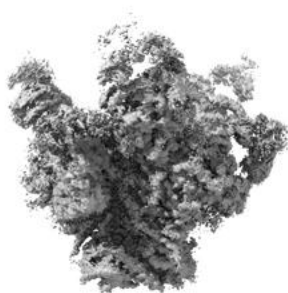
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0115. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

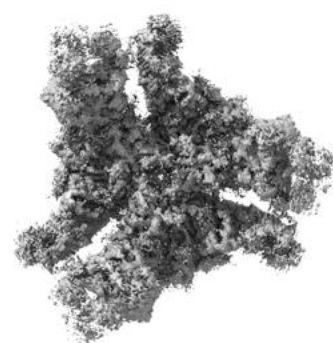
6.5.2 Raw map



X



Y



Z

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

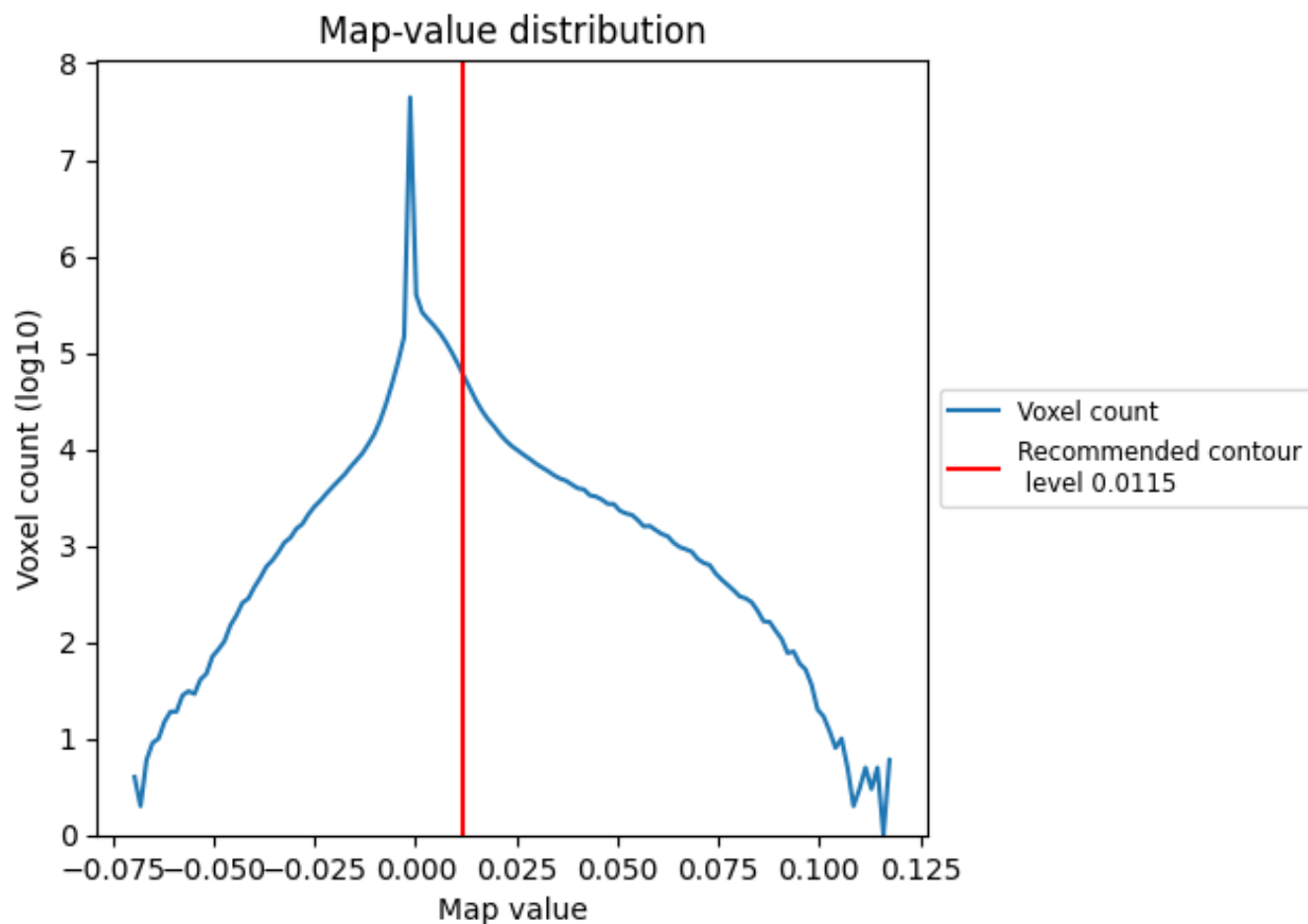
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

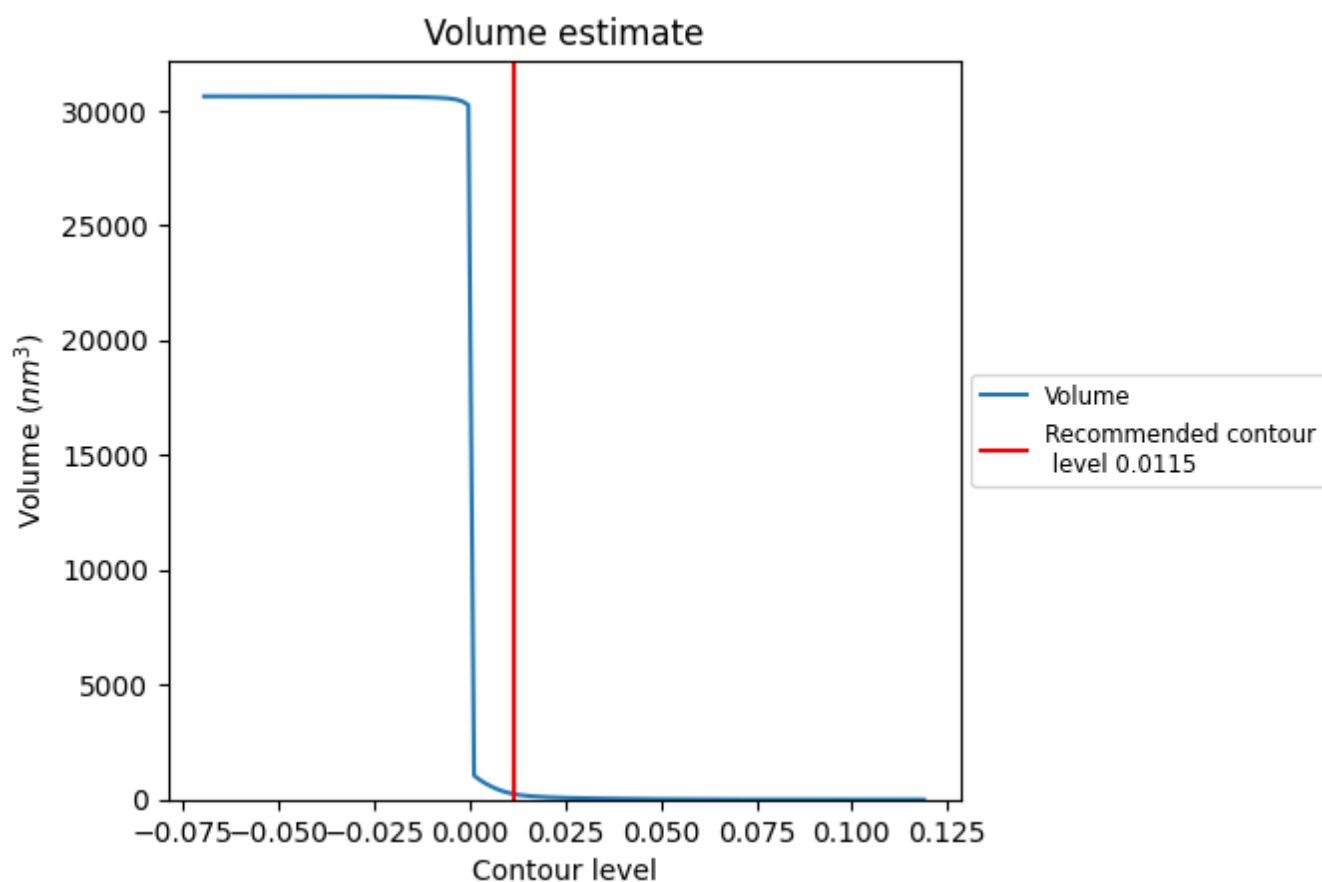
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

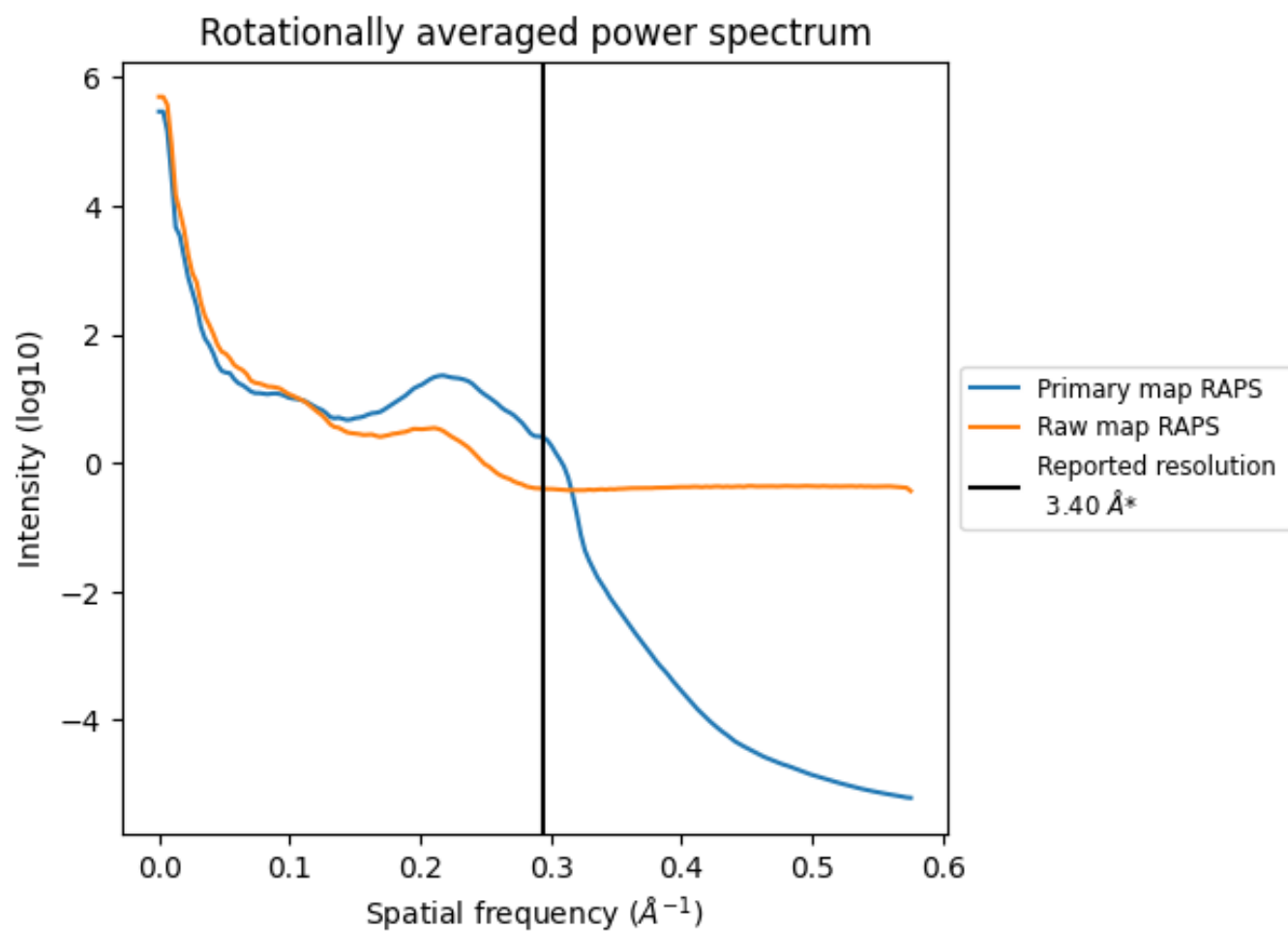
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 243 nm^3 ; this corresponds to an approximate mass of 219 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 ⓘ

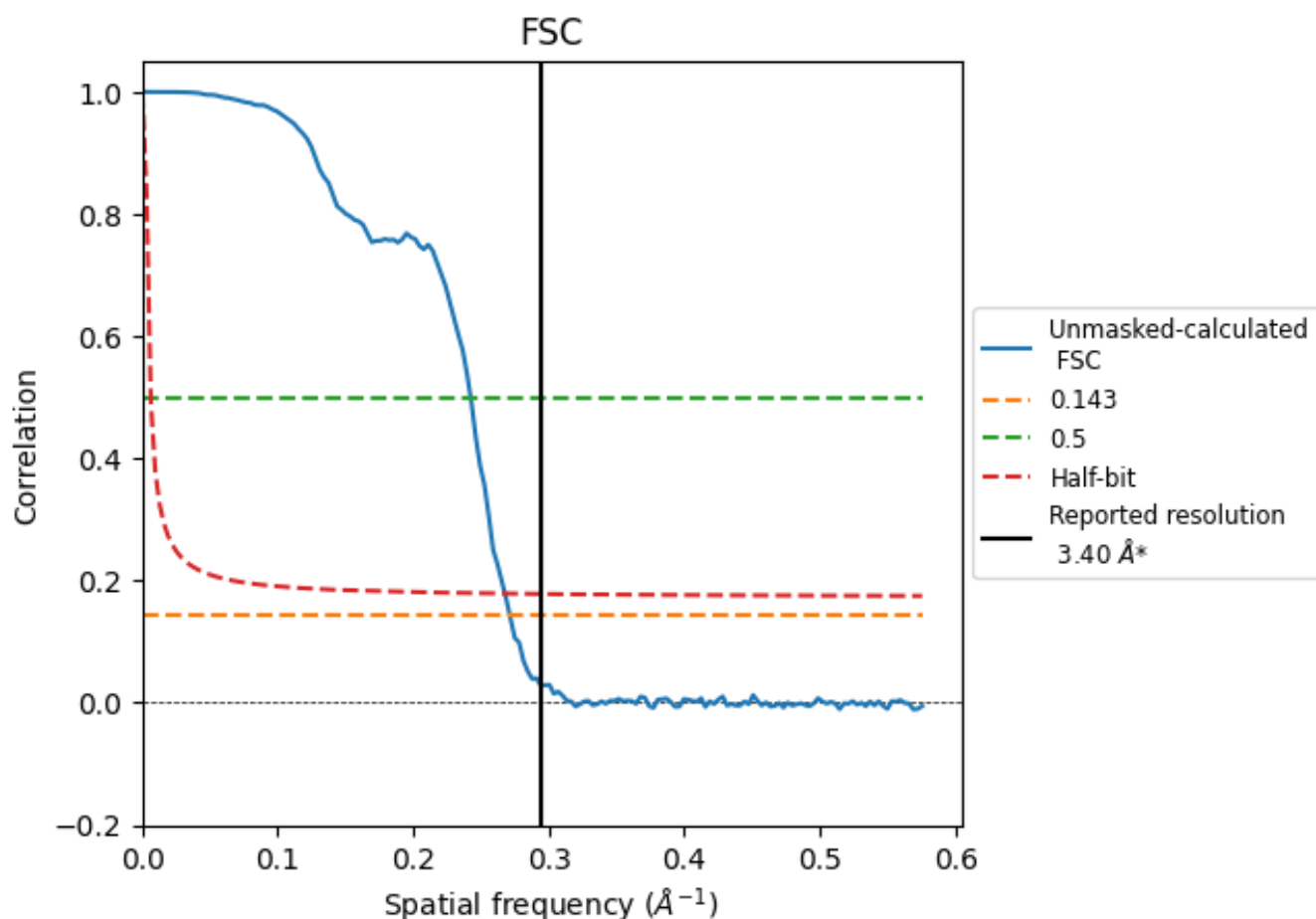


*Reported resolution corresponds to spatial frequency of 0.294 $\text{\AA}^{-1}$

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 0.294 Å⁻¹

8.2 Resolution estimates [i](#)

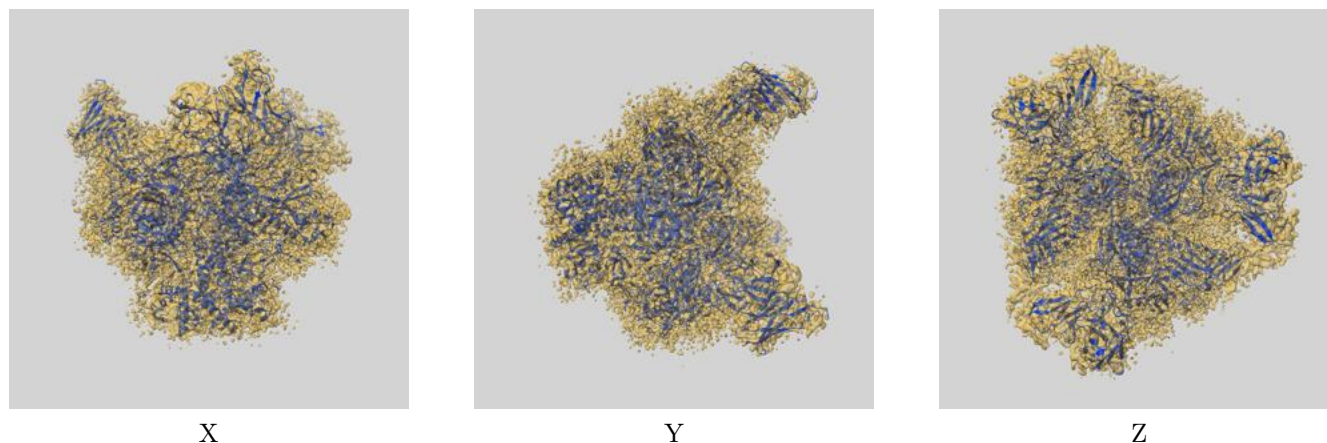
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.69	4.13	3.74

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

9 Map-model fit [i](#)

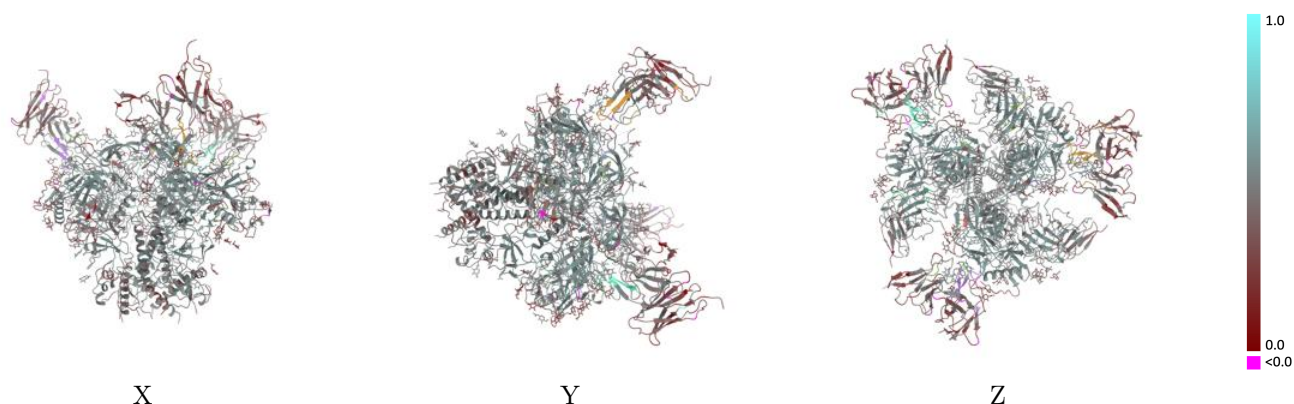
This section contains information regarding the fit between EMDB map EMD-26496 and PDB model 7UGQ. Per-residue inclusion information can be found in section [3](#) on page [15](#).

9.1 Map-model overlay [i](#)



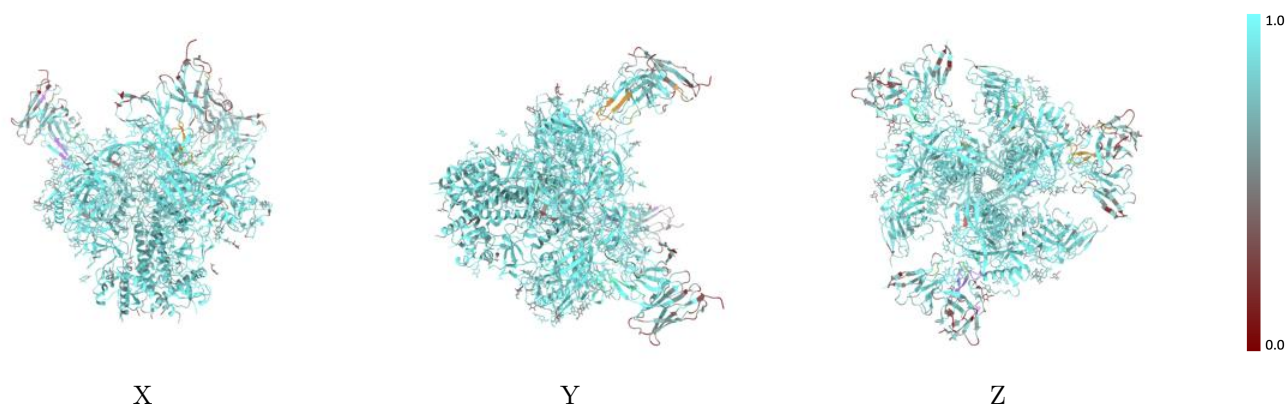
The images above show the 3D surface view of the map at the recommended contour level 0.0115 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



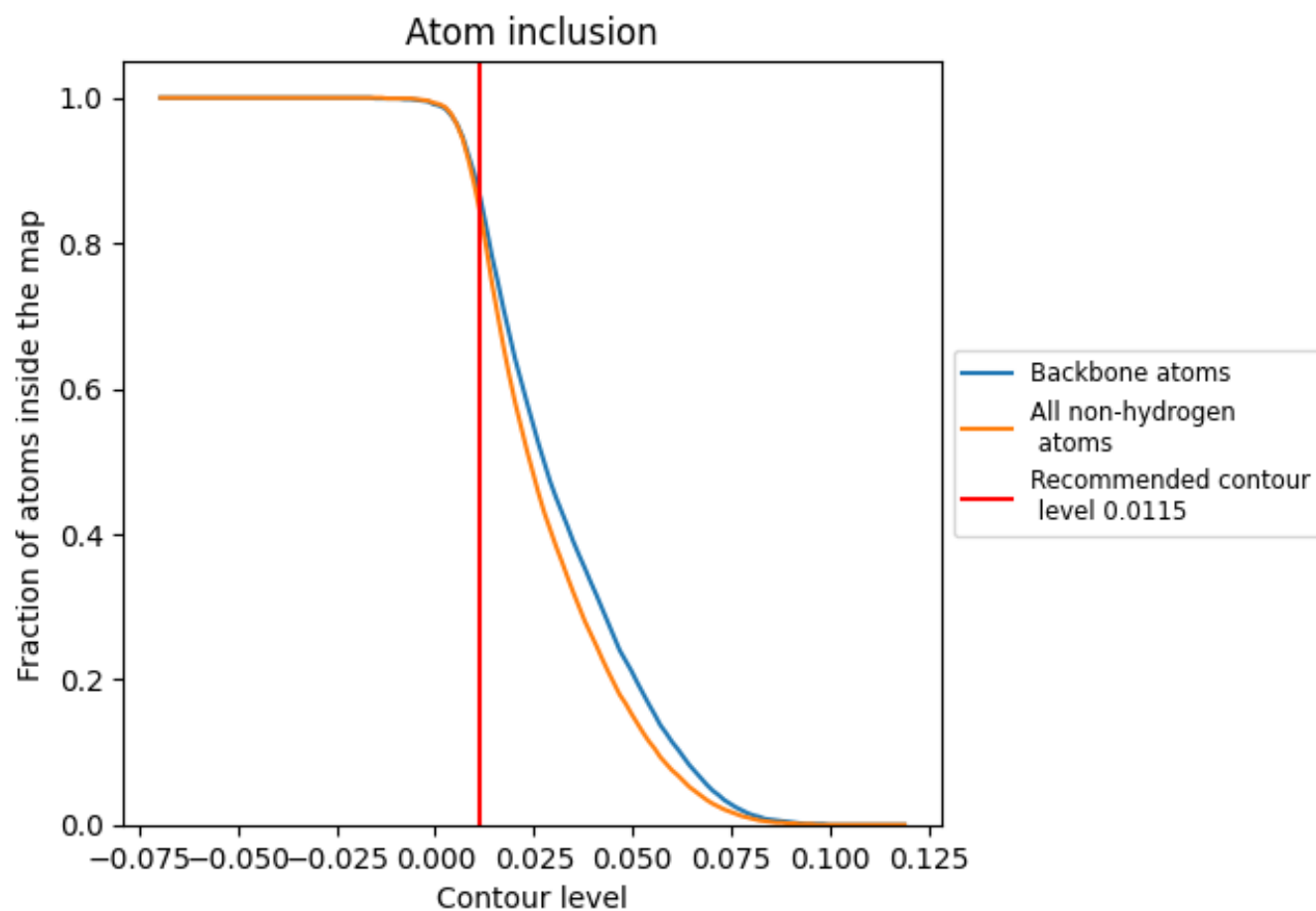
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0115).

9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 84% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.0115) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8390	 0.4700
0	 0.6150	 0.3230
1	 0.8000	 0.4520
A	 0.8890	 0.5090
B	 0.8920	 0.5090
C	 0.8910	 0.5100
D	 0.8520	 0.4660
E	 0.8560	 0.4690
F	 0.8530	 0.4680
G	 0.9000	 0.5280
H	 0.8990	 0.5230
I	 0.8960	 0.5230
J	 0.8350	 0.4460
K	 0.8390	 0.4470
L	 0.8440	 0.4460
M	 0.6580	 0.3690
N	 0.6510	 0.3690
O	 0.6530	 0.3640
P	 0.7900	 0.4220
Q	 0.7930	 0.4170
R	 0.7900	 0.4170
S	 0.8210	 0.4340
T	 0.8570	 0.4680
U	 0.7860	 0.4400
V	 0.6670	 0.3180
W	 0.8200	 0.4350
X	 0.8210	 0.4790
Y	 0.7860	 0.3310
Z	 0.5000	 0.3260
a	 0.7140	 0.3840
b	 0.8930	 0.4870
c	 0.7180	 0.3190
d	 0.8000	 0.4460
e	 0.8210	 0.4350
f	 0.8570	 0.4600



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Chain	Atom inclusion	Q-score
g	 0.7860	 0.4220
h	 0.7440	 0.3850
i	 0.8530	 0.4370
j	 0.7860	 0.4660
k	 0.7500	 0.3370
l	 0.5000	 0.3130
m	 0.8570	 0.4290
n	 0.8570	 0.4750
o	 0.6670	 0.3470
p	 0.8300	 0.4660
q	 0.8210	 0.4400
r	 0.8570	 0.4590
s	 0.7860	 0.4300
t	 0.6670	 0.3360
u	 0.8530	 0.4360
v	 0.7860	 0.4850
w	 0.7500	 0.3560
x	 0.5360	 0.3360
y	 0.7140	 0.3920
z	 0.8570	 0.4750