



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 01:27 PM UTC

PDB ID : 6NF2 / pdb_00006nf2
EMDB ID : EMD-9359
Title : Cryo-EM structure of vaccine-elicited antibody 0PV-c.01 in complex with HIV-1 Env BG505 DS-SOSIP and antibodies VRC03 and PGT122
Authors : Gorman, J.; Kwong, P.D.
Deposited on : 2018-12-18
Resolution : 3.70 Å(reported)
Based on initial model : 6CDI

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

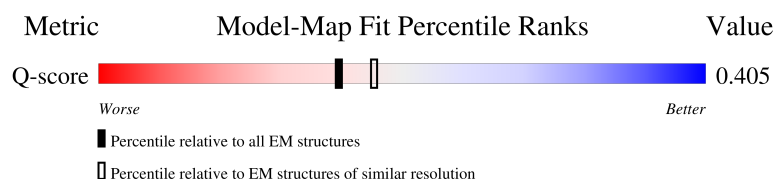
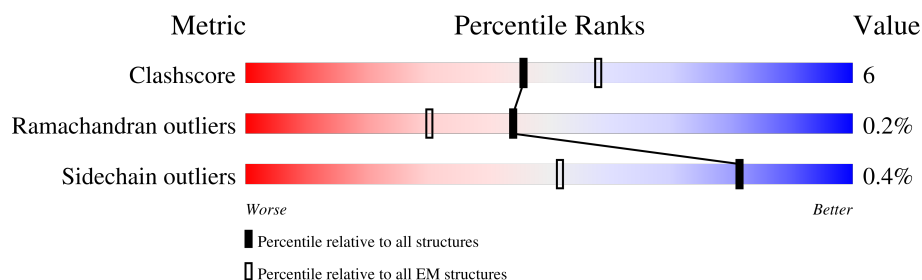
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.70 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.



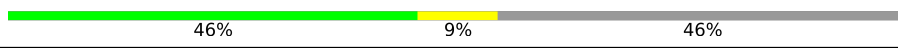
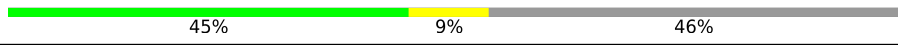
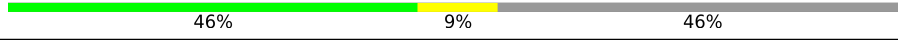
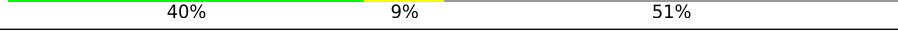
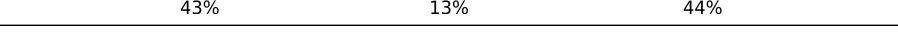
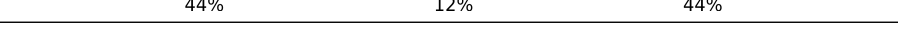
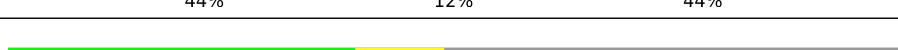
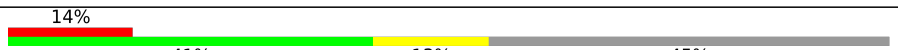
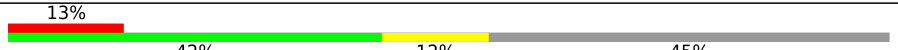
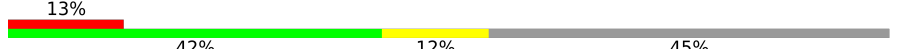
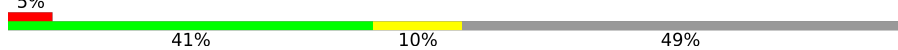
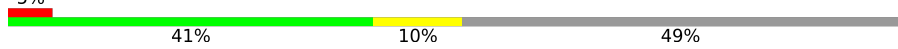
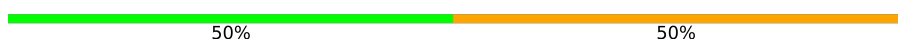
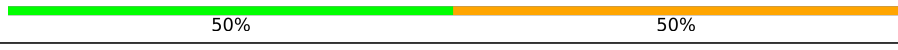
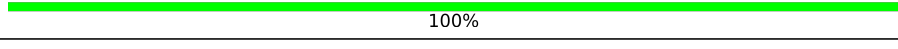
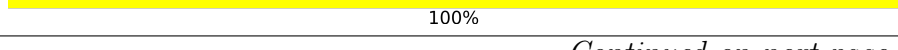
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11569 (3.20 - 4.20)

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	480	
1	G	480	
1	Q	480	
2	B	153	

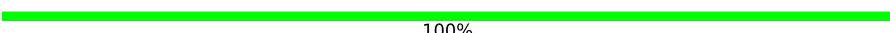
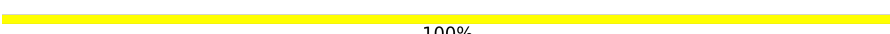
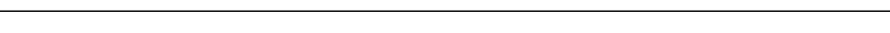
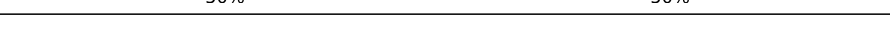
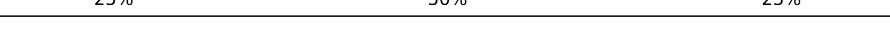
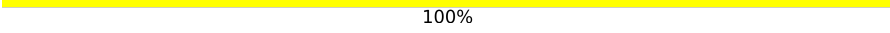
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Mol	Chain	Length	Quality of chain
2	I	153	
2	R	153	
3	C	235	
3	N	235	
3	V	235	
4	D	209	
4	M	209	
4	U	209	
5	E	235	
5	J	235	
5	S	235	
6	F	213	
6	K	213	
6	T	213	
7	H	229	
7	O	229	
7	W	229	
8	L	219	
8	P	219	
8	X	219	
9	0	2	
9	4	2	
9	5	2	
9	CA	2	
9	DA	2	

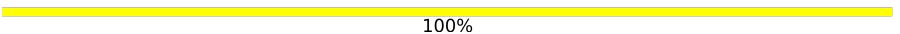
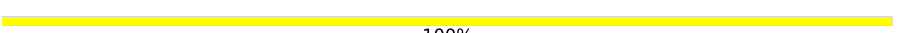
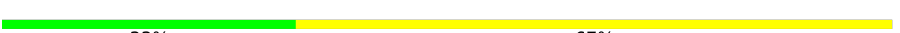
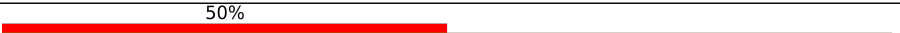
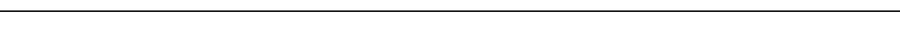
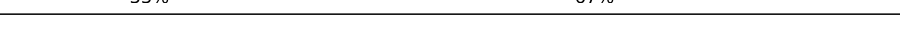
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Mol	Chain	Length	Quality of chain
9	GA	2	 50% 50%
9	Y	2	 50% 50%
9	Z	2	 50% 50%
9	g	2	 100%
9	h	2	 100%
9	k	2	 50% 50%
9	o	2	 50% 50%
9	p	2	 50% 50%
9	w	2	 100%
9	x	2	 100%
10	3	4	 50% 25% 25%
10	6	4	 25% 50% 25%
10	JA	4	 50% 50%
10	a	4	 25% 50% 25%
10	n	4	 50% 50%
10	q	4	 25% 50% 25%
11	7	5	 20% 80%
11	FA	5	 60% 40%
11	b	5	 20% 80%
11	j	5	 60% 40%
11	r	5	 20% 80%
11	z	5	 60% 40%
12	2	3	 100%
12	8	3	 33% 67%
12	AA	3	 67% 33%

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Mol	Chain	Length	Quality of chain
12	EA	3	 67% 33%
12	IA	3	 100%
12	c	3	 33% 67%
12	e	3	 67% 33%
12	i	3	 67% 33%
12	m	3	 100%
12	s	3	 33% 67%
12	u	3	 67% 33%
12	y	3	 100%
13	9	4	 50% 50% 50%
13	d	4	 50% 50% 50%
13	t	4	 50% 50% 50%
14	BA	6	 33% 67%
14	f	6	 33% 67%
14	v	6	 33% 67%
15	1	9	 11% 67% 22%
15	HA	9	 11% 67% 22%
15	l	9	 11% 67% 22%

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 32604 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	453	Total	C	N	O	S	0	0
			3564	2233	630	671	30		
1	G	453	Total	C	N	O	S	0	0
			3564	2233	630	671	30		
1	Q	453	Total	C	N	O	S	0	0
			3564	2233	630	671	30		

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	CYS	ILE	engineered mutation	UNP Q2N0S6
A	332	ASN	THR	engineered mutation	UNP Q2N0S6
A	433	CYS	ALA	engineered mutation	UNP Q2N0S6
A	501	CYS	ALA	engineered mutation	UNP Q2N0S6
A	509	ARG	GLU	engineered mutation	UNP Q2N0S6
A	510	ARG	LYS	engineered mutation	UNP Q2N0S6
A	512	ARG	ALA	engineered mutation	UNP Q2N0S6
G	201	CYS	ILE	engineered mutation	UNP Q2N0S6
G	332	ASN	THR	engineered mutation	UNP Q2N0S6
G	433	CYS	ALA	engineered mutation	UNP Q2N0S6
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	GLU	engineered mutation	UNP Q2N0S6
G	510	ARG	LYS	engineered mutation	UNP Q2N0S6
G	512	ARG	ALA	engineered mutation	UNP Q2N0S6
Q	201	CYS	ILE	engineered mutation	UNP Q2N0S6
Q	332	ASN	THR	engineered mutation	UNP Q2N0S6
Q	433	CYS	ALA	engineered mutation	UNP Q2N0S6
Q	501	CYS	ALA	engineered mutation	UNP Q2N0S6
Q	509	ARG	GLU	engineered mutation	UNP Q2N0S6
Q	510	ARG	LYS	engineered mutation	UNP Q2N0S6
Q	512	ARG	ALA	engineered mutation	UNP Q2N0S6

- Molecule 2 is a protein called Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	132	Total	C	N	O	S	0	0
			1034	654	178	196	6		
2	I	132	Total	C	N	O	S	0	0
			1034	654	178	196	6		
2	R	132	Total	C	N	O	S	0	0
			1034	654	178	196	6		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6
I	559	PRO	ILE	engineered mutation	UNP Q2N0S6
I	605	CYS	THR	engineered mutation	UNP Q2N0S6
R	559	PRO	ILE	engineered mutation	UNP Q2N0S6
R	605	CYS	THR	engineered mutation	UNP Q2N0S6

- Molecule 3 is a protein called VRC03 Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	128	Total	C	N	O	S	0	0
			1023	657	175	185	6		
3	N	128	Total	C	N	O	S	0	0
			1023	657	175	185	6		
3	V	128	Total	C	N	O	S	0	0
			1023	657	175	185	6		

- Molecule 4 is a protein called VRC03 Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	102	Total	C	N	O	S	0	0
			802	510	137	152	3		
4	M	102	Total	C	N	O	S	0	0
			802	510	137	152	3		
4	U	102	Total	C	N	O	S	0	0
			802	510	137	152	3		

- Molecule 5 is a protein called PGT122 Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		

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Mol	Chain	Residues	Atoms					AltConf	Trace
5	J	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		
5	S	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		

- Molecule 6 is a protein called PGT122 Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	105	Total	C	N	O	S	0	0
			805	504	139	160	2		
6	K	105	Total	C	N	O	S	0	0
			805	504	139	160	2		
6	T	105	Total	C	N	O	S	0	0
			805	504	139	160	2		

- Molecule 7 is a protein called 0PV-c.01 Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	H	125	Total	C	N	O	S	0	0
			951	603	164	182	2		
7	O	125	Total	C	N	O	S	0	0
			951	603	164	182	2		
7	W	125	Total	C	N	O	S	0	0
			951	603	164	182	2		

- Molecule 8 is a protein called 0PV-c.01 Light Chain.

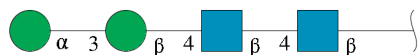
Mol	Chain	Residues	Atoms					AltConf	Trace
8	L	112	Total	C	N	O	S	0	0
			855	538	139	173	5		
8	P	112	Total	C	N	O	S	0	0
			855	538	139	173	5		
8	X	112	Total	C	N	O	S	0	0
			855	538	139	173	5		

- Molecule 9 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
9	Y	2	Total	C	N	O	0	0
			28	16	2	10		
9	Z	2	Total	C	N	O	0	0
			28	16	2	10		
9	g	2	Total	C	N	O	0	0
			28	16	2	10		
9	h	2	Total	C	N	O	0	0
			28	16	2	10		
9	k	2	Total	C	N	O	0	0
			28	16	2	10		
9	o	2	Total	C	N	O	0	0
			28	16	2	10		
9	p	2	Total	C	N	O	0	0
			28	16	2	10		
9	w	2	Total	C	N	O	0	0
			28	16	2	10		
9	x	2	Total	C	N	O	0	0
			28	16	2	10		
9	0	2	Total	C	N	O	0	0
			28	16	2	10		
9	4	2	Total	C	N	O	0	0
			28	16	2	10		
9	5	2	Total	C	N	O	0	0
			28	16	2	10		
9	CA	2	Total	C	N	O	0	0
			28	16	2	10		
9	DA	2	Total	C	N	O	0	0
			28	16	2	10		
9	GA	2	Total	C	N	O	0	0
			28	16	2	10		

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



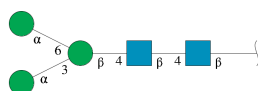
Mol	Chain	Residues	Atoms				AltConf	Trace
10	a	4	Total	C	N	O	0	0
			50	28	2	20		
10	n	4	Total	C	N	O	0	0
			50	28	2	20		

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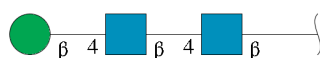
Mol	Chain	Residues	Atoms				AltConf	Trace
10	q	4	Total	C	N	O	0	0
			50	28	2	20		
10	3	4	Total	C	N	O	0	0
			50	28	2	20		
10	6	4	Total	C	N	O	0	0
			50	28	2	20		
10	JA	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 11 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
11	b	5	Total	C	N	O	0	0
			61	34	2	25		
11	j	5	Total	C	N	O	0	0
			61	34	2	25		
11	r	5	Total	C	N	O	0	0
			61	34	2	25		
11	z	5	Total	C	N	O	0	0
			61	34	2	25		
11	7	5	Total	C	N	O	0	0
			61	34	2	25		
11	FA	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 12 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
12	c	3	Total	C	N	O	0	0
			39	22	2	15		

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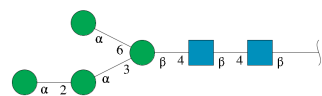
Mol	Chain	Residues	Atoms				AltConf	Trace
12	e	3	Total	C	N	O	0	0
			39	22	2	15		
12	i	3	Total	C	N	O	0	0
			39	22	2	15		
12	m	3	Total	C	N	O	0	0
			39	22	2	15		
12	s	3	Total	C	N	O	0	0
			39	22	2	15		
12	u	3	Total	C	N	O	0	0
			39	22	2	15		
12	y	3	Total	C	N	O	0	0
			39	22	2	15		
12	2	3	Total	C	N	O	0	0
			39	22	2	15		
12	8	3	Total	C	N	O	0	0
			39	22	2	15		
12	AA	3	Total	C	N	O	0	0
			39	22	2	15		
12	EA	3	Total	C	N	O	0	0
			39	22	2	15		
12	IA	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 13 is an oligosaccharide called 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.



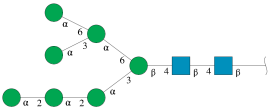
Mol	Chain	Residues	Atoms				AltConf	Trace
13	d	4	Total	C	N	O	0	0
			50	28	2	20		
13	t	4	Total	C	N	O	0	0
			50	28	2	20		
13	9	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 14 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-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
14	f	6	Total	C	N	O	0	0
			72	40	2	30		
14	v	6	Total	C	N	O	0	0
			72	40	2	30		
14	BA	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 15 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
15	1	9	Total	C	N	O	0	0
			105	58	2	45		
15	1	9	Total	C	N	O	0	0
			105	58	2	45		
15	HA	9	Total	C	N	O	0	0
			105	58	2	45		

- Molecule 16 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

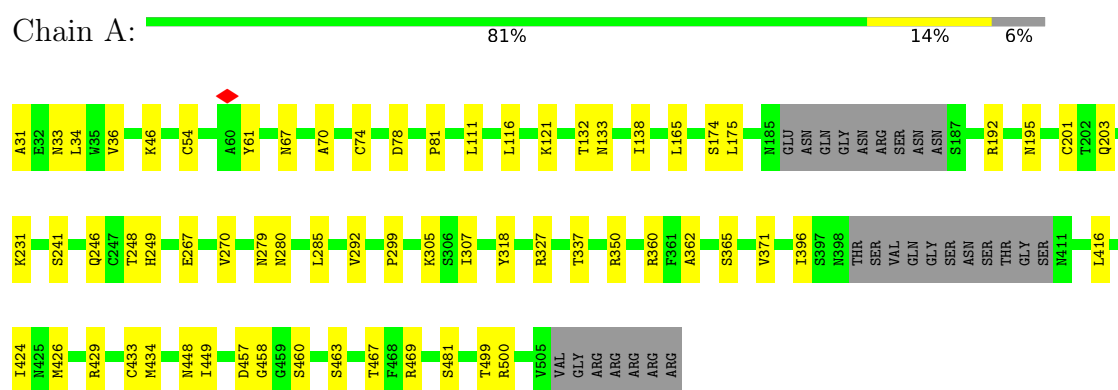


Mol	Chain	Residues	Atoms				AltConf
16	A	1	Total	C	N	O	0
			14	8	1	5	
16	B	1	Total	C	N	O	0
			14	8	1	5	
16	B	1	Total	C	N	O	0
			14	8	1	5	
16	G	1	Total	C	N	O	0
			14	8	1	5	
16	I	1	Total	C	N	O	0
			14	8	1	5	
16	I	1	Total	C	N	O	0
			14	8	1	5	
16	Q	1	Total	C	N	O	0
			14	8	1	5	
16	R	1	Total	C	N	O	0
			14	8	1	5	
16	R	1	Total	C	N	O	0
			14	8	1	5	

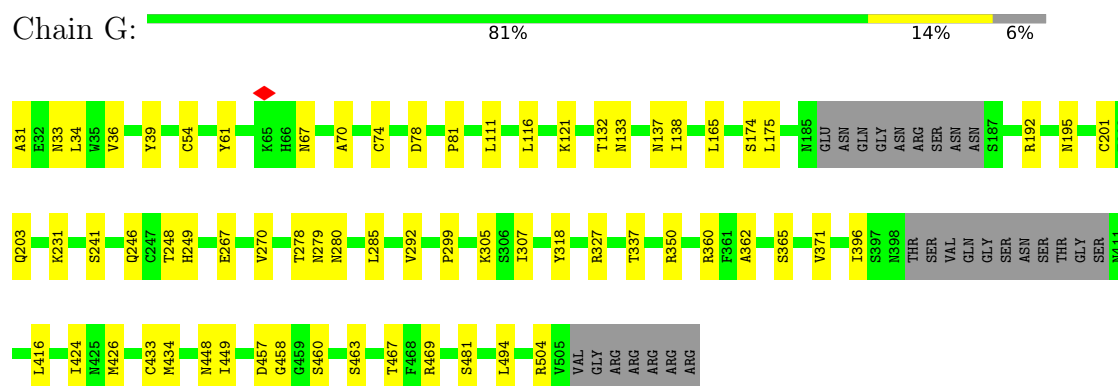
3 Residue-property plots [i](#)

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

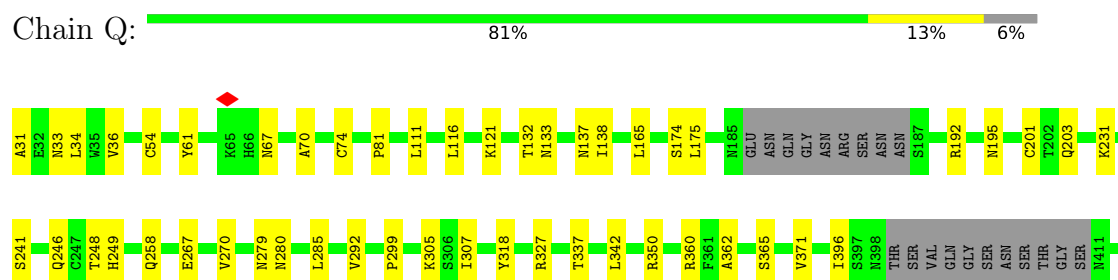
- Molecule 1: Envelope glycoprotein gp120

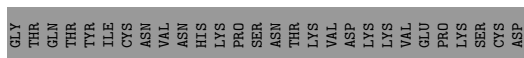


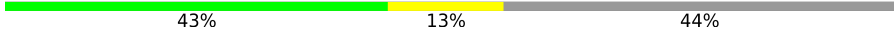
- Molecule 1: Envelope glycoprotein gp120

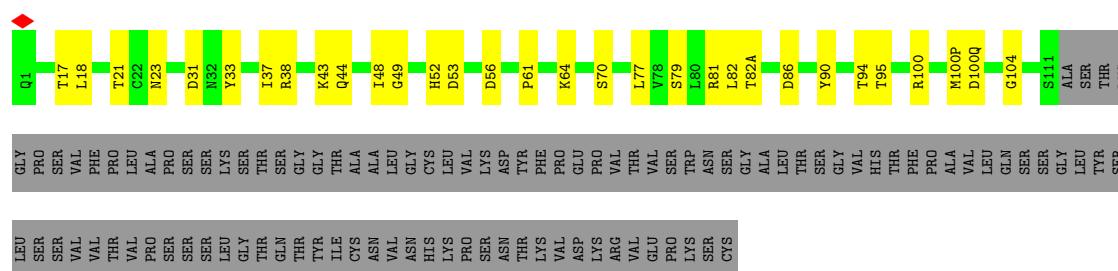


- Molecule 1: Envelope glycoprotein gp120

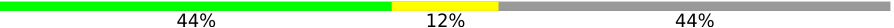


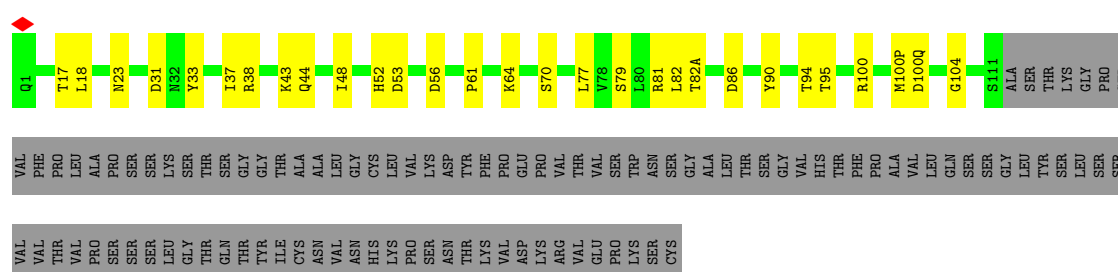


Chain E: 

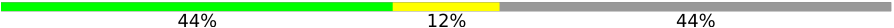


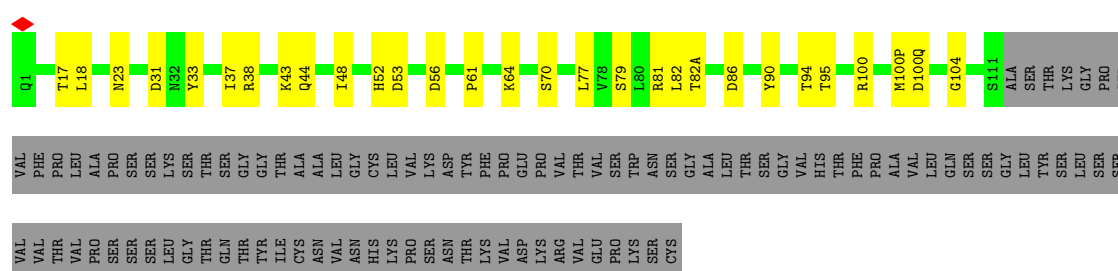
• Molecule 5: PGT122 Heavy Chain

Chain J: 



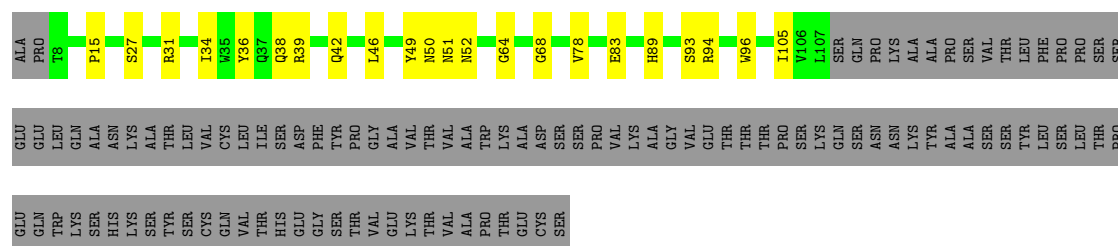
• Molecule 5: PGT122 Heavy Chain

Chain S: 



• Molecule 6: PGT122 Light Chain

Chain F: 

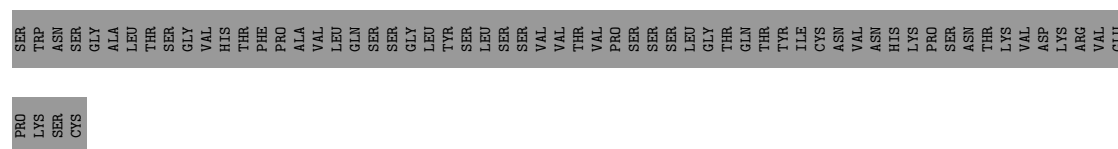


• Molecule 6: PGT122 Light Chain

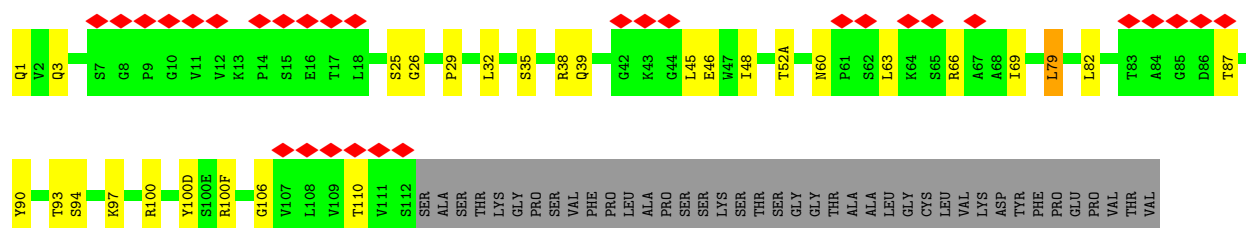
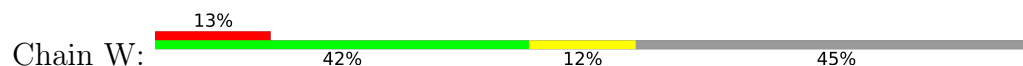
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LYS	LYS	PRO
SER	ALA	T8
TYR	THR	
SER	LEU	P15
CYS	VAL	
GLN	CYS	S27
VAL	LEU	
THR	ILE	R31
HIS	SER	
GLY	ASP	I34
GLU	PHE	
SER	TYR	Q38
THR	PRO	R39
VAL	GLY	
GLU	ALA	Q42
LYS	VAL	
THR	THR	L46
VAL	VAL	
ALA	ALA	Y49
PRO	TRP	N50
THR	LYS	N51
GLY	ALA	N52
CYS	ASP	
SER	SER	G64
	SER	
	PRO	G68
	VAL	
	LYS	V78
	ALA	
	GLY	E83
	VAL	
	GLU	H89
	THR	
	THR	S93
	THR	
	PRO	I105
	SER	V106
	SER	L107
	LYS	
	GLN	SER
	SER	
	ASN	GLN
	ASN	PRO
	LYS	LYS
	LYS	ALA
	TYR	ALA
	ALA	PRO
	ALA	SER
	SER	VAL
	SER	THR
	TYR	LEU
	LEU	PHE
	SER	PRO
	LEU	PRO
	THR	SER
	PRO	SER
	GLU	GLU
	GLN	GLU
	TRP	LEU
	LYS	GLN
SER	SER	ALA

SER	ALA	ALA
HIS	ASN	PRO
LYS	LYS	T8
SER	ALA	P15
TYR	THR	S27
SER	LEU	R31
CYS	VAL	I34
GLN	LEU	Q38
VAL	LEU	R39
THR	IIE	Q42
HIS	SER	L46
GLU	ASP	V49
GLY	PHE	N50
SER	TYR	N51
THR	PRO	N52
VAL	GLY	G64
GLU	ALA	G68
LYS	VAL	V78
THR	THR	E83
VAL	VAL	H89
VAL	LYS	S93
PRO	ALA	R94
THR	THR	I105
THR	PRO	V106
SER	SER	L107
LYS	LYS	SER
GLN	GLN	GLN
SER	SER	GLN
ASN	ASN	PRO
LYS	LYS	LYS
TYR	TYR	ALA
ALA	ALA	PRO
SER	ALA	SER
SER	SER	VAL
TYR	TYR	THR
LEU	LEU	LEU
SER	SER	PHE
PRO	PRO	PRO
SER	SER	PRO
THR	THR	SER
PRO	PRO	SER
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GLU	GLU	GLU
LYS	LYS	LEU
TYR	TYR	GLN

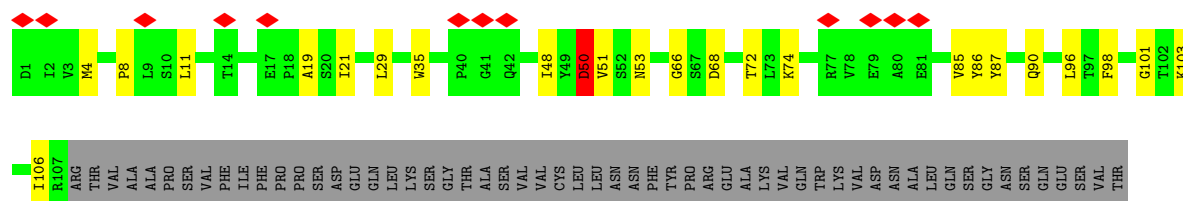
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Y90	T93	S94	R95	A96	K97	R100	Y100B	G106	V107	L108	V109	V111	S112	SER	ALA	SER	THR	LYS	GLY	PRO	SER	VAL	PHE	LEU	PRO	ALA	PRO	SER	SER	LYS	THR	GLY	GLY	THR	ALA	ALA	LEU	GLY	CYS	LEU	VAL	LYS	ASP	TYR	PHE	PRO	GLU	PRO	VAL	THR	VAL



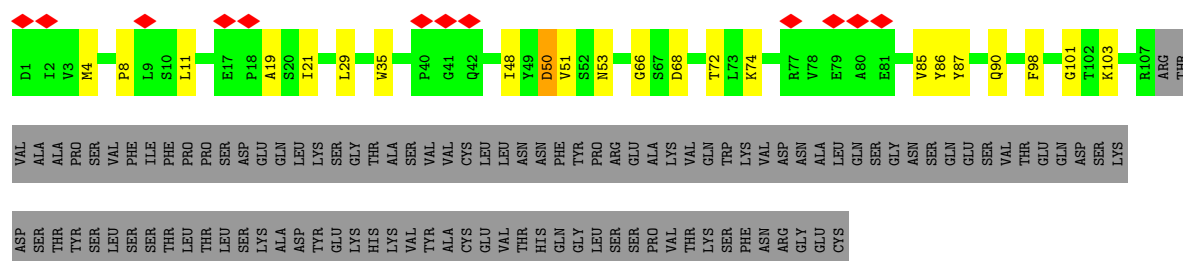
• Molecule 7: 0PV-c.01 Heavy Chain



• Molecule 8: 0PV-c.01 Light Chain

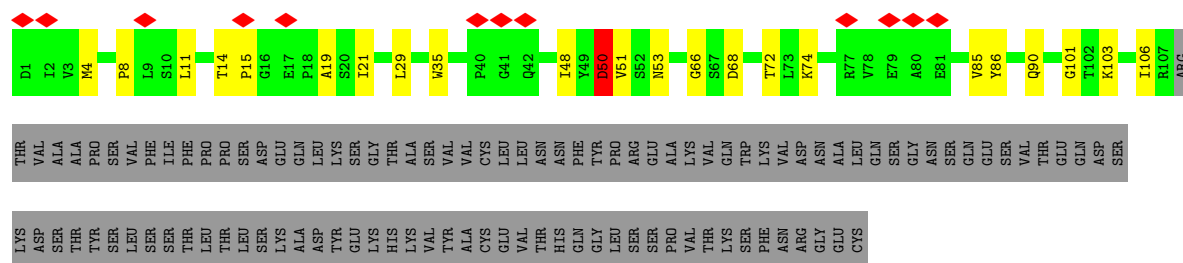


• Molecule 8: 0PV-c.01 Light Chain



• Molecule 8: 0PV-c.01 Light Chain





- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain g: 100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain h: 100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain k: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain o:  50% 50%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain p:  50% 50%

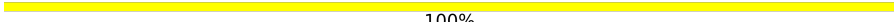

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain w:  100%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain x:  100%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 0:  50% 50%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 4:  50% 50%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 5:  50% 50%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain CA:  100%

NAG1
NAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain DA:  100%

NAG1
NAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain GA:  50% 50%

NAG1
NAG2

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

Chain a:  25% 50% 25%

NAG1
NAG2
BMA3
MAN4

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

Chain n:  50% 50%

NAG1
NAG2
BMA3
MAN4

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

Chain q:  25% 50% 25%

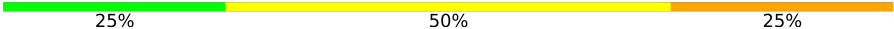
NAG1
NAG2
BMA3
MAN4

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

Chain 3:  50% 25% 25%



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

Chain 6:  25% 50% 25%

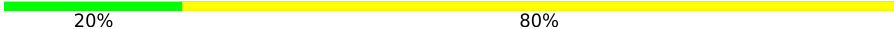


- Molecule 10: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain JA:  50% 50%

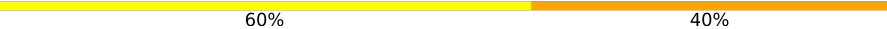


- Molecule 11: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:  20% 80%

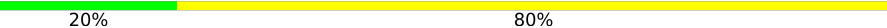


- Molecule 11: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain j:  60% 40%

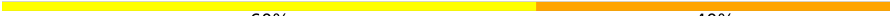


- Molecule 11: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain r:  20% 80%



- Molecule 11: 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 z:  60% 40%



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

Chain 7:  20% 80%



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

Chain FA:  60% 40%



- Molecule 12: β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain c:  33% 67%



- Molecule 12: β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain e:  67% 33%



- Molecule 12: β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain i:  67% 33%



- Molecule 12: β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain m:  100%



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain s:



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain u:



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain y:



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 2:



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 8:



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain AA:

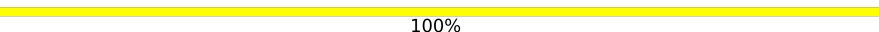


- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain EA:  67% 33%



- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain IA:  100%



- Molecule 13: 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:  50% 50% 50%



- Molecule 13: 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 t:  50% 50% 50%



- Molecule 13: 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 9:  50% 50% 50%

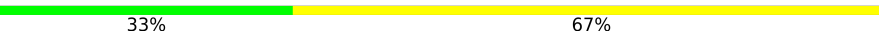


- Molecule 14: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain f:  33% 67%

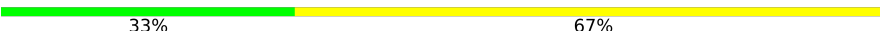


- Molecule 14: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain v:  33% 67%

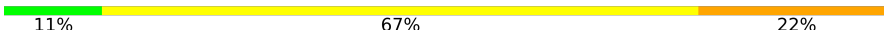


- Molecule 14: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain BA:  33% 67%

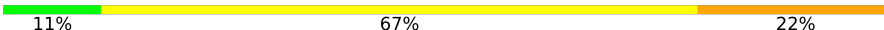


- Molecule 15: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain l:  11% 67% 22%

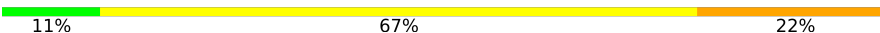


- Molecule 15: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain 1:  11% 67% 22%



- Molecule 15: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain HA:  11% 67% 22%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	98419	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	73.46	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.915	Depositor
Minimum map value	-1.365	Depositor
Average map value	-0.002	Depositor
Map value standard deviation	0.074	Depositor
Recommended contour level	0.4	Depositor
Map size (Å)	429.30002, 429.30002, 429.30002	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07325, 1.07325, 1.07325	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.42	0/3638	0.63	0/4939
1	G	0.42	0/3638	0.63	0/4939
1	Q	0.42	0/3638	0.63	0/4939
2	B	0.35	0/1052	0.70	0/1427
2	I	0.35	0/1052	0.70	0/1427
2	R	0.35	0/1052	0.69	0/1427
3	C	0.43	0/1056	0.52	0/1439
3	N	0.43	0/1056	0.52	0/1439
3	V	0.43	0/1056	0.52	0/1439
4	D	0.35	0/820	0.61	0/1107
4	M	0.35	0/820	0.61	0/1107
4	U	0.35	0/820	0.61	0/1107
5	E	0.31	0/1076	0.57	0/1465
5	J	0.31	0/1076	0.57	0/1465
5	S	0.32	0/1076	0.56	0/1465
6	F	0.33	0/826	0.60	0/1130
6	K	0.34	0/826	0.60	0/1130
6	T	0.33	0/826	0.60	0/1130
7	H	0.36	0/975	0.94	1/1332 (0.1%)
7	O	0.36	0/975	0.94	1/1332 (0.1%)
7	W	0.36	0/975	0.94	1/1332 (0.1%)
8	L	0.36	0/873	1.01	3/1187 (0.3%)
8	P	0.36	0/873	1.00	3/1187 (0.3%)
8	X	0.36	0/873	1.00	3/1187 (0.3%)
All	All	0.38	0/30948	0.69	12/42078 (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
8	L	0	1
8	P	0	1
8	X	0	1
All	All	0	3

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	P	53	ASN	CA-C-N	5.78	130.64	122.09
8	P	53	ASN	C-N-CA	5.78	130.64	122.09
8	X	53	ASN	CA-C-N	5.77	130.64	122.09
8	X	53	ASN	C-N-CA	5.77	130.64	122.09
8	L	53	ASN	CA-C-N	5.76	130.61	122.09
8	L	53	ASN	C-N-CA	5.76	130.61	122.09
7	W	79	LEU	CA-CB-CG	5.57	135.80	116.30
7	O	79	LEU	CA-CB-CG	5.56	135.75	116.30
7	H	79	LEU	CA-CB-CG	5.54	135.71	116.30
8	X	66	GLY	N-CA-C	5.45	118.53	110.63
8	L	66	GLY	N-CA-C	5.43	118.50	110.63
8	P	66	GLY	N-CA-C	5.40	118.47	110.63

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	L	50	ASP	Peptide
8	P	50	ASP	Peptide
8	X	50	ASP	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	3564	0	3489	43	0
1	G	3564	0	3489	45	0
1	Q	3564	0	3489	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1034	0	1013	12	0
2	I	1034	0	1013	14	0
2	R	1034	0	1013	9	0
3	C	1023	0	971	13	0
3	N	1023	0	971	14	0
3	V	1023	0	971	13	0
4	D	802	0	780	13	0
4	M	802	0	780	10	0
4	U	802	0	780	10	0
5	E	1047	0	1026	22	0
5	J	1047	0	1026	19	0
5	S	1047	0	1026	19	0
6	F	805	0	761	15	0
6	K	805	0	761	12	0
6	T	805	0	761	13	0
7	H	951	0	946	18	0
7	O	951	0	946	20	0
7	W	951	0	946	15	0
8	L	855	0	831	14	0
8	P	855	0	831	13	0
8	X	855	0	831	12	0
9	0	28	0	25	1	0
9	4	28	0	25	0	0
9	5	28	0	25	1	0
9	CA	28	0	25	0	0
9	DA	28	0	25	0	0
9	GA	28	0	25	1	0
9	Y	28	0	25	0	0
9	Z	28	0	25	1	0
9	g	28	0	25	0	0
9	h	28	0	25	0	0
9	k	28	0	25	1	0
9	o	28	0	25	0	0
9	p	28	0	25	1	0
9	w	28	0	25	0	0
9	x	28	0	25	0	0
10	3	50	0	43	1	0
10	6	50	0	43	1	0
10	JA	50	0	43	0	0
10	a	50	0	43	1	0
10	n	50	0	43	0	0
10	q	50	0	43	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	7	61	0	52	0	0
11	FA	61	0	52	2	0
11	b	61	0	52	0	0
11	j	61	0	52	2	0
11	r	61	0	52	0	0
11	z	61	0	52	2	0
12	2	39	0	34	0	0
12	8	39	0	34	0	0
12	AA	39	0	34	0	0
12	EA	39	0	34	0	0
12	IA	39	0	34	0	0
12	c	39	0	34	0	0
12	e	39	0	34	0	0
12	i	39	0	34	0	0
12	m	39	0	34	0	0
12	s	39	0	34	0	0
12	u	39	0	34	0	0
12	y	39	0	34	0	0
13	9	50	0	43	1	0
13	d	50	0	43	1	0
13	t	50	0	43	1	0
14	BA	72	0	61	0	0
14	f	72	0	61	0	0
14	v	72	0	61	0	0
15	1	105	0	88	4	0
15	HA	105	0	88	3	0
15	l	105	0	88	3	0
16	A	14	0	13	0	0
16	B	28	0	26	0	0
16	G	14	0	13	0	0
16	I	28	0	26	0	0
16	Q	14	0	13	0	0
16	R	28	0	26	0	0
All	All	32604	0	31497	398	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (398) 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)
5:J:90:TYR:O	5:J:104:GLY:HA2	1.70	0.92
5:S:90:TYR:O	5:S:104:GLY:HA2	1.70	0.91
5:E:90:TYR:O	5:E:104:GLY:HA2	1.70	0.89
8:X:85:VAL:HG22	8:X:103:LYS:HG3	1.72	0.72
8:P:85:VAL:HG22	8:P:103:LYS:HG3	1.72	0.72
8:L:85:VAL:HG22	8:L:103:LYS:HG3	1.72	0.71
7:W:100(F):ARG:NH1	8:X:50:ASP:OD2	2.24	0.70
2:I:570:VAL:O	2:I:573:ILE:HB	1.92	0.70
2:R:570:VAL:O	2:R:573:ILE:HB	1.92	0.70
2:B:570:VAL:O	2:B:573:ILE:HB	1.92	0.69
7:H:39:GLN:NE2	8:L:87:TYR:OH	2.26	0.68
2:B:617:ARG:NH2	2:B:626:MET:SD	2.70	0.65
4:M:36:GLN:HE21	4:M:44:ARG:HD3	1.63	0.64
4:U:36:GLN:HE21	4:U:44:ARG:HD3	1.63	0.64
4:D:36:GLN:HE21	4:D:44:ARG:HD3	1.63	0.63
7:H:46:GLU:OE2	7:H:60:ASN:ND2	2.32	0.63
2:R:617:ARG:NH2	2:R:626:MET:SD	2.70	0.63
7:W:46:GLU:OE2	7:W:60:ASN:ND2	2.32	0.63
1:Q:231:LYS:HD2	1:Q:267:GLU:HG3	1.80	0.63
1:G:231:LYS:HD2	1:G:267:GLU:HG3	1.80	0.62
5:J:18:LEU:HB3	5:J:82:LEU:HB3	1.80	0.62
7:O:46:GLU:OE2	7:O:60:ASN:ND2	2.32	0.62
1:A:195:ASN:ND2	1:A:201:CYS:SG	2.72	0.62
1:G:195:ASN:ND2	1:G:201:CYS:SG	2.72	0.62
2:I:617:ARG:NH2	2:I:626:MET:SD	2.70	0.62
5:J:100:ARG:NH2	15:l:4:MAN:O6	2.33	0.62
5:S:18:LEU:HB3	5:S:82:LEU:HB3	1.80	0.62
5:E:18:LEU:HB3	5:E:82:LEU:HB3	1.81	0.62
1:A:231:LYS:HD2	1:A:267:GLU:HG3	1.80	0.61
1:Q:195:ASN:ND2	1:Q:201:CYS:SG	2.72	0.61
3:N:98:CYS:SG	3:N:100(A):CYS:N	2.73	0.61
3:V:98:CYS:SG	3:V:100(A):CYS:N	2.73	0.61
3:C:98:CYS:SG	3:C:100(A):CYS:N	2.73	0.61
1:G:70:ALA:HB2	1:G:111:LEU:HD11	1.84	0.60
1:Q:70:ALA:HB2	1:Q:111:LEU:HD11	1.84	0.60
7:H:1:GLN:N	7:H:26:GLY:O	2.33	0.60
7:W:1:GLN:N	7:W:26:GLY:O	2.33	0.60
11:j:2:NAG:H82	9:k:1:NAG:H83	1.84	0.60
1:A:81:PRO:HA	1:A:246:GLN:HE21	1.67	0.60
7:O:1:GLN:N	7:O:26:GLY:O	2.33	0.60
11:z:2:NAG:H82	9:0:1:NAG:H83	1.84	0.60
4:D:33:THR:HG22	4:D:48:TYR:HA	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:81:PRO:HA	1:G:246:GLN:HE21	1.67	0.59
1:Q:81:PRO:HA	1:Q:246:GLN:HE21	1.67	0.59
4:U:33:THR:HG22	4:U:48:TYR:HA	1.84	0.58
1:A:70:ALA:HB2	1:A:111:LEU:HD11	1.83	0.58
11:FA:2:NAG:H82	9:GA:1:NAG:H83	1.84	0.58
5:J:52:HIS:HB3	5:J:56:ASP:HB3	1.86	0.58
5:S:52:HIS:HB3	5:S:56:ASP:HB3	1.85	0.58
8:X:19:ALA:O	8:X:74:LYS:HA	2.03	0.58
4:M:33:THR:HG22	4:M:48:TYR:HA	1.84	0.58
8:P:19:ALA:O	8:P:74:LYS:HA	2.03	0.58
1:Q:457:ASP:O	3:V:61:ARG:NH1	2.36	0.58
8:L:19:ALA:O	8:L:74:LYS:HA	2.03	0.58
3:C:61:ARG:NH1	1:G:457:ASP:O	2.37	0.57
5:S:100:ARG:NH2	15:HA:4:MAN:O6	2.37	0.57
3:N:38:ARG:NH1	3:N:46:GLU:OE1	2.38	0.57
3:V:12:LYS:HG3	3:V:18:VAL:HG12	1.86	0.57
3:C:12:LYS:HG3	3:C:18:VAL:HG12	1.87	0.57
5:E:52:HIS:HB3	5:E:56:ASP:HB3	1.85	0.57
1:A:365:SER:OG	1:A:469:ARG:NH1	2.39	0.56
3:V:38:ARG:NH1	3:V:46:GLU:OE1	2.38	0.56
3:N:12:LYS:HG3	3:N:18:VAL:HG12	1.86	0.56
1:Q:365:SER:OG	1:Q:469:ARG:NH1	2.39	0.56
3:C:38:ARG:NH1	3:C:46:GLU:OE1	2.38	0.56
1:G:305:LYS:HE2	1:G:307:ILE:HD11	1.88	0.56
5:E:31:ASP:OD1	5:E:31:ASP:N	2.38	0.56
1:A:305:LYS:HE2	1:A:307:ILE:HD11	1.88	0.56
5:E:100:ARG:NH2	15:1:4:MAN:O6	2.39	0.56
7:O:39:GLN:NE2	8:P:87:TYR:OH	2.39	0.56
5:E:70:SER:OG	5:E:79:SER:OG	2.24	0.55
7:H:60:ASN:HB3	7:H:63:LEU:HD23	1.88	0.55
6:F:94:ARG:O	1:G:137:ASN:N	2.39	0.55
6:K:50:ASN:ND2	15:l:5:MAN:O6	2.39	0.55
5:J:70:SER:OG	5:J:79:SER:OG	2.24	0.55
7:O:60:ASN:HB3	7:O:63:LEU:HD23	1.88	0.55
5:S:31:ASP:OD1	5:S:31:ASP:N	2.38	0.55
2:I:514:GLY:O	7:O:95:ARG:NH2	2.40	0.55
7:O:44:GLY:HA2	8:P:87:TYR:HE1	1.72	0.55
1:Q:305:LYS:HE2	1:Q:307:ILE:HD11	1.88	0.55
4:U:37:LYS:HD2	4:U:43:PRO:HG3	1.89	0.55
7:H:90:TYR:O	7:H:106:GLY:HA2	2.07	0.55
5:S:70:SER:OG	5:S:79:SER:OG	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:W:60:ASN:HB3	7:W:63:LEU:HD23	1.88	0.54
1:A:458:GLY:HA3	3:N:47:TRP:HZ3	1.71	0.54
7:H:100(F):ARG:NH1	8:L:50:ASP:OD2	2.40	0.54
8:L:21:ILE:O	8:L:72:THR:HA	2.08	0.54
1:A:457:ASP:O	3:N:61:ARG:NH1	2.40	0.54
4:D:37:LYS:HD2	4:D:43:PRO:HG3	1.89	0.54
5:E:70:SER:HB3	5:E:81:ARG:HH21	1.73	0.54
7:W:90:TYR:O	7:W:106:GLY:HA2	2.07	0.54
5:J:70:SER:HB3	5:J:81:ARG:HH21	1.73	0.54
8:P:21:ILE:O	8:P:72:THR:HA	2.08	0.54
1:A:195:ASN:ND2	1:A:433:CYS:SG	2.80	0.54
3:C:56:ALA:HB2	1:G:371:VAL:HG11	1.88	0.54
7:O:90:TYR:O	7:O:106:GLY:HA2	2.07	0.54
6:T:51:ASN:ND2	15:HA:5:MAN:O4	2.36	0.54
1:A:371:VAL:HG11	3:N:56:ALA:HB2	1.90	0.54
4:M:37:LYS:HD2	4:M:43:PRO:HG3	1.89	0.54
1:Q:371:VAL:HG11	3:V:56:ALA:HB2	1.90	0.53
8:X:21:ILE:O	8:X:72:THR:HA	2.08	0.53
2:B:546:SER:OG	2:I:588:ARG:NE	2.41	0.53
5:S:70:SER:HB3	5:S:81:ARG:HH21	1.73	0.53
2:B:536:THR:O	2:B:536:THR:OG1	2.26	0.53
6:F:27:SER:OG	6:F:68:GLY:N	2.42	0.53
6:T:83:GLU:HG2	6:T:105:ILE:HA	1.91	0.53
1:G:365:SER:OG	1:G:469:ARG:NH1	2.39	0.52
6:K:83:GLU:HG2	6:K:105:ILE:HA	1.92	0.52
1:G:195:ASN:ND2	1:G:433:CYS:SG	2.80	0.52
5:E:86:ASP:OD1	5:E:86:ASP:N	2.43	0.52
6:T:50:ASN:ND2	15:HA:5:MAN:O6	2.43	0.52
3:C:47:TRP:HZ3	1:G:458:GLY:HA3	1.74	0.52
5:J:31:ASP:OD1	5:J:31:ASP:N	2.38	0.52
5:E:43:LYS:NZ	5:E:44:GLN:O	2.41	0.52
1:Q:67:ASN:OD1	1:Q:67:ASN:N	2.43	0.52
6:T:27:SER:OG	6:T:68:GLY:N	2.42	0.52
7:W:32:LEU:HD23	7:W:94:SER:HB2	1.91	0.52
7:O:32:LEU:HD23	7:O:94:SER:HB2	1.91	0.51
2:B:610:TRP:HB3	1:G:34:LEU:HG	1.93	0.51
6:K:27:SER:OG	6:K:68:GLY:N	2.41	0.51
5:E:37:ILE:HD11	5:E:100(P):MET:HE3	1.93	0.51
6:F:83:GLU:HG2	6:F:105:ILE:HA	1.92	0.51
7:H:32:LEU:HD23	7:H:94:SER:HB2	1.91	0.51
5:J:52:HIS:ND1	5:J:53:ASP:O	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:350:ARG:NH2	1:G:396:ILE:O	2.40	0.51
5:S:52:HIS:ND1	5:S:53:ASP:O	2.44	0.51
1:A:31:ALA:N	1:A:33:ASN:OD1	2.44	0.51
5:S:43:LYS:NZ	5:S:44:GLN:O	2.41	0.51
1:Q:195:ASN:ND2	1:Q:433:CYS:SG	2.80	0.51
1:Q:31:ALA:N	1:Q:33:ASN:OD1	2.44	0.50
4:M:3:VAL:O	4:M:26:SER:OG	2.29	0.50
4:U:3:VAL:O	4:U:26:SER:OG	2.29	0.50
5:J:37:ILE:HD11	5:J:100(P):MET:HE3	1.93	0.50
8:L:86:TYR:O	8:L:101:GLY:HA2	2.12	0.50
1:G:67:ASN:OD1	1:G:67:ASN:N	2.43	0.50
1:Q:174:SER:OG	1:Q:175:LEU:N	2.45	0.50
2:B:617:ARG:NH2	2:B:634:GLU:OE1	2.45	0.50
1:G:174:SER:OG	1:G:175:LEU:N	2.45	0.50
8:X:86:TYR:O	8:X:101:GLY:HA2	2.12	0.50
6:F:93:SER:O	6:F:93:SER:OG	2.30	0.50
1:G:31:ALA:N	1:G:33:ASN:OD1	2.44	0.50
1:A:174:SER:OG	1:A:175:LEU:N	2.45	0.49
2:I:617:ARG:NH2	2:I:634:GLU:OE1	2.45	0.49
6:K:51:ASN:ND2	15:I:5:MAN:O4	2.42	0.49
5:S:37:ILE:HD11	5:S:100(P):MET:HE3	1.93	0.49
4:D:3:VAL:O	4:D:26:SER:OG	2.29	0.49
5:J:43:LYS:NZ	5:J:44:GLN:O	2.41	0.49
3:C:37:VAL:HG23	3:C:47:TRP:HA	1.94	0.49
8:P:86:TYR:O	8:P:101:GLY:HA2	2.12	0.49
1:Q:448:ASN:OD1	1:Q:448:ASN:N	2.44	0.49
2:R:617:ARG:NH2	2:R:634:GLU:OE1	2.45	0.49
5:S:17:THR:HG22	5:S:82(A):THR:HG23	1.94	0.49
5:E:17:THR:HG22	5:E:82(A):THR:HG23	1.94	0.49
1:G:292:VAL:HG23	1:G:449:ILE:HB	1.95	0.49
5:J:38:ARG:HB2	5:J:48:ILE:HD11	1.95	0.49
1:Q:458:GLY:HA3	3:V:47:TRP:HZ3	1.76	0.49
1:A:67:ASN:N	1:A:67:ASN:OD1	2.43	0.49
7:H:47:TRP:CD1	8:L:96:LEU:HD12	2.48	0.49
5:S:38:ARG:HB2	5:S:48:ILE:HD11	1.95	0.49
1:A:292:VAL:HG23	1:A:449:ILE:HB	1.94	0.49
6:F:38:GLN:NE2	6:F:39:ARG:O	2.39	0.49
1:G:61:TYR:OH	1:G:74:CYS:O	2.31	0.49
5:J:17:THR:HG22	5:J:82(A):THR:HG23	1.94	0.49
1:A:36:VAL:HG22	2:I:610:TRP:HE3	1.77	0.48
5:E:38:ARG:HB2	5:E:48:ILE:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:V:37:VAL:HG23	3:V:47:TRP:HA	1.94	0.48
1:A:448:ASN:N	1:A:448:ASN:OD1	2.44	0.48
5:E:52:HIS:ND1	5:E:53:ASP:O	2.44	0.48
1:G:448:ASN:OD1	1:G:448:ASN:N	2.44	0.48
5:S:86:ASP:OD1	5:S:86:ASP:N	2.43	0.48
1:G:318:TYR:OH	3:N:76(B):ASP:OD2	2.25	0.48
1:G:362:ALA:HB3	1:G:469:ARG:HE	1.79	0.48
1:Q:362:ALA:HB3	1:Q:469:ARG:HE	1.79	0.48
1:A:78:ASP:OD1	1:A:78:ASP:N	2.41	0.48
6:K:38:GLN:NE2	6:K:39:ARG:O	2.39	0.48
5:E:56:ASP:OD2	10:q:3:BMA:O2	2.30	0.48
1:G:457:ASP:OD2	1:G:467:THR:OG1	2.29	0.48
7:H:45:LEU:HB2	8:L:98:PHE:CG	2.49	0.48
1:Q:350:ARG:NH2	1:Q:396:ILE:O	2.40	0.48
5:S:56:ASP:OD2	10:6:3:BMA:O2	2.31	0.48
6:T:38:GLN:NE2	6:T:42:GLN:O	2.44	0.48
8:L:29:LEU:HD11	8:L:90:GLN:HB2	1.96	0.48
5:S:94:THR:OG1	5:S:100(Q):ASP:OD1	2.32	0.48
8:P:29:LEU:HD11	8:P:90:GLN:HB2	1.96	0.48
1:Q:61:TYR:OH	1:Q:74:CYS:O	2.31	0.48
1:Q:457:ASP:OD2	1:Q:467:THR:OG1	2.30	0.48
8:X:68:ASP:OD1	8:X:68:ASP:N	2.42	0.48
1:A:61:TYR:OH	1:A:74:CYS:O	2.31	0.48
1:Q:292:VAL:HG23	1:Q:449:ILE:HB	1.95	0.48
2:R:632:ASP:O	2:R:636:SER:HB3	2.14	0.48
5:J:61:PRO:HA	5:J:64:LYS:HB2	1.96	0.47
11:z:3:BMA:H4	11:z:5:MAN:H3	1.95	0.47
3:N:37:VAL:HG23	3:N:47:TRP:HA	1.94	0.47
1:A:138:ILE:HD13	9:Z:1:NAG:HN2	1.80	0.47
5:E:94:THR:OG1	5:E:100(Q):ASP:OD1	2.32	0.47
2:I:515:ILE:HG13	7:O:95:ARG:HB3	1.96	0.47
5:E:100(P):MET:N	6:F:36:TYR:OH	2.42	0.47
5:J:56:ASP:OD2	10:a:3:BMA:O2	2.31	0.47
1:Q:138:ILE:HD13	9:5:1:NAG:HN2	1.80	0.47
1:G:54:CYS:HB3	1:G:74:CYS:HB3	1.57	0.47
6:T:38:GLN:NE2	6:T:39:ARG:O	2.39	0.47
5:E:61:PRO:HA	5:E:64:LYS:HB2	1.96	0.47
1:G:78:ASP:OD1	1:G:78:ASP:N	2.41	0.47
6:K:38:GLN:NE2	6:K:42:GLN:O	2.44	0.47
7:O:69:ILE:HA	7:O:79:LEU:O	2.15	0.47
11:j:3:BMA:H4	11:j:5:MAN:H3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:632:ASP:O	2:B:636:SER:HB3	2.15	0.47
1:G:248:THR:OG1	1:G:249:HIS:N	2.48	0.47
5:J:94:THR:OG1	5:J:100(Q):ASP:OD1	2.32	0.47
1:Q:248:THR:OG1	1:Q:249:HIS:N	2.48	0.47
8:X:35:TRP:HB2	8:X:48:ILE:HB	1.97	0.47
11:FA:3:BMA:H4	11:FA:5:MAN:H3	1.95	0.47
1:G:138:ILE:HD13	9:p:1:NAG:HN2	1.80	0.47
8:P:68:ASP:OD1	8:P:68:ASP:N	2.42	0.47
1:A:362:ALA:HB3	1:A:469:ARG:HE	1.79	0.47
1:A:457:ASP:OD2	1:A:467:THR:OG1	2.29	0.47
1:Q:460:SER:O	1:Q:463:SER:OG	2.33	0.47
1:A:460:SER:O	1:A:463:SER:OG	2.33	0.46
2:B:610:TRP:HE3	1:G:36:VAL:HG22	1.80	0.46
7:W:69:ILE:HA	7:W:79:LEU:O	2.15	0.46
5:S:61:PRO:HA	5:S:64:LYS:HB2	1.96	0.46
1:G:203:GLN:HE22	1:G:318:TYR:H	1.62	0.46
8:P:35:TRP:HB2	8:P:48:ILE:HB	1.97	0.46
1:Q:36:VAL:HG22	2:R:610:TRP:HE3	1.79	0.46
4:D:16:GLY:N	4:D:77:LEU:O	2.48	0.46
2:I:546:SER:OG	2:R:588:ARG:NE	2.47	0.46
4:U:45:LEU:HD22	4:U:54:ALA:HB2	1.98	0.46
7:W:3:GLN:HB2	7:W:25:SER:HB3	1.98	0.46
7:H:69:ILE:HA	7:H:79:LEU:O	2.15	0.46
8:L:35:TRP:HB2	8:L:48:ILE:HB	1.97	0.46
8:X:29:LEU:HD11	8:X:90:GLN:HB2	1.95	0.46
6:F:15:PRO:HA	6:F:78:VAL:HG13	1.98	0.46
2:I:632:ASP:O	2:I:636:SER:HB3	2.14	0.46
1:Q:203:GLN:HE22	1:Q:318:TYR:H	1.62	0.46
1:A:165:LEU:HD11	1:Q:192:ARG:HD3	1.97	0.46
1:A:350:ARG:NH2	1:A:396:ILE:O	2.40	0.46
6:K:27:SER:HA	6:K:31:ARG:HH21	1.81	0.46
6:K:34:ILE:HG13	6:K:89:HIS:HB3	1.98	0.46
6:T:15:PRO:HA	6:T:78:VAL:HG13	1.98	0.46
6:T:27:SER:HA	6:T:31:ARG:HH21	1.81	0.46
1:Q:285:LEU:HD23	1:Q:481:SER:HB2	1.98	0.46
1:A:203:GLN:HE22	1:A:318:TYR:H	1.62	0.46
4:D:45:LEU:HD22	4:D:54:ALA:HB2	1.97	0.46
7:H:3:GLN:HB2	7:H:25:SER:HB3	1.98	0.46
5:J:100(Q):ASP:OD1	5:J:100(Q):ASP:N	2.40	0.46
6:K:93:SER:O	6:K:93:SER:OG	2.30	0.46
4:M:45:LEU:HD22	4:M:54:ALA:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:O:3:GLN:HB2	7:O:25:SER:HB3	1.98	0.46
6:F:38:GLN:NE2	6:F:42:GLN:O	2.44	0.45
1:G:285:LEU:HD23	1:G:481:SER:HB2	1.98	0.45
6:T:34:ILE:HG13	6:T:89:HIS:HB3	1.98	0.45
4:U:38:ARG:NH1	4:U:80:GLU:O	2.50	0.45
1:G:201:CYS:HB2	1:G:433:CYS:HB2	1.70	0.45
1:A:248:THR:OG1	1:A:249:HIS:N	2.48	0.45
13:9:3:BMA:H62	13:9:4:MAN:H2	1.74	0.45
4:D:38:ARG:NH1	4:D:80:GLU:O	2.50	0.45
5:E:100(Q):ASP:OD1	5:E:100(Q):ASP:N	2.40	0.45
6:F:27:SER:HA	6:F:31:ARG:HH21	1.81	0.45
1:G:121:LYS:HE2	1:G:121:LYS:HB2	1.76	0.45
7:H:39:GLN:HB2	7:H:45:LEU:HD23	1.99	0.45
3:N:47:TRP:HZ2	3:N:50:TRP:HD1	1.65	0.45
1:A:285:LEU:HD23	1:A:481:SER:HB2	1.98	0.45
1:G:460:SER:O	1:G:463:SER:OG	2.33	0.45
1:G:504:ARG:HE	2:I:661:LEU:HD11	1.82	0.45
7:O:39:GLN:HB2	7:O:45:LEU:HD23	1.99	0.45
6:T:93:SER:O	6:T:93:SER:OG	2.30	0.45
6:K:15:PRO:HA	6:K:78:VAL:HG13	1.98	0.45
3:C:47:TRP:HZ2	3:C:50:TRP:HD1	1.65	0.44
5:J:23:ASN:HA	5:J:77:LEU:HG	1.99	0.44
4:M:38:ARG:NH1	4:M:80:GLU:O	2.50	0.44
1:Q:360:ARG:O	1:Q:467:THR:HA	2.18	0.44
3:V:47:TRP:HZ2	3:V:50:TRP:HD1	1.65	0.44
1:A:360:ARG:O	1:A:467:THR:HA	2.17	0.44
4:D:49:ASP:OD2	4:D:52:ARG:NH1	2.51	0.44
1:G:360:ARG:O	1:G:467:THR:HA	2.18	0.44
5:E:23:ASN:HA	5:E:77:LEU:HG	1.99	0.44
6:F:50:ASN:ND2	15:1:5:MAN:O6	2.49	0.44
5:S:23:ASN:HA	5:S:77:LEU:HG	1.99	0.44
3:V:94:ARG:NH1	3:V:95:ARG:O	2.51	0.44
15:1:5:MAN:O4	15:1:5:MAN:O6	2.34	0.44
7:W:35:SER:HB2	7:W:93:THR:HG22	1.99	0.44
7:W:39:GLN:HB2	7:W:45:LEU:HD23	1.98	0.44
3:C:94:ARG:NH1	3:C:95:ARG:O	2.50	0.44
6:F:34:ILE:HG13	6:F:89:HIS:HB3	1.98	0.44
7:H:35:SER:HB2	7:H:93:THR:HG22	1.99	0.44
4:M:49:ASP:OD2	4:M:52:ARG:NH1	2.51	0.44
3:N:94:ARG:NH1	3:N:95:ARG:O	2.50	0.44
1:Q:132:THR:OG1	1:Q:133:ASN:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:ARG:HD3	1:G:165:LEU:HD11	1.99	0.44
4:D:33:THR:OG1	4:D:88:GLN:O	2.33	0.44
1:A:54:CYS:HB3	1:A:74:CYS:HB3	1.57	0.43
2:I:536:THR:O	2:I:536:THR:OG1	2.26	0.43
6:K:52:ASN:ND2	6:K:64:GLY:O	2.51	0.43
3:C:13:THR:O	3:C:16:SER:OG	2.36	0.43
1:G:132:THR:OG1	1:G:133:ASN:N	2.51	0.43
7:W:29:PRO:HB2	7:W:100:ARG:NH2	2.33	0.43
1:A:121:LYS:HE2	1:A:121:LYS:HB2	1.76	0.43
2:B:606:THR:OG1	1:G:36:VAL:O	2.34	0.43
1:G:279:ASN:OD1	1:G:280:ASN:N	2.52	0.43
4:M:16:GLY:N	4:M:77:LEU:O	2.48	0.43
7:O:29:PRO:HB2	7:O:100:ARG:NH2	2.33	0.43
7:O:44:GLY:HA2	8:P:87:TYR:CE1	2.53	0.43
6:T:52:ASN:ND2	6:T:64:GLY:O	2.52	0.43
1:A:132:THR:OG1	1:A:133:ASN:N	2.51	0.43
2:B:588:ARG:NE	2:R:546:SER:OG	2.50	0.43
1:G:241:SER:O	1:G:241:SER:OG	2.35	0.43
3:N:13:THR:O	3:N:16:SER:OG	2.36	0.43
4:U:49:ASP:OD2	4:U:52:ARG:NH1	2.51	0.43
1:A:279:ASN:OD1	1:A:280:ASN:N	2.52	0.43
8:X:4:MET:HE1	8:X:90:GLN:HB3	2.01	0.43
4:D:27:GLN:HE22	1:G:278:THR:HG21	1.83	0.43
1:G:116:LEU:HD21	1:G:434:MET:HE2	2.00	0.43
6:T:46:LEU:HD21	6:T:49:TYR:HB3	2.01	0.43
4:D:45:LEU:HD21	4:D:48:TYR:HB3	2.01	0.43
7:O:35:SER:HB2	7:O:93:THR:HG22	1.99	0.43
1:A:201:CYS:HB2	1:A:433:CYS:HB2	1.70	0.43
6:F:52:ASN:ND2	6:F:64:GLY:O	2.51	0.43
7:H:29:PRO:HB2	7:H:100:ARG:NH2	2.34	0.43
1:A:116:LEU:HD21	1:A:434:MET:HE2	2.00	0.42
1:Q:34:LEU:HG	2:R:610:TRP:HB3	2.01	0.42
1:Q:137:ASN:N	6:T:94:ARG:O	2.52	0.42
1:Q:116:LEU:HD21	1:Q:434:MET:HE2	2.00	0.42
13:d:3:BMA:H62	13:d:4:MAN:H2	1.74	0.42
7:O:97:LYS:HB3	7:O:100(D):TYR:HE1	1.84	0.42
1:Q:424:ILE:HD12	1:Q:426:MET:HB2	2.01	0.42
5:J:86:ASP:N	5:J:86:ASP:OD1	2.43	0.42
6:K:46:LEU:HD21	6:K:49:TYR:HB3	2.01	0.42
1:Q:279:ASN:OD1	1:Q:280:ASN:N	2.52	0.42
1:G:192:ARG:HD3	1:Q:165:LEU:HD11	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:H:97:LYS:HB3	7:H:100(D):TYR:HE1	1.84	0.42
8:L:68:ASP:OD1	8:L:68:ASP:N	2.42	0.42
2:R:571:TRP:O	2:R:574:LYS:HG3	2.20	0.42
1:A:46:LYS:HA	1:A:46:LYS:HD3	1.88	0.42
8:L:4:MET:HE1	8:L:90:GLN:HB3	2.01	0.42
4:M:32:MET:HE2	4:M:87:CYS:HB2	2.01	0.42
1:Q:241:SER:O	1:Q:241:SER:OG	2.35	0.42
5:S:100(Q):ASP:OD1	5:S:100(Q):ASP:N	2.40	0.42
4:U:16:GLY:N	4:U:77:LEU:O	2.48	0.42
1:A:34:LEU:HG	2:I:610:TRP:HB3	2.02	0.42
1:A:36:VAL:O	2:I:606:THR:OG1	2.37	0.42
5:E:49:GLY:HA2	6:F:96:TRP:HZ3	1.84	0.42
8:P:4:MET:HE1	8:P:90:GLN:HB3	2.00	0.42
4:M:45:LEU:HD21	4:M:48:TYR:HB3	2.01	0.42
8:X:8:PRO:HG2	8:X:11:LEU:HB2	2.02	0.42
4:D:32:MET:HE2	4:D:87:CYS:HB2	2.01	0.42
2:I:571:TRP:O	2:I:574:LYS:HG3	2.20	0.42
8:L:8:PRO:HG2	8:L:11:LEU:HB2	2.02	0.42
6:F:51:ASN:ND2	15:1:5:MAN:O4	2.45	0.41
5:J:33:TYR:HB2	5:J:95:THR:HG23	2.02	0.41
4:U:45:LEU:HD21	4:U:48:TYR:HB3	2.01	0.41
6:F:46:LEU:HD21	6:F:49:TYR:HB3	2.00	0.41
1:Q:121:LYS:HE2	1:Q:121:LYS:HB2	1.76	0.41
8:X:14:THR:HA	8:X:15:PRO:HD3	1.95	0.41
1:A:429:ARG:HE	1:A:429:ARG:HB2	1.75	0.41
5:S:33:TYR:HB2	5:S:95:THR:HG23	2.02	0.41
1:Q:54:CYS:HB3	1:Q:74:CYS:HB3	1.57	0.41
1:Q:258:GLN:H	1:Q:258:GLN:HG2	1.78	0.41
4:U:32:MET:HE2	4:U:87:CYS:HB2	2.01	0.41
1:A:424:ILE:HD12	1:A:426:MET:HB2	2.02	0.41
2:B:571:TRP:O	2:B:574:LYS:HG3	2.20	0.41
3:V:87:THR:HG23	3:V:110:VAL:HA	2.03	0.41
7:W:97:LYS:HB3	7:W:100(D):TYR:HE1	1.84	0.41
4:D:90:PHE:CE1	10:3:1:NAG:H82	2.56	0.41
3:V:13:THR:O	3:V:16:SER:OG	2.36	0.41
8:P:8:PRO:HG2	8:P:11:LEU:HB2	2.02	0.41
7:W:87:THR:HG23	7:W:110:THR:HA	2.03	0.41
3:N:87:THR:HG23	3:N:110:VAL:HA	2.03	0.41
3:C:66:ARG:NH2	3:C:86:ASP:OD2	2.54	0.41
3:C:87:THR:HG23	3:C:110:VAL:HA	2.03	0.41
7:H:66:ARG:HE	7:H:82:LEU:HD21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:L:106:ILE:HD13	8:L:106:ILE:HA	1.94	0.41
3:N:29:PHE:CD1	3:N:76(G):GLY:HA3	2.56	0.41
1:A:299:PRO:HG2	1:A:327:ARG:HB2	2.03	0.41
3:C:29:PHE:CD1	3:C:76(G):GLY:HA3	2.56	0.41
1:G:39:TYR:O	1:G:494:LEU:HA	2.21	0.41
1:G:424:ILE:HD12	1:G:426:MET:HB2	2.02	0.41
1:Q:201:CYS:HB2	1:Q:433:CYS:HB2	1.70	0.41
8:X:106:ILE:HD13	8:X:106:ILE:HA	1.94	0.41
1:A:499:THR:OG1	1:A:500:ARG:N	2.54	0.40
5:E:21:THR:HB	5:E:77:LEU:HD23	2.03	0.40
7:H:87:THR:HG23	7:H:110:THR:HA	2.02	0.40
7:O:28:THR:HA	7:O:29:PRO:HD3	1.80	0.40
1:Q:299:PRO:HG2	1:Q:327:ARG:HB2	2.03	0.40
3:V:6:GLN:HE22	3:V:91:PHE:HA	1.87	0.40
3:V:29:PHE:CD1	3:V:76(G):GLY:HA3	2.56	0.40
7:W:66:ARG:HE	7:W:82:LEU:HD21	1.86	0.40
5:E:33:TYR:HB2	5:E:95:THR:HG23	2.02	0.40
7:H:38:ARG:HD3	7:H:48:ILE:HD11	2.03	0.40
7:O:45:LEU:O	8:P:98:PHE:HB2	2.22	0.40
7:W:38:ARG:HD3	7:W:48:ILE:HD11	2.03	0.40
1:G:299:PRO:HG2	1:G:327:ARG:HB2	2.03	0.40
3:N:66:ARG:NH2	3:N:86:ASP:OD2	2.54	0.40
7:O:38:ARG:HD3	7:O:48:ILE:HD11	2.03	0.40
13:t:3:BMA:H62	13:t:4:MAN:H2	1.74	0.40
1:A:241:SER:O	1:A:241:SER:OG	2.35	0.40
2:B:578:ALA:O	2:B:581:LEU:HB3	2.22	0.40
7:O:64:LYS:HA	7:O:64:LYS:HD3	1.81	0.40
1:Q:342:LEU:HD23	1:Q:342:LEU:HA	1.89	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	447/480 (93%)	416 (93%)	31 (7%)	0	100	100
1	G	447/480 (93%)	416 (93%)	31 (7%)	0	100	100
1	Q	447/480 (93%)	416 (93%)	31 (7%)	0	100	100
2	B	128/153 (84%)	118 (92%)	10 (8%)	0	100	100
2	I	128/153 (84%)	118 (92%)	10 (8%)	0	100	100
2	R	128/153 (84%)	118 (92%)	10 (8%)	0	100	100
3	C	126/235 (54%)	122 (97%)	4 (3%)	0	100	100
3	N	126/235 (54%)	122 (97%)	4 (3%)	0	100	100
3	V	126/235 (54%)	122 (97%)	4 (3%)	0	100	100
4	D	100/209 (48%)	93 (93%)	7 (7%)	0	100	100
4	M	100/209 (48%)	93 (93%)	7 (7%)	0	100	100
4	U	100/209 (48%)	93 (93%)	7 (7%)	0	100	100
5	E	130/235 (55%)	121 (93%)	9 (7%)	0	100	100
5	J	130/235 (55%)	121 (93%)	9 (7%)	0	100	100
5	S	130/235 (55%)	121 (93%)	9 (7%)	0	100	100
6	F	103/213 (48%)	94 (91%)	9 (9%)	0	100	100
6	K	103/213 (48%)	94 (91%)	9 (9%)	0	100	100
6	T	103/213 (48%)	94 (91%)	9 (9%)	0	100	100
7	H	123/229 (54%)	118 (96%)	5 (4%)	0	100	100
7	O	123/229 (54%)	118 (96%)	5 (4%)	0	100	100
7	W	123/229 (54%)	118 (96%)	5 (4%)	0	100	100
8	L	110/219 (50%)	105 (96%)	3 (3%)	2 (2%)	6	34
8	P	110/219 (50%)	105 (96%)	3 (3%)	2 (2%)	6	34
8	X	110/219 (50%)	105 (96%)	3 (3%)	2 (2%)	6	34
All	All	3801/5919 (64%)	3561 (94%)	234 (6%)	6 (0%)	44	72

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
8	L	51	VAL
8	P	51	VAL
8	X	51	VAL
8	L	50	ASP
8	P	50	ASP
8	X	50	ASP

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	405/428 (95%)	402 (99%)	3 (1%)	76	77
1	G	405/428 (95%)	402 (99%)	3 (1%)	76	77
1	Q	405/428 (95%)	402 (99%)	3 (1%)	76	77
2	B	110/129 (85%)	110 (100%)	0	100	100
2	I	110/129 (85%)	110 (100%)	0	100	100
2	R	110/129 (85%)	110 (100%)	0	100	100
3	C	108/201 (54%)	108 (100%)	0	100	100
3	N	108/201 (54%)	108 (100%)	0	100	100
3	V	108/201 (54%)	108 (100%)	0	100	100
4	D	86/182 (47%)	85 (99%)	1 (1%)	63	72
4	M	86/182 (47%)	85 (99%)	1 (1%)	63	72
4	U	86/182 (47%)	85 (99%)	1 (1%)	63	72
5	E	116/205 (57%)	116 (100%)	0	100	100
5	J	116/205 (57%)	116 (100%)	0	100	100
5	S	116/205 (57%)	116 (100%)	0	100	100
6	F	88/181 (49%)	88 (100%)	0	100	100
6	K	88/181 (49%)	88 (100%)	0	100	100
6	T	88/181 (49%)	88 (100%)	0	100	100
7	H	105/195 (54%)	104 (99%)	1 (1%)	68	74
7	O	105/195 (54%)	104 (99%)	1 (1%)	68	74
7	W	105/195 (54%)	104 (99%)	1 (1%)	68	74
8	L	100/196 (51%)	100 (100%)	0	100	100
8	P	100/196 (51%)	100 (100%)	0	100	100
8	X	100/196 (51%)	100 (100%)	0	100	100
All	All	3354/5151 (65%)	3339 (100%)	15 (0%)	81	81

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

Mol	Chain	Res	Type
1	A	270	VAL
1	A	337	THR
1	A	416	LEU
4	D	50	THR
1	G	270	VAL
1	G	337	THR
1	G	416	LEU
7	H	52(A)	THR
4	M	50	THR
7	O	52(A)	THR
1	Q	270	VAL
1	Q	337	THR
1	Q	416	LEU
4	U	50	THR
7	W	52(A)	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (61) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	94	ASN
1	A	130	GLN
1	A	183	GLN
1	A	203	GLN
1	A	246	GLN
1	A	283	ASN
1	A	411	ASN
2	B	577	GLN
3	C	64	GLN
3	C	72	GLN
4	D	27	GLN
4	D	36	GLN
5	E	58	ASN
6	F	52	ASN
1	G	82	GLN
1	G	94	ASN
1	G	130	GLN
1	G	183	GLN
1	G	203	GLN
1	G	246	GLN
1	G	283	ASN
1	G	328	GLN
1	G	411	ASN

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Mol	Chain	Res	Type
7	H	39	GLN
2	I	577	GLN
2	I	618	ASN
5	J	1	GLN
5	J	3	HIS
5	J	5	GLN
5	J	58	ASN
6	K	52	ASN
8	L	53	ASN
4	M	27	GLN
4	M	36	GLN
3	N	64	GLN
3	N	72	GLN
7	O	39	GLN
8	P	53	ASN
1	Q	82	GLN
1	Q	94	ASN
1	Q	183	GLN
1	Q	203	GLN
1	Q	246	GLN
1	Q	283	ASN
1	Q	328	GLN
1	Q	411	ASN
2	R	577	GLN
2	R	607	ASN
2	R	618	ASN
5	S	1	GLN
5	S	3	HIS
5	S	5	GLN
5	S	58	ASN
6	T	52	ASN
4	U	27	GLN
4	U	36	GLN
3	V	64	GLN
3	V	72	GLN
7	W	39	GLN
8	X	53	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 ⓘ

177 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
9	NAG	0	1	9,1	14,14,15	0.64	1 (7%)	17,19,21	0.89	1 (5%)
9	NAG	0	2	9	14,14,15	0.38	0	17,19,21	0.41	0
15	NAG	1	1	1,15	14,14,15	0.44	0	17,19,21	0.53	0
15	NAG	1	2	15	14,14,15	0.33	0	17,19,21	1.27	3 (17%)
15	BMA	1	3	15	11,11,12	0.91	0	15,15,17	1.34	4 (26%)
15	MAN	1	4	15	11,11,12	0.56	0	15,15,17	1.36	1 (6%)
15	MAN	1	5	15	11,11,12	0.69	0	15,15,17	1.54	2 (13%)
15	MAN	1	6	15	11,11,12	0.95	1 (9%)	15,15,17	0.87	1 (6%)
15	MAN	1	7	15	11,11,12	0.62	0	15,15,17	1.13	2 (13%)
15	MAN	1	8	15	11,11,12	0.76	0	15,15,17	0.96	2 (13%)
15	MAN	1	9	15	11,11,12	0.71	0	15,15,17	1.13	2 (13%)
12	NAG	2	1	1,12	14,14,15	0.88	1 (7%)	17,19,21	1.57	3 (17%)
12	NAG	2	2	12	14,14,15	0.95	1 (7%)	17,19,21	1.25	1 (5%)
12	BMA	2	3	12	11,11,12	0.75	0	15,15,17	0.89	1 (6%)
10	NAG	3	1	10,1	14,14,15	0.19	0	17,19,21	0.69	1 (5%)
10	NAG	3	2	10	14,14,15	0.18	0	17,19,21	0.52	0
10	BMA	3	3	10	11,11,12	0.64	0	15,15,17	0.88	0
10	MAN	3	4	10	11,11,12	0.95	0	15,15,17	1.43	2 (13%)
9	NAG	4	1	9,1	14,14,15	0.42	0	17,19,21	0.48	0
9	NAG	4	2	9	14,14,15	0.84	1 (7%)	17,19,21	2.40	3 (17%)
9	NAG	5	1	9,1	14,14,15	1.00	1 (7%)	17,19,21	1.56	2 (11%)
9	NAG	5	2	9	14,14,15	0.41	0	17,19,21	0.47	0
10	NAG	6	1	10,1	14,14,15	0.32	0	17,19,21	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	NAG	6	2	10	14,14,15	0.19	0	17,19,21	0.97	2 (11%)
10	BMA	6	3	10	11,11,12	1.09	2 (18%)	15,15,17	1.36	1 (6%)
10	MAN	6	4	10	11,11,12	1.62	3 (27%)	15,15,17	2.29	2 (13%)
11	NAG	7	1	1,11	14,14,15	0.63	1 (7%)	17,19,21	0.59	0
11	NAG	7	2	11	14,14,15	0.22	0	17,19,21	0.47	0
11	BMA	7	3	11	11,11,12	0.55	0	15,15,17	1.02	1 (6%)
11	MAN	7	4	11	11,11,12	0.85	1 (9%)	15,15,17	1.13	2 (13%)
11	MAN	7	5	11	11,11,12	1.42	3 (27%)	15,15,17	1.86	3 (20%)
12	NAG	8	1	1,12	14,14,15	0.66	0	17,19,21	1.20	2 (11%)
12	NAG	8	2	12	14,14,15	0.22	0	17,19,21	0.71	1 (5%)
12	BMA	8	3	12	11,11,12	0.77	0	15,15,17	0.80	0
13	NAG	9	1	1,13	14,14,15	0.55	0	17,19,21	0.78	0
13	NAG	9	2	13	14,14,15	0.31	0	17,19,21	0.67	0
13	BMA	9	3	13	11,11,12	0.90	0	15,15,17	1.25	1 (6%)
13	MAN	9	4	13	11,11,12	0.96	1 (9%)	15,15,17	0.95	1 (6%)
12	NAG	AA	1	1,12	14,14,15	0.32	0	17,19,21	0.52	0
12	NAG	AA	2	12	14,14,15	0.22	0	17,19,21	0.54	0
12	BMA	AA	3	12	11,11,12	0.64	0	15,15,17	1.01	1 (6%)
14	NAG	BA	1	1,14	14,14,15	0.37	0	17,19,21	0.47	0
14	NAG	BA	2	14	14,14,15	0.20	0	17,19,21	0.46	0
14	BMA	BA	3	14	11,11,12	0.57	0	15,15,17	1.05	2 (13%)
14	MAN	BA	4	14	11,11,12	0.67	0	15,15,17	1.14	2 (13%)
14	MAN	BA	5	14	11,11,12	0.67	0	15,15,17	1.17	2 (13%)
14	MAN	BA	6	14	11,11,12	0.84	0	15,15,17	0.93	2 (13%)
9	NAG	CA	1	9,1	14,14,15	0.21	0	17,19,21	0.58	0
9	NAG	CA	2	9	14,14,15	0.26	0	17,19,21	0.48	0
9	NAG	DA	1	9,1	14,14,15	0.45	0	17,19,21	1.08	1 (5%)
9	NAG	DA	2	9	14,14,15	0.39	0	17,19,21	0.61	1 (5%)
12	NAG	EA	1	1,12	14,14,15	0.24	0	17,19,21	0.58	0
12	NAG	EA	2	12	14,14,15	0.27	0	17,19,21	0.61	0
12	BMA	EA	3	12	11,11,12	0.76	0	15,15,17	0.72	1 (6%)
11	NAG	FA	1	1,11	14,14,15	0.34	0	17,19,21	0.81	1 (5%)
11	NAG	FA	2	11	14,14,15	0.45	0	17,19,21	0.69	0
11	BMA	FA	3	11	11,11,12	0.93	0	15,15,17	0.91	1 (6%)
11	MAN	FA	4	11	11,11,12	0.88	0	15,15,17	1.21	2 (13%)
11	MAN	FA	5	11	11,11,12	0.83	0	15,15,17	0.97	2 (13%)
9	NAG	GA	1	9,1	14,14,15	0.65	1 (7%)	17,19,21	0.89	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	NAG	GA	2	9	14,14,15	0.38	0	17,19,21	0.41	0
15	NAG	HA	1	1,15	14,14,15	0.46	0	17,19,21	0.53	0
15	NAG	HA	2	15	14,14,15	0.34	0	17,19,21	1.28	3 (17%)
15	BMA	HA	3	15	11,11,12	0.91	0	15,15,17	1.34	4 (26%)
15	MAN	HA	4	15	11,11,12	0.54	0	15,15,17	1.37	1 (6%)
15	MAN	HA	5	15	11,11,12	0.70	0	15,15,17	1.55	2 (13%)
15	MAN	HA	6	15	11,11,12	0.96	1 (9%)	15,15,17	0.86	1 (6%)
15	MAN	HA	7	15	11,11,12	0.64	0	15,15,17	1.14	2 (13%)
15	MAN	HA	8	15	11,11,12	0.77	0	15,15,17	0.94	2 (13%)
15	MAN	HA	9	15	11,11,12	0.69	0	15,15,17	1.13	2 (13%)
12	NAG	IA	1	1,12	14,14,15	0.88	1 (7%)	17,19,21	1.58	3 (17%)
12	NAG	IA	2	12	14,14,15	0.95	1 (7%)	17,19,21	1.25	1 (5%)
12	BMA	IA	3	12	11,11,12	0.76	0	15,15,17	0.91	1 (6%)
10	NAG	JA	1	10,1	14,14,15	0.18	0	17,19,21	0.70	1 (5%)
10	NAG	JA	2	10	14,14,15	0.22	0	17,19,21	0.52	0
10	BMA	JA	3	10	11,11,12	0.63	0	15,15,17	0.87	0
10	MAN	JA	4	10	11,11,12	0.96	0	15,15,17	1.41	2 (13%)
9	NAG	Y	1	9,1	14,14,15	0.41	0	17,19,21	0.49	0
9	NAG	Y	2	9	14,14,15	0.82	1 (7%)	17,19,21	2.38	3 (17%)
9	NAG	Z	1	9,1	14,14,15	1.00	1 (7%)	17,19,21	1.56	2 (11%)
9	NAG	Z	2	9	14,14,15	0.43	0	17,19,21	0.47	0
10	NAG	a	1	10,1	14,14,15	0.32	0	17,19,21	0.54	0
10	NAG	a	2	10	14,14,15	0.17	0	17,19,21	0.98	2 (11%)
10	BMA	a	3	10	11,11,12	1.08	2 (18%)	15,15,17	1.35	1 (6%)
10	MAN	a	4	10	11,11,12	1.62	3 (27%)	15,15,17	2.29	3 (20%)
11	NAG	b	1	1,11	14,14,15	0.63	1 (7%)	17,19,21	0.58	0
11	NAG	b	2	11	14,14,15	0.23	0	17,19,21	0.47	0
11	BMA	b	3	11	11,11,12	0.54	0	15,15,17	1.01	1 (6%)
11	MAN	b	4	11	11,11,12	0.86	1 (9%)	15,15,17	1.13	2 (13%)
11	MAN	b	5	11	11,11,12	1.43	3 (27%)	15,15,17	1.85	3 (20%)
12	NAG	c	1	1,12	14,14,15	0.68	1 (7%)	17,19,21	1.20	1 (5%)
12	NAG	c	2	12	14,14,15	0.22	0	17,19,21	0.72	1 (5%)
12	BMA	c	3	12	11,11,12	0.77	0	15,15,17	0.80	0
13	NAG	d	1	1,13	14,14,15	0.54	0	17,19,21	0.78	0
13	NAG	d	2	13	14,14,15	0.31	0	17,19,21	0.67	0
13	BMA	d	3	13	11,11,12	0.92	0	15,15,17	1.25	1 (6%)
13	MAN	d	4	13	11,11,12	0.95	1 (9%)	15,15,17	0.95	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
12	NAG	e	1	1,12	14,14,15	0.33	0	17,19,21	0.51	0
12	NAG	e	2	12	14,14,15	0.22	0	17,19,21	0.56	0
12	BMA	e	3	12	11,11,12	0.64	0	15,15,17	1.03	1 (6%)
14	NAG	f	1	1,14	14,14,15	0.38	0	17,19,21	0.47	0
14	NAG	f	2	14	14,14,15	0.20	0	17,19,21	0.47	0
14	BMA	f	3	14	11,11,12	0.58	0	15,15,17	1.05	2 (13%)
14	MAN	f	4	14	11,11,12	0.67	0	15,15,17	1.13	2 (13%)
14	MAN	f	5	14	11,11,12	0.66	0	15,15,17	1.16	2 (13%)
14	MAN	f	6	14	11,11,12	0.85	0	15,15,17	0.92	2 (13%)
9	NAG	g	1	9,1	14,14,15	0.23	0	17,19,21	0.56	0
9	NAG	g	2	9	14,14,15	0.25	0	17,19,21	0.47	0
9	NAG	h	1	9,1	14,14,15	0.48	0	17,19,21	1.06	1 (5%)
9	NAG	h	2	9	14,14,15	0.40	0	17,19,21	0.61	1 (5%)
12	NAG	i	1	1,12	14,14,15	0.25	0	17,19,21	0.56	0
12	NAG	i	2	12	14,14,15	0.30	0	17,19,21	0.62	0
12	BMA	i	3	12	11,11,12	0.75	0	15,15,17	0.74	1 (6%)
11	NAG	j	1	1,11	14,14,15	0.33	0	17,19,21	0.81	1 (5%)
11	NAG	j	2	11	14,14,15	0.45	0	17,19,21	0.68	0
11	BMA	j	3	11	11,11,12	0.93	0	15,15,17	0.90	1 (6%)
11	MAN	j	4	11	11,11,12	0.87	0	15,15,17	1.21	2 (13%)
11	MAN	j	5	11	11,11,12	0.82	0	15,15,17	0.96	2 (13%)
9	NAG	k	1	9,1	14,14,15	0.66	1 (7%)	17,19,21	0.90	1 (5%)
9	NAG	k	2	9	14,14,15	0.40	0	17,19,21	0.41	0
15	NAG	l	1	1,15	14,14,15	0.44	0	17,19,21	0.52	0
15	NAG	l	2	15	14,14,15	0.33	0	17,19,21	1.27	3 (17%)
15	BMA	l	3	15	11,11,12	0.92	0	15,15,17	1.34	3 (20%)
15	MAN	l	4	15	11,11,12	0.56	0	15,15,17	1.37	1 (6%)
15	MAN	l	5	15	11,11,12	0.70	0	15,15,17	1.55	2 (13%)
15	MAN	l	6	15	11,11,12	0.95	1 (9%)	15,15,17	0.87	1 (6%)
15	MAN	l	7	15	11,11,12	0.64	0	15,15,17	1.12	2 (13%)
15	MAN	l	8	15	11,11,12	0.76	0	15,15,17	0.96	2 (13%)
15	MAN	l	9	15	11,11,12	0.71	0	15,15,17	1.12	2 (13%)
12	NAG	m	1	1,12	14,14,15	0.87	1 (7%)	17,19,21	1.57	3 (17%)
12	NAG	m	2	12	14,14,15	0.98	1 (7%)	17,19,21	1.24	1 (5%)
12	BMA	m	3	12	11,11,12	0.75	0	15,15,17	0.90	1 (6%)
10	NAG	n	1	10,1	14,14,15	0.19	0	17,19,21	0.69	1 (5%)
10	NAG	n	2	10	14,14,15	0.21	0	17,19,21	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	BMA	n	3	10	11,11,12	0.63	0	15,15,17	0.88	0
10	MAN	n	4	10	11,11,12	0.97	0	15,15,17	1.43	2 (13%)
9	NAG	o	1	9,1	14,14,15	0.42	0	17,19,21	0.47	0
9	NAG	o	2	9	14,14,15	0.82	1 (7%)	17,19,21	2.38	3 (17%)
9	NAG	p	1	9,1	14,14,15	1.01	1 (7%)	17,19,21	1.56	2 (11%)
9	NAG	p	2	9	14,14,15	0.41	0	17,19,21	0.47	0
10	NAG	q	1	10,1	14,14,15	0.33	0	17,19,21	0.54	0
10	NAG	q	2	10	14,14,15	0.19	0	17,19,21	0.97	2 (11%)
10	BMA	q	3	10	11,11,12	1.08	2 (18%)	15,15,17	1.36	1 (6%)
10	MAN	q	4	10	11,11,12	1.61	3 (27%)	15,15,17	2.28	3 (20%)
11	NAG	r	1	1,11	14,14,15	0.64	1 (7%)	17,19,21	0.59	0
11	NAG	r	2	11	14,14,15	0.23	0	17,19,21	0.47	0
11	BMA	r	3	11	11,11,12	0.52	0	15,15,17	1.01	1 (6%)
11	MAN	r	4	11	11,11,12	0.86	1 (9%)	15,15,17	1.13	2 (13%)
11	MAN	r	5	11	11,11,12	1.41	3 (27%)	15,15,17	1.85	3 (20%)
12	NAG	s	1	1,12	14,14,15	0.65	0	17,19,21	1.20	1 (5%)
12	NAG	s	2	12	14,14,15	0.21	0	17,19,21	0.72	1 (5%)
12	BMA	s	3	12	11,11,12	0.77	0	15,15,17	0.81	0
13	NAG	t	1	1,13	14,14,15	0.57	0	17,19,21	0.78	0
13	NAG	t	2	13	14,14,15	0.30	0	17,19,21	0.66	0
13	BMA	t	3	13	11,11,12	0.91	0	15,15,17	1.25	1 (6%)
13	MAN	t	4	13	11,11,12	0.95	1 (9%)	15,15,17	0.96	1 (6%)
12	NAG	u	1	1,12	14,14,15	0.32	0	17,19,21	0.51	0
12	NAG	u	2	12	14,14,15	0.22	0	17,19,21	0.56	0
12	BMA	u	3	12	11,11,12	0.64	0	15,15,17	1.02	1 (6%)
14	NAG	v	1	1,14	14,14,15	0.38	0	17,19,21	0.47	0
14	NAG	v	2	14	14,14,15	0.21	0	17,19,21	0.47	0
14	BMA	v	3	14	11,11,12	0.59	0	15,15,17	1.05	1 (6%)
14	MAN	v	4	14	11,11,12	0.68	0	15,15,17	1.13	2 (13%)
14	MAN	v	5	14	11,11,12	0.66	0	15,15,17	1.17	2 (13%)
14	MAN	v	6	14	11,11,12	0.85	0	15,15,17	0.92	2 (13%)
9	NAG	w	1	9,1	14,14,15	0.21	0	17,19,21	0.58	0
9	NAG	w	2	9	14,14,15	0.25	0	17,19,21	0.46	0
9	NAG	x	1	9,1	14,14,15	0.50	0	17,19,21	1.08	1 (5%)
9	NAG	x	2	9	14,14,15	0.37	0	17,19,21	0.60	1 (5%)
12	NAG	y	1	1,12	14,14,15	0.26	0	17,19,21	0.57	0
12	NAG	y	2	12	14,14,15	0.26	0	17,19,21	0.62	0
12	BMA	y	3	12	11,11,12	0.76	0	15,15,17	0.73	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	NAG	z	1	1,11	14,14,15	0.34	0	17,19,21	0.80	1 (5%)
11	NAG	z	2	11	14,14,15	0.43	0	17,19,21	0.69	0
11	BMA	z	3	11	11,11,12	0.92	0	15,15,17	0.90	1 (6%)
11	MAN	z	4	11	11,11,12	0.90	0	15,15,17	1.22	2 (13%)
11	MAN	z	5	11	11,11,12	0.83	0	15,15,17	0.97	2 (13%)

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
9	NAG	0	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	0	2	9	-	2/6/23/26	0/1/1/1
15	NAG	1	1	1,15	-	2/6/23/26	0/1/1/1
15	NAG	1	2	15	-	4/6/23/26	0/1/1/1
15	BMA	1	3	15	-	0/2/19/22	0/1/1/1
15	MAN	1	4	15	-	2/2/19/22	0/1/1/1
15	MAN	1	5	15	-	2/2/19/22	0/1/1/1
15	MAN	1	6	15	-	1/2/19/22	0/1/1/1
15	MAN	1	7	15	-	2/2/19/22	0/1/1/1
15	MAN	1	8	15	-	2/2/19/22	0/1/1/1
15	MAN	1	9	15	-	0/2/19/22	0/1/1/1
12	NAG	2	1	1,12	-	4/6/23/26	0/1/1/1
12	NAG	2	2	12	-	2/6/23/26	0/1/1/1
12	BMA	2	3	12	-	2/2/19/22	0/1/1/1
10	NAG	3	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	3	2	10	-	0/6/23/26	0/1/1/1
10	BMA	3	3	10	-	2/2/19/22	0/1/1/1
10	MAN	3	4	10	-	2/2/19/22	1/1/1/1
9	NAG	4	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	4	2	9	-	5/6/23/26	0/1/1/1
9	NAG	5	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	5	2	9	-	0/6/23/26	0/1/1/1
10	NAG	6	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	6	2	10	-	2/6/23/26	0/1/1/1
10	BMA	6	3	10	-	0/2/19/22	0/1/1/1
10	MAN	6	4	10	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	NAG	7	1	1,11	-	0/6/23/26	0/1/1/1
11	NAG	7	2	11	-	2/6/23/26	0/1/1/1
11	BMA	7	3	11	-	0/2/19/22	0/1/1/1
11	MAN	7	4	11	-	2/2/19/22	0/1/1/1
11	MAN	7	5	11	-	0/2/19/22	0/1/1/1
12	NAG	8	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	8	2	12	-	1/6/23/26	0/1/1/1
12	BMA	8	3	12	-	2/2/19/22	0/1/1/1
13	NAG	9	1	1,13	-	0/6/23/26	0/1/1/1
13	NAG	9	2	13	-	2/6/23/26	0/1/1/1
13	BMA	9	3	13	-	0/2/19/22	0/1/1/1
13	MAN	9	4	13	-	1/2/19/22	0/1/1/1
12	NAG	AA	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	AA	2	12	-	2/6/23/26	0/1/1/1
12	BMA	AA	3	12	-	2/2/19/22	0/1/1/1
14	NAG	BA	1	1,14	-	1/6/23/26	0/1/1/1
14	NAG	BA	2	14	-	2/6/23/26	0/1/1/1
14	BMA	BA	3	14	-	0/2/19/22	0/1/1/1
14	MAN	BA	4	14	-	2/2/19/22	0/1/1/1
14	MAN	BA	5	14	-	0/2/19/22	0/1/1/1
14	MAN	BA	6	14	-	2/2/19/22	0/1/1/1
9	NAG	CA	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	CA	2	9	-	0/6/23/26	0/1/1/1
9	NAG	DA	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	DA	2	9	-	0/6/23/26	0/1/1/1
12	NAG	EA	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	EA	2	12	-	0/6/23/26	0/1/1/1
12	BMA	EA	3	12	-	0/2/19/22	0/1/1/1
11	NAG	FA	1	1,11	-	1/6/23/26	0/1/1/1
11	NAG	FA	2	11	-	2/6/23/26	0/1/1/1
11	BMA	FA	3	11	-	2/2/19/22	0/1/1/1
11	MAN	FA	4	11	-	1/2/19/22	1/1/1/1
11	MAN	FA	5	11	-	0/2/19/22	0/1/1/1
9	NAG	GA	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	GA	2	9	-	2/6/23/26	0/1/1/1
15	NAG	HA	1	1,15	-	2/6/23/26	0/1/1/1
15	NAG	HA	2	15	-	4/6/23/26	0/1/1/1
15	BMA	HA	3	15	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
15	MAN	HA	4	15	-	2/2/19/22	0/1/1/1
15	MAN	HA	5	15	-	2/2/19/22	0/1/1/1
15	MAN	HA	6	15	-	1/2/19/22	0/1/1/1
15	MAN	HA	7	15	-	2/2/19/22	0/1/1/1
15	MAN	HA	8	15	-	2/2/19/22	0/1/1/1
15	MAN	HA	9	15	-	0/2/19/22	0/1/1/1
12	NAG	IA	1	1,12	-	4/6/23/26	0/1/1/1
12	NAG	IA	2	12	-	2/6/23/26	0/1/1/1
12	BMA	IA	3	12	-	2/2/19/22	0/1/1/1
10	NAG	JA	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	JA	2	10	-	0/6/23/26	0/1/1/1
10	BMA	JA	3	10	-	2/2/19/22	0/1/1/1
10	MAN	JA	4	10	-	2/2/19/22	1/1/1/1
9	NAG	Y	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	Y	2	9	-	5/6/23/26	0/1/1/1
9	NAG	Z	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	Z	2	9	-	0/6/23/26	0/1/1/1
10	NAG	a	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	a	2	10	-	2/6/23/26	0/1/1/1
10	BMA	a	3	10	-	0/2/19/22	0/1/1/1
10	MAN	a	4	10	-	2/2/19/22	0/1/1/1
11	NAG	b	1	1,11	-	0/6/23/26	0/1/1/1
11	NAG	b	2	11	-	2/6/23/26	0/1/1/1
11	BMA	b	3	11	-	0/2/19/22	0/1/1/1
11	MAN	b	4	11	-	2/2/19/22	0/1/1/1
11	MAN	b	5	11	-	0/2/19/22	0/1/1/1
12	NAG	c	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	c	2	12	-	1/6/23/26	0/1/1/1
12	BMA	c	3	12	-	2/2/19/22	0/1/1/1
13	NAG	d	1	1,13	-	0/6/23/26	0/1/1/1
13	NAG	d	2	13	-	2/6/23/26	0/1/1/1
13	BMA	d	3	13	-	0/2/19/22	0/1/1/1
13	MAN	d	4	13	-	1/2/19/22	0/1/1/1
12	NAG	e	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	e	2	12	-	2/6/23/26	0/1/1/1
12	BMA	e	3	12	-	2/2/19/22	0/1/1/1
14	NAG	f	1	1,14	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	NAG	f	2	14	-	2/6/23/26	0/1/1/1
14	BMA	f	3	14	-	0/2/19/22	0/1/1/1
14	MAN	f	4	14	-	2/2/19/22	0/1/1/1
14	MAN	f	5	14	-	0/2/19/22	0/1/1/1
14	MAN	f	6	14	-	2/2/19/22	0/1/1/1
9	NAG	g	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	g	2	9	-	0/6/23/26	0/1/1/1
9	NAG	h	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	h	2	9	-	0/6/23/26	0/1/1/1
12	NAG	i	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	i	2	12	-	0/6/23/26	0/1/1/1
12	BMA	i	3	12	-	0/2/19/22	0/1/1/1
11	NAG	j	1	1,11	-	1/6/23/26	0/1/1/1
11	NAG	j	2	11	-	2/6/23/26	0/1/1/1
11	BMA	j	3	11	-	2/2/19/22	0/1/1/1
11	MAN	j	4	11	-	1/2/19/22	1/1/1/1
11	MAN	j	5	11	-	0/2/19/22	0/1/1/1
9	NAG	k	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	k	2	9	-	2/6/23/26	0/1/1/1
15	NAG	l	1	1,15	-	2/6/23/26	0/1/1/1
15	NAG	l	2	15	-	4/6/23/26	0/1/1/1
15	BMA	l	3	15	-	0/2/19/22	0/1/1/1
15	MAN	l	4	15	-	2/2/19/22	0/1/1/1
15	MAN	l	5	15	-	2/2/19/22	0/1/1/1
15	MAN	l	6	15	-	1/2/19/22	0/1/1/1
15	MAN	l	7	15	-	2/2/19/22	0/1/1/1
15	MAN	l	8	15	-	2/2/19/22	0/1/1/1
15	MAN	l	9	15	-	0/2/19/22	0/1/1/1
12	NAG	m	1	1,12	-	4/6/23/26	0/1/1/1
12	NAG	m	2	12	-	2/6/23/26	0/1/1/1
12	BMA	m	3	12	-	2/2/19/22	0/1/1/1
10	NAG	n	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	n	2	10	-	0/6/23/26	0/1/1/1
10	BMA	n	3	10	-	2/2/19/22	0/1/1/1
10	MAN	n	4	10	-	2/2/19/22	1/1/1/1
9	NAG	o	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	o	2	9	-	5/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	p	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	p	2	9	-	0/6/23/26	0/1/1/1
10	NAG	q	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	q	2	10	-	2/6/23/26	0/1/1/1
10	BMA	q	3	10	-	0/2/19/22	0/1/1/1
10	MAN	q	4	10	-	2/2/19/22	0/1/1/1
11	NAG	r	1	1,11	-	0/6/23/26	0/1/1/1
11	NAG	r	2	11	-	2/6/23/26	0/1/1/1
11	BMA	r	3	11	-	0/2/19/22	0/1/1/1
11	MAN	r	4	11	-	2/2/19/22	0/1/1/1
11	MAN	r	5	11	-	0/2/19/22	0/1/1/1
12	NAG	s	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	s	2	12	-	1/6/23/26	0/1/1/1
12	BMA	s	3	12	-	2/2/19/22	0/1/1/1
13	NAG	t	1	1,13	-	0/6/23/26	0/1/1/1
13	NAG	t	2	13	-	2/6/23/26	0/1/1/1
13	BMA	t	3	13	-	0/2/19/22	0/1/1/1
13	MAN	t	4	13	-	1/2/19/22	0/1/1/1
12	NAG	u	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	u	2	12	-	2/6/23/26	0/1/1/1
12	BMA	u	3	12	-	2/2/19/22	0/1/1/1
14	NAG	v	1	1,14	-	1/6/23/26	0/1/1/1
14	NAG	v	2	14	-	2/6/23/26	0/1/1/1
14	BMA	v	3	14	-	0/2/19/22	0/1/1/1
14	MAN	v	4	14	-	2/2/19/22	0/1/1/1
14	MAN	v	5	14	-	0/2/19/22	0/1/1/1
14	MAN	v	6	14	-	2/2/19/22	0/1/1/1
9	NAG	w	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	w	2	9	-	0/6/23/26	0/1/1/1
9	NAG	x	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	x	2	9	-	0/6/23/26	0/1/1/1
12	NAG	y	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	y	2	12	-	0/6/23/26	0/1/1/1
12	BMA	y	3	12	-	0/2/19/22	0/1/1/1
11	NAG	z	1	1,11	-	1/6/23/26	0/1/1/1
11	NAG	z	2	11	-	2/6/23/26	0/1/1/1
11	BMA	z	3	11	-	2/2/19/22	0/1/1/1
11	MAN	z	4	11	-	1/2/19/22	1/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	MAN	z	5	11	-	0/2/19/22	0/1/1/1

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	5	1	NAG	O5-C1	3.11	1.48	1.43
9	p	1	NAG	O5-C1	3.11	1.48	1.43
9	Z	1	NAG	O5-C1	3.10	1.48	1.43
10	6	4	MAN	C1-C2	3.09	1.59	1.52
10	a	4	MAN	C1-C2	3.08	1.59	1.52
10	q	4	MAN	C1-C2	3.05	1.59	1.52
10	a	4	MAN	O5-C5	3.00	1.49	1.43
10	6	4	MAN	O5-C5	2.98	1.49	1.43
10	q	4	MAN	O5-C5	2.97	1.49	1.43
12	m	2	NAG	O5-C1	2.92	1.48	1.43
10	6	4	MAN	O5-C1	2.89	1.48	1.43
10	q	4	MAN	O5-C1	2.87	1.48	1.43
12	2	2	NAG	O5-C1	2.87	1.48	1.43
12	IA	2	NAG	O5-C1	2.86	1.48	1.43
10	a	4	MAN	O5-C1	2.84	1.48	1.43
11	b	5	MAN	C2-C3	2.65	1.56	1.52
11	r	5	MAN	C2-C3	2.63	1.56	1.52
11	7	5	MAN	C2-C3	2.61	1.56	1.52
12	2	1	NAG	O5-C1	2.59	1.48	1.43
12	IA	1	NAG	O5-C1	2.59	1.48	1.43
12	m	1	NAG	O5-C1	2.58	1.48	1.43
11	7	5	MAN	C1-C2	2.57	1.58	1.52
11	b	5	MAN	C1-C2	2.57	1.58	1.52
11	r	5	MAN	C1-C2	2.56	1.58	1.52
9	4	2	NAG	C1-C2	2.50	1.55	1.52
9	o	2	NAG	C1-C2	2.46	1.55	1.52
10	6	3	BMA	O5-C5	2.39	1.48	1.43
9	Y	2	NAG	C1-C2	2.39	1.55	1.52
10	q	3	BMA	O5-C5	2.36	1.48	1.43
10	a	3	BMA	O5-C5	2.33	1.48	1.43
11	b	5	MAN	O5-C5	2.28	1.47	1.43
11	r	1	NAG	O5-C1	-2.28	1.39	1.43
11	7	1	NAG	O5-C1	-2.27	1.39	1.43
11	r	5	MAN	O5-C5	2.26	1.47	1.43
11	b	1	NAG	O5-C1	-2.24	1.39	1.43
11	7	5	MAN	O5-C5	2.22	1.47	1.43
13	9	4	MAN	C1-C2	2.16	1.57	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	b	4	MAN	C1-C2	2.16	1.57	1.52
15	l	6	MAN	O5-C1	-2.15	1.40	1.43
10	a	3	BMA	C2-C3	-2.15	1.49	1.52
11	r	4	MAN	C1-C2	2.14	1.57	1.52
15	HA	6	MAN	O5-C1	-2.14	1.40	1.43
10	6	3	BMA	C2-C3	-2.13	1.49	1.52
15	1	6	MAN	O5-C1	-2.13	1.40	1.43
13	t	4	MAN	C1-C2	2.13	1.57	1.52
9	GA	1	NAG	O5-C1	-2.12	1.40	1.43
11	7	4	MAN	C1-C2	2.12	1.57	1.52
13	d	4	MAN	C1-C2	2.11	1.57	1.52
9	k	1	NAG	O5-C1	-2.11	1.40	1.43
10	q	3	BMA	C2-C3	-2.09	1.49	1.52
9	0	1	NAG	O5-C1	-2.07	1.40	1.43
12	c	1	NAG	C1-C2	2.03	1.55	1.52

All (192) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	4	2	NAG	C2-N2-C7	8.37	134.12	122.90
9	Y	2	NAG	C2-N2-C7	8.25	133.96	122.90
9	o	2	NAG	C2-N2-C7	8.24	133.94	122.90
10	a	4	MAN	C1-O5-C5	7.64	122.43	112.19
10	6	4	MAN	C1-O5-C5	7.64	122.43	112.19
10	q	4	MAN	C1-O5-C5	7.60	122.37	112.19
11	7	5	MAN	C1-O5-C5	5.56	119.63	112.19
11	b	5	MAN	C1-O5-C5	5.51	119.58	112.19
11	r	5	MAN	C1-O5-C5	5.50	119.56	112.19
9	p	1	NAG	C1-O5-C5	5.18	119.13	112.19
9	5	1	NAG	C1-O5-C5	5.18	119.12	112.19
9	Z	1	NAG	C1-O5-C5	5.17	119.11	112.19
12	2	2	NAG	C1-O5-C5	4.94	118.81	112.19
12	IA	2	NAG	C1-O5-C5	4.93	118.79	112.19
12	m	2	NAG	C1-O5-C5	4.90	118.76	112.19
10	n	4	MAN	C1-O5-C5	4.65	118.41	112.19
10	3	4	MAN	C1-O5-C5	4.65	118.41	112.19
10	6	3	BMA	C1-O5-C5	4.60	118.36	112.19
15	l	4	MAN	C1-O5-C5	4.59	118.34	112.19
15	HA	4	MAN	C1-O5-C5	4.58	118.33	112.19
10	q	3	BMA	C1-O5-C5	4.57	118.31	112.19
10	JA	4	MAN	C1-O5-C5	4.57	118.31	112.19
15	1	4	MAN	C1-O5-C5	4.57	118.31	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	a	3	BMA	C1-O5-C5	4.55	118.29	112.19
15	HA	5	MAN	C1-O5-C5	4.29	117.94	112.19
15	l	5	MAN	C1-O5-C5	4.28	117.92	112.19
15	1	5	MAN	C1-O5-C5	4.24	117.87	112.19
9	Y	2	NAG	C1-C2-N2	4.07	116.85	110.43
9	4	2	NAG	C1-C2-N2	4.07	116.85	110.43
12	m	1	NAG	C1-O5-C5	4.07	117.64	112.19
12	IA	1	NAG	C1-O5-C5	4.07	117.63	112.19
12	2	1	NAG	C1-O5-C5	4.06	117.63	112.19
9	o	2	NAG	C1-C2-N2	4.05	116.81	110.43
9	DA	1	NAG	C1-O5-C5	3.95	117.48	112.19
9	x	1	NAG	C1-O5-C5	3.93	117.45	112.19
9	h	1	NAG	C1-O5-C5	3.85	117.34	112.19
13	t	3	BMA	C1-O5-C5	3.78	117.26	112.19
13	9	3	BMA	C1-O5-C5	3.77	117.24	112.19
13	d	3	BMA	C1-O5-C5	3.76	117.23	112.19
11	z	4	MAN	C1-O5-C5	3.75	117.21	112.19
11	j	4	MAN	C1-O5-C5	3.69	117.13	112.19
11	FA	4	MAN	C1-O5-C5	3.69	117.12	112.19
14	v	5	MAN	C1-O5-C5	3.40	116.74	112.19
14	BA	5	MAN	C1-O5-C5	3.39	116.73	112.19
14	f	5	MAN	C1-O5-C5	3.37	116.70	112.19
15	l	5	MAN	O2-C2-C3	-3.34	103.22	110.15
15	HA	5	MAN	O2-C2-C3	-3.33	103.26	110.15
15	1	9	MAN	C1-O5-C5	3.31	116.62	112.19
15	1	5	MAN	O2-C2-C3	-3.31	103.30	110.15
15	HA	9	MAN	C1-O5-C5	3.31	116.62	112.19
15	l	9	MAN	C1-O5-C5	3.30	116.61	112.19
14	BA	4	MAN	C1-O5-C5	3.25	116.54	112.19
12	c	1	NAG	C2-N2-C7	3.25	127.25	122.90
15	l	2	NAG	C2-N2-C7	3.23	127.23	122.90
12	s	1	NAG	C2-N2-C7	3.22	127.22	122.90
15	HA	2	NAG	C2-N2-C7	3.22	127.21	122.90
12	8	1	NAG	C2-N2-C7	3.21	127.21	122.90
15	1	2	NAG	C2-N2-C7	3.19	127.17	122.90
14	v	4	MAN	C1-O5-C5	3.18	116.45	112.19
14	f	4	MAN	C1-O5-C5	3.17	116.43	112.19
9	p	1	NAG	C2-N2-C7	3.14	127.11	122.90
9	5	1	NAG	C2-N2-C7	3.12	127.08	122.90
9	Z	1	NAG	C2-N2-C7	3.12	127.08	122.90
12	IA	1	NAG	C2-N2-C7	3.04	126.97	122.90
12	2	1	NAG	C2-N2-C7	3.04	126.97	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	m	1	NAG	C2-N2-C7	2.98	126.90	122.90
9	0	1	NAG	C1-O5-C5	2.97	116.16	112.19
9	k	1	NAG	C1-O5-C5	2.96	116.16	112.19
9	GA	1	NAG	C1-O5-C5	2.96	116.15	112.19
12	e	3	BMA	C1-O5-C5	2.91	116.09	112.19
15	HA	7	MAN	C1-O5-C5	2.90	116.08	112.19
12	u	3	BMA	C1-O5-C5	2.87	116.03	112.19
15	1	7	MAN	C1-O5-C5	2.87	116.03	112.19
11	j	1	NAG	C1-O5-C5	2.86	116.02	112.19
15	l	7	MAN	C1-O5-C5	2.84	116.00	112.19
11	FA	1	NAG	C1-O5-C5	2.84	115.99	112.19
12	AA	3	BMA	C1-O5-C5	2.83	115.98	112.19
11	z	1	NAG	C1-O5-C5	2.82	115.97	112.19
11	7	3	BMA	C1-O5-C5	2.77	115.90	112.19
15	l	3	BMA	C1-O5-C5	2.75	115.88	112.19
14	v	3	BMA	C1-O5-C5	2.74	115.86	112.19
14	f	3	BMA	C1-O5-C5	2.73	115.84	112.19
11	b	3	BMA	C1-O5-C5	2.72	115.83	112.19
15	1	3	BMA	C1-O5-C5	2.71	115.82	112.19
14	BA	3	BMA	C1-O5-C5	2.71	115.81	112.19
11	r	3	BMA	C1-O5-C5	2.70	115.81	112.19
15	HA	3	BMA	C1-O5-C5	2.69	115.79	112.19
11	r	4	MAN	C1-O5-C5	2.68	115.78	112.19
11	b	4	MAN	C1-O5-C5	2.67	115.76	112.19
11	7	4	MAN	C1-O5-C5	2.65	115.74	112.19
11	7	5	MAN	C1-C2-C3	2.64	113.48	109.64
11	b	5	MAN	C1-C2-C3	2.62	113.46	109.64
11	r	5	MAN	C1-C2-C3	2.60	113.44	109.64
10	a	2	NAG	C1-O5-C5	2.59	115.65	112.19
12	IA	1	NAG	O4-C4-C5	2.58	115.67	109.32
12	m	1	NAG	O4-C4-C5	2.57	115.65	109.32
10	6	2	NAG	C1-O5-C5	2.56	115.62	112.19
10	q	2	NAG	C1-O5-C5	2.56	115.61	112.19
12	2	1	NAG	O4-C4-C5	2.55	115.61	109.32
15	1	7	MAN	O2-C2-C3	-2.49	105.00	110.15
15	l	7	MAN	O2-C2-C3	-2.47	105.03	110.15
15	HA	7	MAN	O2-C2-C3	-2.47	105.03	110.15
15	1	8	MAN	O2-C2-C3	-2.42	105.13	110.15
12	s	2	NAG	C1-O5-C5	2.42	115.43	112.19
15	l	8	MAN	O2-C2-C3	-2.42	105.14	110.15
12	c	2	NAG	C1-O5-C5	2.39	115.39	112.19
15	HA	8	MAN	O2-C2-C3	-2.39	105.19	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	a	4	MAN	O2-C2-C3	-2.39	105.20	110.15
10	6	4	MAN	O2-C2-C3	-2.39	105.20	110.15
11	r	4	MAN	O2-C2-C3	-2.37	105.23	110.15
11	b	4	MAN	O2-C2-C3	-2.37	105.24	110.15
12	8	2	NAG	C1-O5-C5	2.35	115.33	112.19
10	q	4	MAN	O2-C2-C3	-2.35	105.29	110.15
15	HA	9	MAN	O2-C2-C3	-2.34	105.31	110.15
11	7	4	MAN	O2-C2-C3	-2.33	105.33	110.15
15	1	9	MAN	O2-C2-C3	-2.33	105.33	110.15
15	1	6	MAN	O2-C2-C3	-2.32	105.35	110.15
15	HA	6	MAN	O2-C2-C3	-2.32	105.35	110.15
11	FA	5	MAN	C1-O5-C5	2.31	115.28	112.19
10	JA	1	NAG	C1-O5-C5	2.31	115.28	112.19
11	b	5	MAN	O2-C2-C3	-2.31	105.37	110.15
11	z	5	MAN	C1-O5-C5	2.31	115.28	112.19
15	l	9	MAN	O2-C2-C3	-2.30	105.39	110.15
15	l	6	MAN	O2-C2-C3	-2.30	105.39	110.15
11	r	5	MAN	O2-C2-C3	-2.29	105.41	110.15
10	n	1	NAG	C1-O5-C5	2.28	115.24	112.19
11	7	5	MAN	O2-C2-C3	-2.28	105.43	110.15
11	j	4	MAN	O2-C2-C3	-2.28	105.44	110.15
12	IA	3	BMA	C1-O5-C5	2.27	115.23	112.19
10	3	1	NAG	C1-O5-C5	2.27	115.23	112.19
15	HA	2	NAG	C1-O5-C5	2.27	115.22	112.19
11	FA	4	MAN	O2-C2-C3	-2.26	105.47	110.15
11	z	4	MAN	O2-C2-C3	-2.26	105.47	110.15
11	j	5	MAN	O2-C2-C3	-2.25	105.48	110.15
11	j	5	MAN	C1-O5-C5	2.25	115.20	112.19
11	FA	5	MAN	O2-C2-C3	-2.24	105.50	110.15
11	z	5	MAN	O2-C2-C3	-2.24	105.50	110.15
12	m	3	BMA	C1-O5-C5	2.23	115.18	112.19
13	9	4	MAN	O2-C2-C3	-2.23	105.53	110.15
13	t	4	MAN	O2-C2-C3	-2.22	105.55	110.15
14	BA	6	MAN	O2-C2-C3	-2.22	105.55	110.15
15	l	3	BMA	O2-C2-C3	-2.22	105.55	110.15
14	f	6	MAN	O2-C2-C3	-2.22	105.56	110.15
13	d	4	MAN	O2-C2-C3	-2.22	105.56	110.15
14	v	5	MAN	O2-C2-C3	-2.21	105.57	110.15
15	HA	3	BMA	O2-C2-C3	-2.21	105.57	110.15
11	z	3	BMA	C1-O5-C5	2.21	115.15	112.19
15	l	2	NAG	C1-O5-C5	2.21	115.15	112.19
14	v	6	MAN	O2-C2-C3	-2.21	105.58	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	1	3	BMA	O2-C2-C3	-2.21	105.58	110.15
14	BA	5	MAN	O2-C2-C3	-2.21	105.58	110.15
14	f	5	MAN	O2-C2-C3	-2.21	105.58	110.15
11	FA	3	BMA	C1-O5-C5	2.20	115.14	112.19
10	n	4	MAN	O2-C2-C3	-2.19	105.61	110.15
10	3	4	MAN	O2-C2-C3	-2.19	105.61	110.15
10	JA	4	MAN	O2-C2-C3	-2.19	105.61	110.15
15	1	2	NAG	C1-C2-N2	2.18	113.87	110.43
15	l	2	NAG	C1-C2-N2	2.18	113.87	110.43
15	HA	2	NAG	C1-C2-N2	2.17	113.86	110.43
9	Y	2	NAG	C8-C7-N2	2.17	119.72	116.12
15	1	2	NAG	C1-O5-C5	2.16	115.08	112.19
9	o	2	NAG	C8-C7-N2	2.16	119.70	116.12
12	2	3	BMA	C1-O5-C5	2.16	115.08	112.19
11	j	3	BMA	C1-O5-C5	2.15	115.06	112.19
9	4	2	NAG	C8-C7-N2	2.11	119.62	116.12
14	f	6	MAN	C1-O5-C5	2.08	114.98	112.19
9	h	2	NAG	C1-O5-C5	2.08	114.97	112.19
14	BA	6	MAN	C1-O5-C5	2.08	114.97	112.19
15	l	3	BMA	O5-C5-C4	-2.08	105.78	110.83
9	x	2	NAG	C1-O5-C5	2.08	114.97	112.19
9	DA	2	NAG	C1-O5-C5	2.07	114.97	112.19
15	l	8	MAN	C1-O5-C5	2.07	114.97	112.19
14	v	6	MAN	C1-O5-C5	2.07	114.96	112.19
15	1	3	BMA	O5-C5-C4	-2.07	105.80	110.83
15	HA	3	BMA	O5-C5-C4	-2.06	105.82	110.83
15	HA	3	BMA	C3-C4-C5	-2.06	106.50	110.23
15	1	3	BMA	C3-C4-C5	-2.05	106.51	110.23
15	HA	8	MAN	C1-O5-C5	2.05	114.94	112.19
15	1	8	MAN	C1-O5-C5	2.05	114.93	112.19
10	a	4	MAN	C1-C2-C3	2.03	112.60	109.64
14	f	3	BMA	O2-C2-C3	-2.03	105.95	110.15
10	a	2	NAG	O4-C4-C3	2.03	115.15	110.38
10	q	4	MAN	C1-C2-C3	2.02	112.59	109.64
10	6	2	NAG	O4-C4-C3	2.02	115.14	110.38
14	f	4	MAN	O2-C2-C3	-2.02	105.97	110.15
10	q	2	NAG	O4-C4-C3	2.01	115.12	110.38
12	EA	3	BMA	O2-C2-C3	-2.01	105.99	110.15
14	BA	3	BMA	O2-C2-C3	-2.01	105.99	110.15
14	v	4	MAN	O2-C2-C3	-2.01	105.99	110.15
12	8	1	NAG	C1-C2-N2	2.00	113.59	110.43
12	i	3	BMA	O2-C2-C3	-2.00	106.00	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	BA	4	MAN	O2-C2-C3	-2.00	106.00	110.15

There are no chirality outliers.

All (237) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
15	l	2	NAG	C4-C5-C6-O6
15	1	2	NAG	C4-C5-C6-O6
15	HA	2	NAG	C4-C5-C6-O6
12	e	2	NAG	O5-C5-C6-O6
12	u	2	NAG	O5-C5-C6-O6
12	AA	2	NAG	O5-C5-C6-O6
14	f	2	NAG	O5-C5-C6-O6
14	v	2	NAG	O5-C5-C6-O6
14	BA	2	NAG	O5-C5-C6-O6
15	l	2	NAG	O5-C5-C6-O6
15	1	2	NAG	O5-C5-C6-O6
15	HA	2	NAG	O5-C5-C6-O6
10	n	4	MAN	O5-C5-C6-O6
10	3	4	MAN	O5-C5-C6-O6
10	JA	4	MAN	O5-C5-C6-O6
11	j	3	BMA	C4-C5-C6-O6
11	z	3	BMA	C4-C5-C6-O6
11	FA	3	BMA	C4-C5-C6-O6
11	b	2	NAG	O5-C5-C6-O6
11	r	2	NAG	O5-C5-C6-O6
11	7	2	NAG	O5-C5-C6-O6
12	e	1	NAG	O5-C5-C6-O6
12	u	1	NAG	O5-C5-C6-O6
12	AA	1	NAG	O5-C5-C6-O6
10	n	3	BMA	C4-C5-C6-O6
10	3	3	BMA	C4-C5-C6-O6
10	JA	3	BMA	C4-C5-C6-O6
12	e	2	NAG	C4-C5-C6-O6
12	m	2	NAG	C4-C5-C6-O6
12	u	2	NAG	C4-C5-C6-O6
12	2	2	NAG	C4-C5-C6-O6
12	AA	2	NAG	C4-C5-C6-O6
12	IA	2	NAG	C4-C5-C6-O6
12	m	1	NAG	O5-C5-C6-O6
10	a	4	MAN	O5-C5-C6-O6
10	q	4	MAN	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
10	6	4	MAN	O5-C5-C6-O6
11	j	3	BMA	O5-C5-C6-O6
11	z	3	BMA	O5-C5-C6-O6
11	FA	3	BMA	O5-C5-C6-O6
12	2	1	NAG	O5-C5-C6-O6
12	IA	1	NAG	O5-C5-C6-O6
11	b	4	MAN	O5-C5-C6-O6
15	l	8	MAN	O5-C5-C6-O6
15	1	8	MAN	O5-C5-C6-O6
15	HA	8	MAN	O5-C5-C6-O6
11	r	4	MAN	O5-C5-C6-O6
11	7	4	MAN	O5-C5-C6-O6
10	n	4	MAN	C4-C5-C6-O6
10	3	4	MAN	C4-C5-C6-O6
10	JA	4	MAN	C4-C5-C6-O6
15	l	5	MAN	O5-C5-C6-O6
15	1	5	MAN	O5-C5-C6-O6
15	HA	5	MAN	O5-C5-C6-O6
14	f	2	NAG	C4-C5-C6-O6
14	v	2	NAG	C4-C5-C6-O6
14	BA	2	NAG	C4-C5-C6-O6
10	n	3	BMA	O5-C5-C6-O6
10	3	3	BMA	O5-C5-C6-O6
10	JA	3	BMA	O5-C5-C6-O6
15	l	1	NAG	O5-C5-C6-O6
15	1	1	NAG	O5-C5-C6-O6
15	HA	1	NAG	O5-C5-C6-O6
15	l	8	MAN	C4-C5-C6-O6
15	1	8	MAN	C4-C5-C6-O6
15	HA	8	MAN	C4-C5-C6-O6
12	e	1	NAG	C4-C5-C6-O6
12	u	1	NAG	C4-C5-C6-O6
12	AA	1	NAG	C4-C5-C6-O6
13	d	2	NAG	O5-C5-C6-O6
13	t	2	NAG	O5-C5-C6-O6
13	9	2	NAG	O5-C5-C6-O6
11	b	4	MAN	C4-C5-C6-O6
11	j	2	NAG	C4-C5-C6-O6
11	r	4	MAN	C4-C5-C6-O6
11	z	2	NAG	C4-C5-C6-O6
11	7	4	MAN	C4-C5-C6-O6
11	FA	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
15	l	1	NAG	C4-C5-C6-O6
15	1	1	NAG	C4-C5-C6-O6
15	HA	1	NAG	C4-C5-C6-O6
12	m	1	NAG	C4-C5-C6-O6
12	2	1	NAG	C4-C5-C6-O6
12	IA	1	NAG	C4-C5-C6-O6
10	q	2	NAG	O5-C5-C6-O6
10	6	2	NAG	O5-C5-C6-O6
11	j	2	NAG	O5-C5-C6-O6
11	z	2	NAG	O5-C5-C6-O6
11	FA	2	NAG	O5-C5-C6-O6
12	m	2	NAG	O5-C5-C6-O6
12	2	2	NAG	O5-C5-C6-O6
12	IA	2	NAG	O5-C5-C6-O6
14	f	4	MAN	O5-C5-C6-O6
14	v	4	MAN	O5-C5-C6-O6
14	BA	4	MAN	O5-C5-C6-O6
13	d	2	NAG	C4-C5-C6-O6
10	a	2	NAG	O5-C5-C6-O6
13	t	2	NAG	C4-C5-C6-O6
13	9	2	NAG	C4-C5-C6-O6
9	Y	1	NAG	C8-C7-N2-C2
9	Y	1	NAG	O7-C7-N2-C2
9	Y	2	NAG	C8-C7-N2-C2
9	Y	2	NAG	O7-C7-N2-C2
9	o	1	NAG	C8-C7-N2-C2
9	o	1	NAG	O7-C7-N2-C2
9	o	2	NAG	C8-C7-N2-C2
9	o	2	NAG	O7-C7-N2-C2
9	4	1	NAG	C8-C7-N2-C2
9	4	1	NAG	O7-C7-N2-C2
9	4	2	NAG	C8-C7-N2-C2
9	4	2	NAG	O7-C7-N2-C2
9	Y	1	NAG	O5-C5-C6-O6
9	o	1	NAG	O5-C5-C6-O6
9	4	1	NAG	O5-C5-C6-O6
12	i	1	NAG	O5-C5-C6-O6
12	y	1	NAG	O5-C5-C6-O6
12	EA	1	NAG	O5-C5-C6-O6
11	b	2	NAG	C4-C5-C6-O6
11	r	2	NAG	C4-C5-C6-O6
11	7	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
10	a	2	NAG	C4-C5-C6-O6
10	q	2	NAG	C4-C5-C6-O6
10	6	2	NAG	C4-C5-C6-O6
9	Y	1	NAG	C4-C5-C6-O6
9	o	1	NAG	C4-C5-C6-O6
9	4	1	NAG	C4-C5-C6-O6
12	i	1	NAG	C4-C5-C6-O6
12	y	1	NAG	C4-C5-C6-O6
12	EA	1	NAG	C4-C5-C6-O6
15	l	4	MAN	O5-C5-C6-O6
15	1	4	MAN	O5-C5-C6-O6
15	HA	4	MAN	O5-C5-C6-O6
10	a	4	MAN	C4-C5-C6-O6
10	q	4	MAN	C4-C5-C6-O6
10	6	4	MAN	C4-C5-C6-O6
15	HA	4	MAN	C4-C5-C6-O6
15	l	4	MAN	C4-C5-C6-O6
15	1	4	MAN	C4-C5-C6-O6
9	0	2	NAG	O5-C5-C6-O6
9	GA	2	NAG	O5-C5-C6-O6
9	k	2	NAG	O5-C5-C6-O6
15	l	7	MAN	C4-C5-C6-O6
15	1	7	MAN	C4-C5-C6-O6
15	HA	7	MAN	C4-C5-C6-O6
15	1	7	MAN	O5-C5-C6-O6
15	l	7	MAN	O5-C5-C6-O6
15	HA	7	MAN	O5-C5-C6-O6
9	4	2	NAG	O5-C5-C6-O6
9	Y	2	NAG	O5-C5-C6-O6
9	o	2	NAG	O5-C5-C6-O6
11	j	1	NAG	O5-C5-C6-O6
11	z	1	NAG	O5-C5-C6-O6
11	FA	1	NAG	O5-C5-C6-O6
11	j	4	MAN	O5-C5-C6-O6
11	z	4	MAN	O5-C5-C6-O6
11	FA	4	MAN	O5-C5-C6-O6
12	c	3	BMA	C4-C5-C6-O6
12	s	3	BMA	C4-C5-C6-O6
12	8	3	BMA	C4-C5-C6-O6
14	v	1	NAG	O5-C5-C6-O6
14	BA	1	NAG	O5-C5-C6-O6
14	f	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
13	d	4	MAN	O5-C5-C6-O6
13	t	4	MAN	O5-C5-C6-O6
13	9	4	MAN	O5-C5-C6-O6
12	c	3	BMA	O5-C5-C6-O6
12	s	3	BMA	O5-C5-C6-O6
12	8	3	BMA	O5-C5-C6-O6
15	l	5	MAN	C4-C5-C6-O6
15	1	5	MAN	C4-C5-C6-O6
15	HA	5	MAN	C4-C5-C6-O6
12	c	1	NAG	C3-C2-N2-C7
12	s	1	NAG	C3-C2-N2-C7
12	8	1	NAG	C3-C2-N2-C7
15	l	2	NAG	C3-C2-N2-C7
15	1	2	NAG	C3-C2-N2-C7
15	HA	2	NAG	C3-C2-N2-C7
14	v	4	MAN	C4-C5-C6-O6
9	p	1	NAG	C4-C5-C6-O6
14	f	4	MAN	C4-C5-C6-O6
14	BA	4	MAN	C4-C5-C6-O6
9	5	1	NAG	C4-C5-C6-O6
9	p	1	NAG	O5-C5-C6-O6
9	Z	1	NAG	C4-C5-C6-O6
9	0	2	NAG	C4-C5-C6-O6
9	GA	2	NAG	C4-C5-C6-O6
9	k	2	NAG	C4-C5-C6-O6
9	Z	1	NAG	O5-C5-C6-O6
9	5	1	NAG	O5-C5-C6-O6
14	v	6	MAN	C4-C5-C6-O6
14	BA	6	MAN	C4-C5-C6-O6
14	f	6	MAN	C4-C5-C6-O6
15	l	6	MAN	C4-C5-C6-O6
15	1	6	MAN	C4-C5-C6-O6
15	HA	6	MAN	C4-C5-C6-O6
9	Y	2	NAG	C1-C2-N2-C7
9	Z	1	NAG	C1-C2-N2-C7
9	o	2	NAG	C1-C2-N2-C7
9	p	1	NAG	C1-C2-N2-C7
9	4	2	NAG	C1-C2-N2-C7
9	5	1	NAG	C1-C2-N2-C7
12	c	1	NAG	C1-C2-N2-C7
12	m	1	NAG	C1-C2-N2-C7
12	s	1	NAG	C1-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
12	2	1	NAG	C1-C2-N2-C7
12	8	1	NAG	C1-C2-N2-C7
12	IA	1	NAG	C1-C2-N2-C7
15	l	2	NAG	C1-C2-N2-C7
15	1	2	NAG	C1-C2-N2-C7
15	HA	2	NAG	C1-C2-N2-C7
14	f	6	MAN	O5-C5-C6-O6
14	BA	6	MAN	O5-C5-C6-O6
14	v	6	MAN	O5-C5-C6-O6
9	Y	2	NAG	C3-C2-N2-C7
9	Z	1	NAG	C3-C2-N2-C7
9	o	2	NAG	C3-C2-N2-C7
9	p	1	NAG	C3-C2-N2-C7
9	4	2	NAG	C3-C2-N2-C7
9	5	1	NAG	C3-C2-N2-C7
12	m	1	NAG	C3-C2-N2-C7
12	2	1	NAG	C3-C2-N2-C7
12	IA	1	NAG	C3-C2-N2-C7
12	AA	3	BMA	O5-C5-C6-O6
12	e	3	BMA	O5-C5-C6-O6
12	u	3	BMA	O5-C5-C6-O6
12	2	3	BMA	C4-C5-C6-O6
12	m	3	BMA	C4-C5-C6-O6
12	IA	3	BMA	C4-C5-C6-O6
12	2	3	BMA	O5-C5-C6-O6
12	m	3	BMA	O5-C5-C6-O6
12	IA	3	BMA	O5-C5-C6-O6
12	c	2	NAG	C4-C5-C6-O6
12	s	2	NAG	C4-C5-C6-O6
12	8	2	NAG	C4-C5-C6-O6
12	AA	3	BMA	C4-C5-C6-O6
12	e	3	BMA	C4-C5-C6-O6
12	u	3	BMA	C4-C5-C6-O6

All (6) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	3	4	MAN	C1-C2-C3-C4-C5-O5
10	n	4	MAN	C1-C2-C3-C4-C5-O5
10	JA	4	MAN	C1-C2-C3-C4-C5-O5
11	FA	4	MAN	C1-C2-C3-C4-C5-O5
11	j	4	MAN	C1-C2-C3-C4-C5-O5

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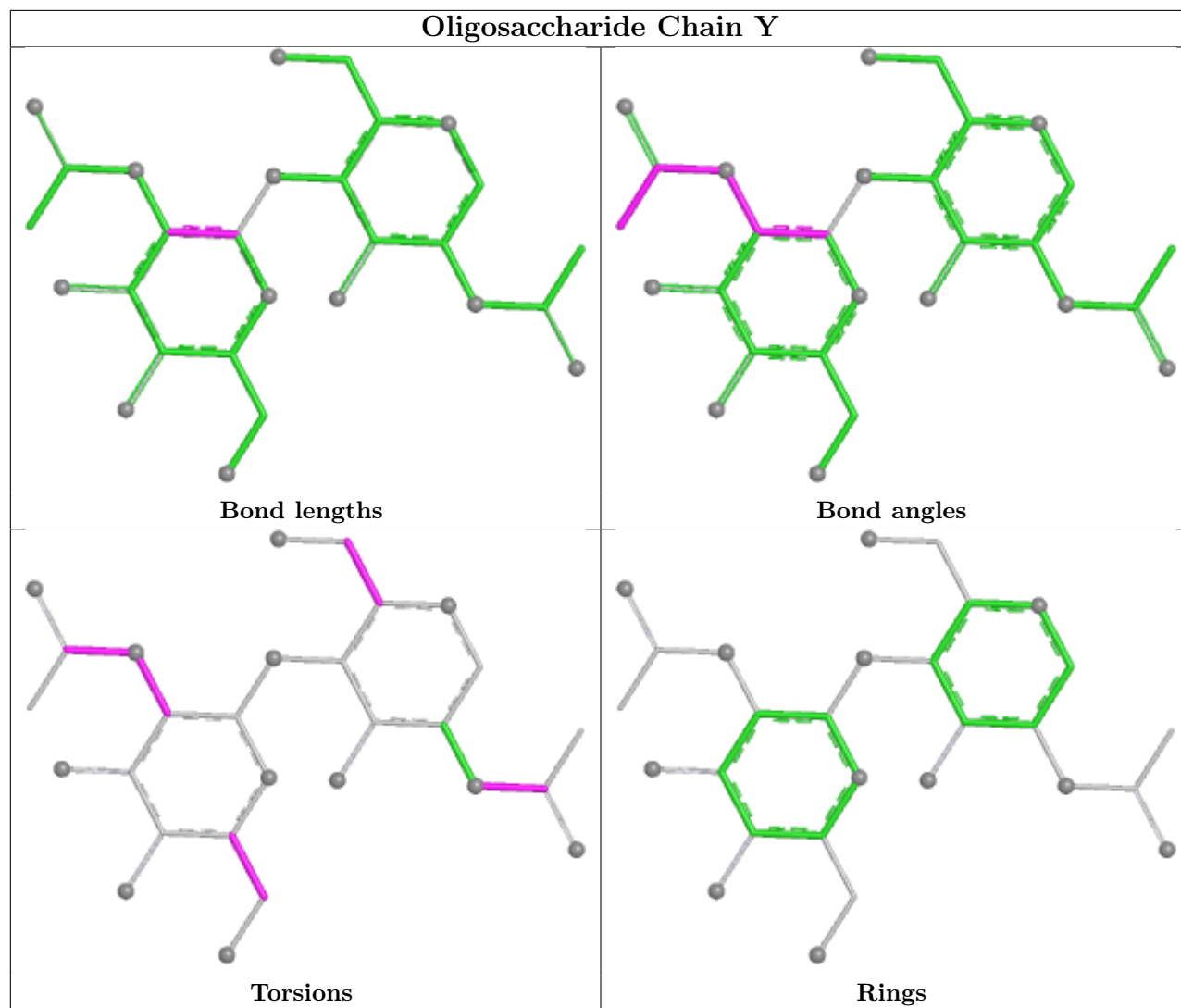
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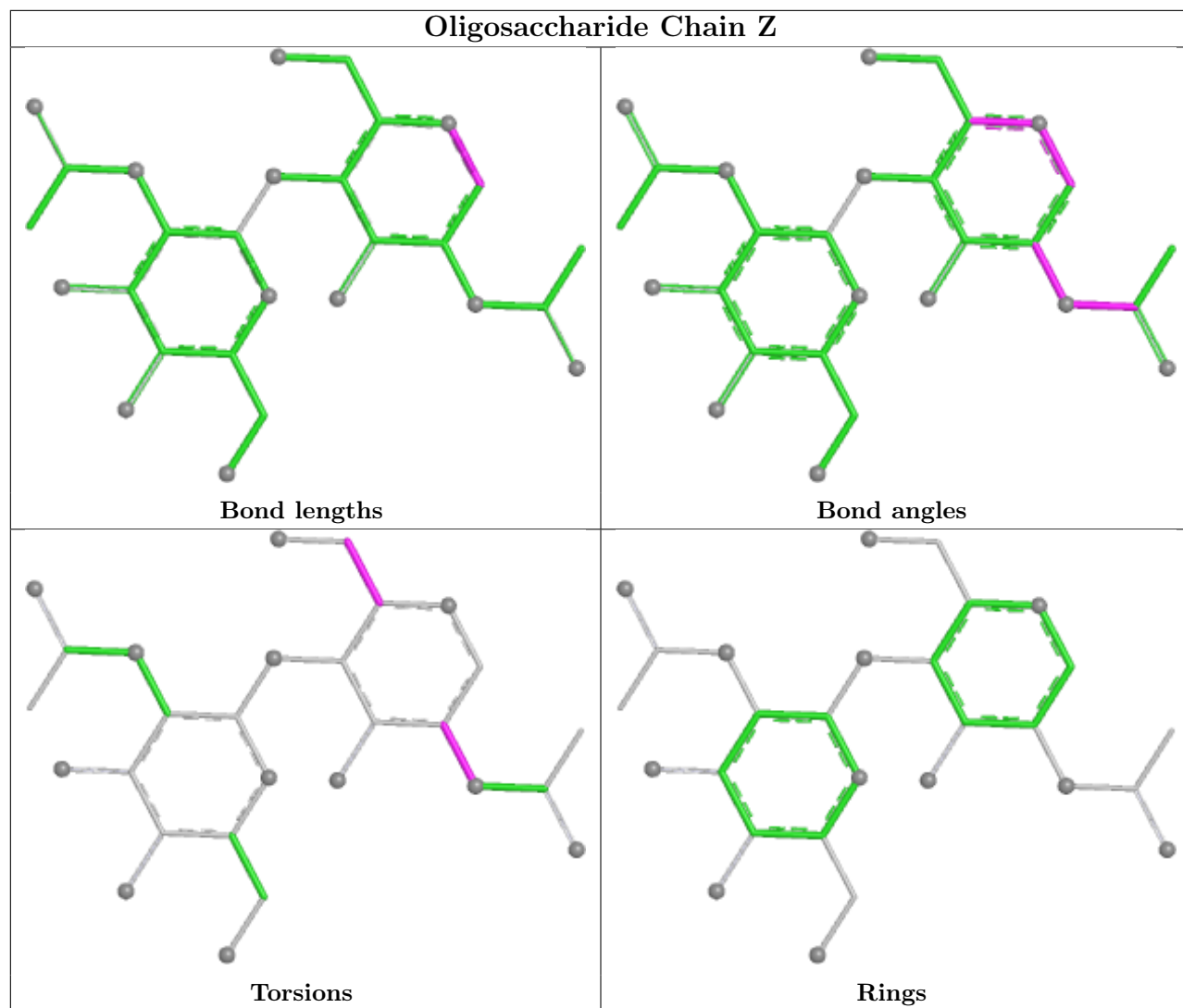
Mol	Chain	Res	Type	Atoms
11	z	4	MAN	C1-C2-C3-C4-C5-O5

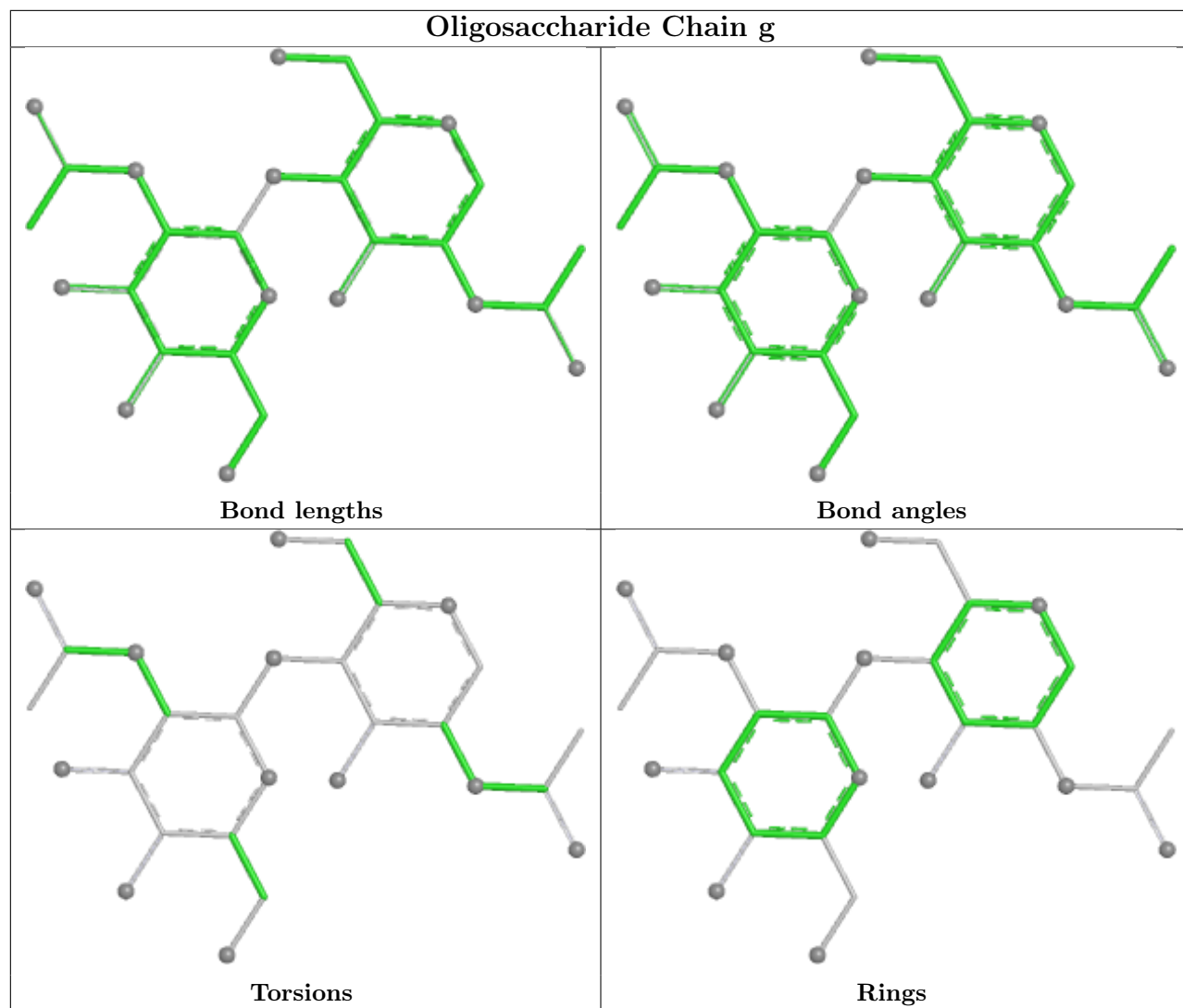
31 monomers are involved in 26 short contacts:

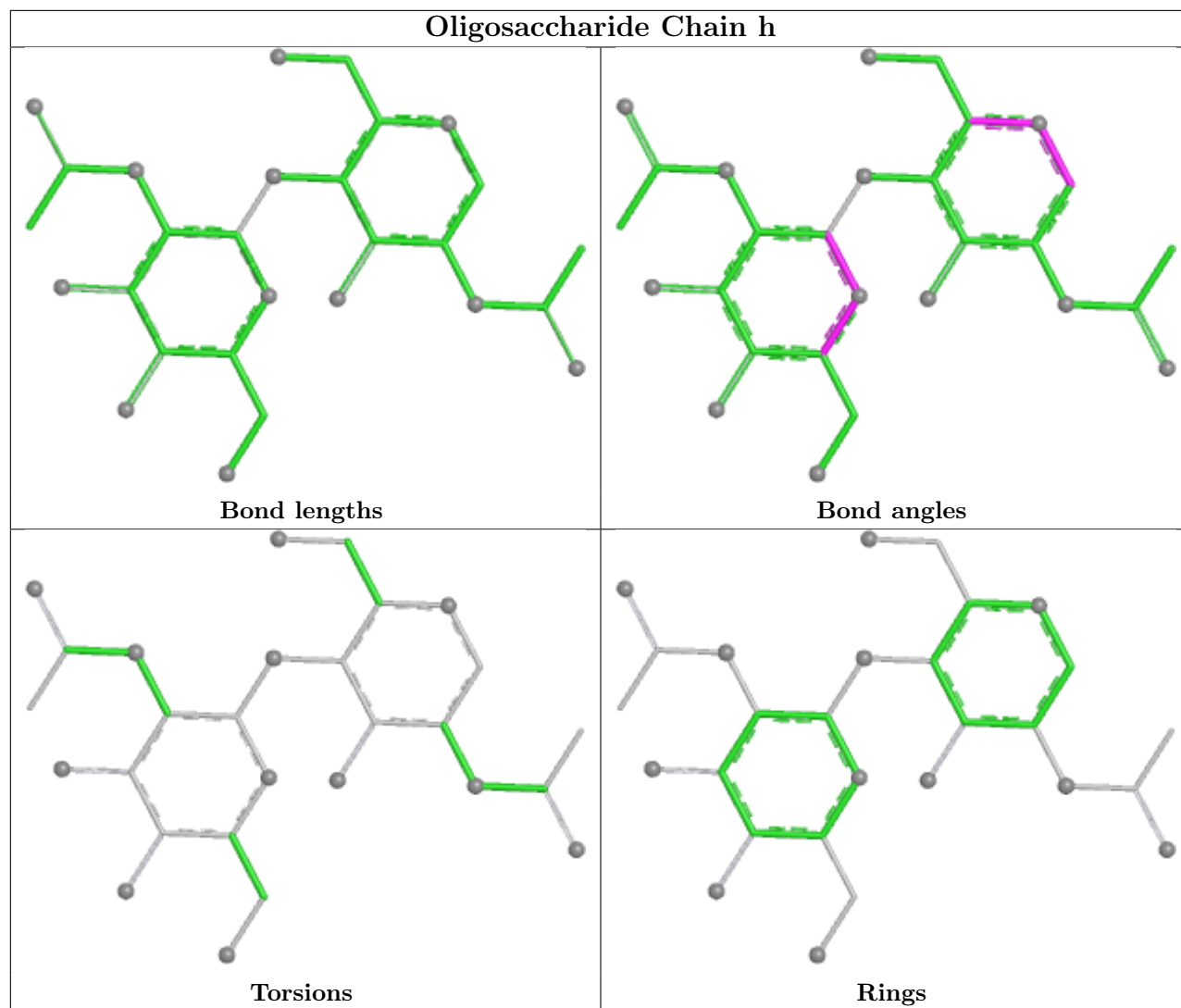
Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	q	3	BMA	1	0
11	FA	5	MAN	1	0
15	1	4	MAN	1	0
11	z	2	NAG	1	0
15	HA	4	MAN	1	0
13	t	3	BMA	1	0
15	HA	5	MAN	2	0
9	5	1	NAG	1	0
11	FA	3	BMA	1	0
9	k	1	NAG	1	0
15	l	4	MAN	1	0
15	l	5	MAN	2	0
9	0	1	NAG	1	0
11	z	3	BMA	1	0
15	1	5	MAN	3	0
11	z	5	MAN	1	0
13	9	4	MAN	1	0
10	3	1	NAG	1	0
11	j	2	NAG	1	0
10	a	3	BMA	1	0
9	p	1	NAG	1	0
13	9	3	BMA	1	0
9	Z	1	NAG	1	0
11	FA	2	NAG	1	0
13	d	4	MAN	1	0
11	j	3	BMA	1	0
9	GA	1	NAG	1	0
11	j	5	MAN	1	0
13	t	4	MAN	1	0
13	d	3	BMA	1	0
10	6	3	BMA	1	0

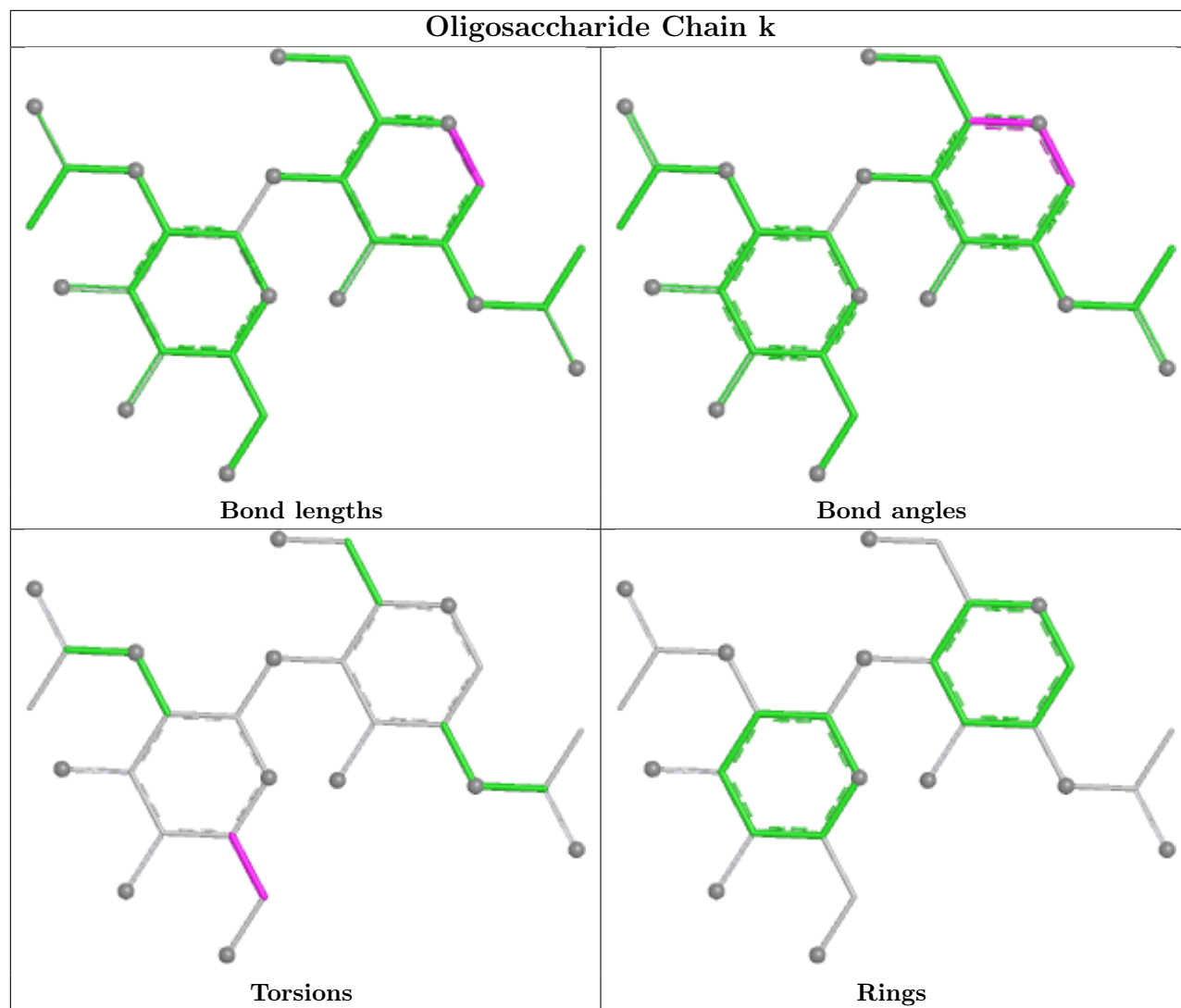
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

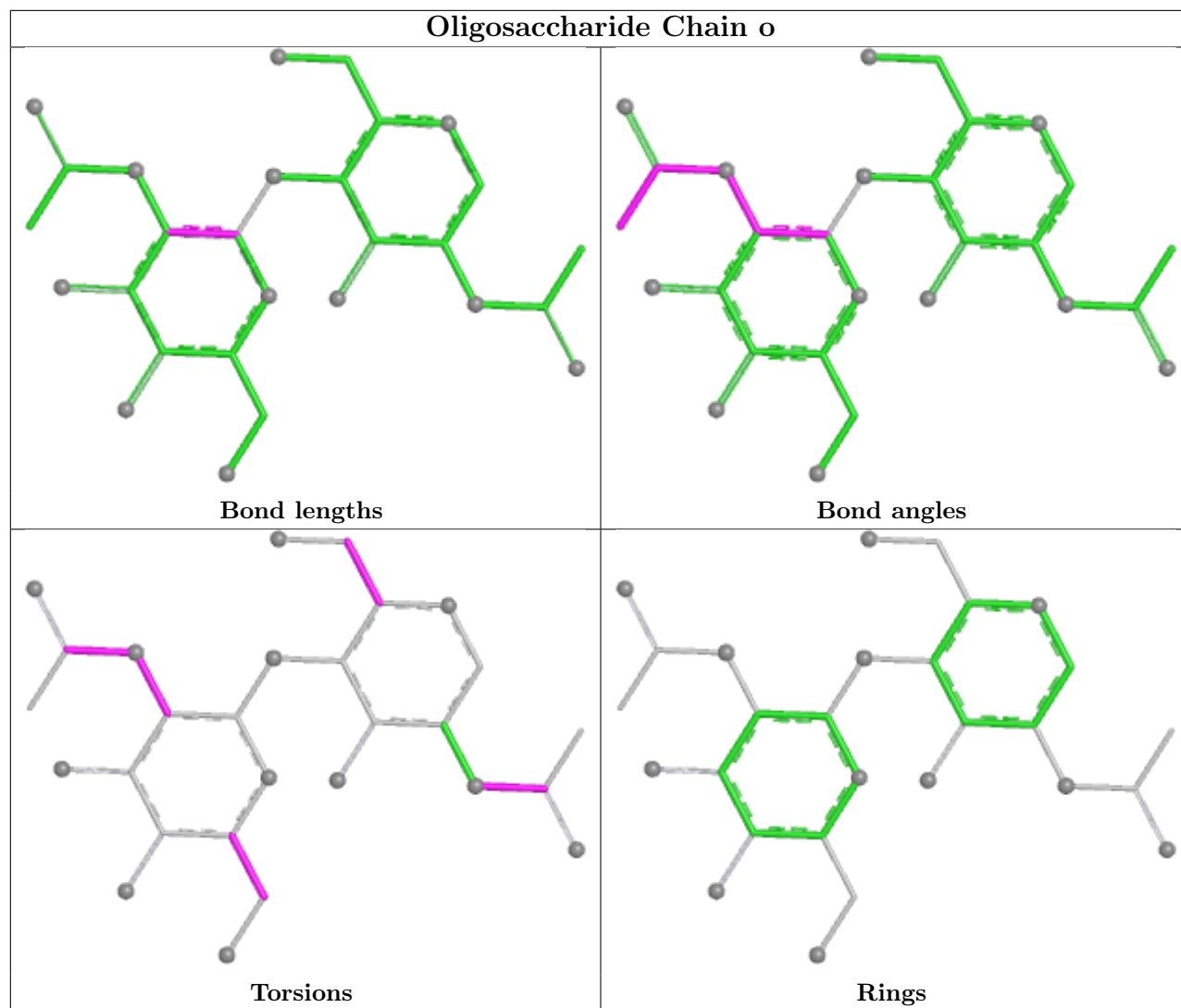


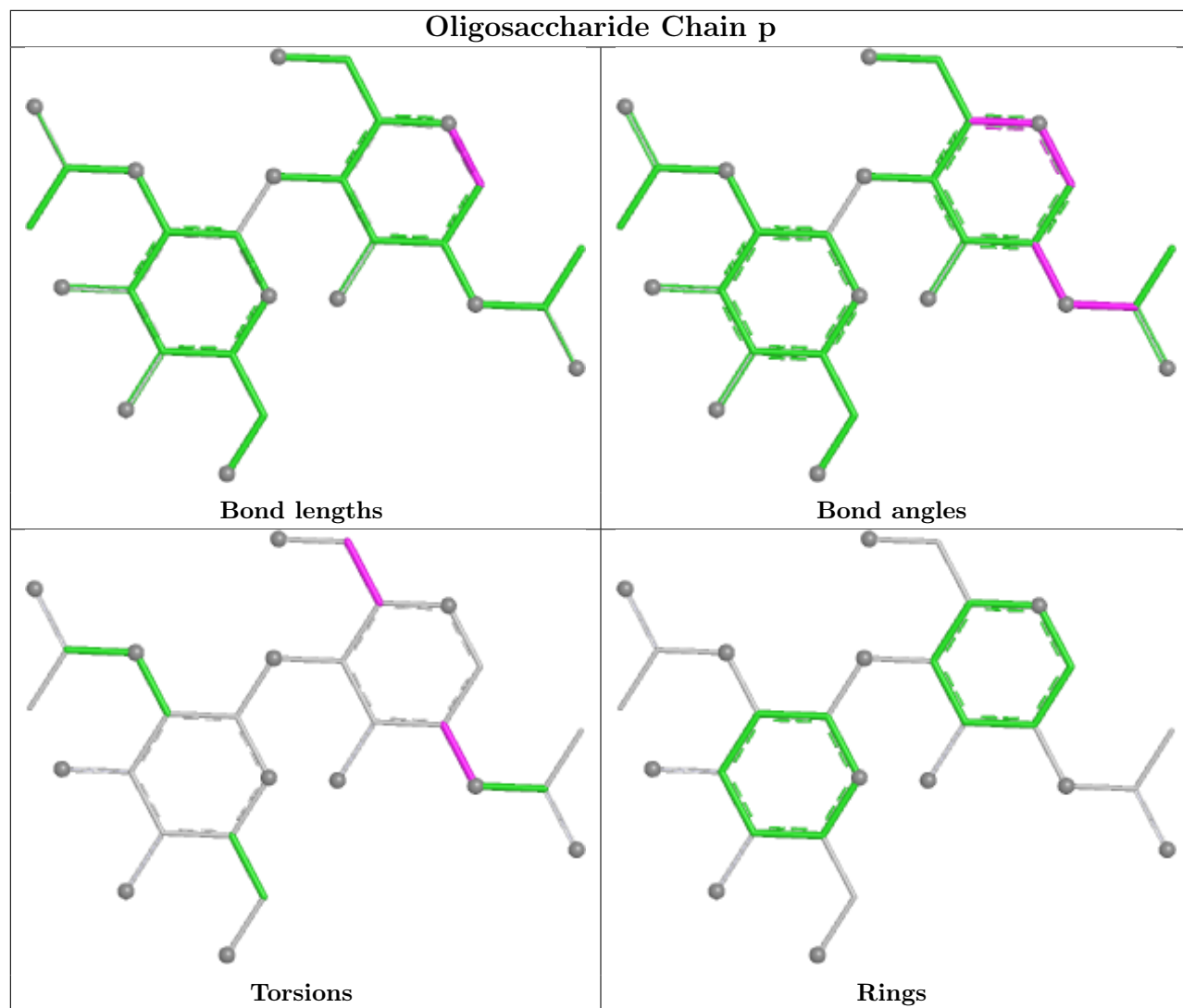


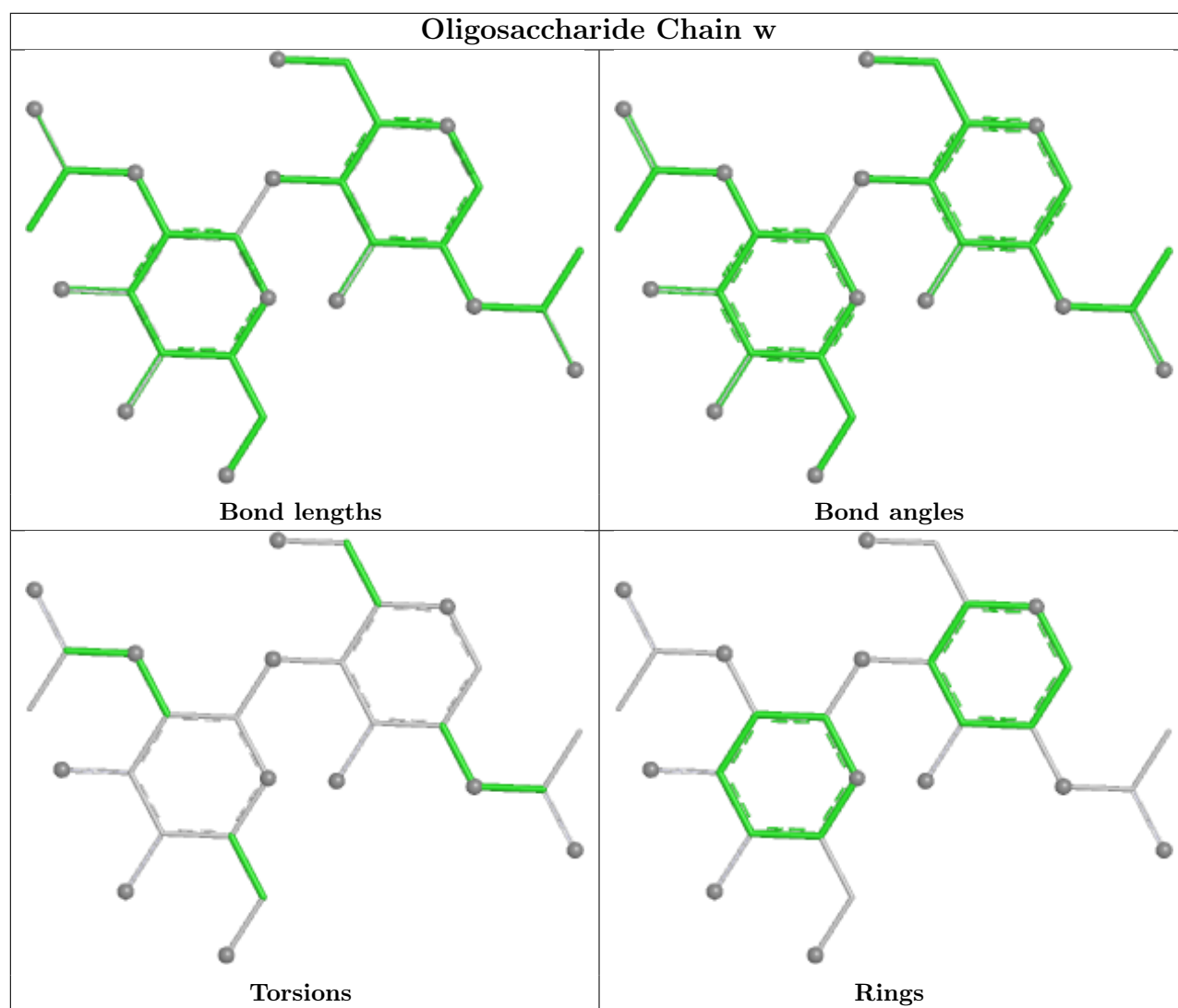


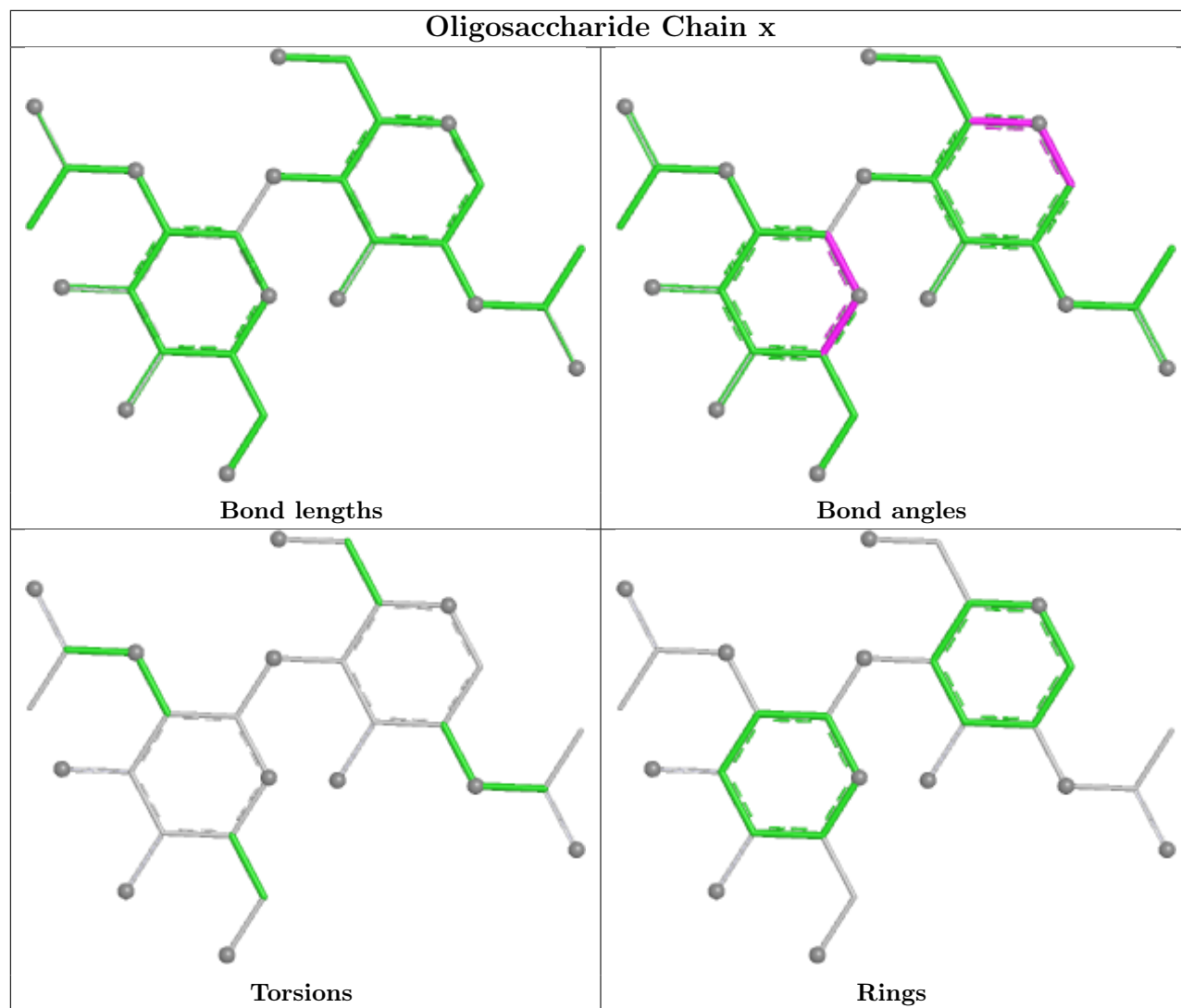


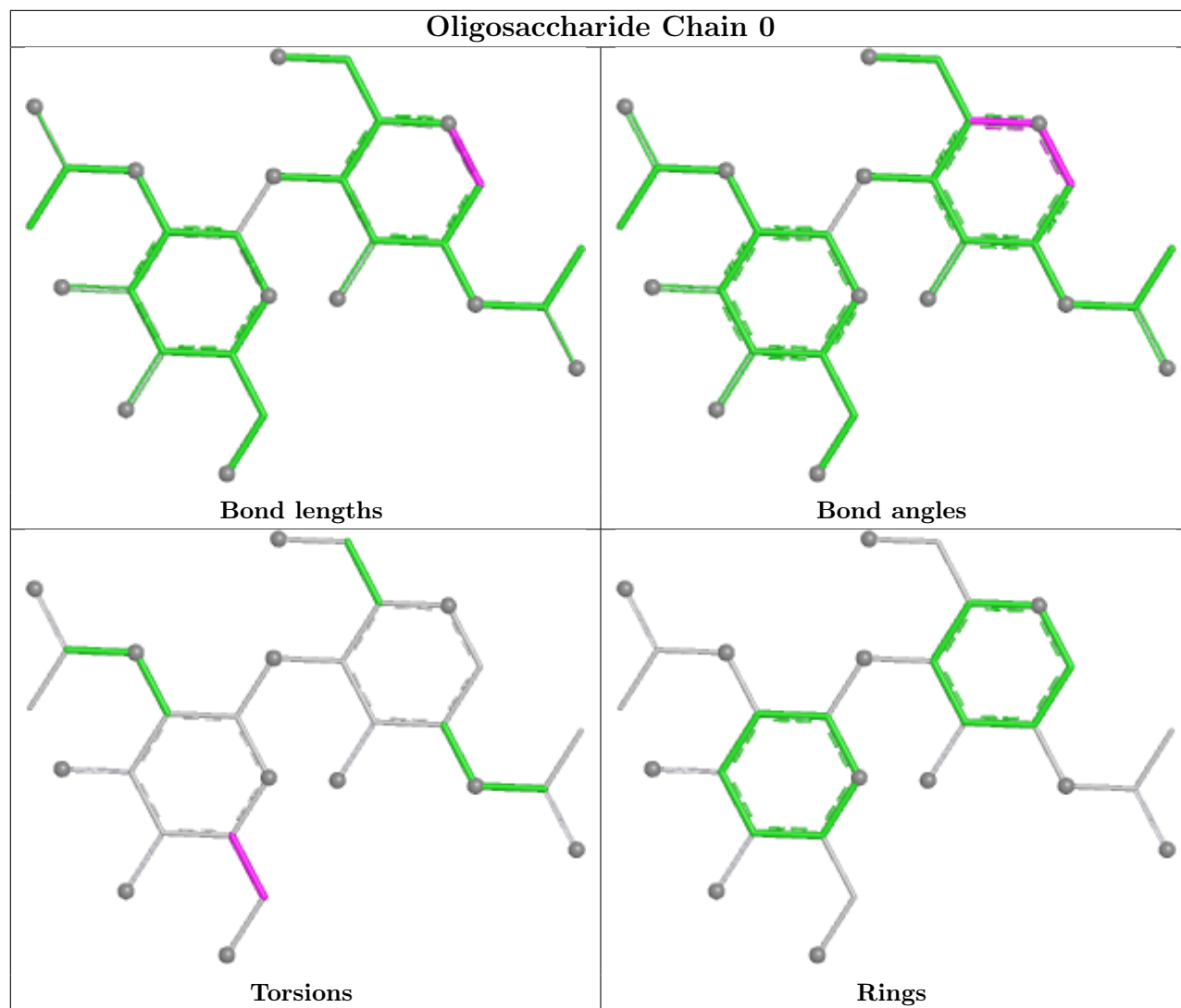


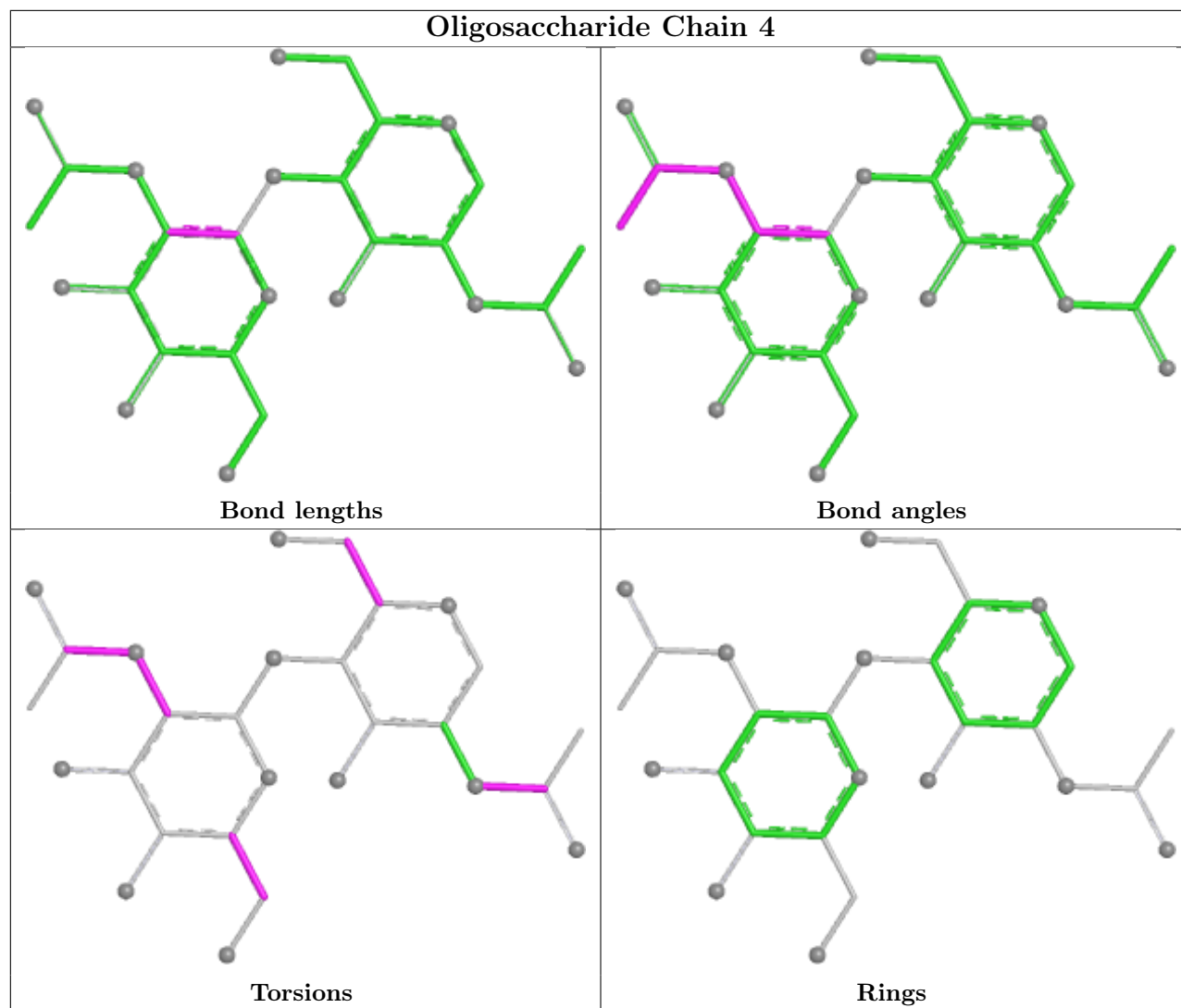


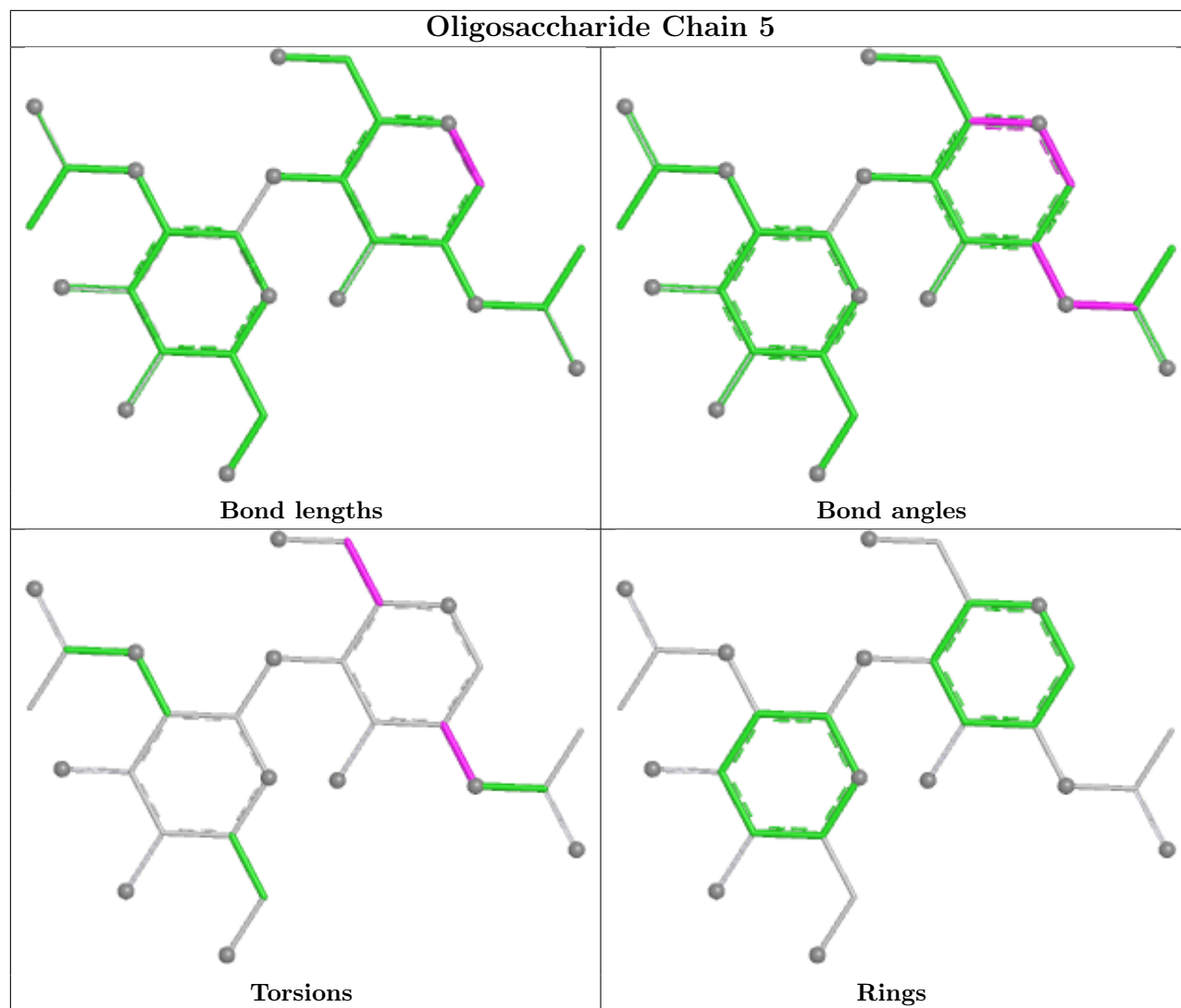


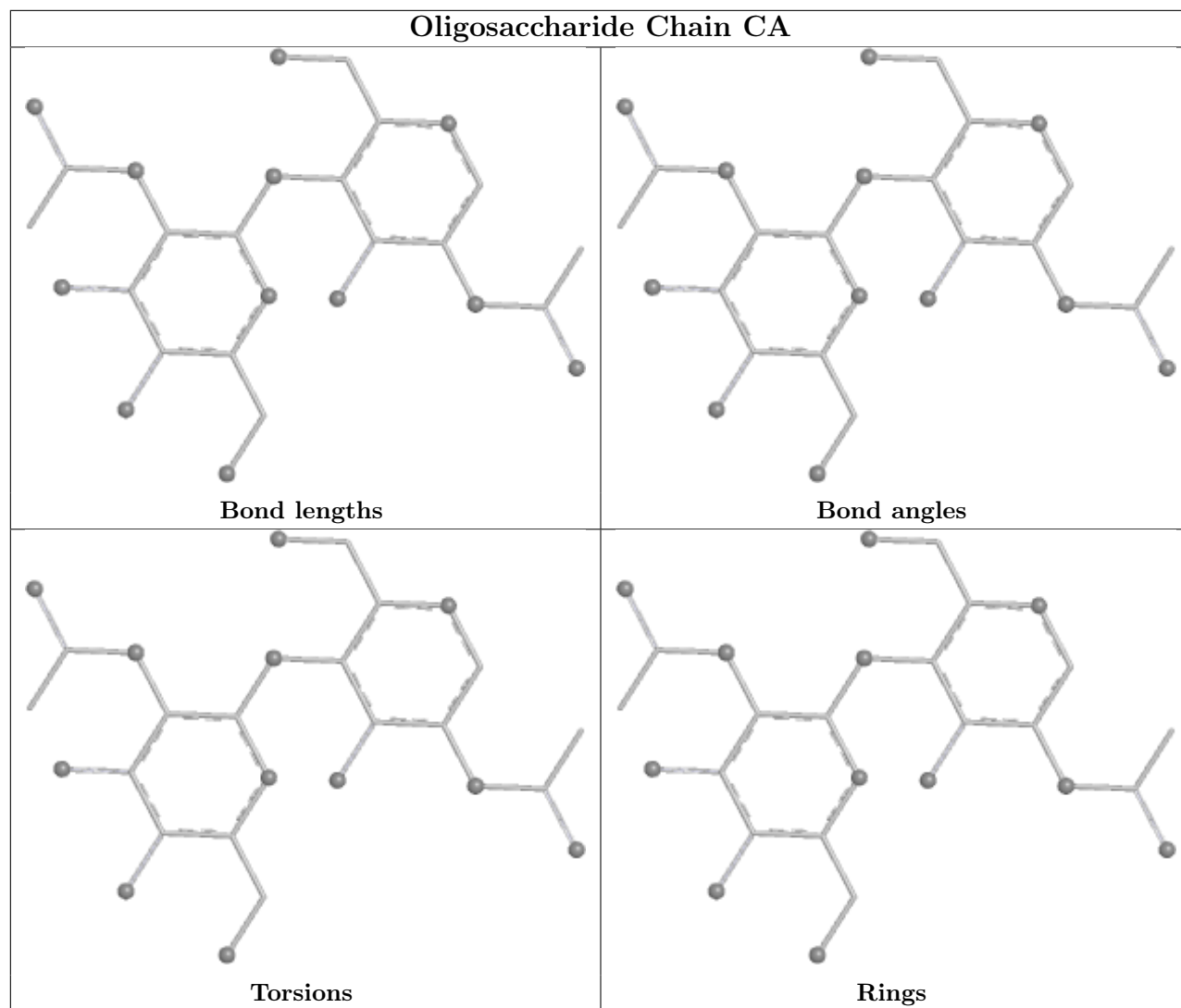


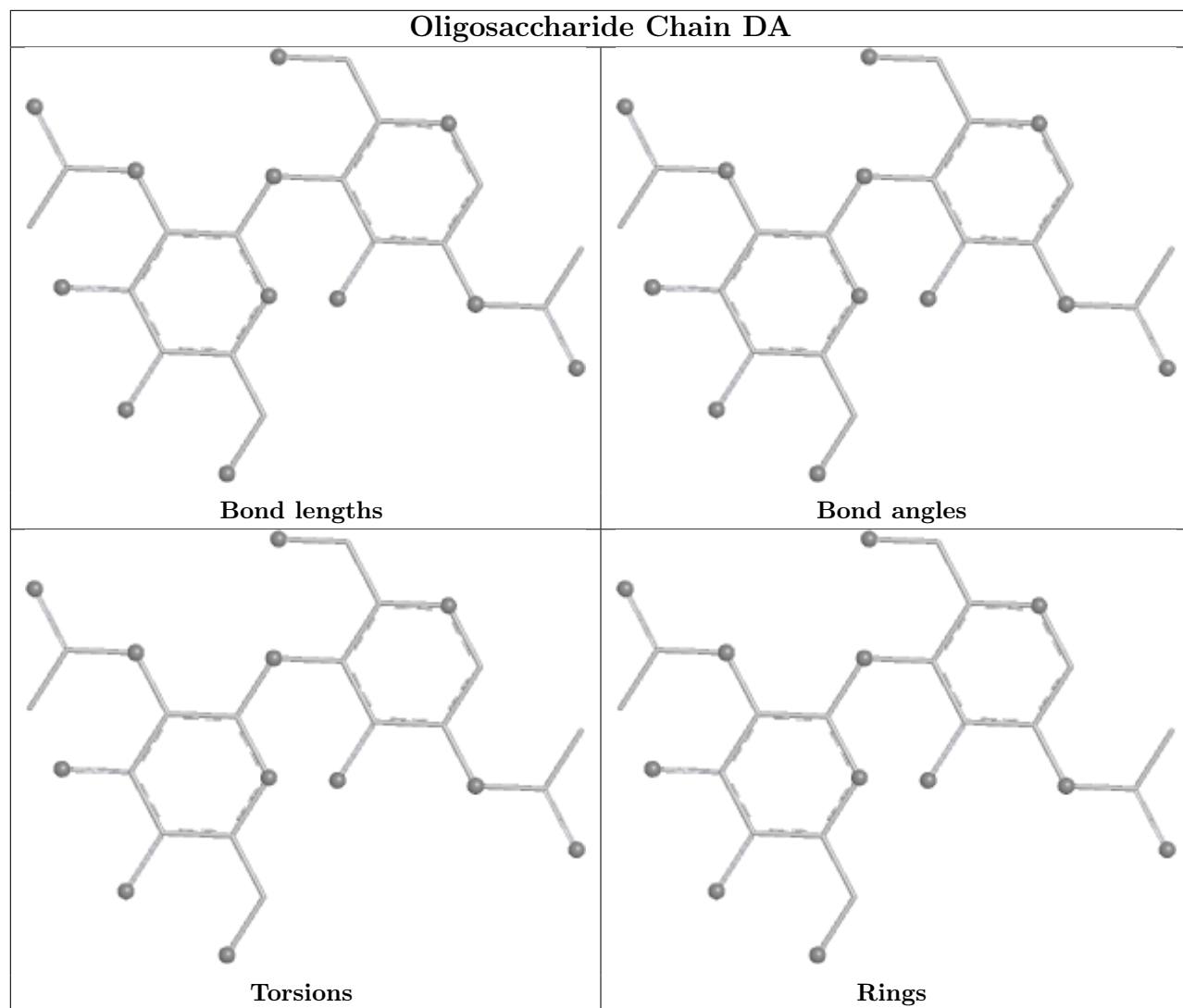


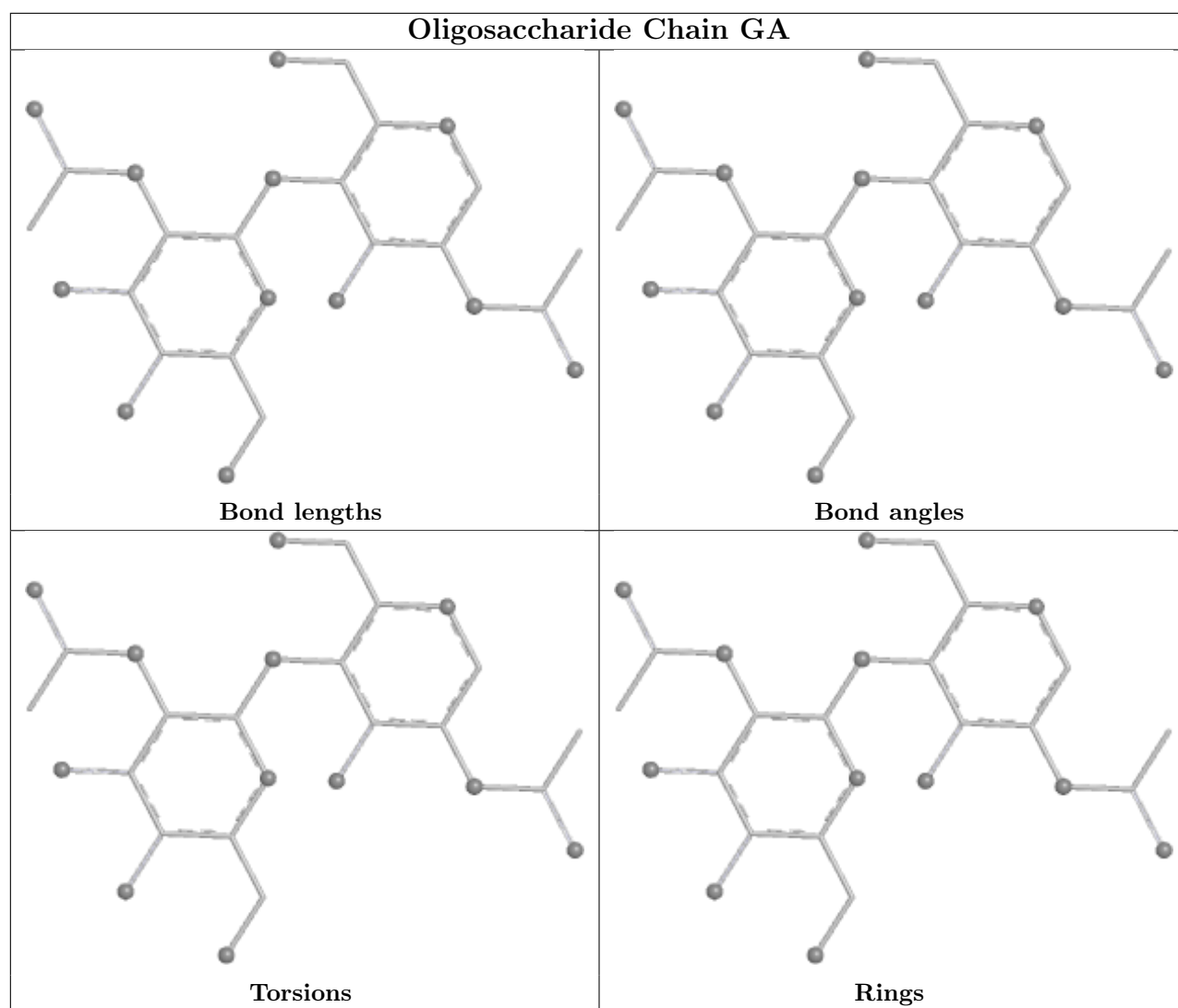


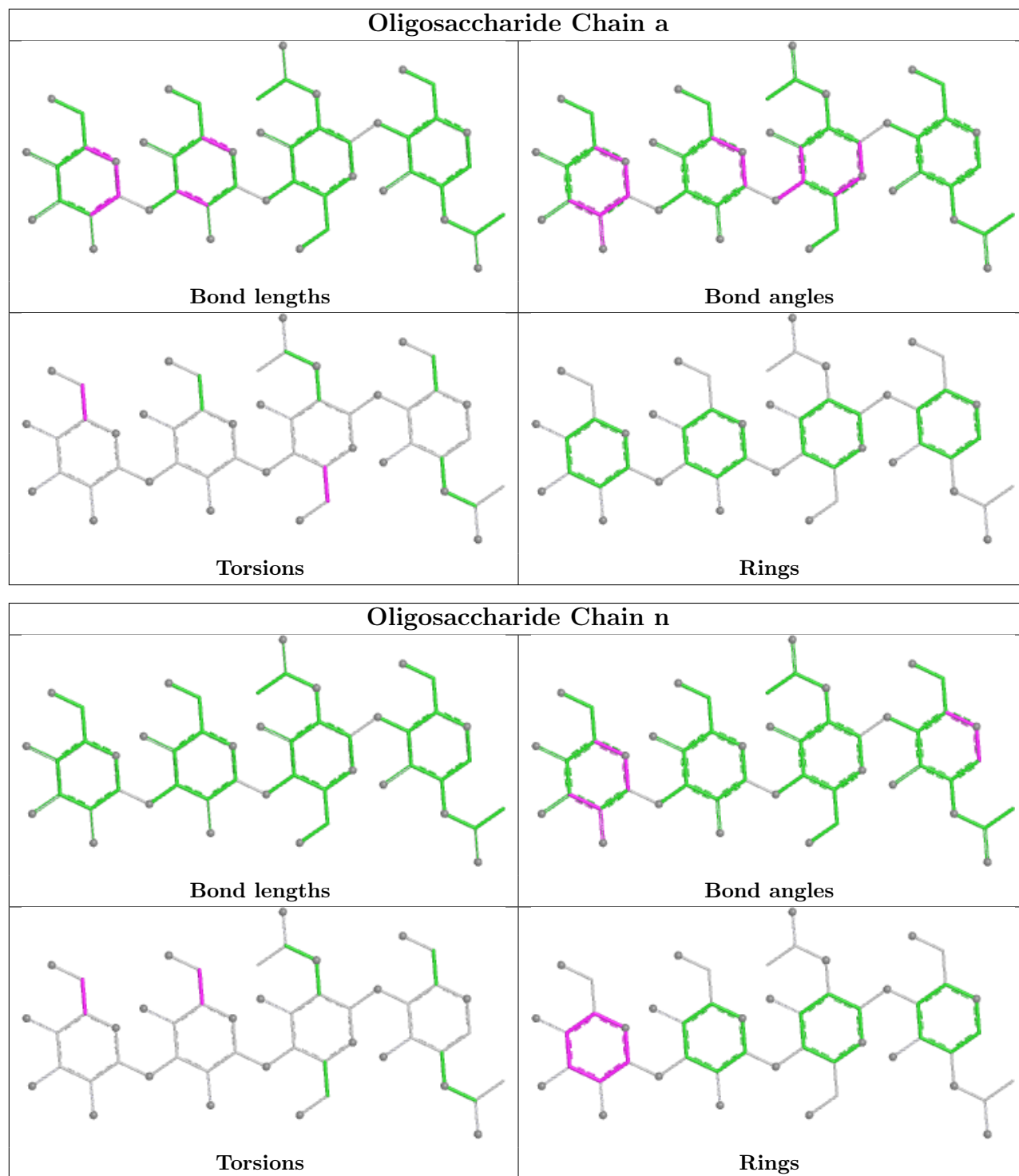


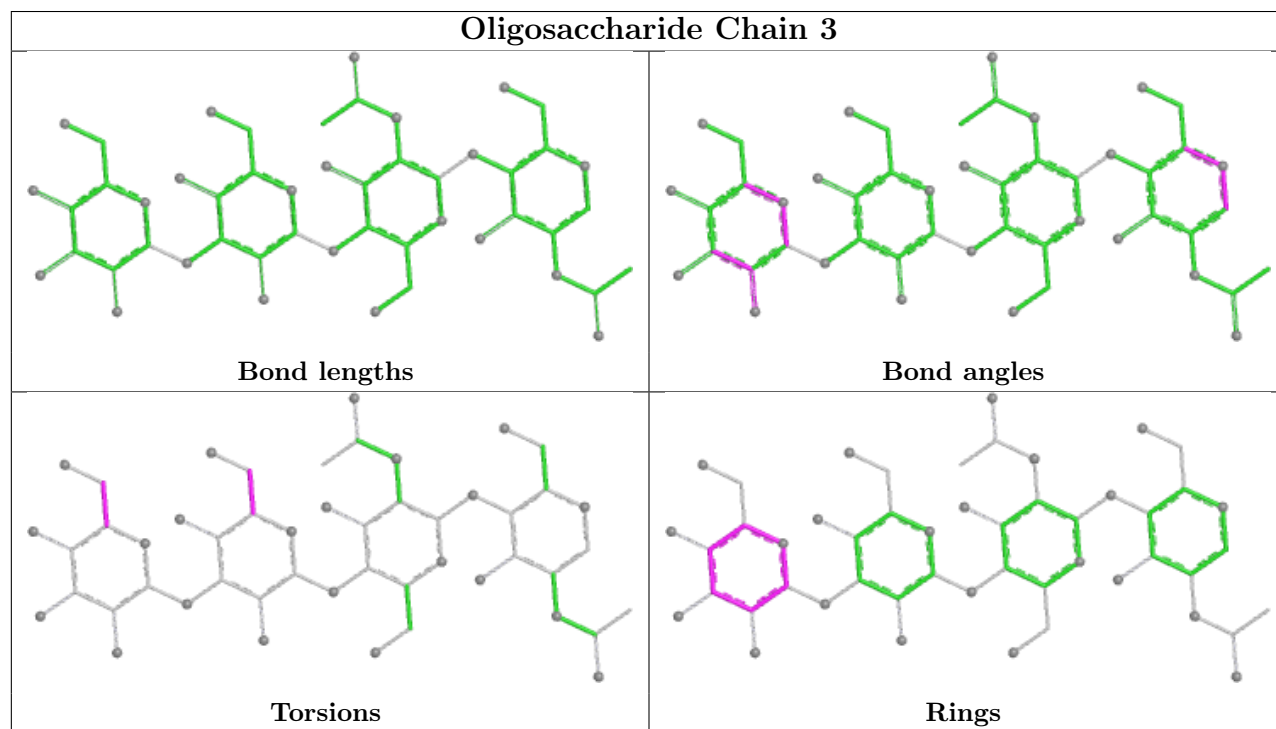
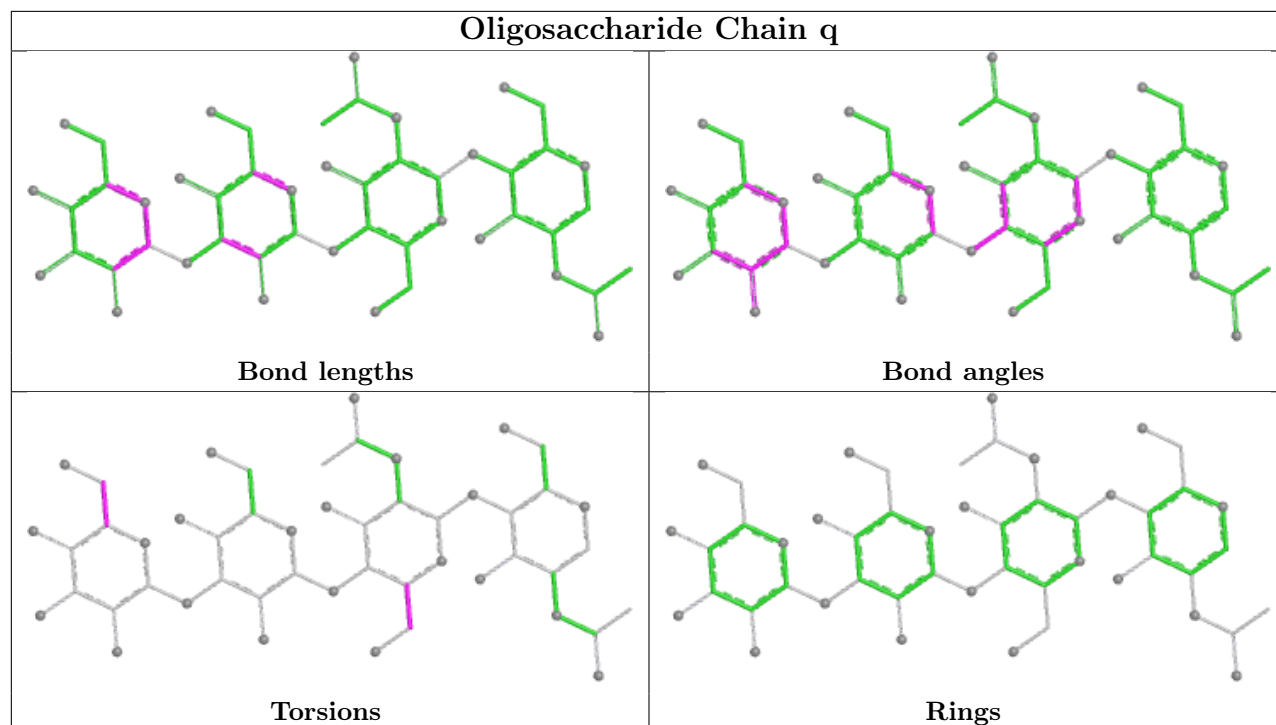


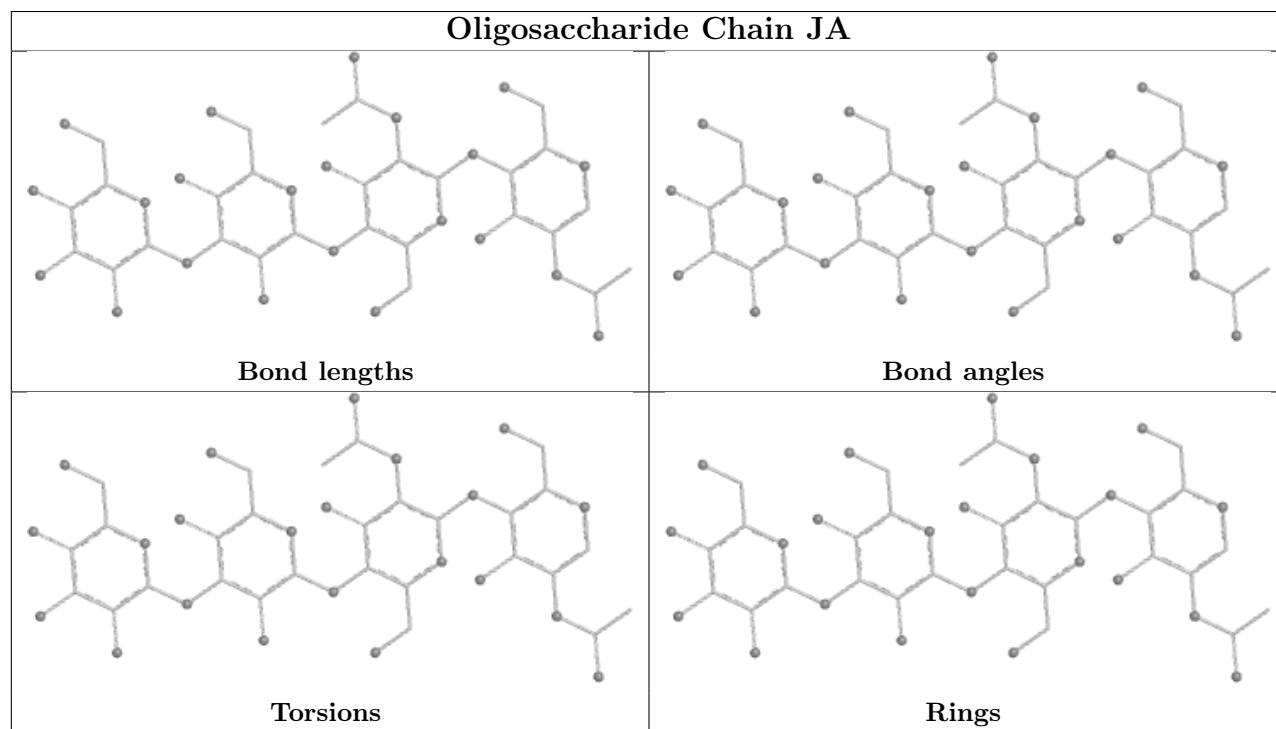
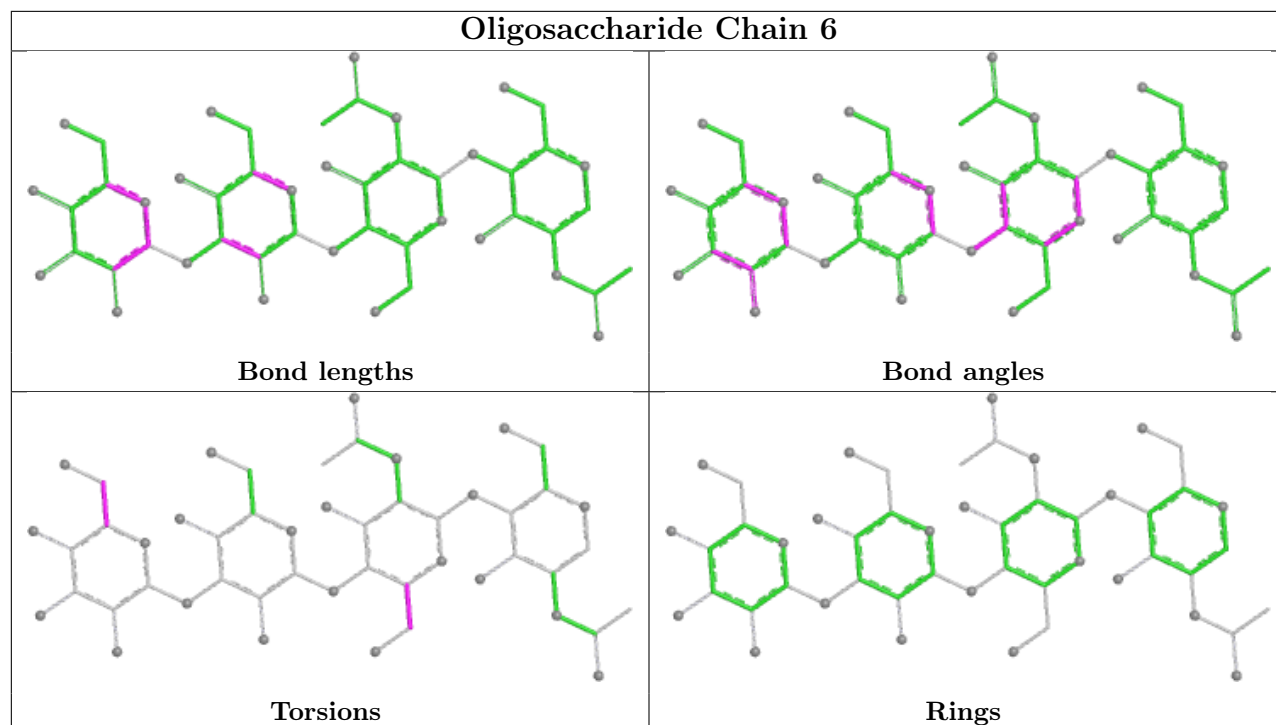


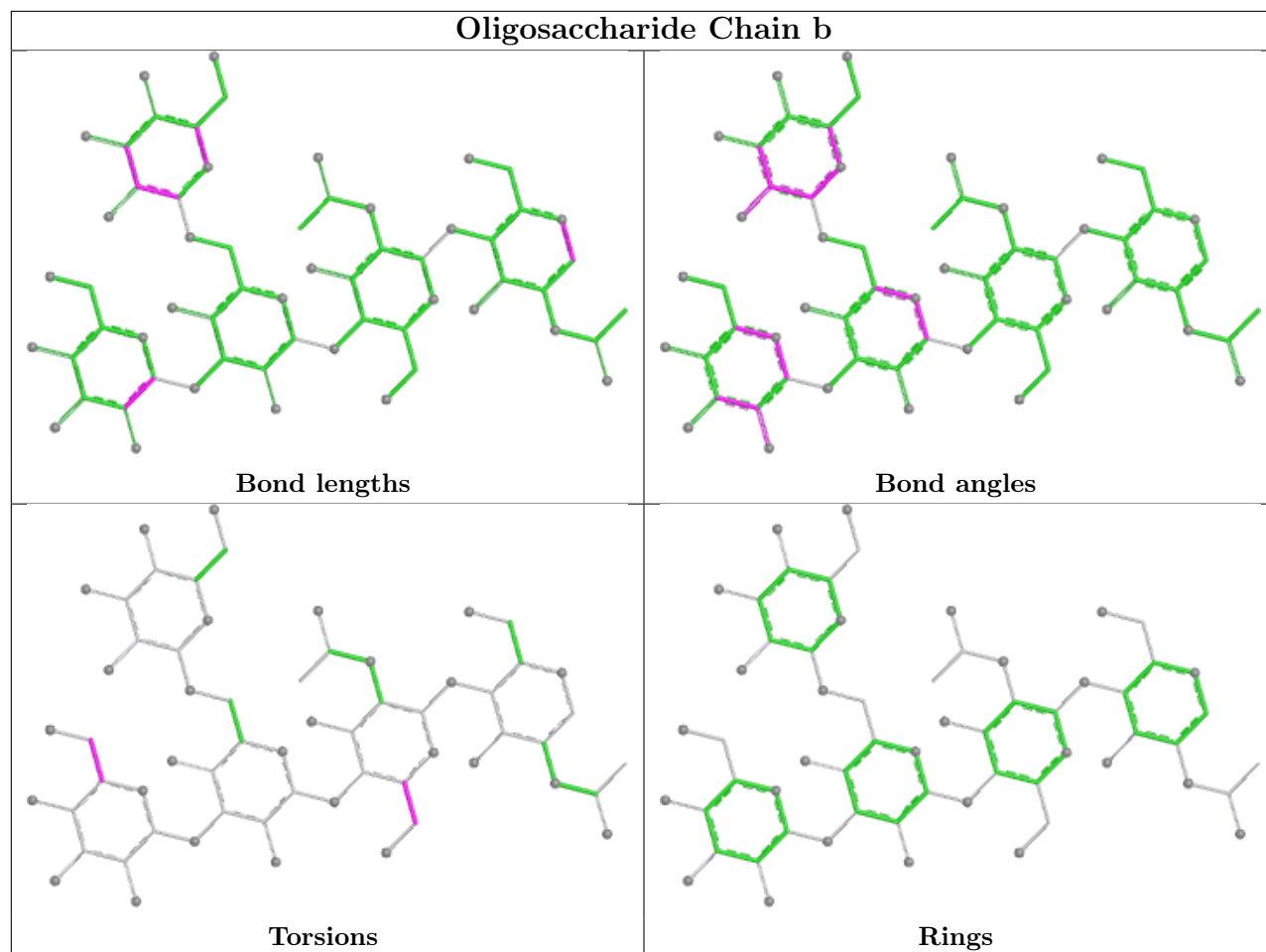


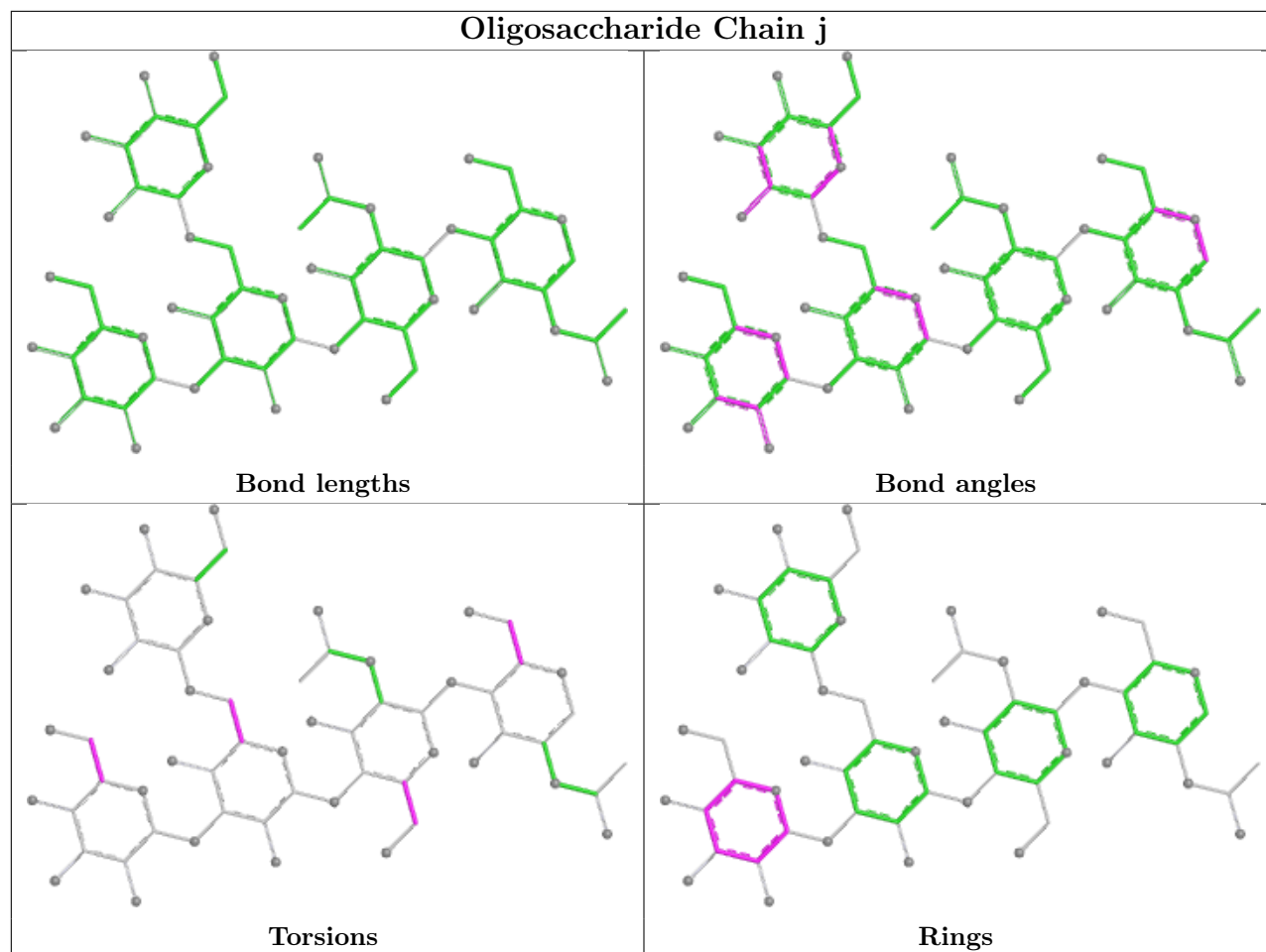


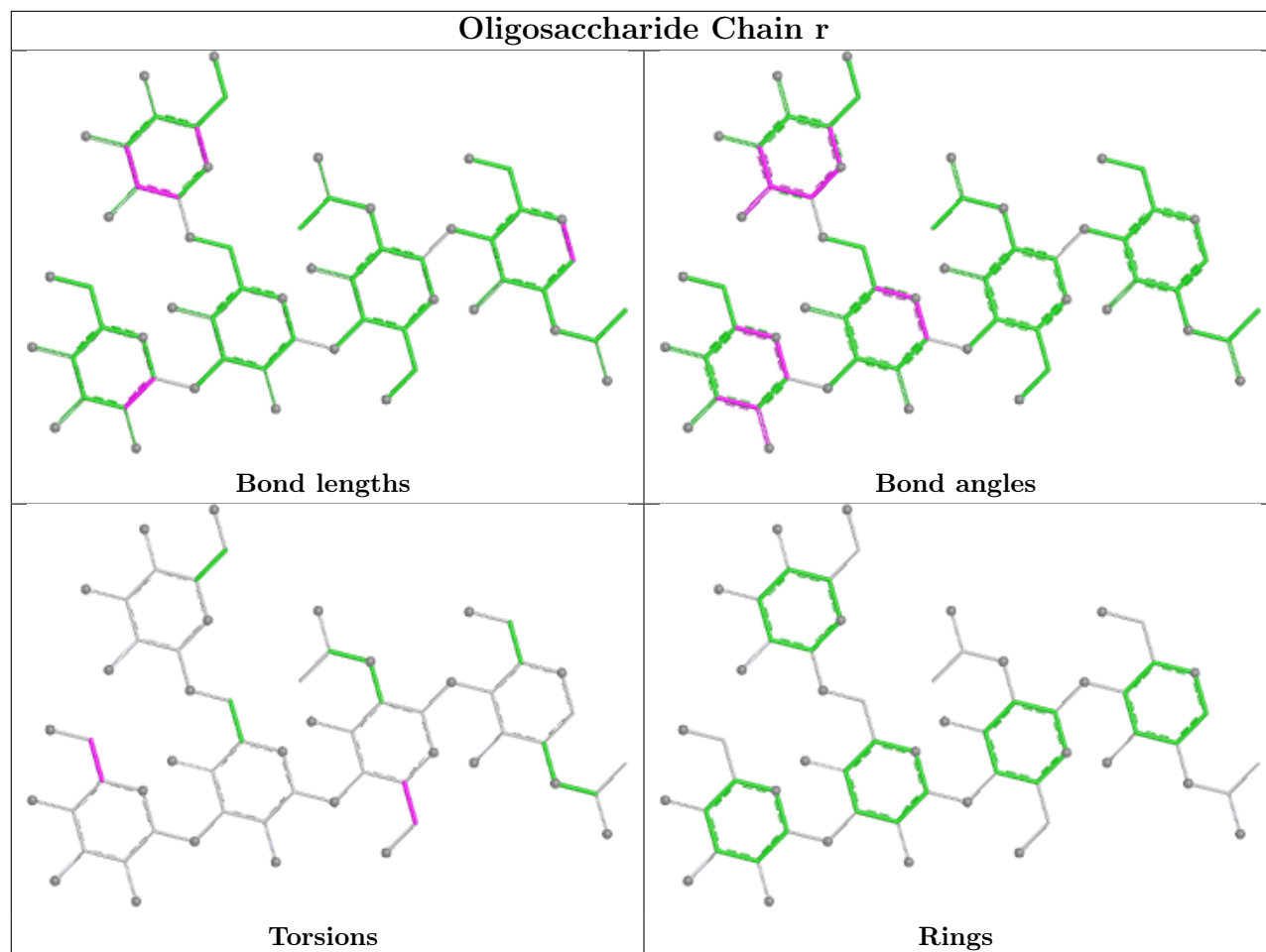


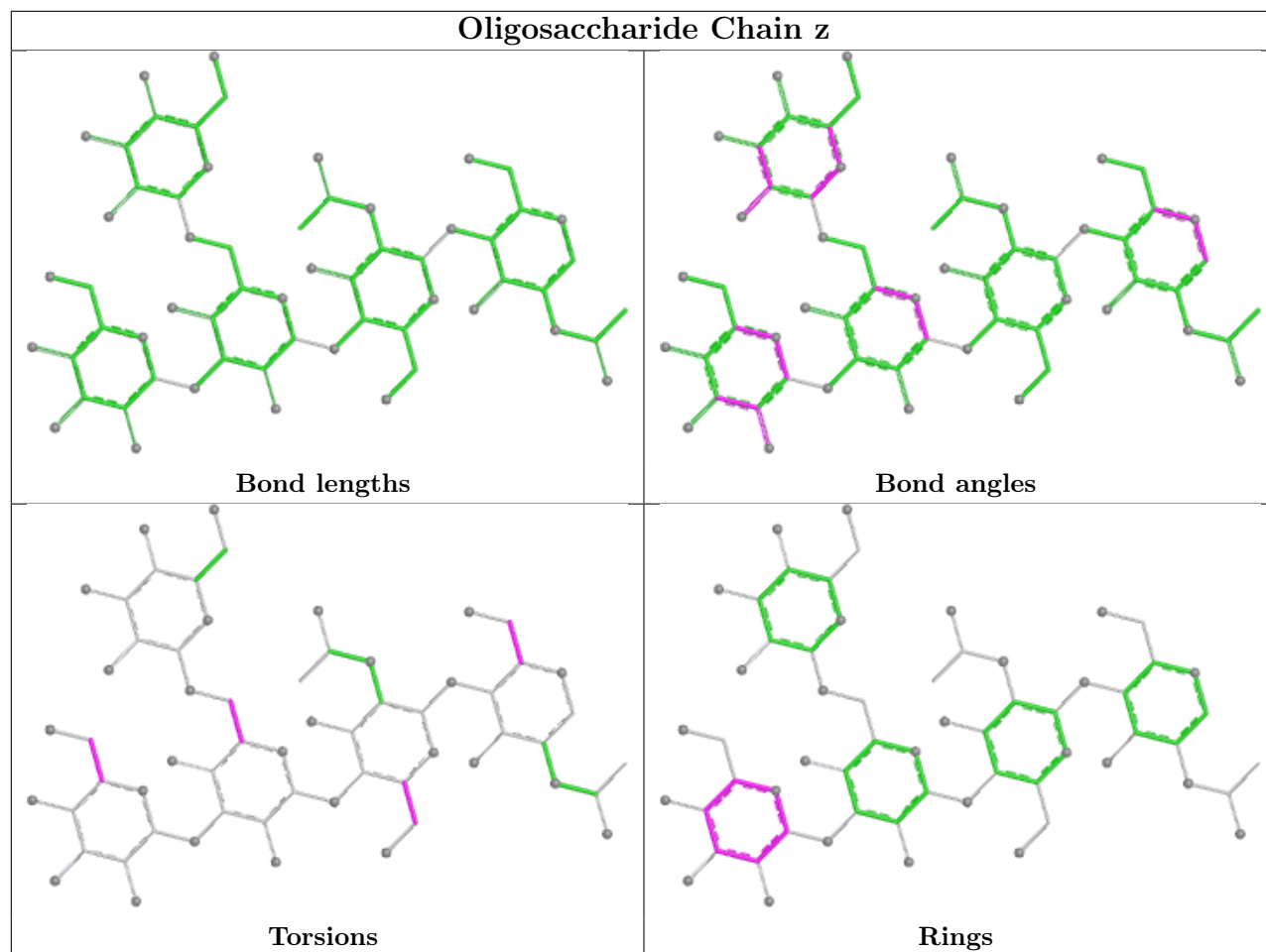


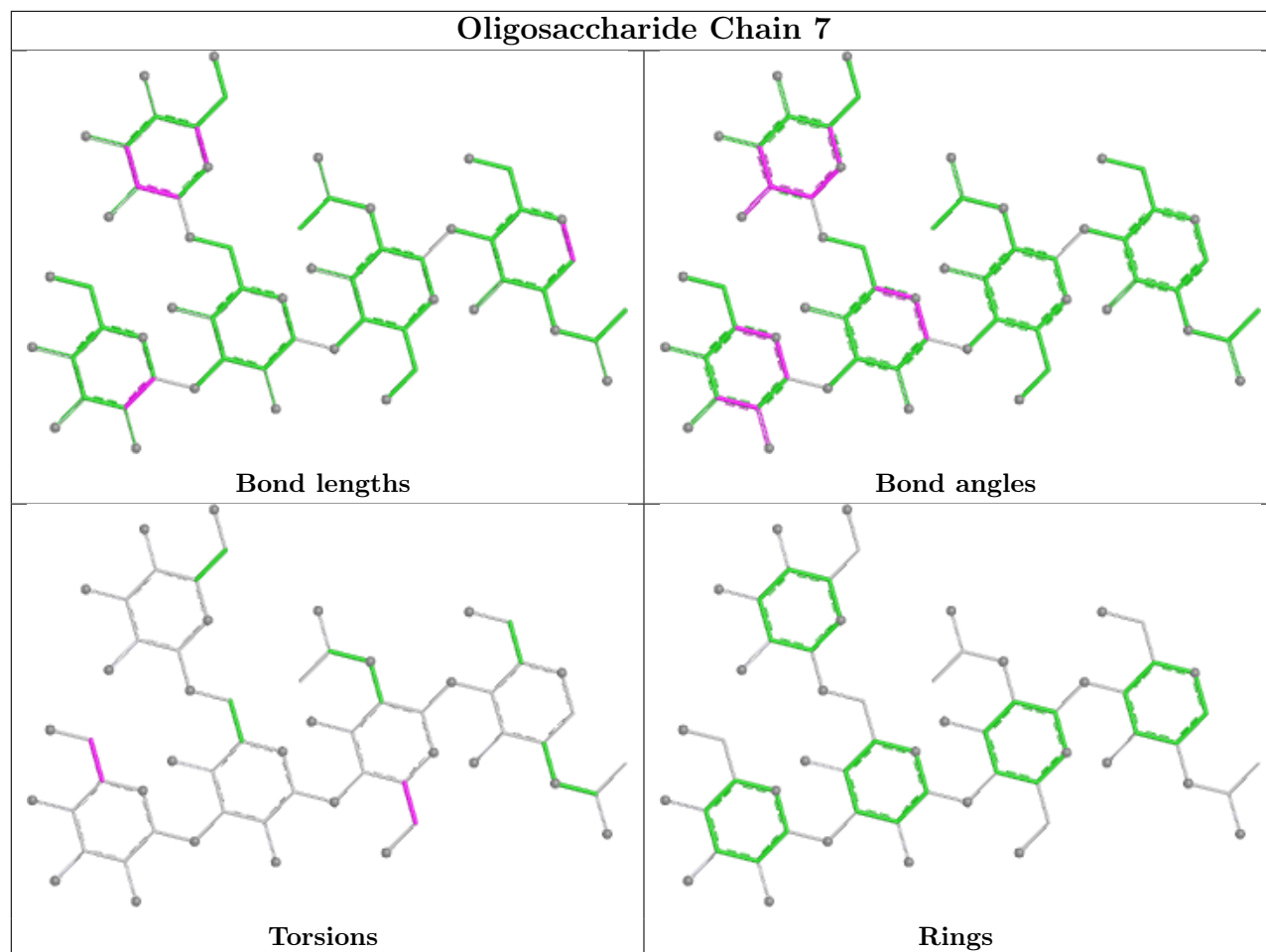




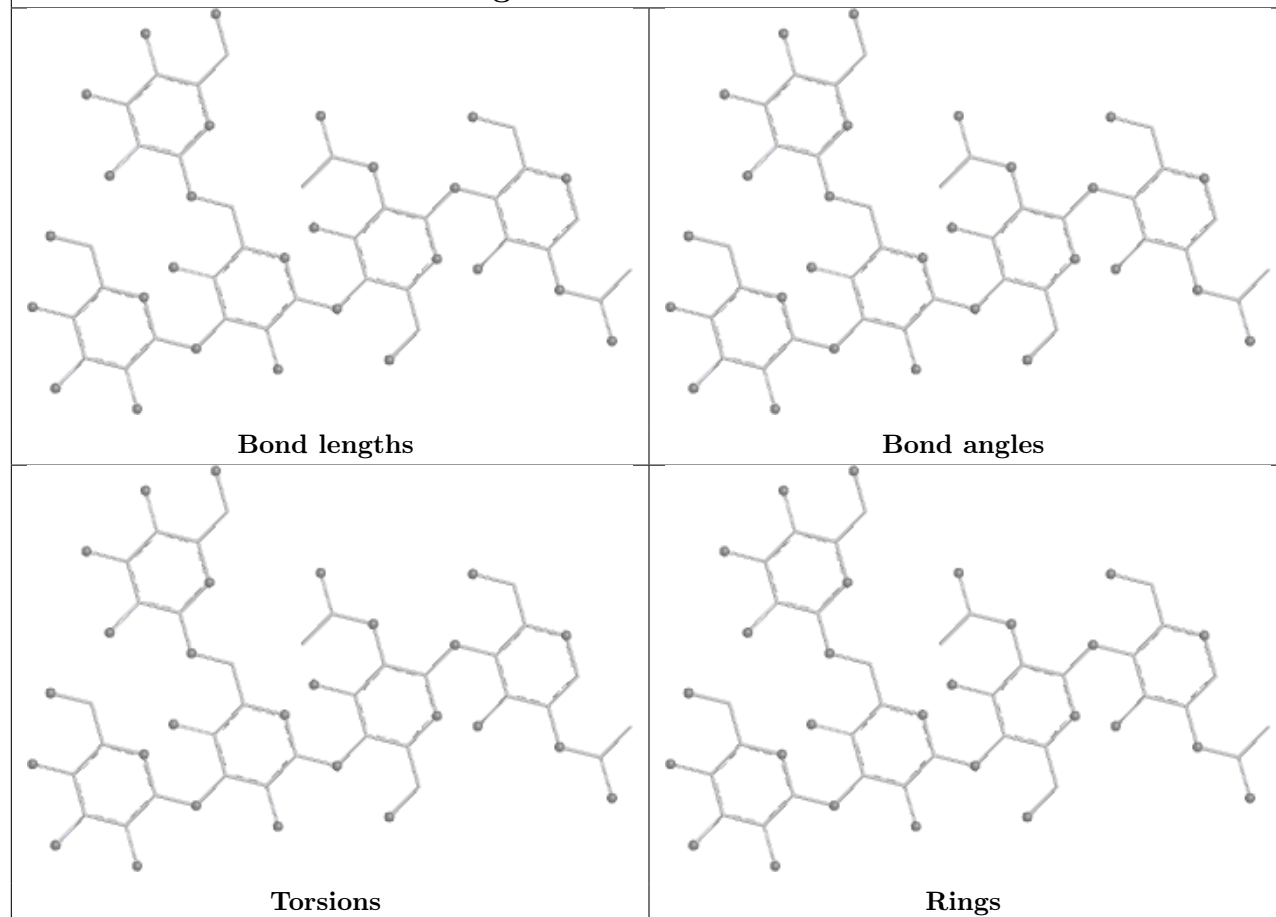




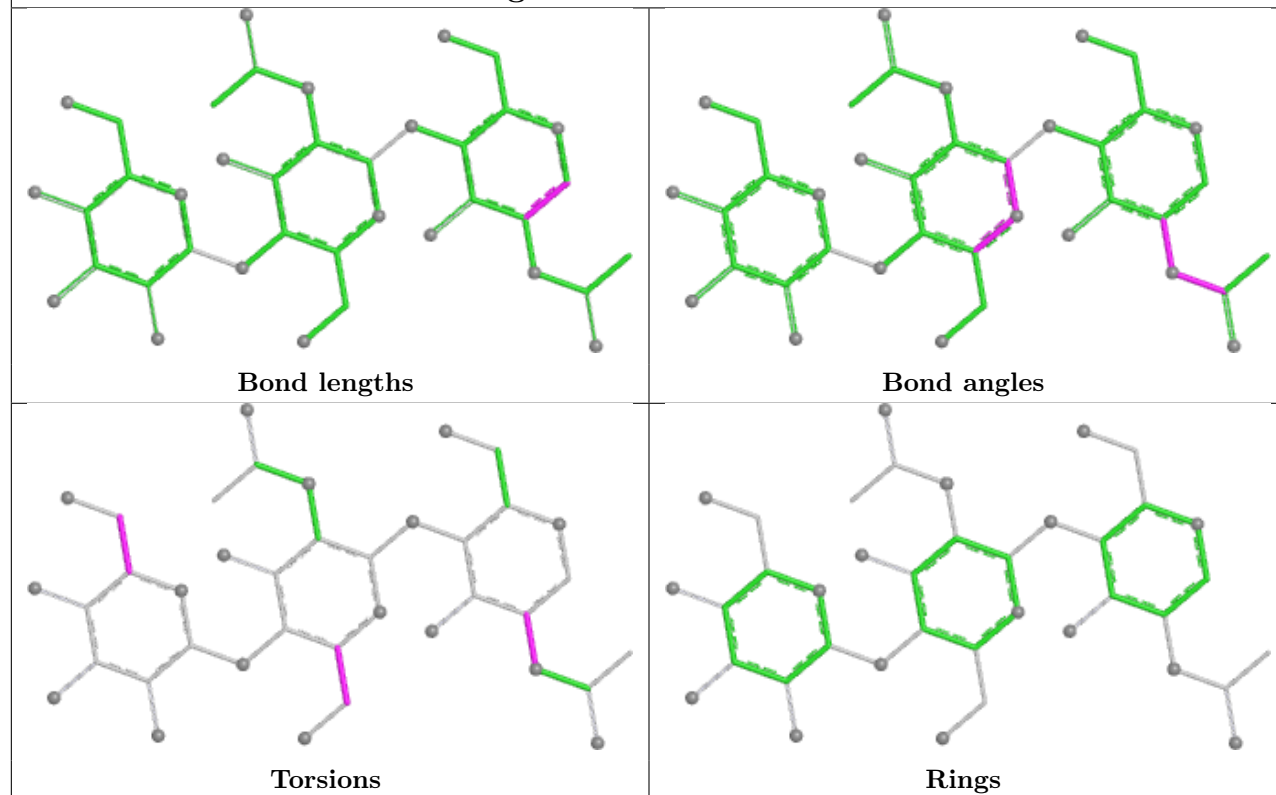


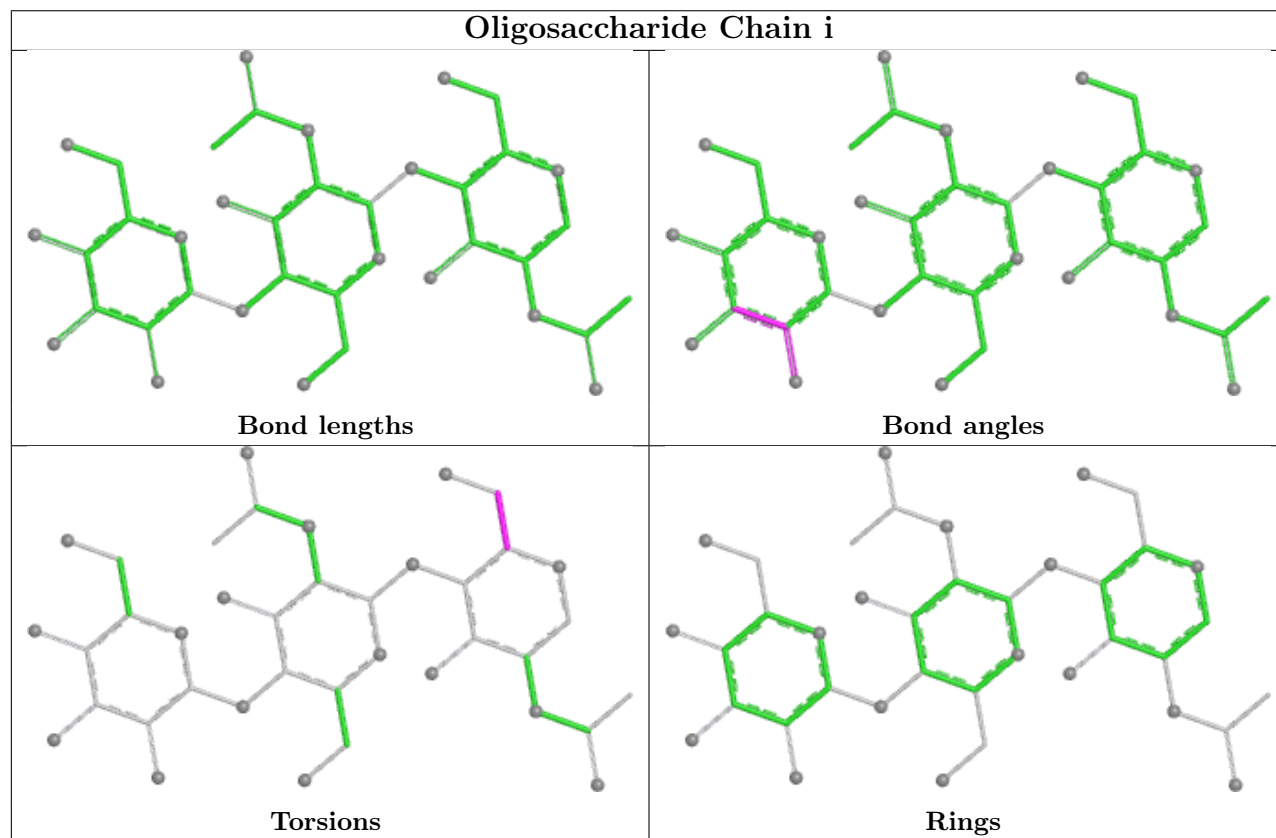
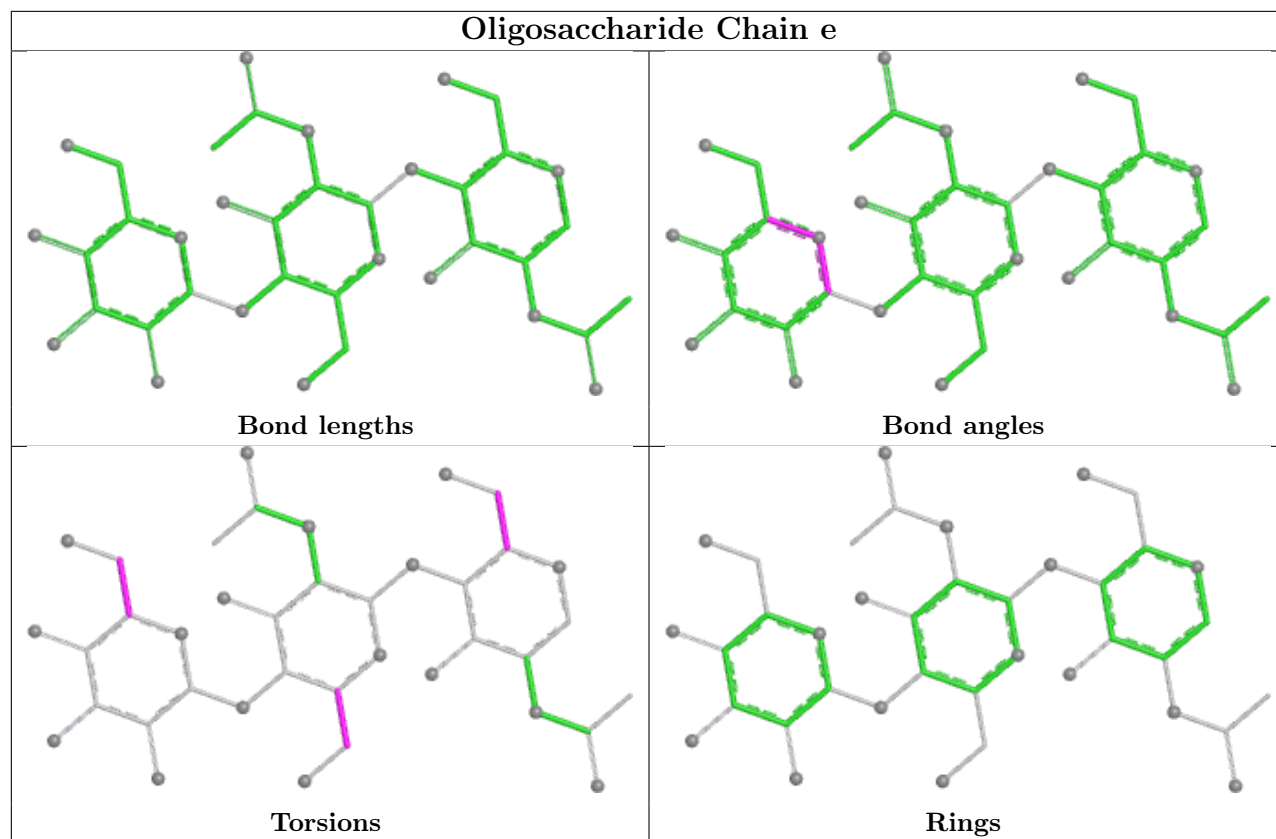


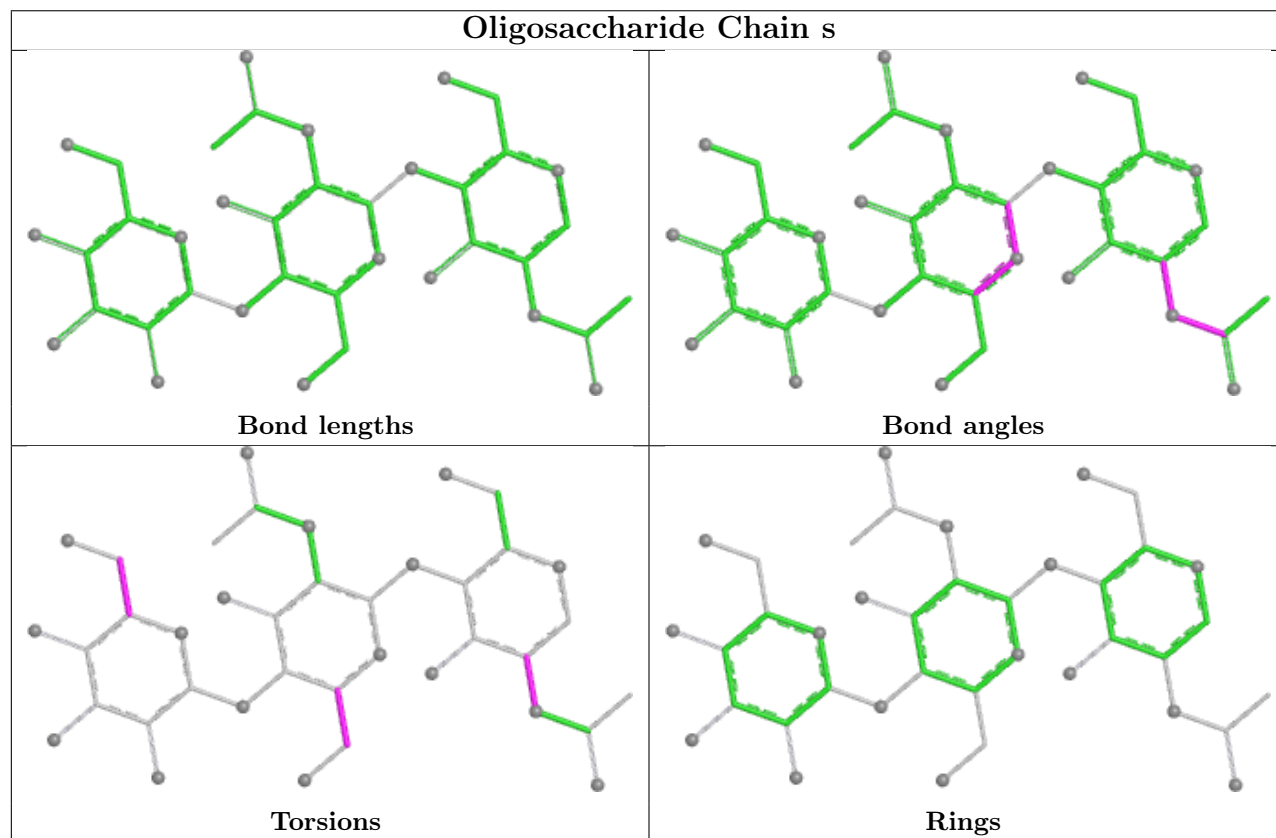
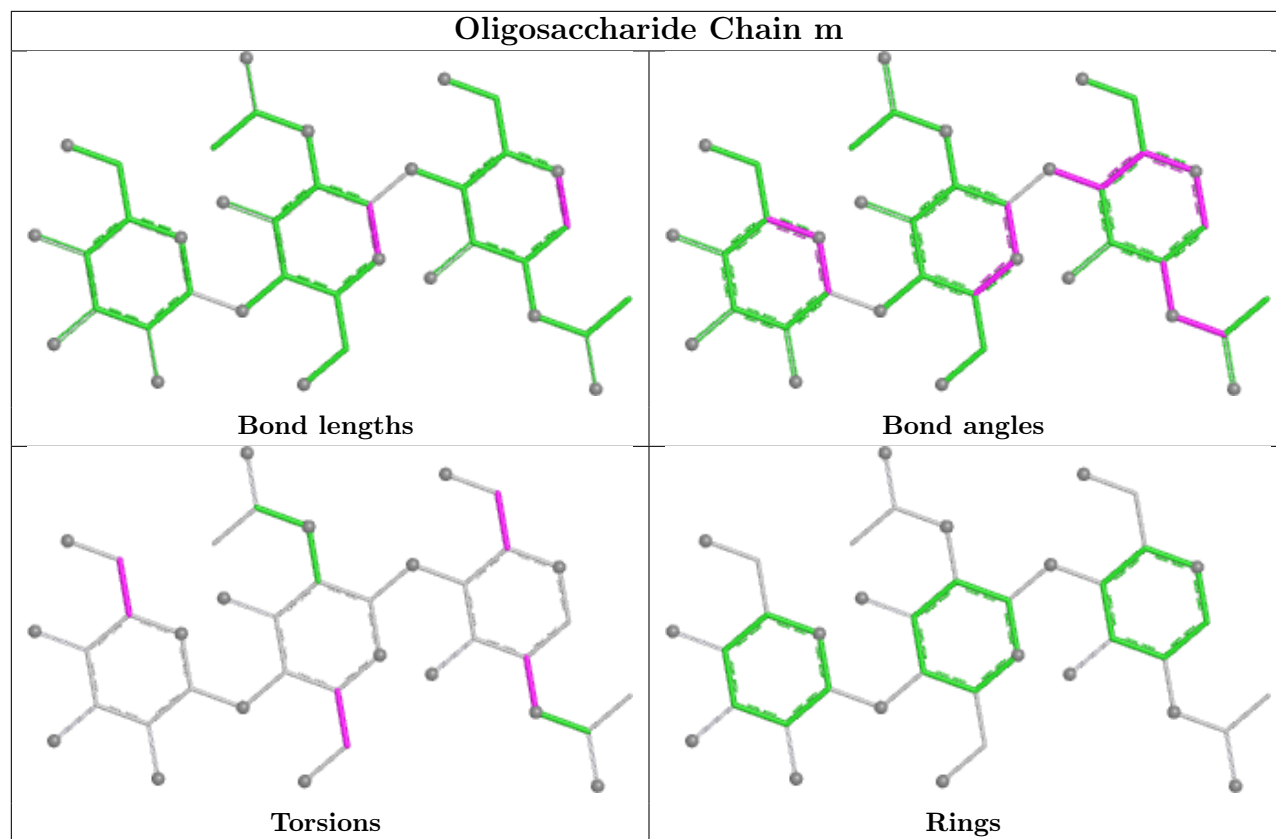
Oligosaccharide Chain FA

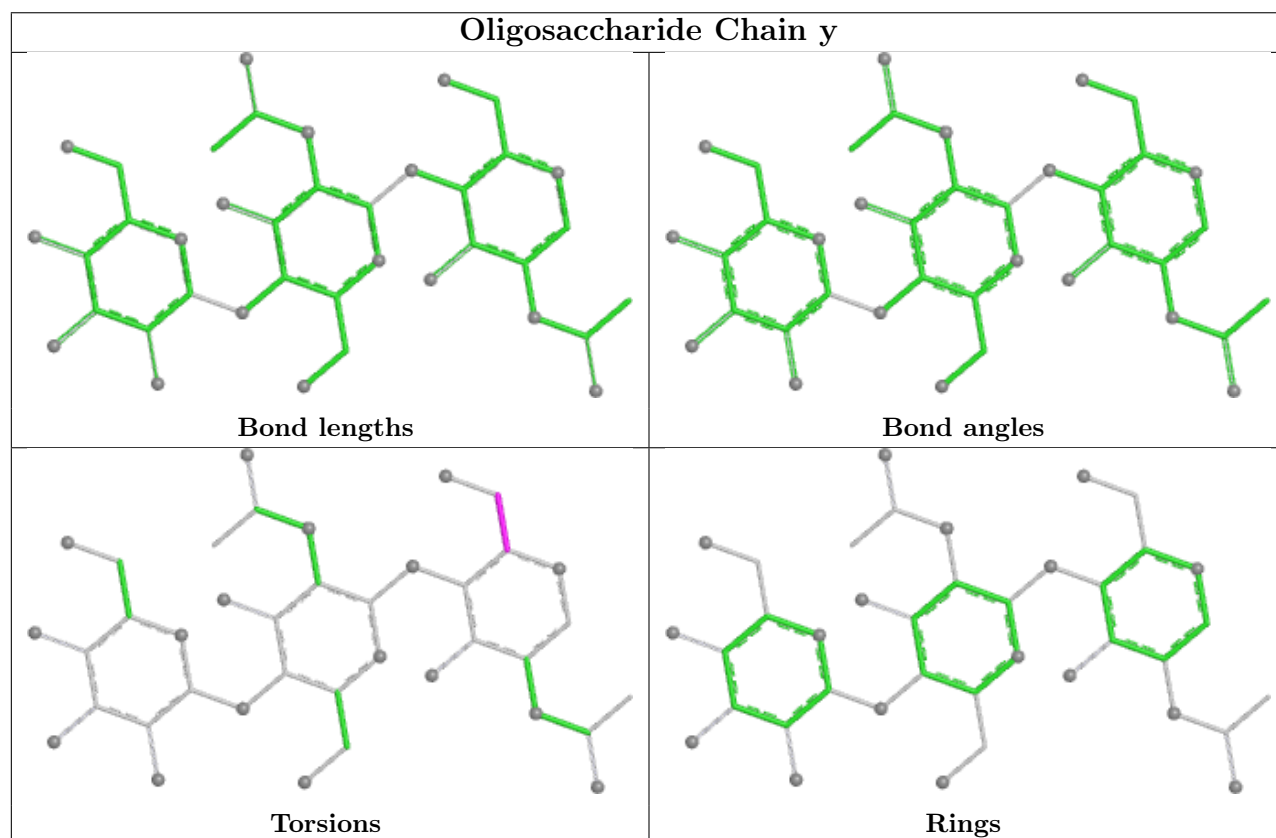
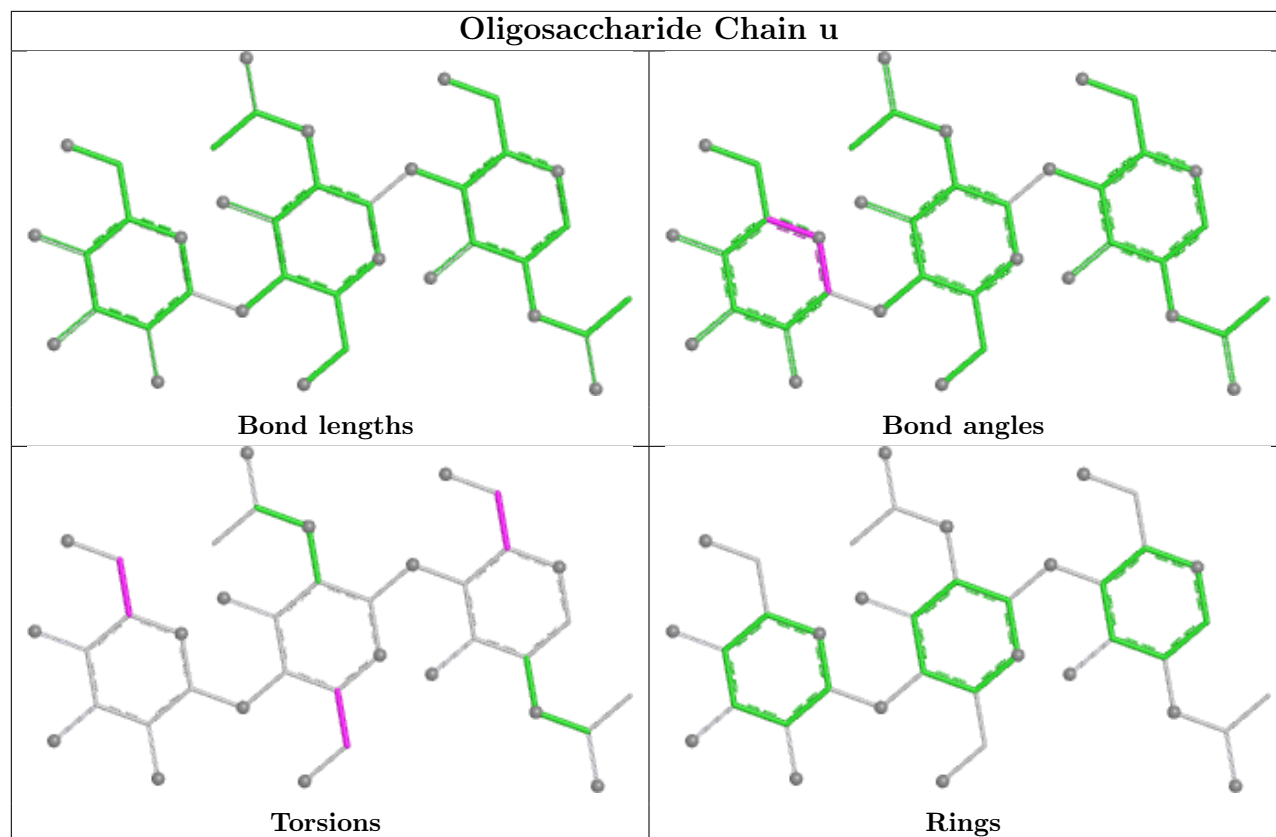


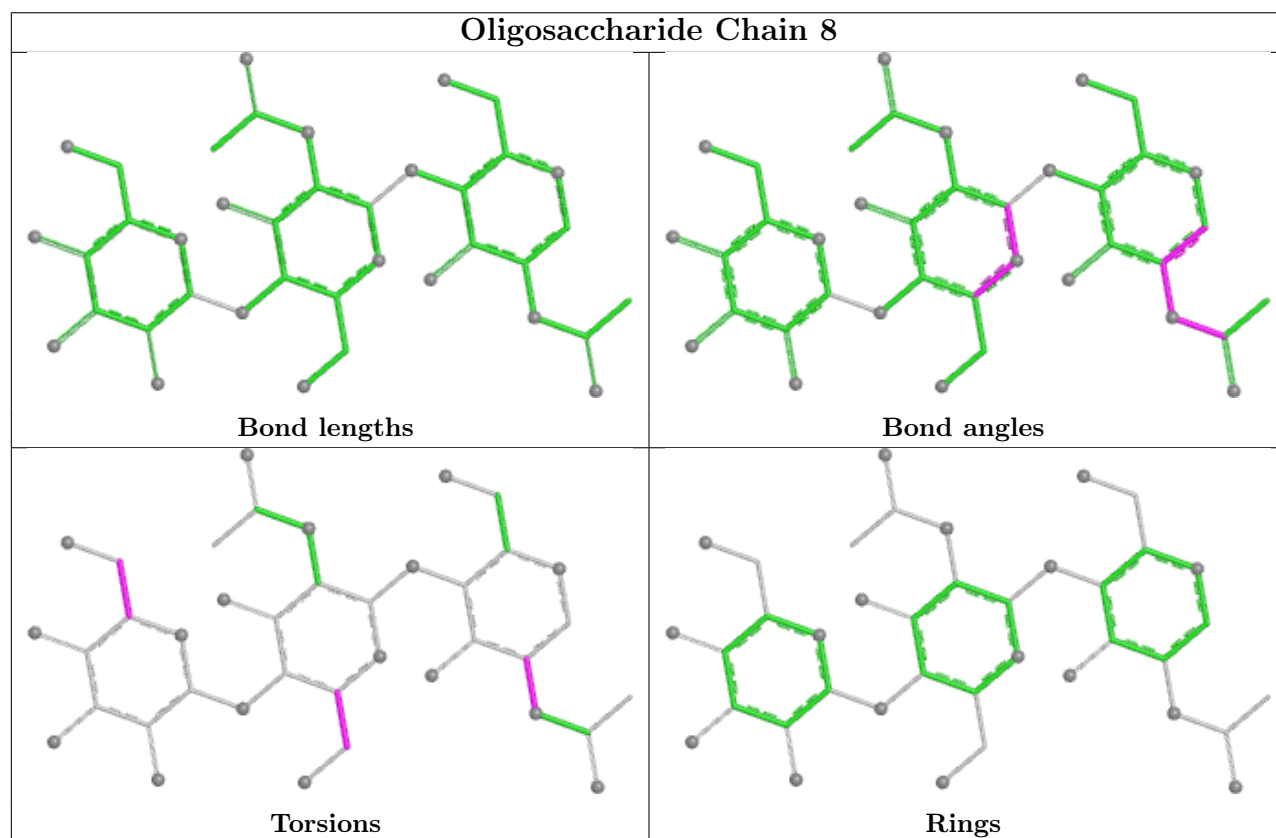
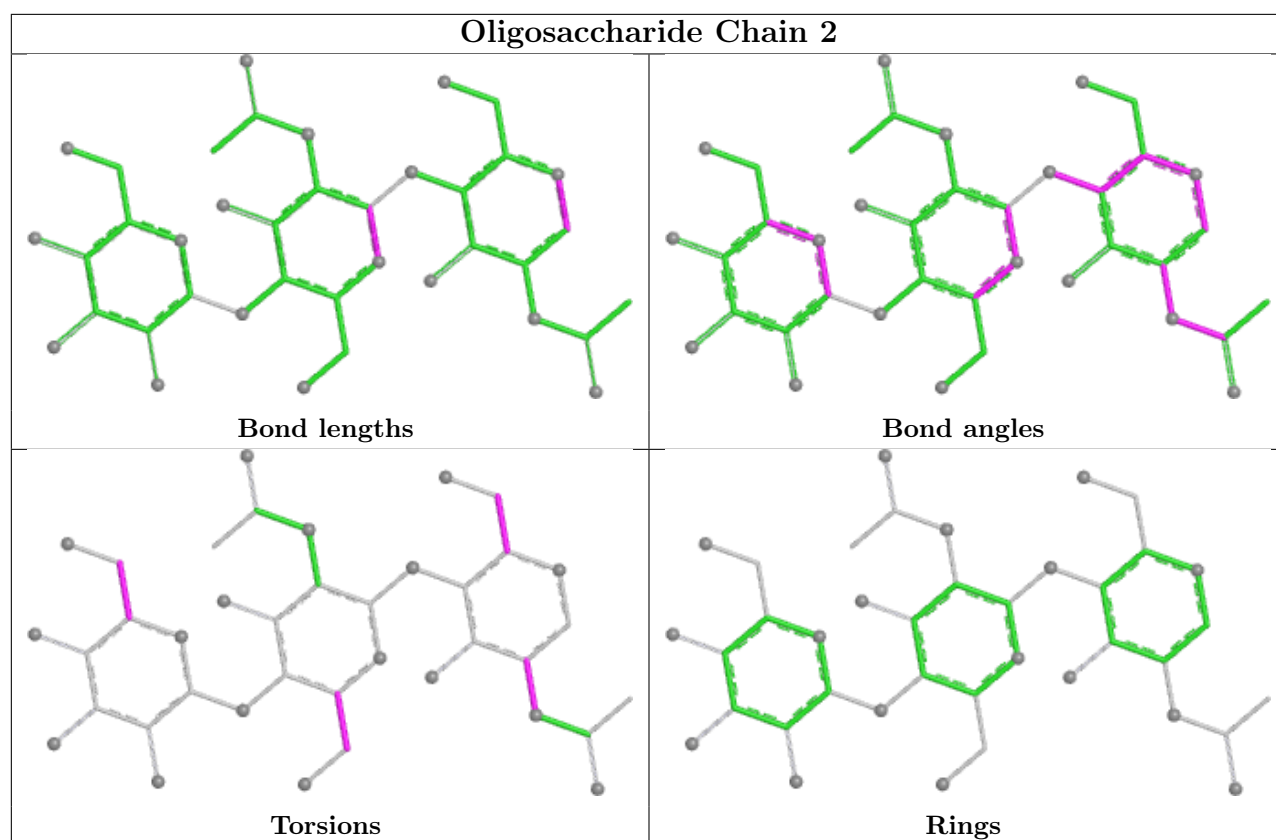
Oligosaccharide Chain c



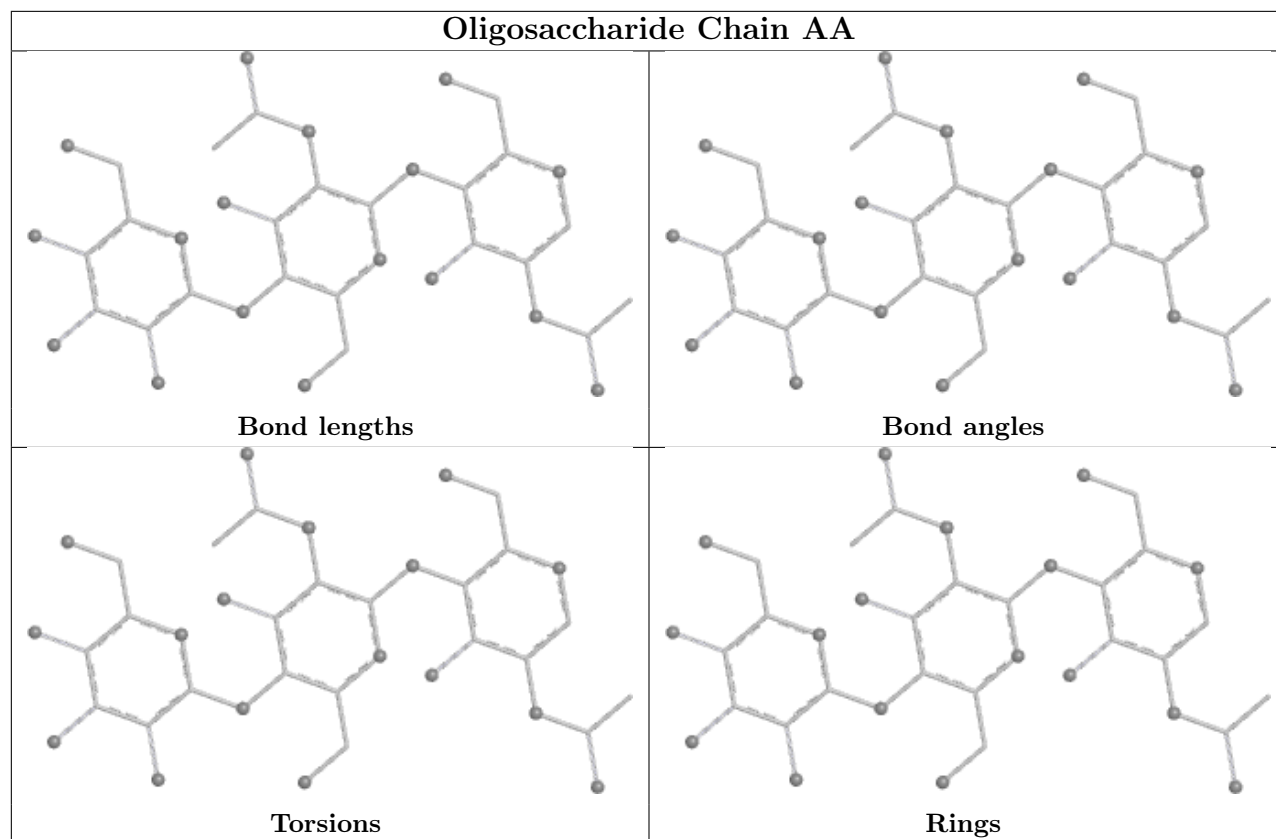




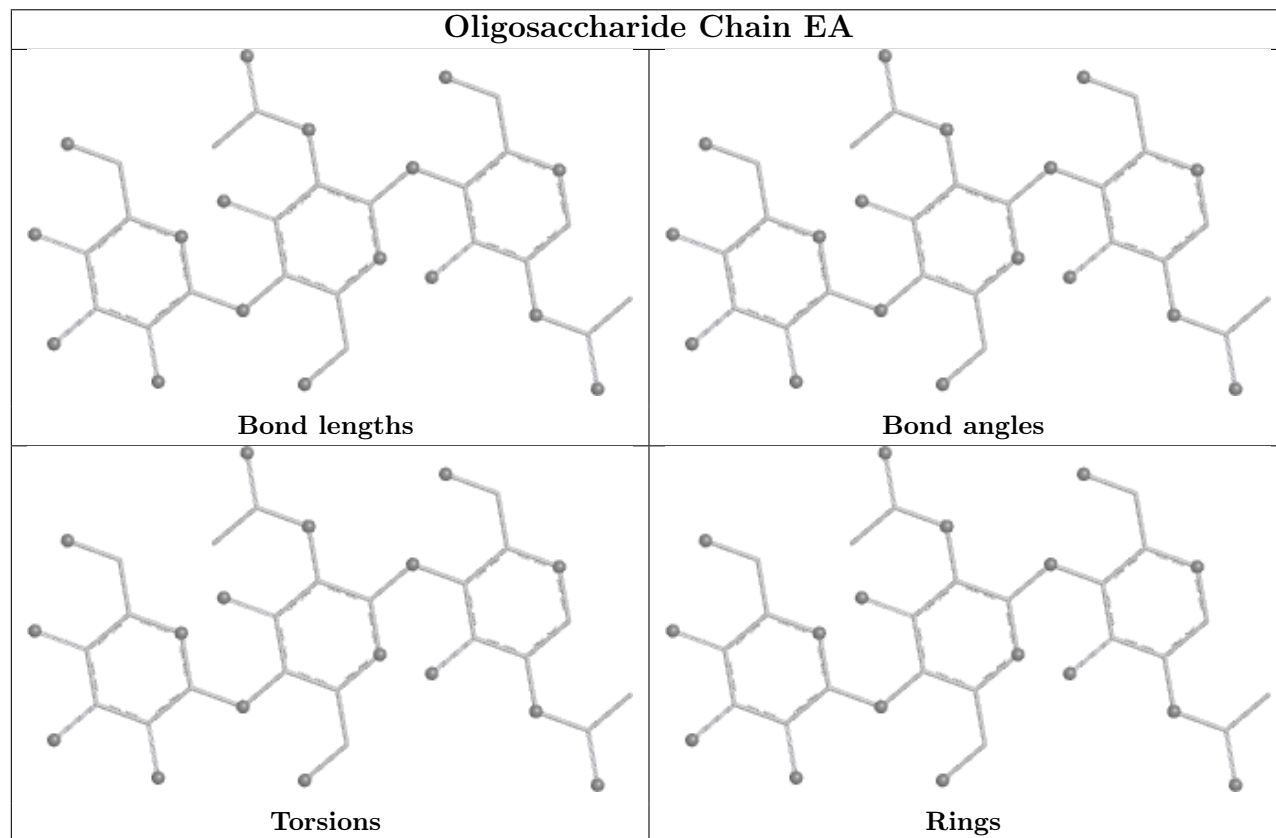


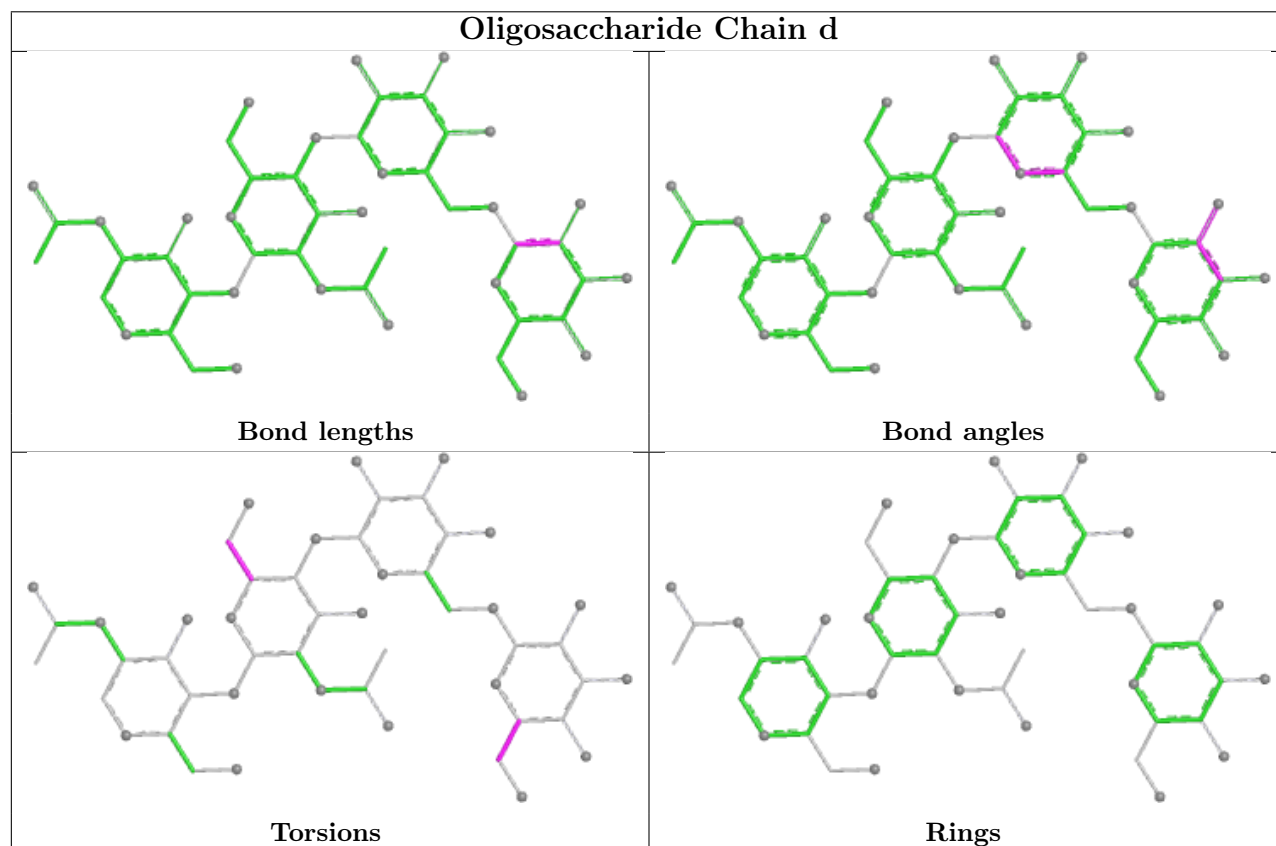
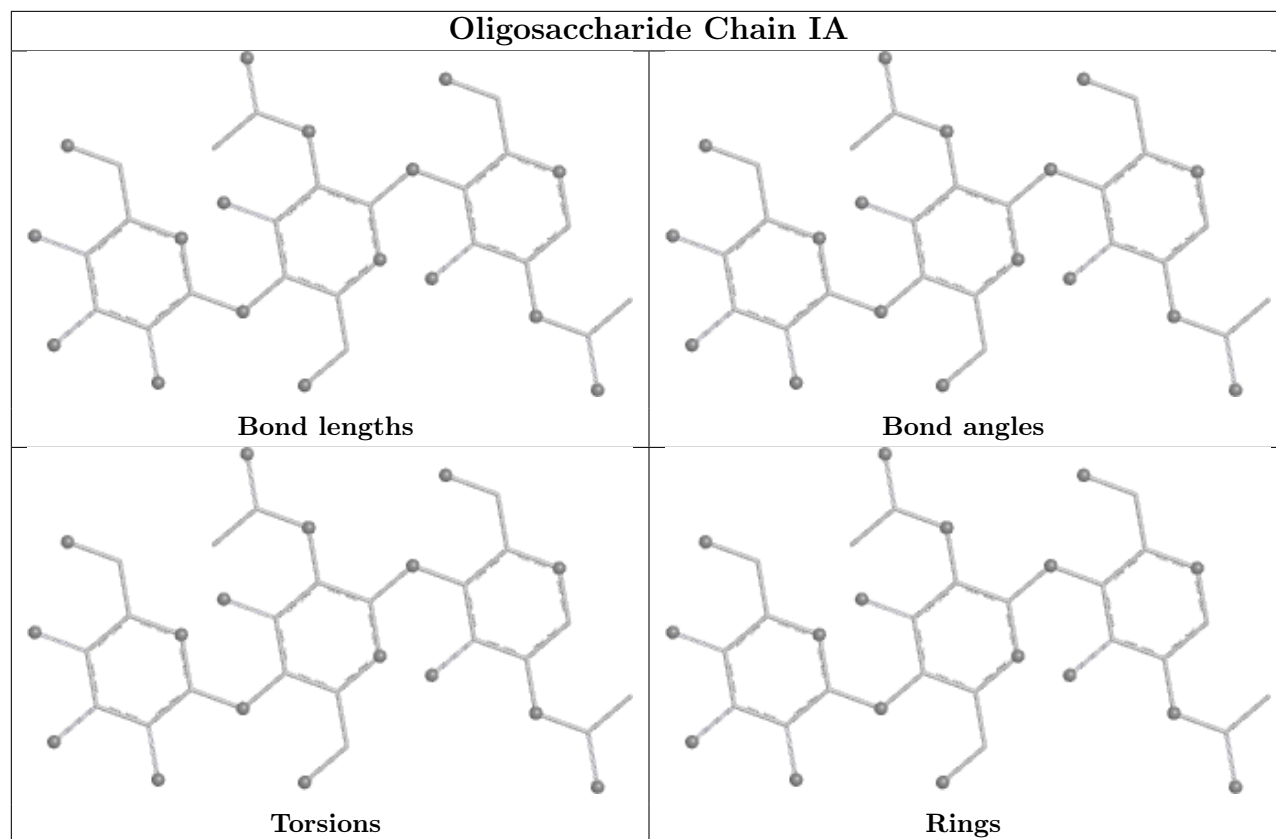


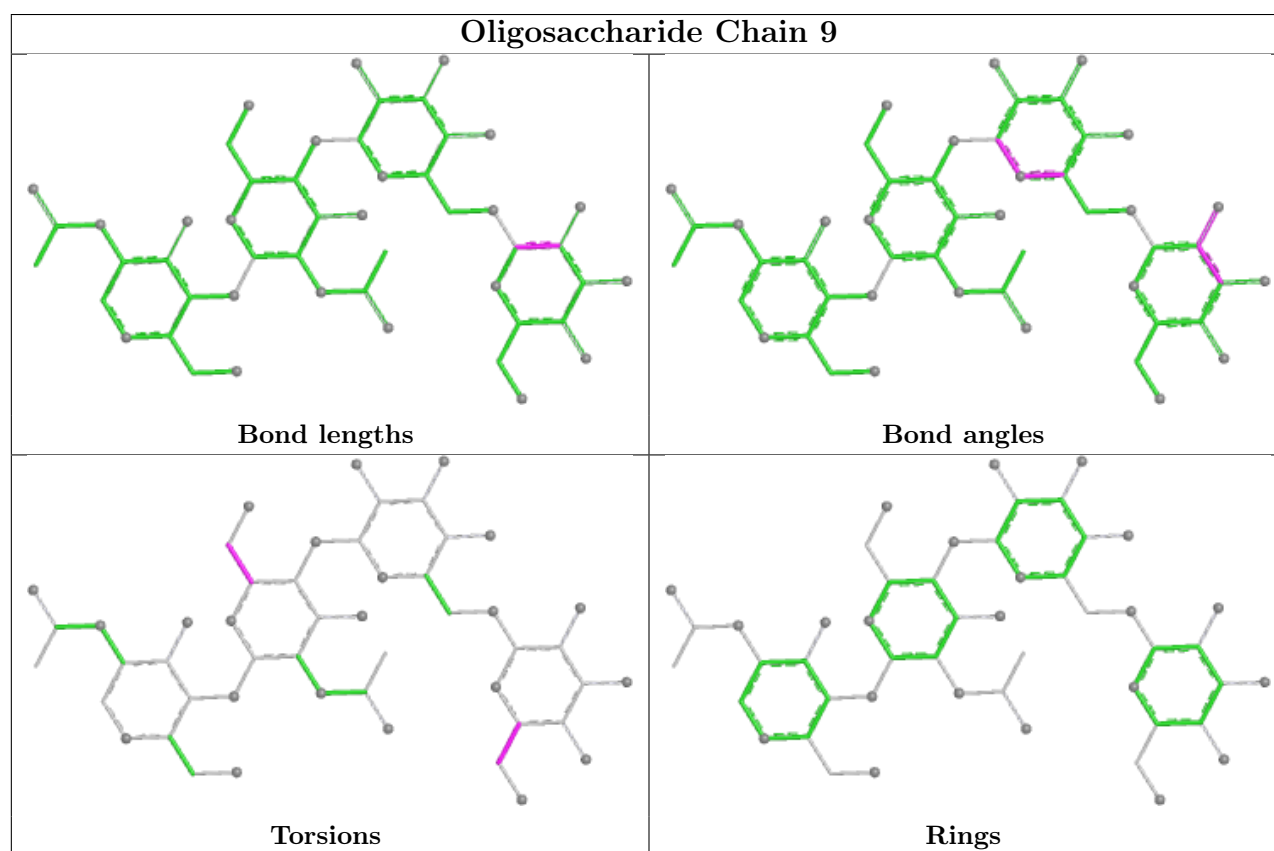
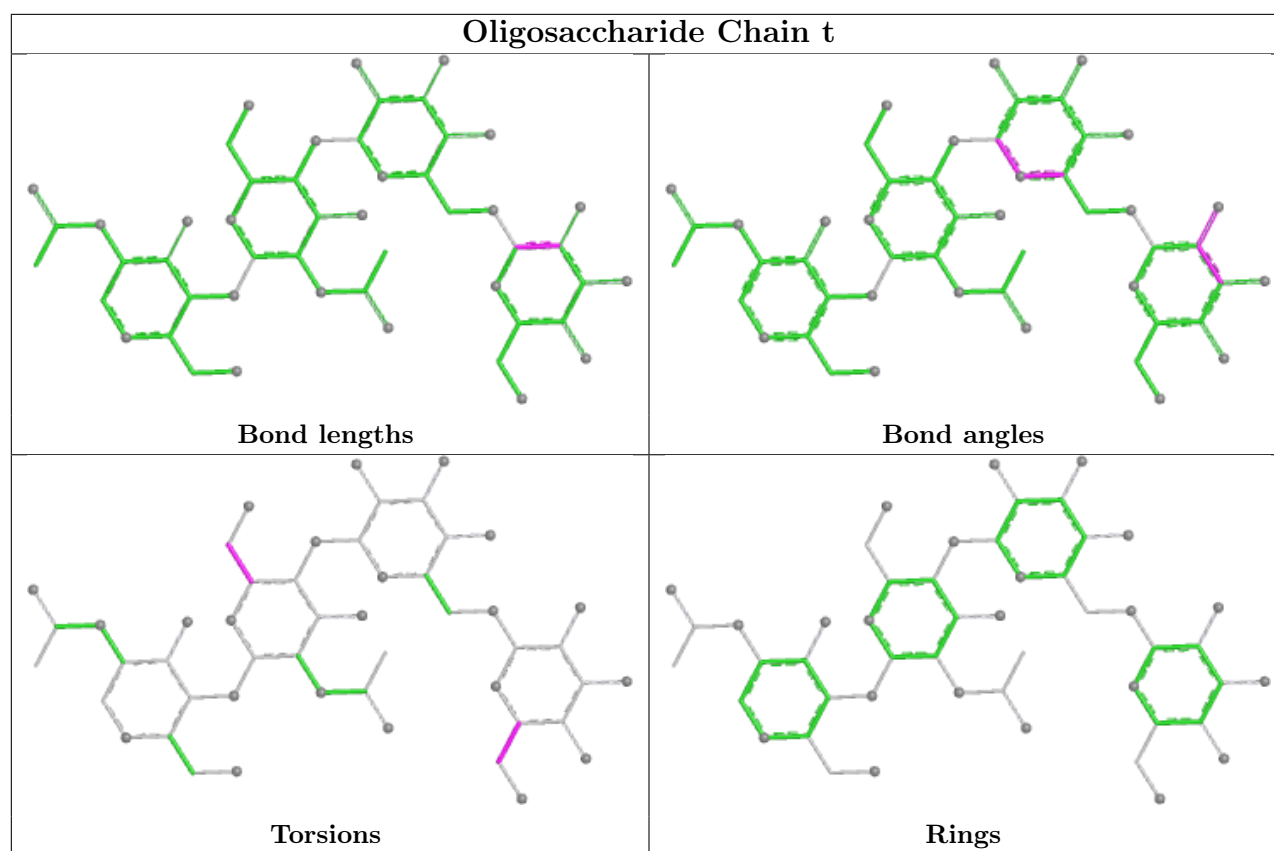
Oligosaccharide Chain AA

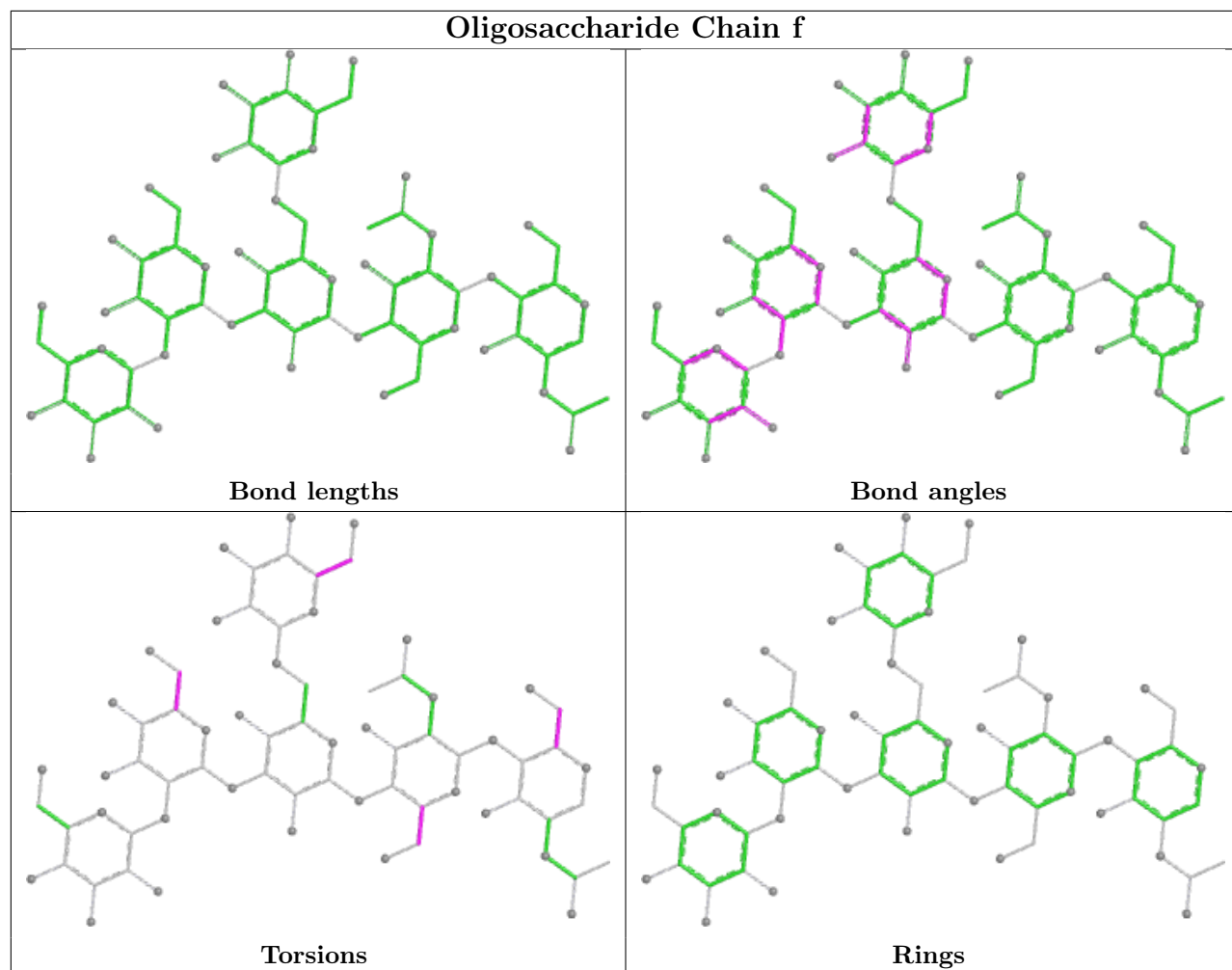


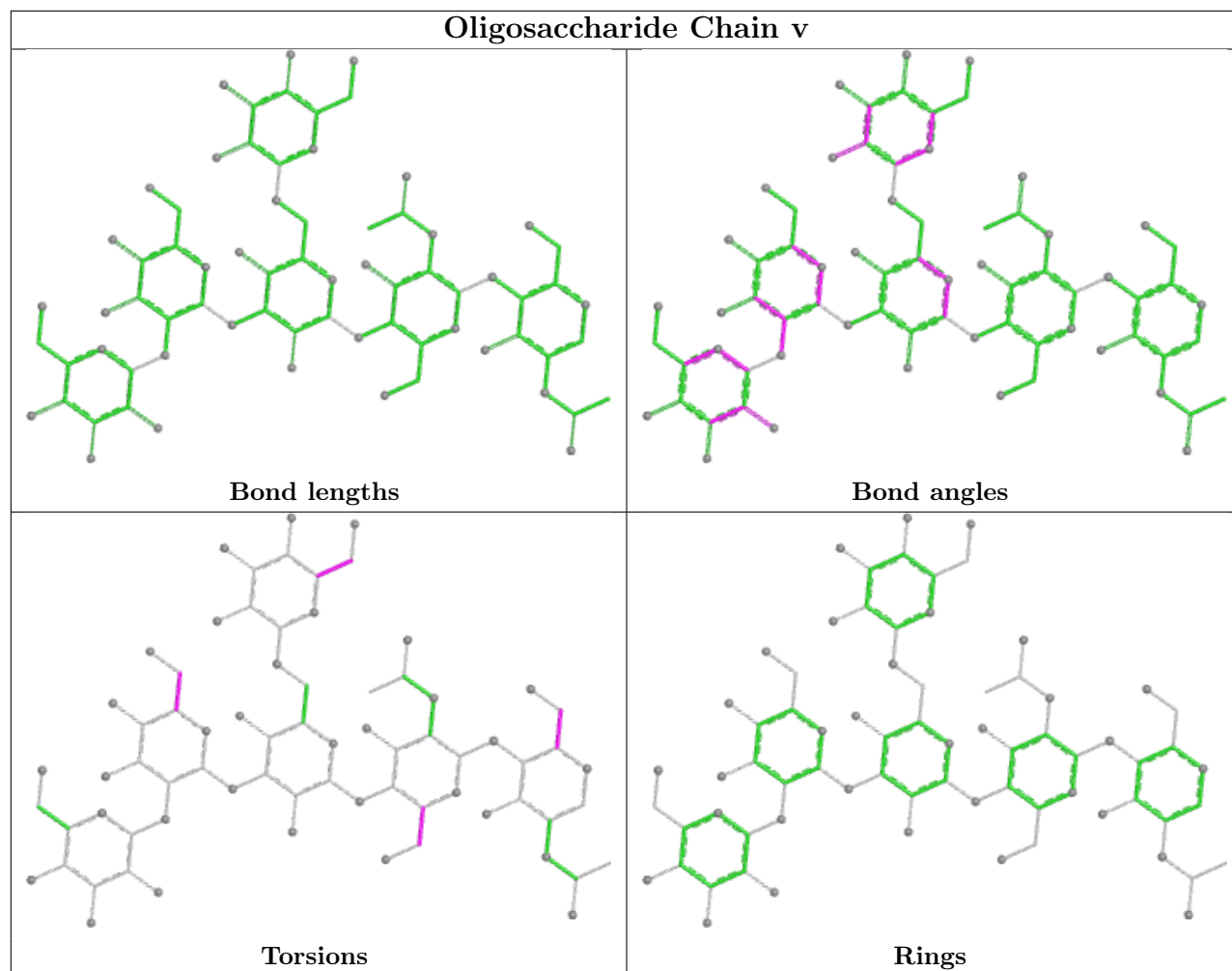
Oligosaccharide Chain EA

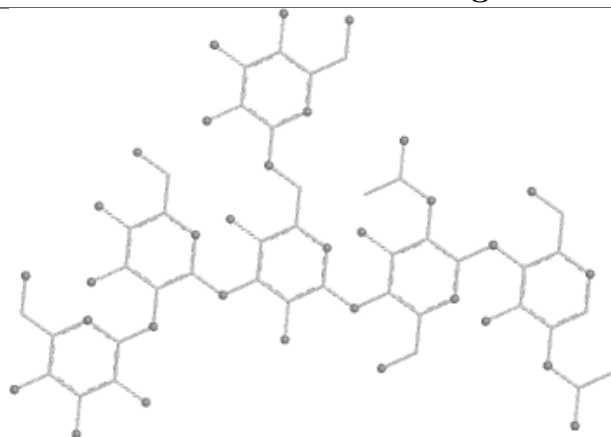
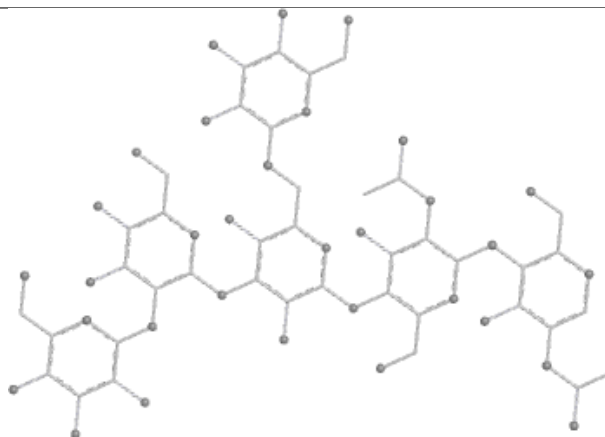
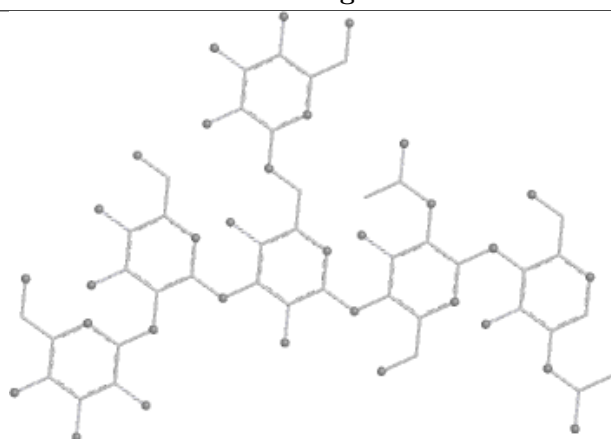
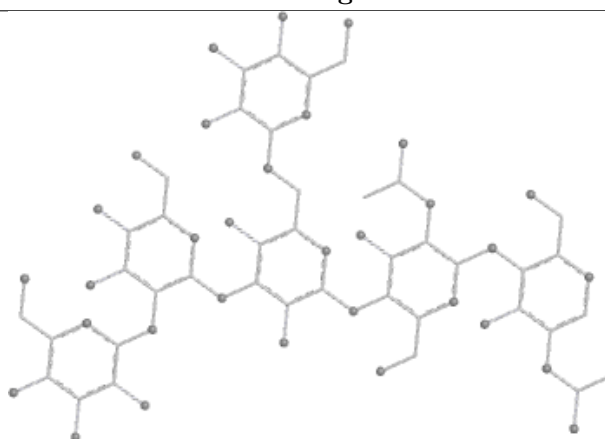


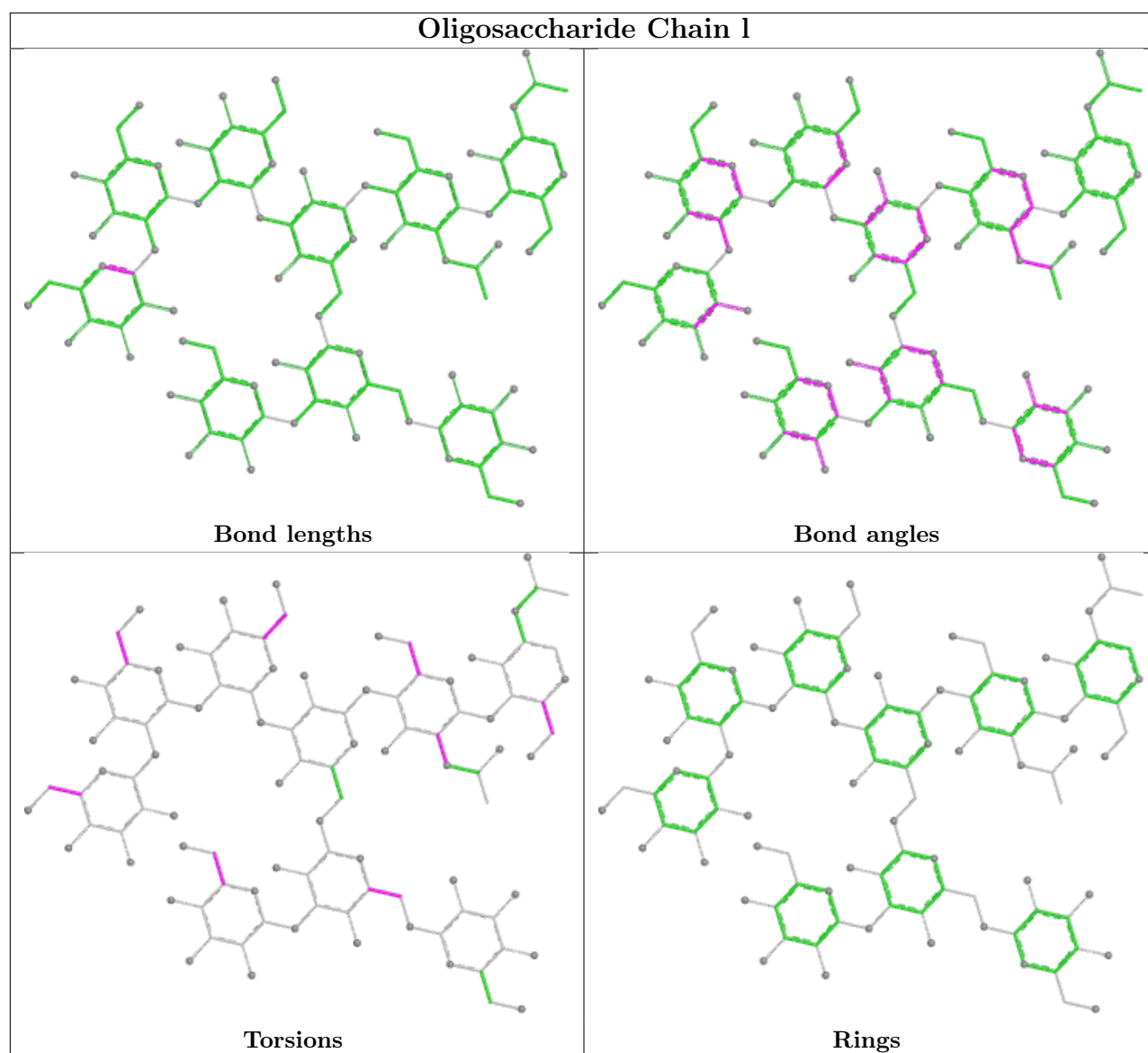


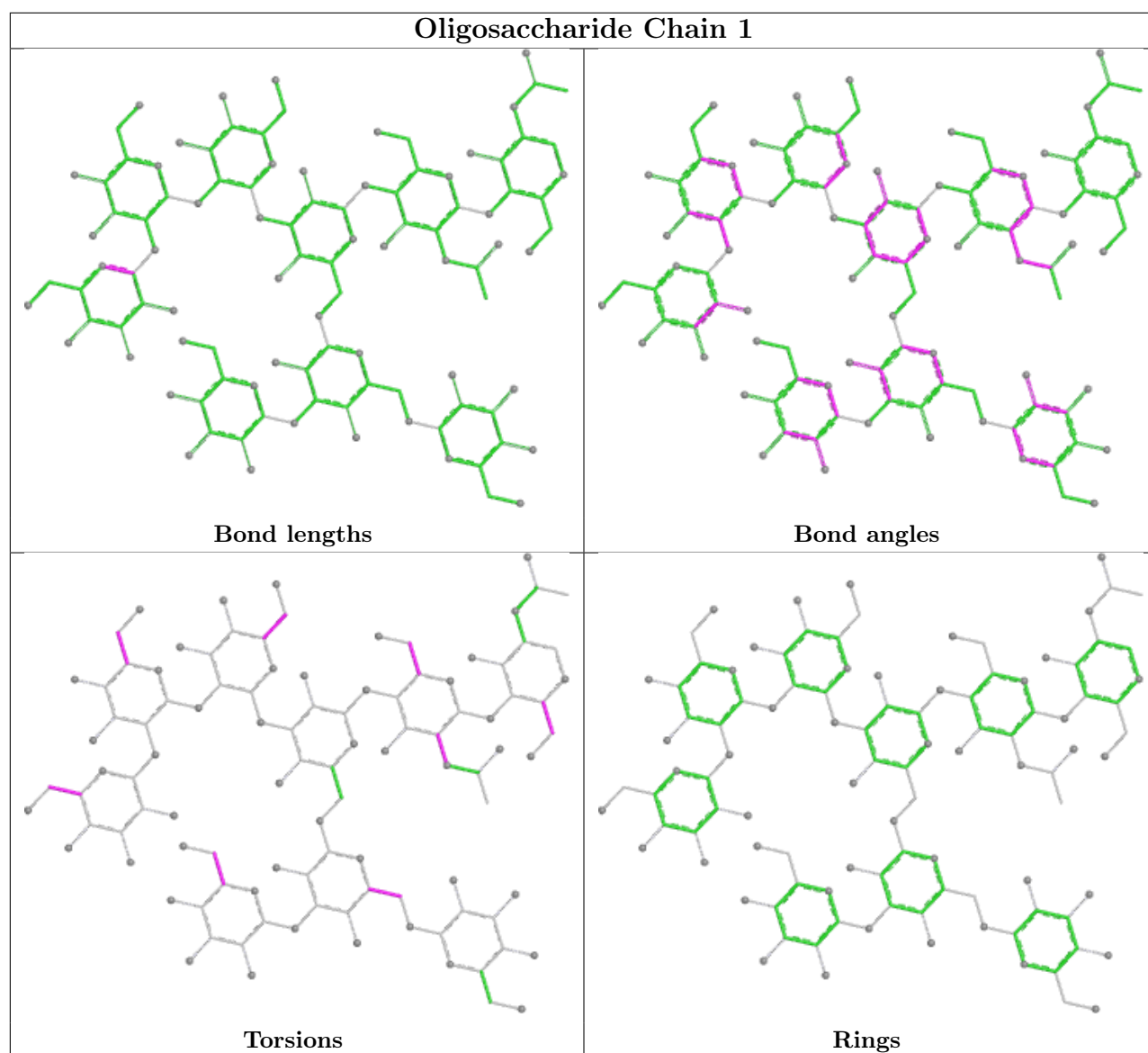


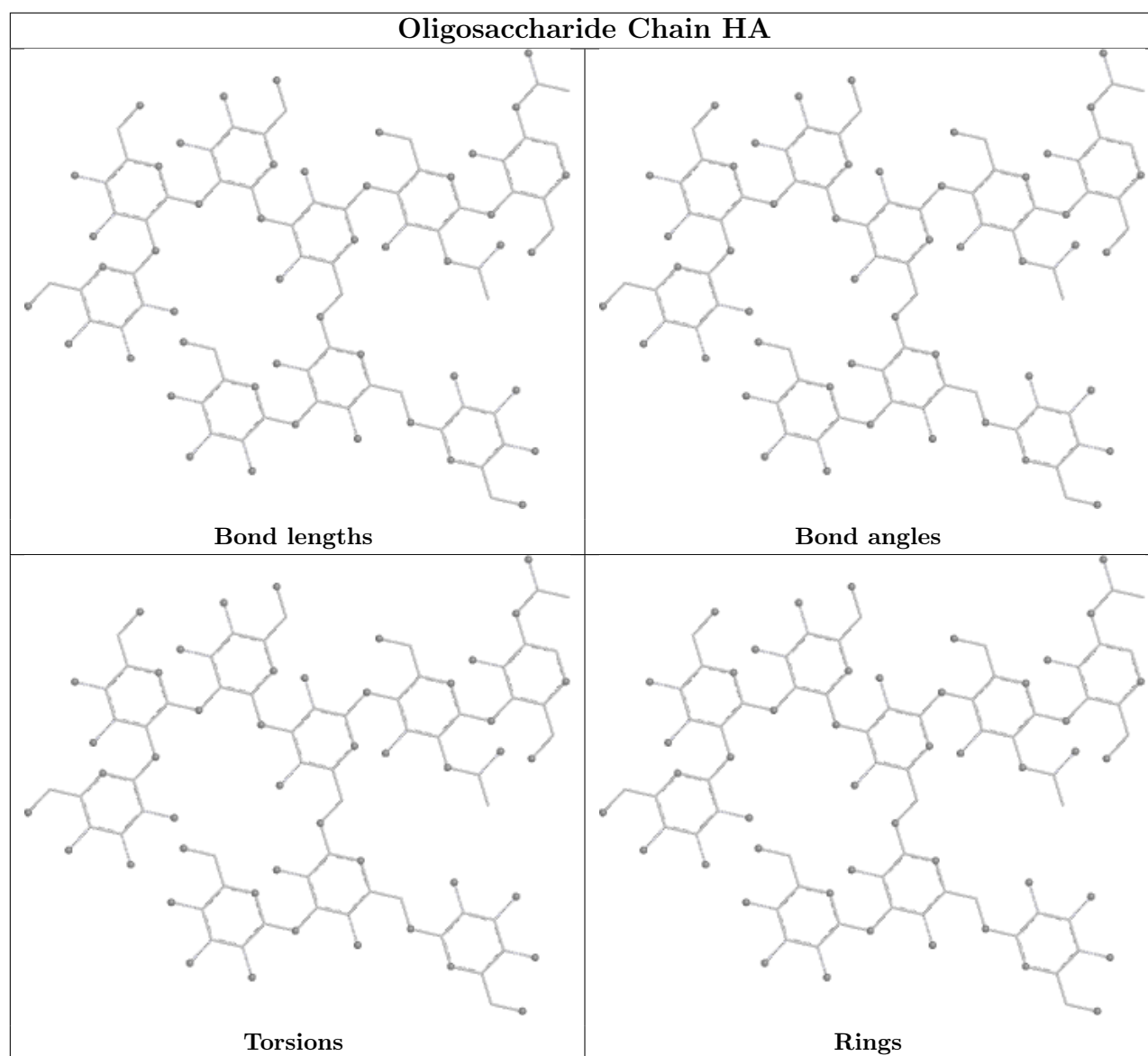




Oligosaccharide Chain BA**Bond lengths****Bond angles****Torsions****Rings**







5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
16	NAG	G	634	1	14,14,15	0.34	0	17,19,21	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
16	NAG	I	701	2	14,14,15	0.73	1 (7%)	17,19,21	0.90	1 (5%)
16	NAG	R	702	2	14,14,15	0.34	0	17,19,21	0.60	1 (5%)
16	NAG	R	701	2	14,14,15	0.74	1 (7%)	17,19,21	0.91	1 (5%)
16	NAG	B	702	2	14,14,15	0.33	0	17,19,21	0.58	0
16	NAG	I	702	2	14,14,15	0.34	0	17,19,21	0.59	0
16	NAG	Q	634	1	14,14,15	0.35	0	17,19,21	0.43	0
16	NAG	B	701	2	14,14,15	0.76	1 (7%)	17,19,21	0.91	1 (5%)
16	NAG	A	634	1	14,14,15	0.34	0	17,19,21	0.42	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
16	NAG	G	634	1	-	2/6/23/26	0/1/1/1
16	NAG	I	701	2	-	2/6/23/26	0/1/1/1
16	NAG	R	702	2	-	3/6/23/26	0/1/1/1
16	NAG	R	701	2	-	2/6/23/26	0/1/1/1
16	NAG	B	702	2	-	3/6/23/26	0/1/1/1
16	NAG	I	702	2	-	3/6/23/26	0/1/1/1
16	NAG	Q	634	1	-	2/6/23/26	0/1/1/1
16	NAG	B	701	2	-	2/6/23/26	0/1/1/1
16	NAG	A	634	1	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	B	701	NAG	C1-C2	2.50	1.55	1.52
16	R	701	NAG	C1-C2	2.40	1.55	1.52
16	I	701	NAG	C1-C2	2.33	1.55	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	R	701	NAG	C1-O5-C5	3.35	116.67	112.19
16	B	701	NAG	C1-O5-C5	3.34	116.66	112.19
16	I	701	NAG	C1-O5-C5	3.29	116.60	112.19
16	R	702	NAG	C1-O5-C5	2.04	114.93	112.19

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
16	A	634	NAG	O5-C5-C6-O6
16	G	634	NAG	O5-C5-C6-O6
16	Q	634	NAG	O5-C5-C6-O6
16	B	701	NAG	O5-C5-C6-O6
16	I	701	NAG	O5-C5-C6-O6
16	R	701	NAG	O5-C5-C6-O6
16	B	701	NAG	C4-C5-C6-O6
16	R	701	NAG	C4-C5-C6-O6
16	I	701	NAG	C4-C5-C6-O6
16	B	702	NAG	O5-C5-C6-O6
16	I	702	NAG	O5-C5-C6-O6
16	R	702	NAG	O5-C5-C6-O6
16	B	702	NAG	C4-C5-C6-O6
16	R	702	NAG	C4-C5-C6-O6
16	I	702	NAG	C4-C5-C6-O6
16	A	634	NAG	C4-C5-C6-O6
16	G	634	NAG	C4-C5-C6-O6
16	Q	634	NAG	C4-C5-C6-O6
16	B	702	NAG	C1-C2-N2-C7
16	I	702	NAG	C1-C2-N2-C7
16	R	702	NAG	C1-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

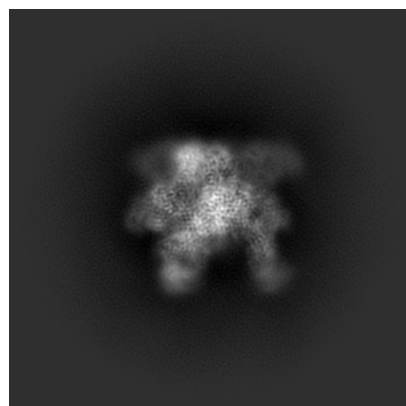
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9359. These allow visual inspection of the internal detail of the map and identification of artifacts.

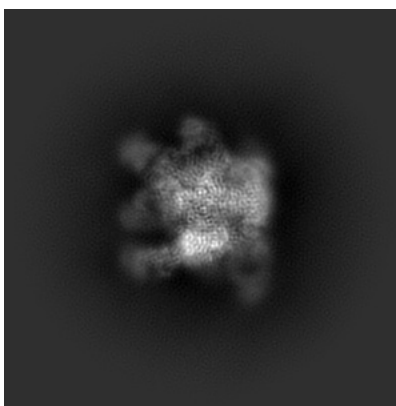
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

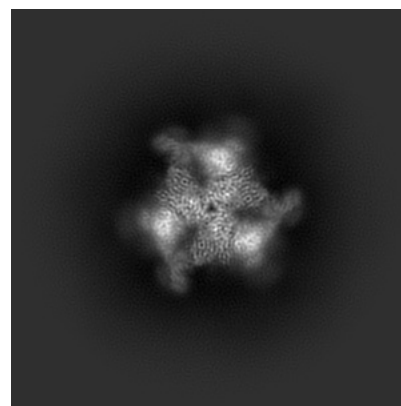
6.1.1 Primary map



X

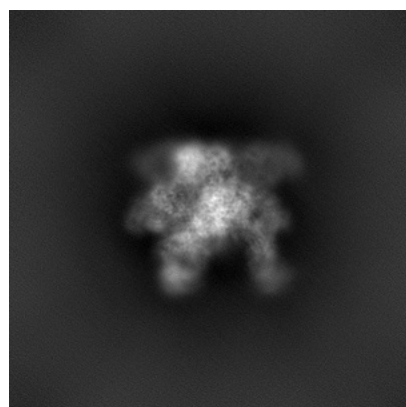


Y

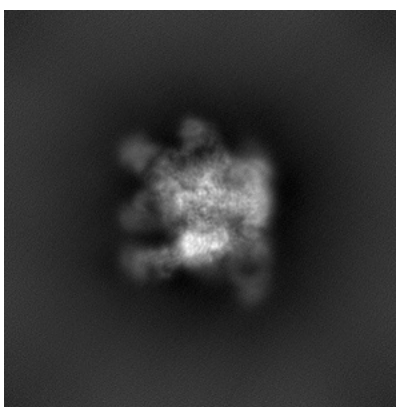


Z

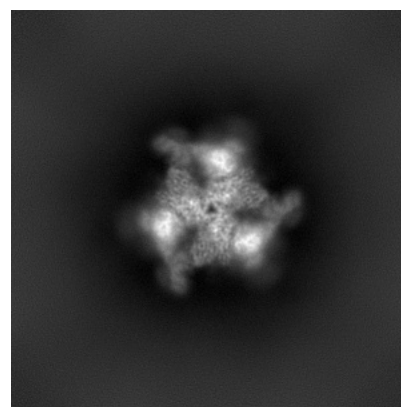
6.1.2 Raw map



X



Y

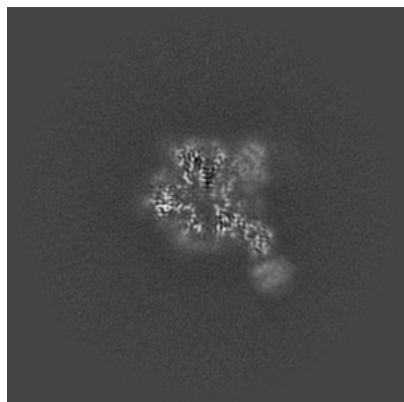


Z

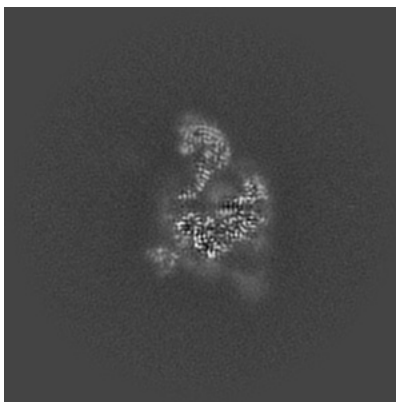
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

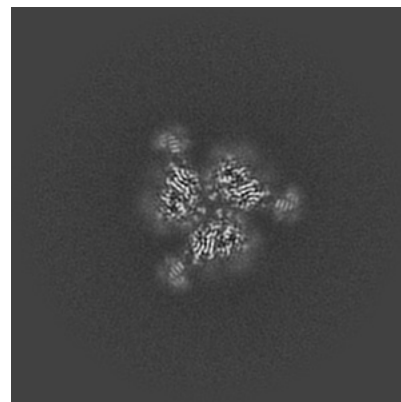
6.2.1 Primary map



X Index: 200

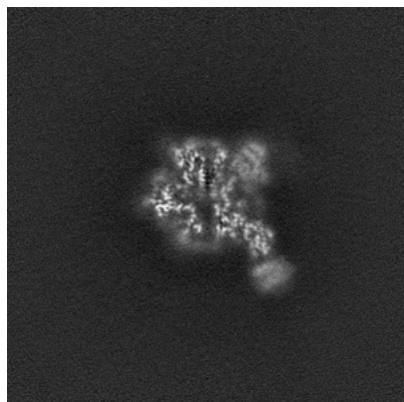


Y Index: 200

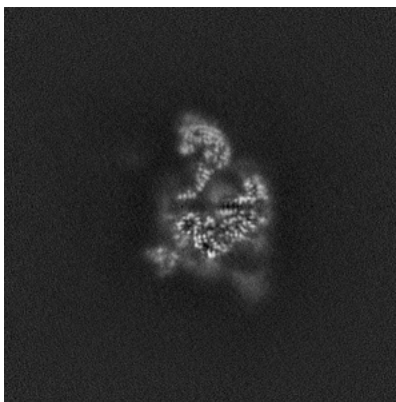


Z Index: 200

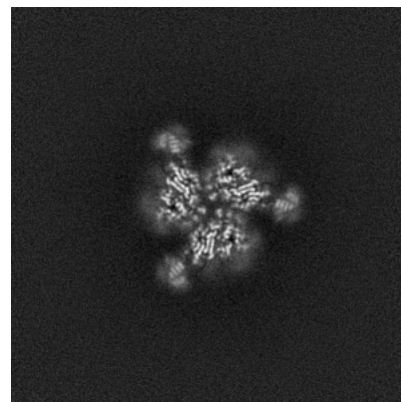
6.2.2 Raw map



X Index: 200



Y Index: 200

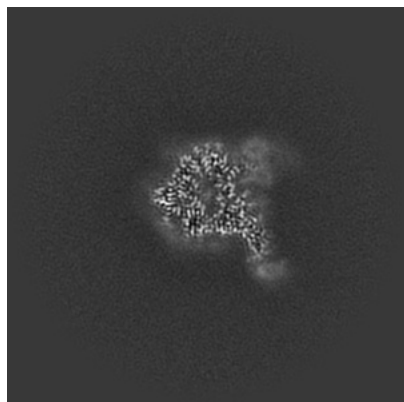


Z Index: 200

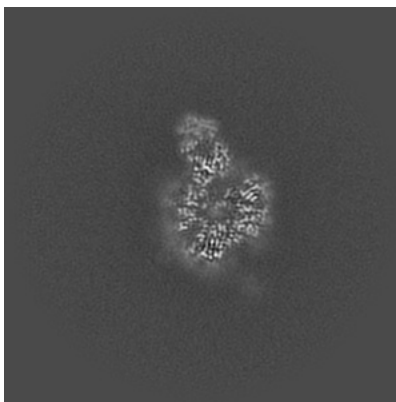
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

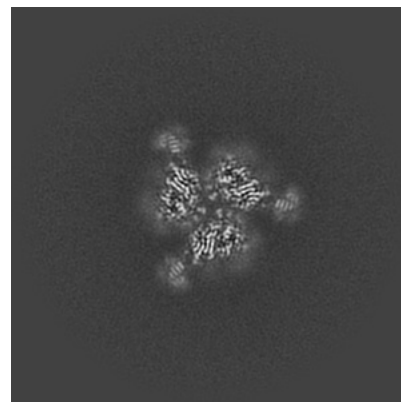
6.3.1 Primary map



X Index: 209

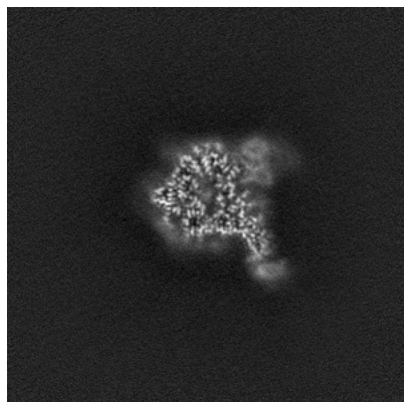


Y Index: 208

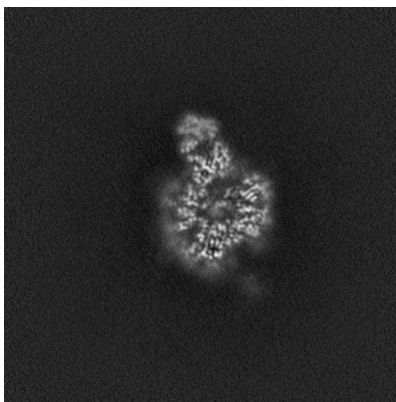


Z Index: 200

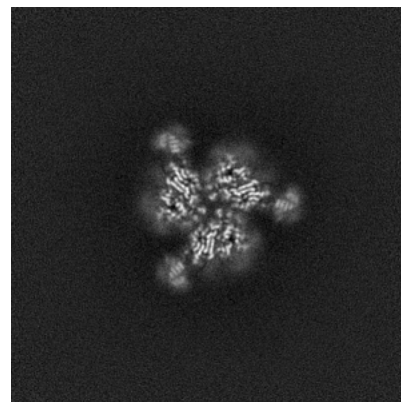
6.3.2 Raw map



X Index: 209



Y Index: 208

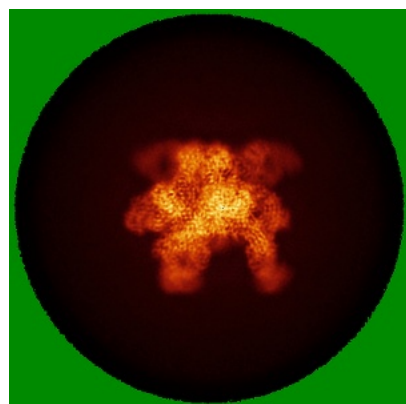


Z Index: 200

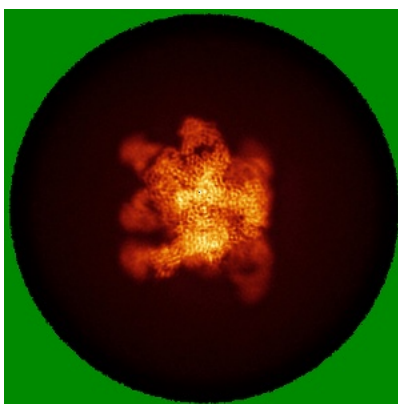
The images above show the largest variance slices of the map in three orthogonal directions.

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

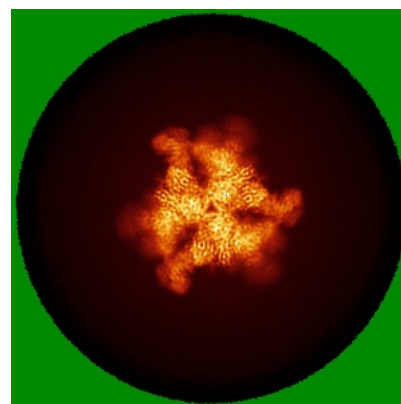
6.4.1 Primary map



X

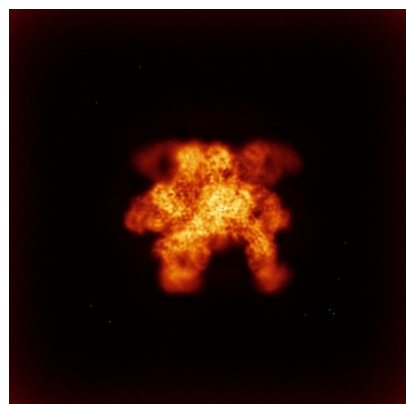


Y

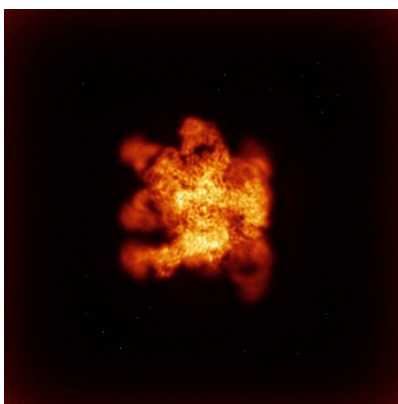


Z

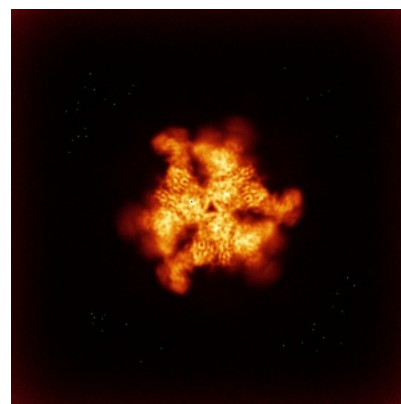
6.4.2 Raw map



X



Y

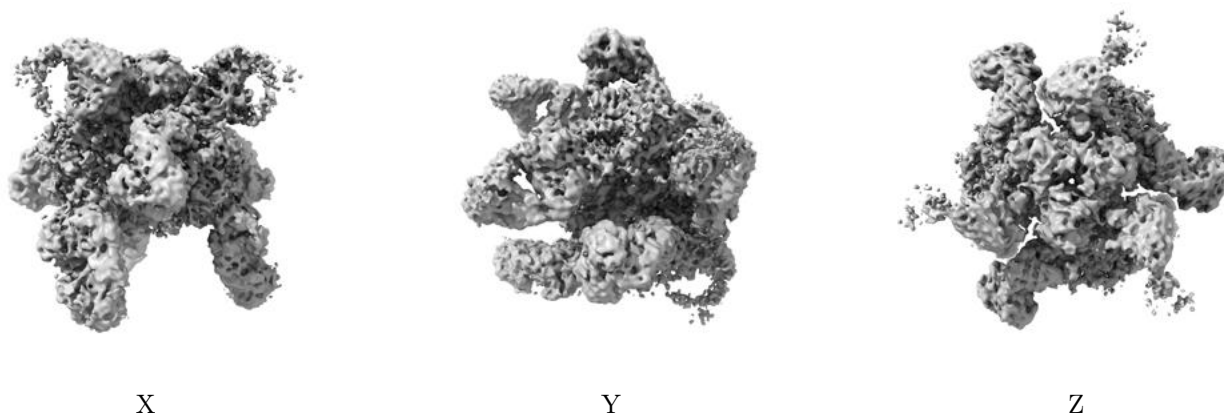


Z

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

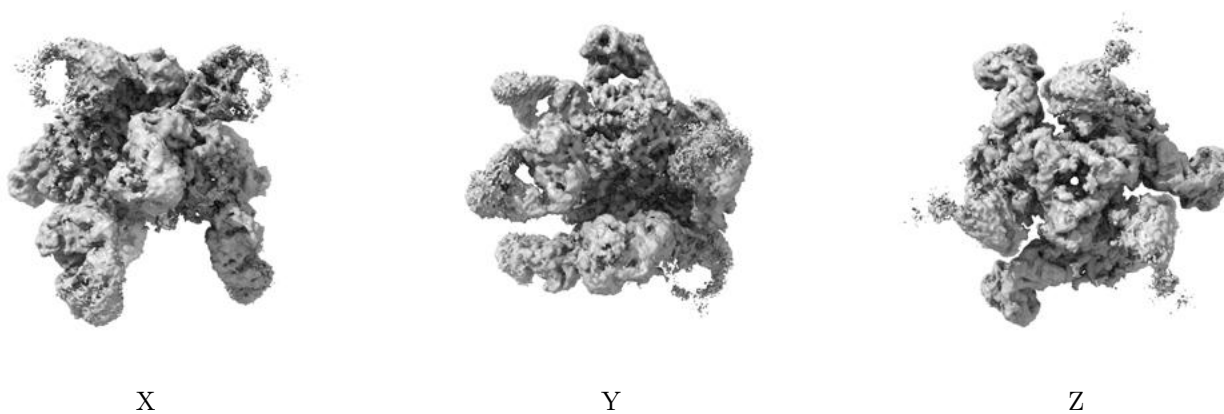
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



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

6.5.2 Raw map



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

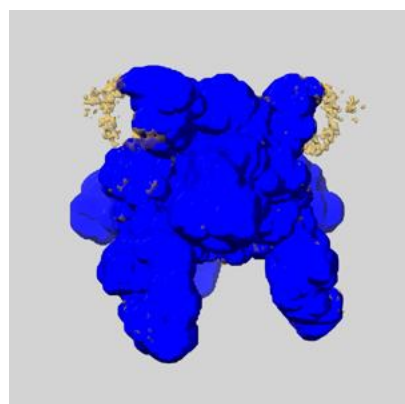
6.6 Mask visualisation [i](#)

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

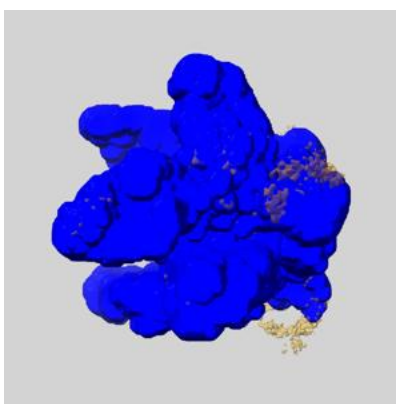
A mask typically either:

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

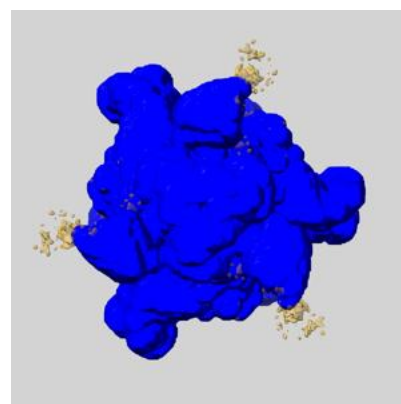
6.6.1 emd_9359_msk_1.map [i](#)



X



Y

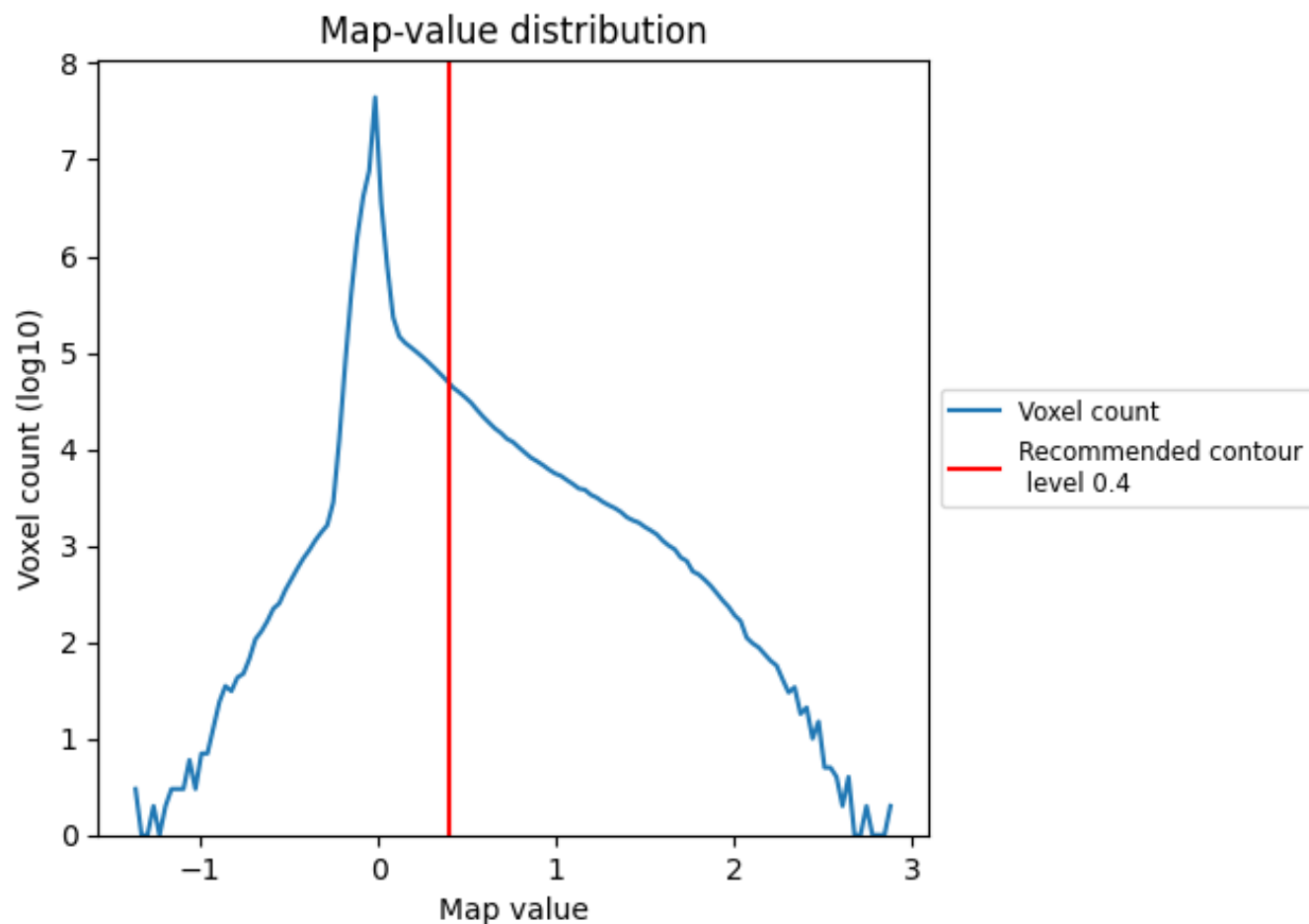


Z

7 Map analysis [i](#)

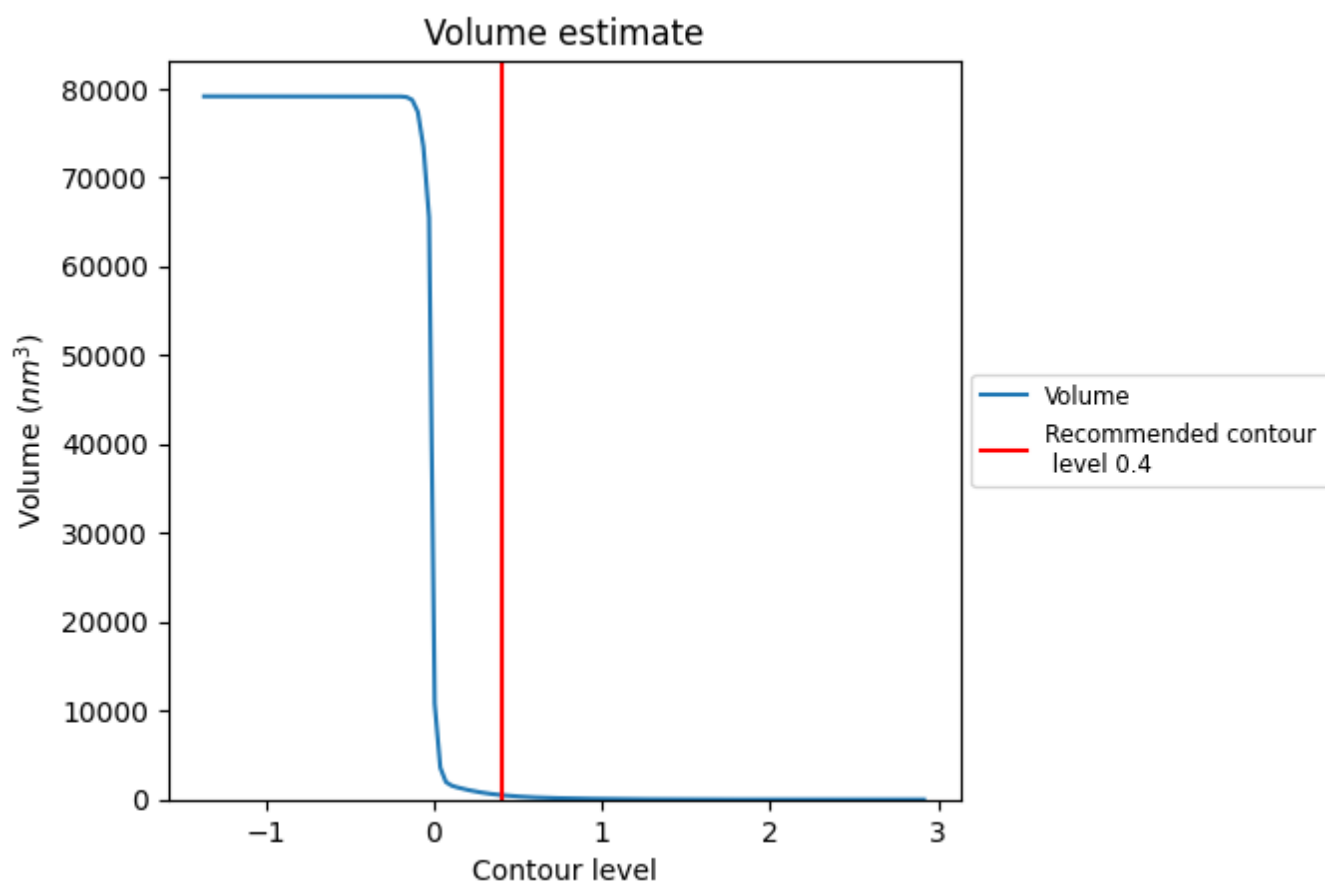
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

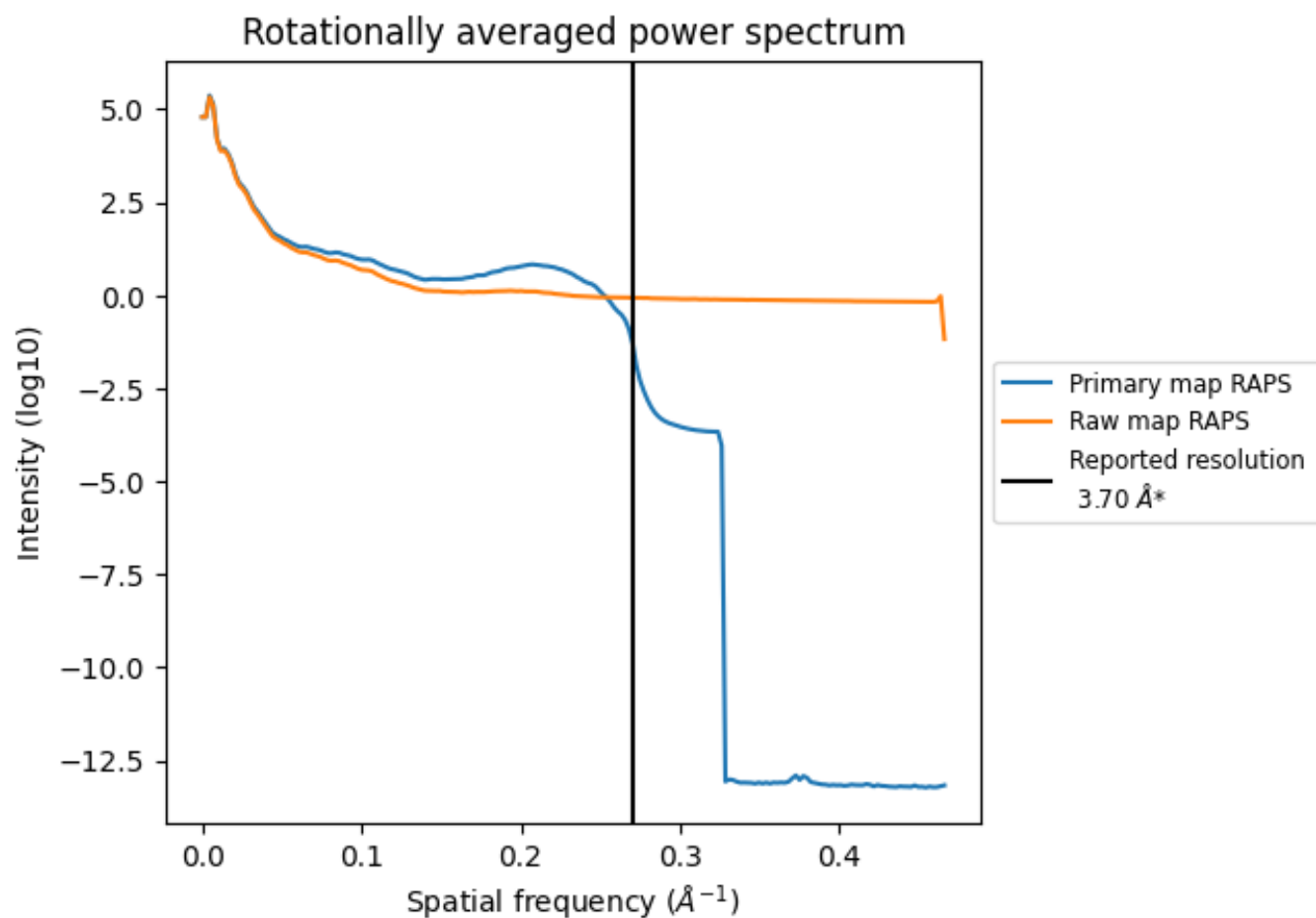
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 508 nm^3 ; this corresponds to an approximate mass of 459 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 ⓘ

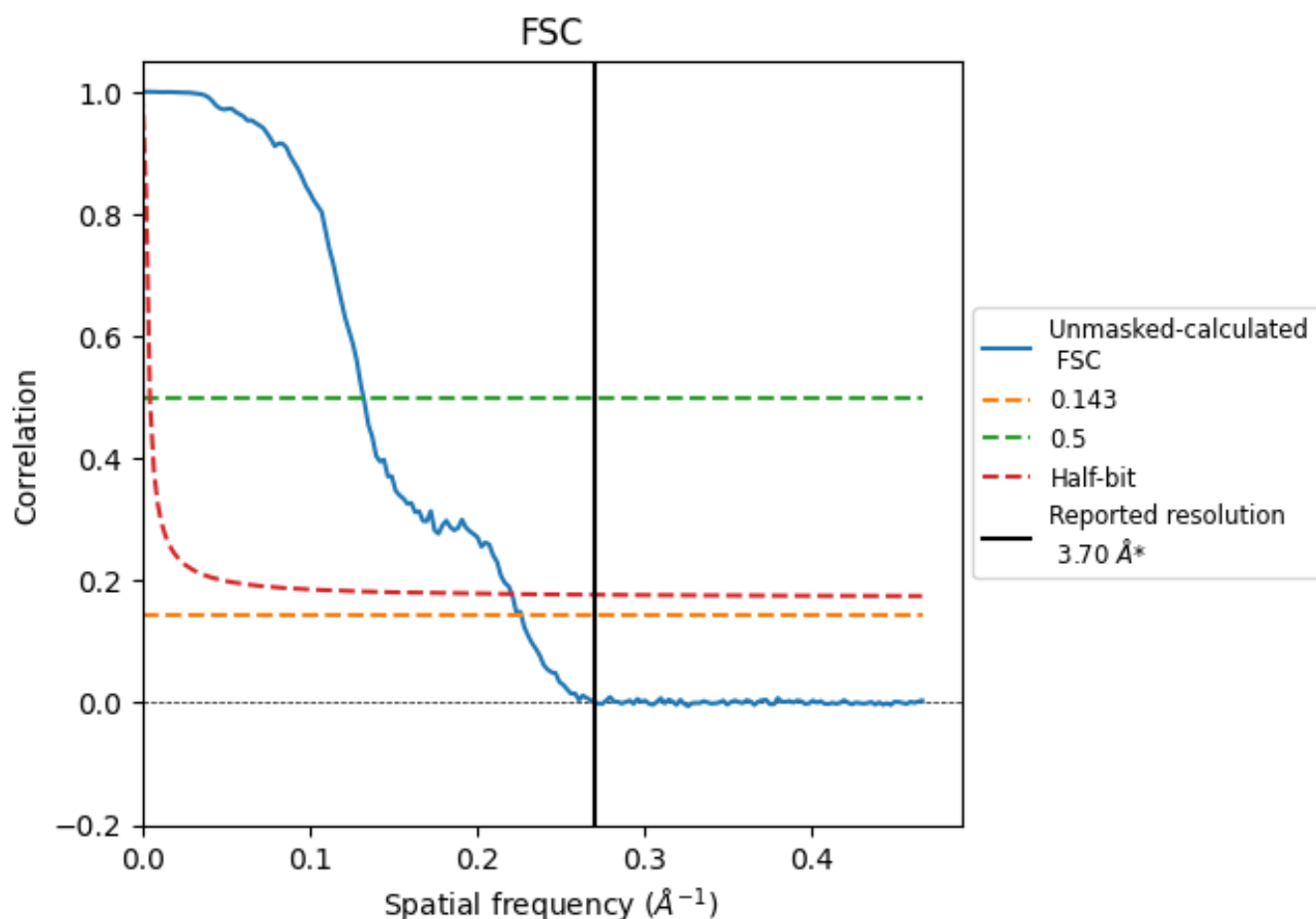


*Reported resolution corresponds to spatial frequency of 0.270 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

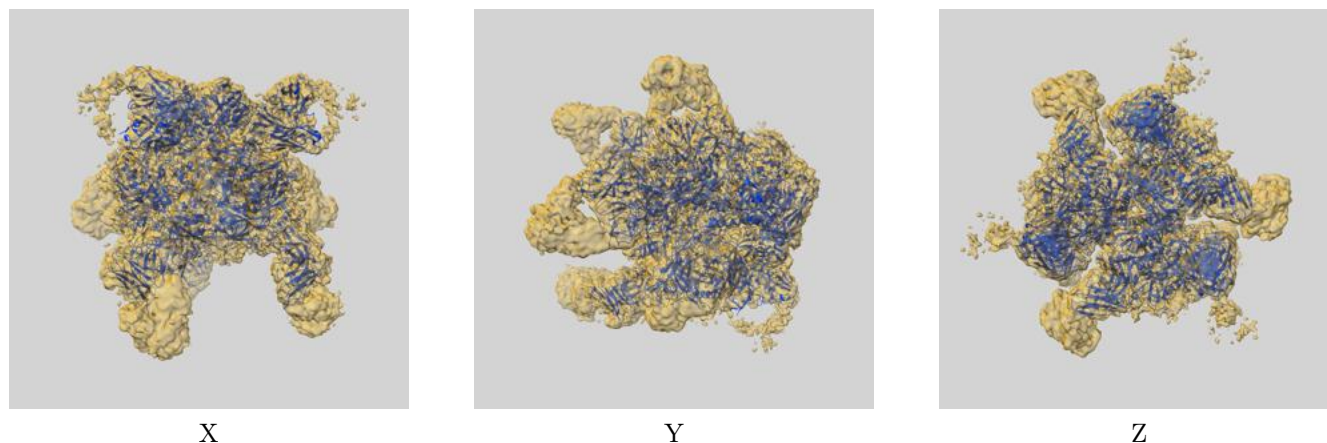
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.42	7.57	4.53

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.42 differs from the reported value 3.7 by more than 10 %

9 Map-model fit [i](#)

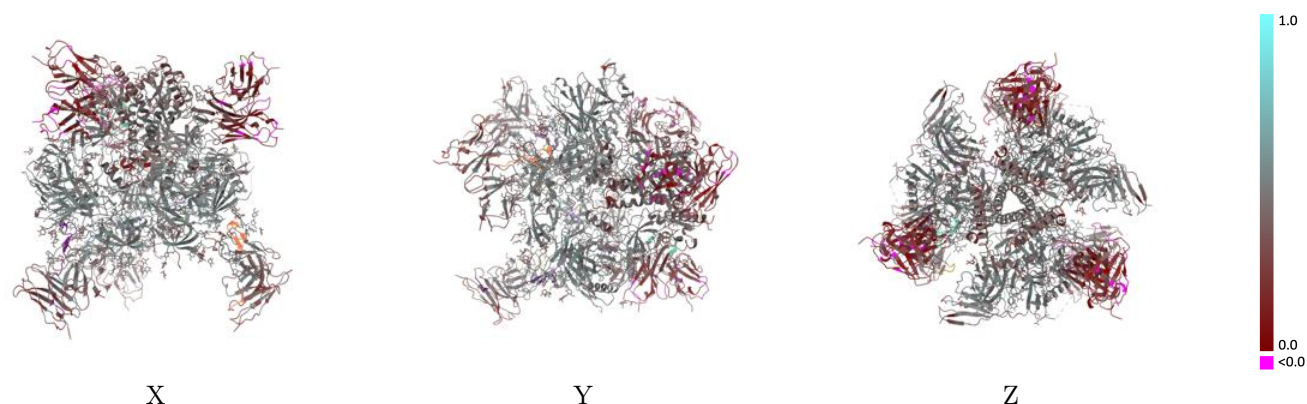
This section contains information regarding the fit between EMDB map EMD-9359 and PDB model 6NF2. Per-residue inclusion information can be found in [section 3](#) on [page 14](#).

9.1 Map-model overlay [i](#)



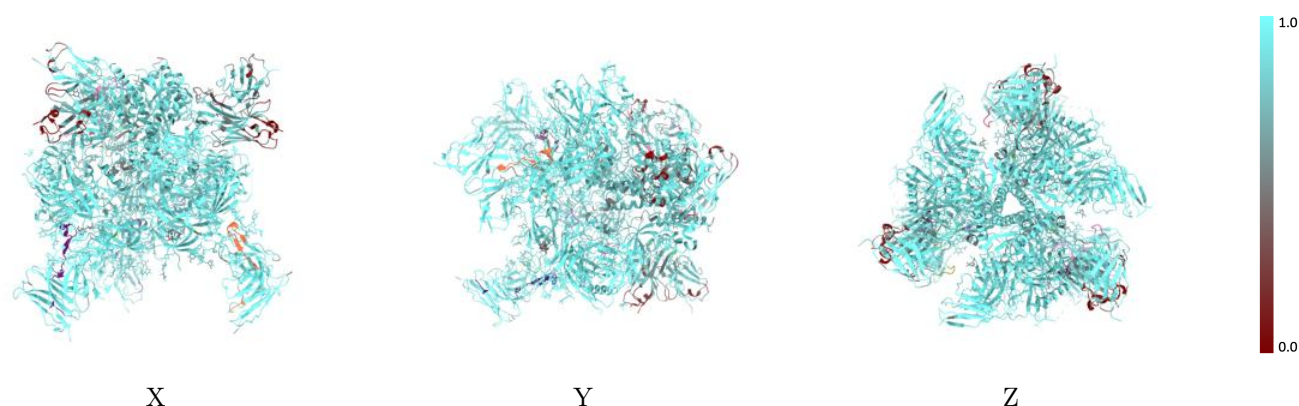
The images above show the 3D surface view of the map at the recommended contour level 0.4 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



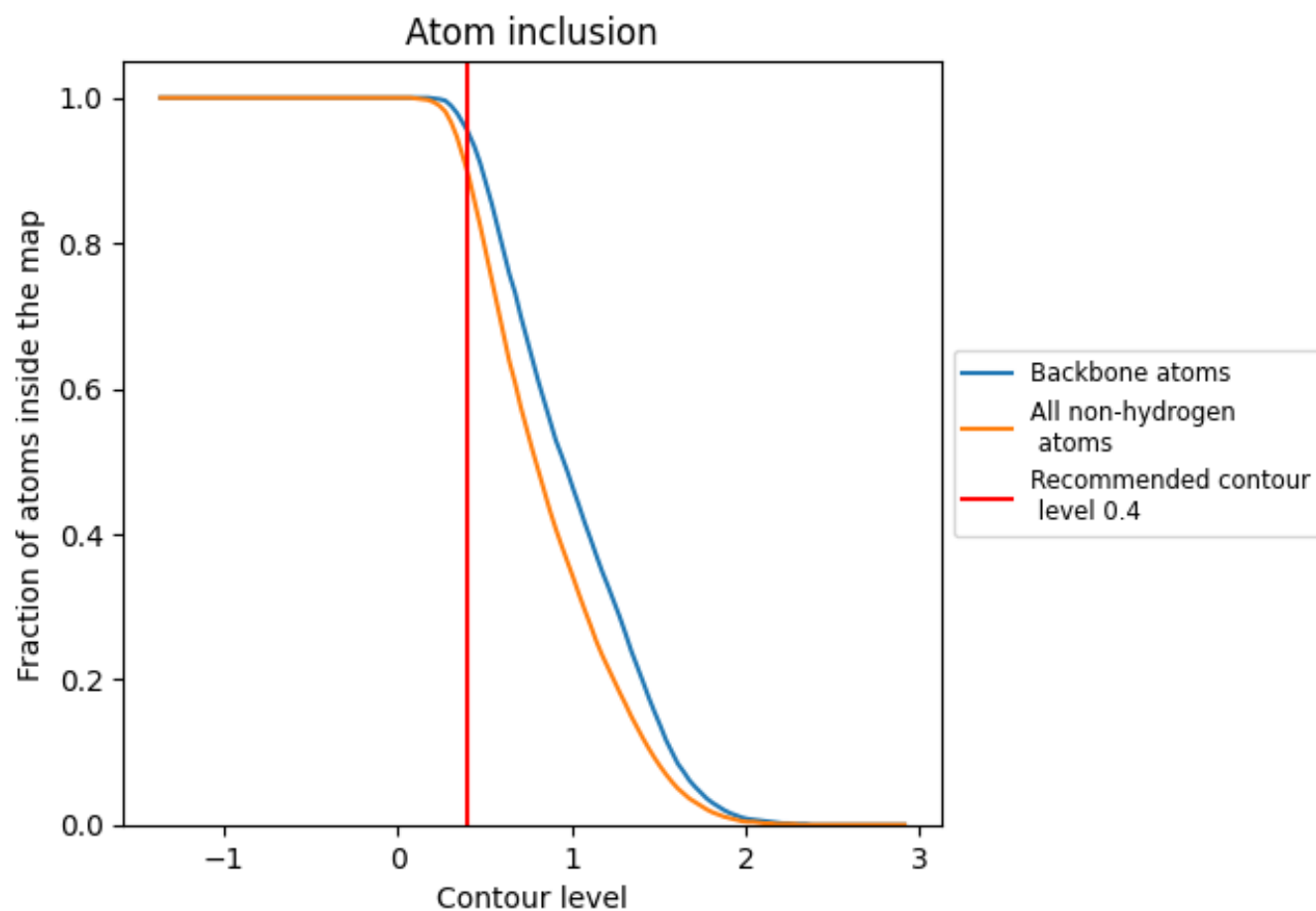
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.4).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 90% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8990	 0.4050
0	 0.8930	 0.3950
1	 0.9810	 0.4750
2	 0.7440	 0.3840
3	 0.9800	 0.4460
4	 0.6790	 0.3810
5	 0.8930	 0.3050
6	 0.9800	 0.4470
7	 0.8690	 0.4670
8	 0.8210	 0.3940
9	 0.5400	 0.3810
A	 0.9410	 0.4700
AA	 0.9740	 0.4700
B	 0.9170	 0.3980
BA	 0.9580	 0.4780
C	 0.9680	 0.4990
CA	 0.8570	 0.4610
D	 0.9520	 0.4610
DA	 0.8930	 0.4450
E	 0.9470	 0.3810
EA	 0.9740	 0.4950
F	 0.9520	 0.4150
FA	 0.9510	 0.4060
G	 0.9410	 0.4700
GA	 0.8930	 0.3850
H	 0.6240	 0.2030
HA	 0.9810	 0.4720
I	 0.9160	 0.4020
IA	 0.7440	 0.3670
J	 0.9450	 0.3810
JA	 0.9800	 0.4440
K	 0.9480	 0.4170
L	 0.7720	 0.1930
M	 0.9530	 0.4610
N	 0.9650	 0.4980



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Chain	Atom inclusion	Q-score
O	 0.6180	 0.1910
P	 0.7760	 0.1890
Q	 0.9410	 0.4710
R	 0.9150	 0.3990
S	 0.9460	 0.3800
T	 0.9520	 0.4170
U	 0.9520	 0.4570
V	 0.9700	 0.5000
W	 0.6310	 0.2110
X	 0.7780	 0.2010
Y	 0.6790	 0.3870
Z	 0.8930	 0.3140
a	 0.9800	 0.4530
b	 0.8690	 0.4650
c	 0.8210	 0.3970
d	 0.5400	 0.3840
e	 1.0000	 0.4670
f	 0.9580	 0.4730
g	 0.8570	 0.4830
h	 0.8930	 0.4460
i	 0.9740	 0.4860
j	 0.9340	 0.4040
k	 0.8930	 0.3750
l	 0.9810	 0.4690
m	 0.7440	 0.3670
n	 0.9800	 0.4460
o	 0.6790	 0.3800
p	 0.8930	 0.2940
q	 0.9800	 0.4530
r	 0.8530	 0.4640
s	 0.7950	 0.3940
t	 0.5400	 0.3950
u	 0.9740	 0.4570
v	 0.9580	 0.4780
w	 0.8570	 0.4640
x	 0.8570	 0.4490
y	 0.9740	 0.4960
z	 0.9340	 0.3990