



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 04:26 PM UTC

PDB ID : 8G6U / pdb_00008g6u
EMDB ID : EMD-29783
Title : Cryo-EM structure of T/F100 SOSIP.664 HIV-1 Env trimer with LMHS mutations in complex with 8ANC195 and 10-1074
Authors : Chen, Y.; Zhou, F.; Huang, R.; Tolbert, W.; Pazgier, M.
Deposited on : 2023-02-16
Resolution : 3.16 Å(reported)
Based on initial model : 8DOK

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:

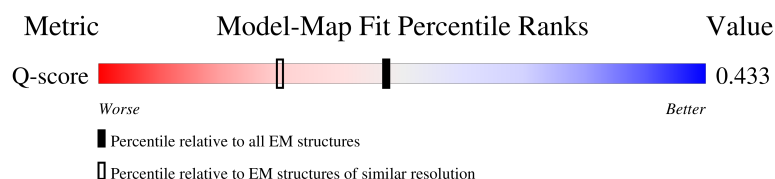
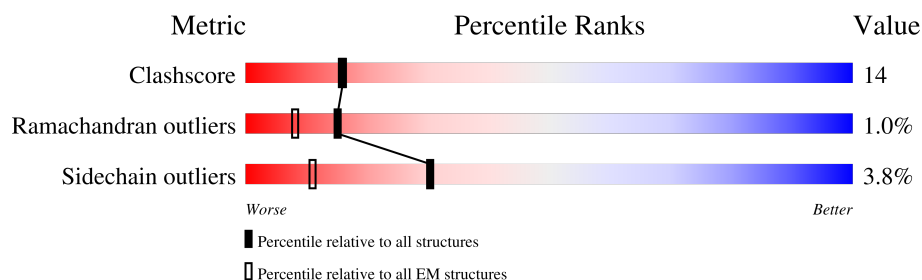
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 3.16 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.



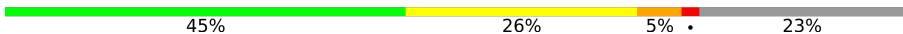
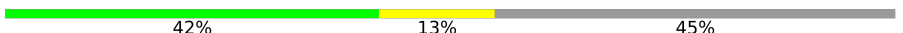
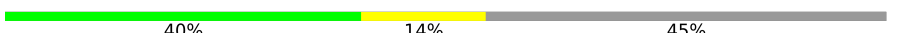
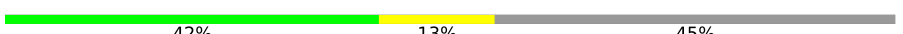
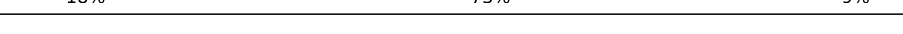
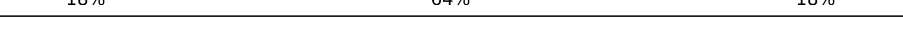
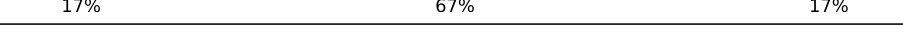
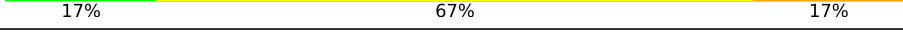
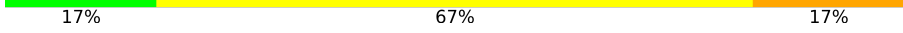
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14474 (2.66 - 3.66)

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	485	
1	E	485	
1	I	485	
2	B	155	

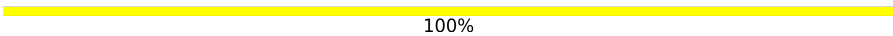
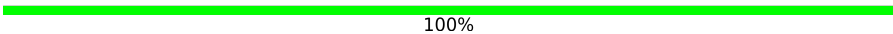
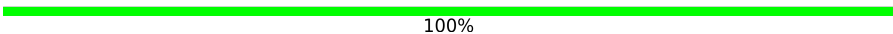
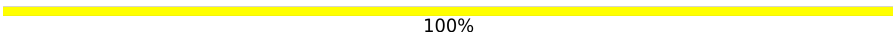
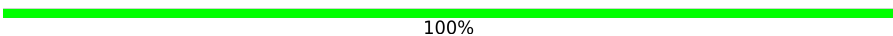
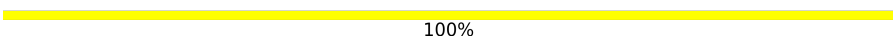
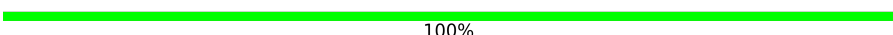
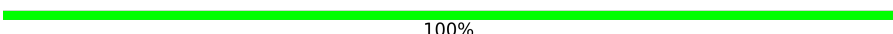
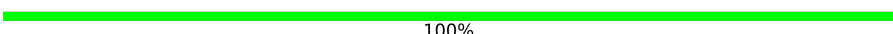
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	155	
2	J	155	
3	C	238	
3	G	238	
3	K	238	
4	D	215	
4	H	215	
4	L	215	
5	M	238	
5	O	238	
5	Q	238	
6	N	214	
6	P	214	
6	R	214	
7	S	11	
7	a	11	
7	j	11	
8	T	6	
8	Y	6	
8	b	6	
8	h	6	
8	k	6	
8	r	6	
9	U	10	
9	d	10	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
9	n	10	
10	0	2	
10	1	2	
10	2	2	
10	V	2	
10	X	2	
10	c	2	
10	e	2	
10	f	2	
10	g	2	
10	l	2	
10	m	2	
10	o	2	
10	p	2	
10	q	2	
10	t	2	
10	u	2	
10	v	2	
10	w	2	
10	x	2	
10	y	2	
10	z	2	
11	Z	5	
11	i	5	
11	s	5	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 26871 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 CRF01_AE T/F100 HIV-1 gp120.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	443	Total	C	N	O	S	0	0
			3510	2212	620	650	28		
1	E	443	Total	C	N	O	S	0	0
			3510	2212	620	650	28		
1	I	443	Total	C	N	O	S	0	0
			3510	2212	620	650	28		

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	TYR	HIS	engineered mutation	UNP A0A6C0ZY47
A	105	HIS	GLN	engineered mutation	UNP A0A6C0ZY47
A	108	ILE	VAL	engineered mutation	UNP A0A6C0ZY47
A	375	SER	HIS	conflict	UNP A0A6C0ZY47
A	?	-	SER	deletion	UNP A0A6C0ZY47
A	474	ASP	ASN	engineered mutation	UNP A0A6C0ZY47
A	475	MET	ILE	engineered mutation	UNP A0A6C0ZY47
A	476	ARG	LYS	engineered mutation	UNP A0A6C0ZY47
A	501	CYS	ALA	conflict	UNP A0A6C0ZY47
A	508	ARG	-	expression tag	UNP A0A6C0ZY47
A	509	ARG	-	expression tag	UNP A0A6C0ZY47
A	510	ARG	-	expression tag	UNP A0A6C0ZY47
A	511	ARG	-	expression tag	UNP A0A6C0ZY47
A	512	ARG	-	expression tag	UNP A0A6C0ZY47
A	513	ARG	-	expression tag	UNP A0A6C0ZY47
E	61	TYR	HIS	engineered mutation	UNP A0A6C0ZY47
E	105	HIS	GLN	engineered mutation	UNP A0A6C0ZY47
E	108	ILE	VAL	engineered mutation	UNP A0A6C0ZY47
E	375	SER	HIS	conflict	UNP A0A6C0ZY47
E	?	-	SER	deletion	UNP A0A6C0ZY47
E	474	ASP	ASN	engineered mutation	UNP A0A6C0ZY47
E	475	MET	ILE	engineered mutation	UNP A0A6C0ZY47
E	476	ARG	LYS	engineered mutation	UNP A0A6C0ZY47
E	501	CYS	ALA	conflict	UNP A0A6C0ZY47

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	508	ARG	-	expression tag	UNP A0A6C0ZY47
E	509	ARG	-	expression tag	UNP A0A6C0ZY47
E	510	ARG	-	expression tag	UNP A0A6C0ZY47
E	511	ARG	-	expression tag	UNP A0A6C0ZY47
E	512	ARG	-	expression tag	UNP A0A6C0ZY47
E	513	ARG	-	expression tag	UNP A0A6C0ZY47
I	61	TYR	HIS	engineered mutation	UNP A0A6C0ZY47
I	105	HIS	GLN	engineered mutation	UNP A0A6C0ZY47
I	108	ILE	VAL	engineered mutation	UNP A0A6C0ZY47
I	375	SER	HIS	conflict	UNP A0A6C0ZY47
I	?	-	SER	deletion	UNP A0A6C0ZY47
I	474	ASP	ASN	engineered mutation	UNP A0A6C0ZY47
I	475	MET	ILE	engineered mutation	UNP A0A6C0ZY47
I	476	ARG	LYS	engineered mutation	UNP A0A6C0ZY47
I	501	CYS	ALA	conflict	UNP A0A6C0ZY47
I	508	ARG	-	expression tag	UNP A0A6C0ZY47
I	509	ARG	-	expression tag	UNP A0A6C0ZY47
I	510	ARG	-	expression tag	UNP A0A6C0ZY47
I	511	ARG	-	expression tag	UNP A0A6C0ZY47
I	512	ARG	-	expression tag	UNP A0A6C0ZY47
I	513	ARG	-	expression tag	UNP A0A6C0ZY47

- Molecule 2 is a protein called CRF-1_AE T/F100 HIV-1 gp41.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	120	Total	C	N	O	S	0	0
			964	616	161	182	5		
2	F	120	Total	C	N	O	S	0	0
			964	616	161	182	5		
2	J	120	Total	C	N	O	S	0	0
			964	616	161	182	5		

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	conflict	UNP A0A6C0ZY47
B	605	CYS	THR	conflict	UNP A0A6C0ZY47
B	665	ALA	-	expression tag	UNP A0A6C0ZY47
B	666	ALA	-	expression tag	UNP A0A6C0ZY47
F	559	PRO	ILE	conflict	UNP A0A6C0ZY47
F	605	CYS	THR	conflict	UNP A0A6C0ZY47
F	665	ALA	-	expression tag	UNP A0A6C0ZY47

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
F	666	ALA	-	expression tag	UNP A0A6C0ZY47
J	559	PRO	ILE	conflict	UNP A0A6C0ZY47
J	605	CYS	THR	conflict	UNP A0A6C0ZY47
J	665	ALA	-	expression tag	UNP A0A6C0ZY47
J	666	ALA	-	expression tag	UNP A0A6C0ZY47

- Molecule 3 is a protein called Heavy chain of 8ANC195.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	130	Total	C	N	O	S	0	0
			1003	636	174	190	3		
3	G	130	Total	C	N	O	S	0	0
			1003	636	174	190	3		
3	K	130	Total	C	N	O	S	0	0
			1003	636	174	190	3		

- Molecule 4 is a protein called Light chain of 8ANC195.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	107	Total	C	N	O	S	0	0
			814	510	143	158	3		
4	H	107	Total	C	N	O	S	0	0
			814	510	143	158	3		
4	L	107	Total	C	N	O	S	0	0
			814	510	143	158	3		

- Molecule 5 is a protein called Heavy chain of 10-1074.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	M	131	Total	C	N	O	S	0	0
			1030	651	173	202	4		
5	O	131	Total	C	N	O	S	0	0
			1030	651	173	202	4		
5	Q	131	Total	C	N	O	S	0	0
			1030	651	173	202	4		

- Molecule 6 is a protein called Light chain of 10-1074.

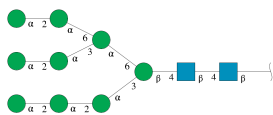
Mol	Chain	Residues	Atoms					AltConf	Trace
6	N	107	Total	C	N	O	S	0	0
			824	515	152	154	3		

Continued on next page...

Continued from previous page...

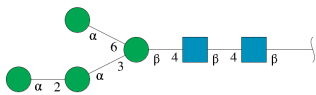
Mol	Chain	Residues	Atoms					AltConf	Trace
6	P	107	Total	C	N	O	S	0	0
			824	515	152	154	3		
6	R	107	Total	C	N	O	S	0	0
			824	515	152	154	3		

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-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
7	S	11	Total	C	N	O		0	0
			127	70	2	55			
7	a	11	Total	C	N	O		0	0
			127	70	2	55			
7	j	11	Total	C	N	O		0	0
			127	70	2	55			

- Molecule 8 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.



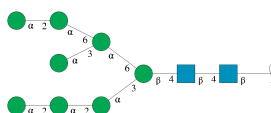
Mol	Chain	Residues	Atoms					AltConf	Trace
8	T	6	Total	C	N	O		0	0
			72	40	2	30			
8	Y	6	Total	C	N	O		0	0
			72	40	2	30			
8	b	6	Total	C	N	O		0	0
			72	40	2	30			
8	h	6	Total	C	N	O		0	0
			72	40	2	30			
8	k	6	Total	C	N	O		0	0
			72	40	2	30			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				AltConf	Trace
8	r	6	Total	C	N	O	0	0
			72	40	2	30		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
9	U	10	Total	C	N	O	0	0
			116	64	2	50		
9	d	10	Total	C	N	O	0	0
			116	64	2	50		
9	n	10	Total	C	N	O	0	0
			116	64	2	50		

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



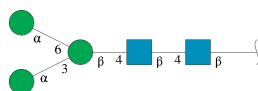
Mol	Chain	Residues	Atoms				AltConf	Trace
10	V	2	Total	C	N	O	0	0
			28	16	2	10		
10	X	2	Total	C	N	O	0	0
			28	16	2	10		
10	c	2	Total	C	N	O	0	0
			28	16	2	10		
10	f	2	Total	C	N	O	0	0
			28	16	2	10		
10	e	2	Total	C	N	O	0	0
			28	16	2	10		
10	g	2	Total	C	N	O	0	0
			28	16	2	10		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				AltConf	Trace
10	l	2	Total	C	N	O	0	0
			28	16	2	10		
10	m	2	Total	C	N	O	0	0
			28	16	2	10		
10	o	2	Total	C	N	O	0	0
			28	16	2	10		
10	p	2	Total	C	N	O	0	0
			28	16	2	10		
10	q	2	Total	C	N	O	0	0
			28	16	2	10		
10	t	2	Total	C	N	O	0	0
			28	16	2	10		
10	u	2	Total	C	N	O	0	0
			28	16	2	10		
10	v	2	Total	C	N	O	0	0
			28	16	2	10		
10	w	2	Total	C	N	O	0	0
			28	16	2	10		
10	x	2	Total	C	N	O	0	0
			28	16	2	10		
10	y	2	Total	C	N	O	0	0
			28	16	2	10		
10	z	2	Total	C	N	O	0	0
			28	16	2	10		
10	0	2	Total	C	N	O	0	0
			28	16	2	10		
10	1	2	Total	C	N	O	0	0
			28	16	2	10		
10	2	2	Total	C	N	O	0	0
			28	16	2	10		

- 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.



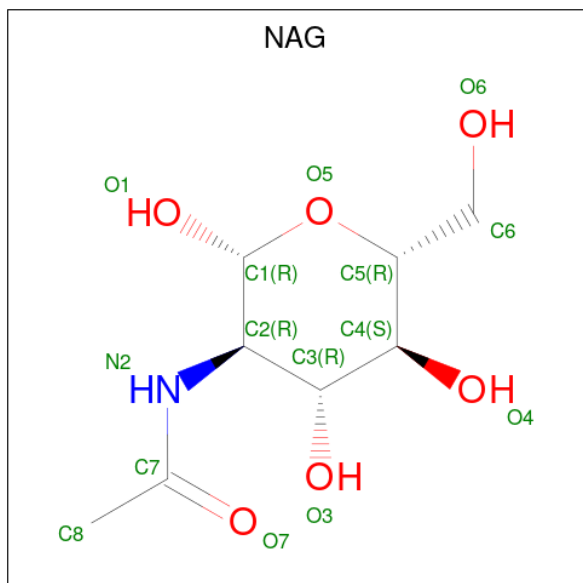
Mol	Chain	Residues	Atoms				AltConf	Trace
11	Z	5	Total	C	N	O	0	0
			61	34	2	25		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				AltConf	Trace
11	i	5	Total	C	N	O	0	0
			61	34	2	25		
11	s	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

Continued on next page...

Continued from previous page...

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

Continued on next page...

Continued from previous page...

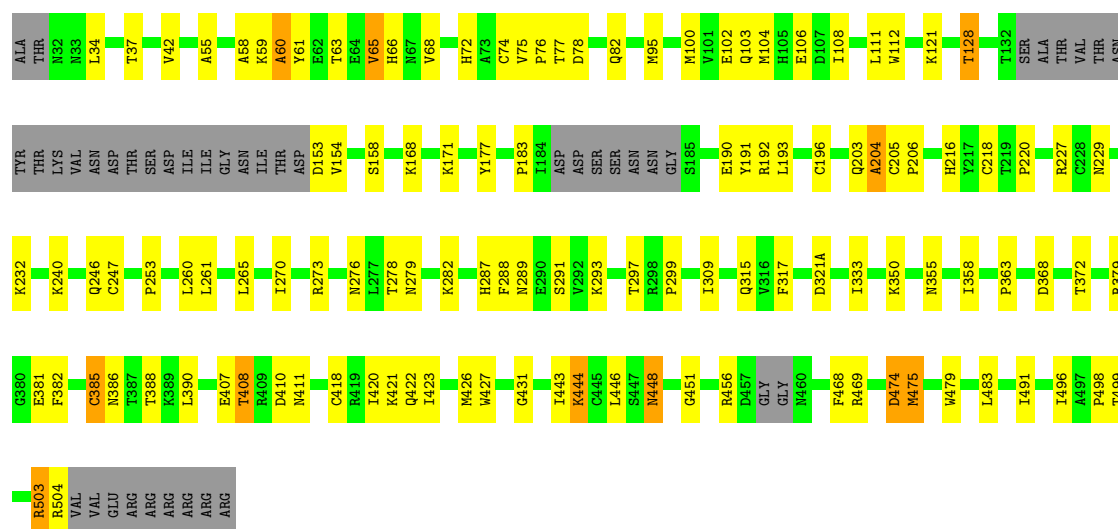
Mol	Chain	Residues	Atoms				AltConf
12	I	1	Total	C	N	O	0
			14	8	1	5	
12	I	1	Total	C	N	O	0
			14	8	1	5	
12	I	1	Total	C	N	O	0
			14	8	1	5	
12	J	1	Total	C	N	O	0
			14	8	1	5	
12	J	1	Total	C	N	O	0
			14	8	1	5	
12	J	1	Total	C	N	O	0
			14	8	1	5	

3 Residue-property plots

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

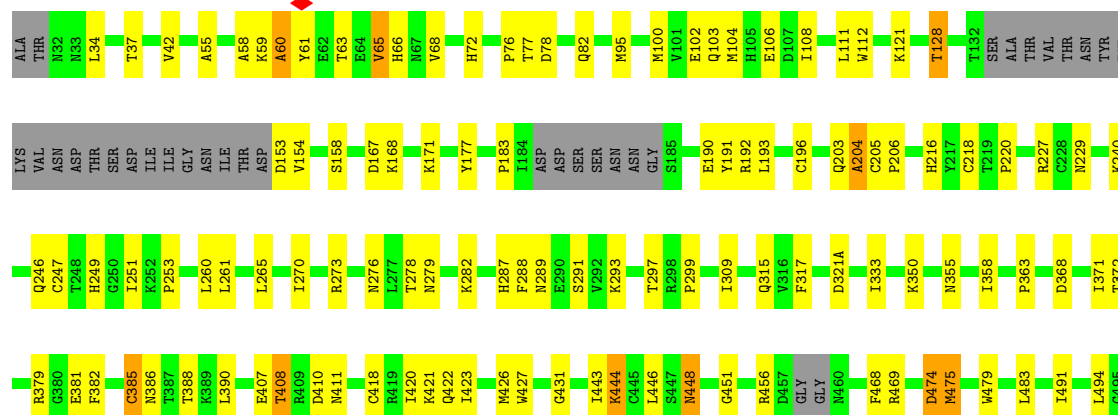
• Molecule 1: CRF01_AE T/F100 HIV-1 gp120

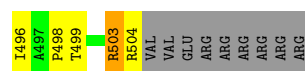
Chain A: 



• Molecule 1: CRF01_AE T/F100 HIV-1 gp120

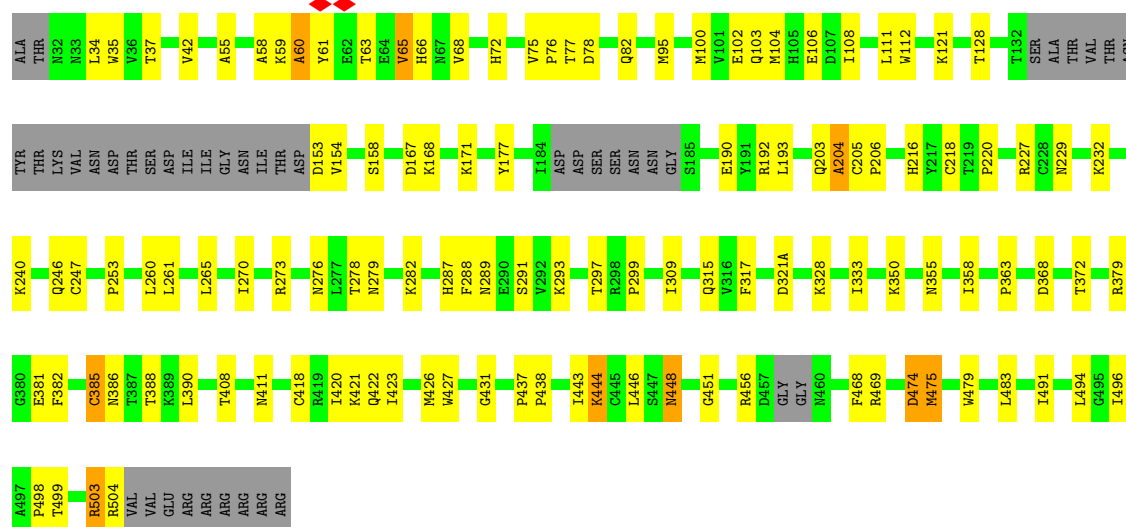
Chain E: 





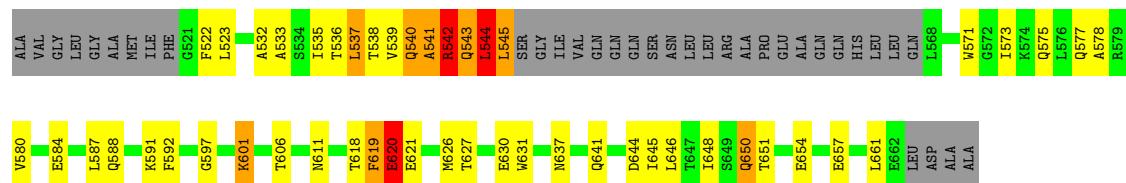
• Molecule 1: CRF01_AE T/F100 HIV-1 gp120

Chain I: 67% 23% 9%



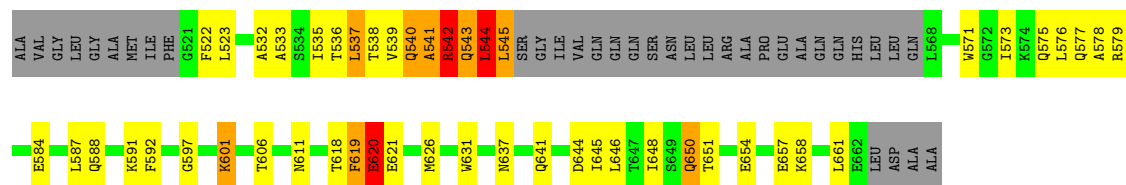
• Molecule 2: CRF-1_AE T/F100 HIV-1 gp41

Chain B: 46% 25% 5% 23%



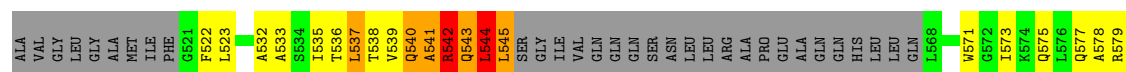
• Molecule 2: CRF-1_AE T/F100 HIV-1 gp41

Chain F: 46% 25% 5% 23%



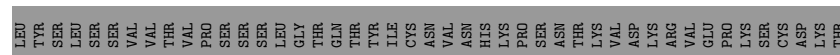
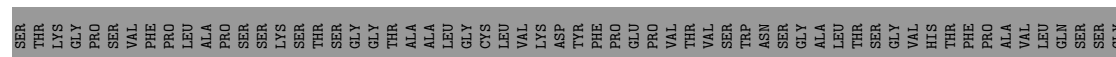
• Molecule 2: CRF-1_AE T/F100 HIV-1 gp41

Chain J: 45% 26% 5% 23%

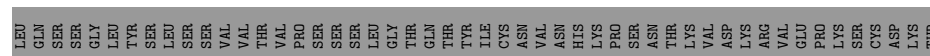
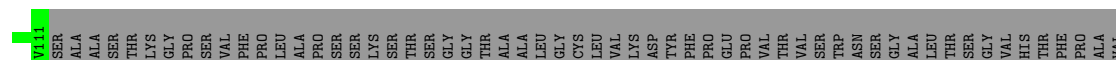




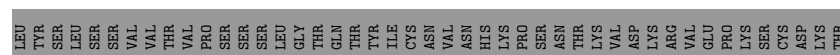
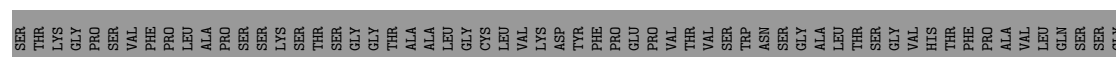
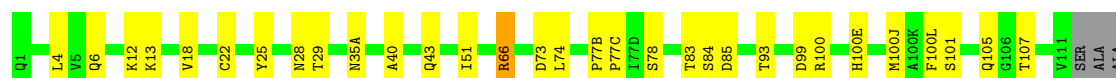
- Molecule 3: Heavy chain of 8ANC195



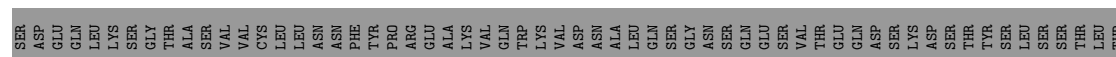
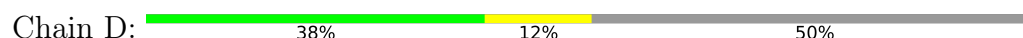
- Molecule 3: Heavy chain of 8ANC195



- Molecule 3: Heavy chain of 8ANC195



- Molecule 4: Light chain of 8ANC195



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

• Molecule 4: Light chain of 8ANC195

Chain H:  39% 11% 50%

D1 M4 T5 Q6 S14 C23 W32 W35 Y36 Q37 Q38 R39 P40 G41 A42 A43 P44 L47 G51 S65 T72 Q79 A80 E81 Y87 C88 Q89 Q90 Y91 D92 V106 LYS ARG THR VAL ALA SER PRO SER VAL PHE ILE PHE PRO PRO ASP SER GLU

GLN LEU LYS SER GLY THR ALA SER VAL VAL CYS VAL LEU ASN ASN PHE GLN TYR PRO ARG GLU ALA LYS VAL THR LYS TRP LYS VAL ASN ASP GLY ALA LEU GLN SER GLY ASN SER GLN GLU SER C88 Q89 Q90 Y91 D92 V106 LYS ASP SER LYS ASP THR TYR PRO SER SER LEU SER SER LEU THR THR LEU ASP SER LYS

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

• Molecule 4: Light chain of 8ANC195

Chain L:  37% 13% 50%

D1 M4 T5 Q6 S14 C23 W32 W35 Y36 Q37 Q38 R39 P40 G41 A42 A43 P44 L47 R50 G51 A52 S65 T72 Q79 A80 E81 Y87 C88 Q89 Q90 Y91 D92 T93 Y94 P95 V106 LYS ARG THR VAL ALA SER PRO SER VAL PHE ILE PHE

PRO PRO SER ASP GLN LEU LYS SER THR ALA SER VAL VAL THR HIS GLN TYR PRO ARG GLU ALA LYS VAL THR LYS TRP LYS VAL ASP ASN ALA GLN SER GLY ASN GLY ALA LEU GLN SER C88 Q89 Q90 Y91 D92 T93 Y94 P95 V106 LYS ASP SER LYS ASP THR TYR PRO SER SER LEU THR THR LEU SER THR

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

• Molecule 5: Heavy chain of 10-1074

Chain M:  10% 30% 24% 45%

Q1 V2 Q3 L4 Q5 E6 S7 G8 P9 G10 L11 V12 K13 P14 S15 E16 T17 L18 S19 V20 V21 T21 C22 V23 V24 D27 S28 M29 N30 Y32 Y33 V34 T35 W36 I37 R38 Q39 S40 P41 G42 K43 E46 W47 I48 G49 Y50 I51 S52 D53 R54 E55 S56 A57 N60 S65 R66

V67 S70 R71 D72 T73 S74 K75 N76 Q77 L78 S79 L80 F81 L82 V82C T87 A88 V89 Y90 C92 A95 R100 I100A F100J Y100M M100P W103 G106 T107 T108 V109 T110 V111 S112 SER ALA SER THR LYS GLY PRO VAL THR VAL PHE PRO LEU LEU ALA SER THR Y50 I51 S52 D53 R54 E55 S56 A57 N60 S65 R66

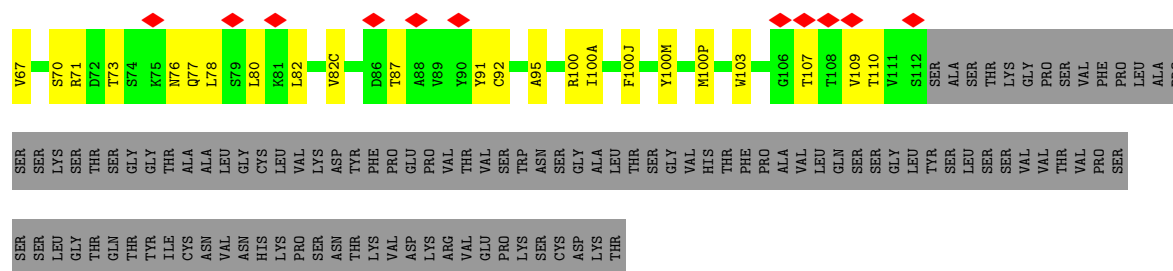
SER LYS SER THR GLY THR GLY THR ALA LEU CYS LEU VAL LYS ASP PHE PRO GLU PRO VAL VAL SER TRP ASN SER GLY CYS ASP LYS LEU THR SER VAL HIS THR PHE PRO ALA VAL LEU GLN SER SER GLY THR TYR VAL LEU SER VAL THR VAL PRO THR VAL LEU LEU ALA SER

SER LEU THR GLN TYR ILE ASN VAL ASN HIS LYS PRO SER ASN THR LYS VAL ASP GLU PRO ARG VAL VAL GLU PRO LYS SER CYS ASP LYS THR

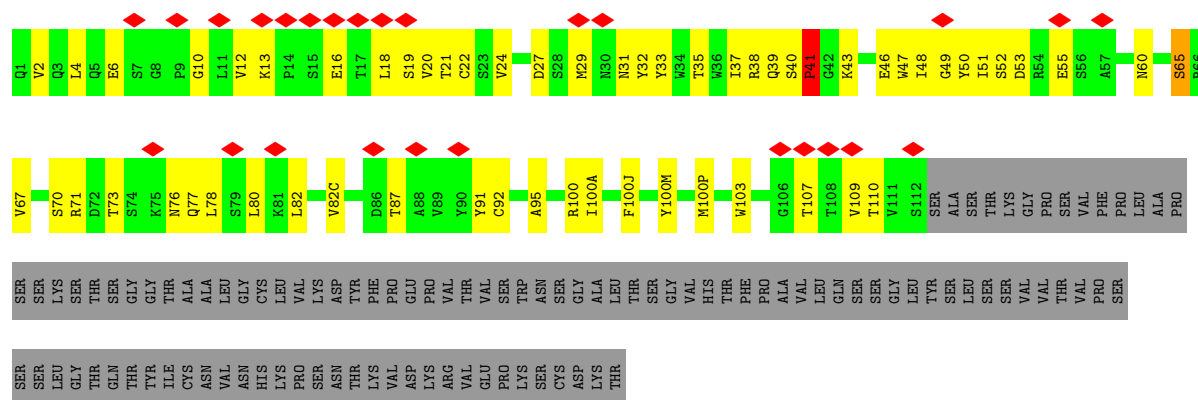
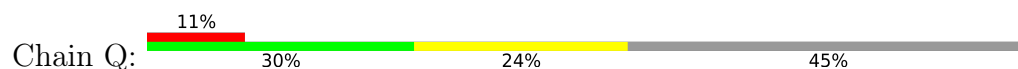
• Molecule 5: Heavy chain of 10-1074

Chain O:  11% 29% 25% 45%

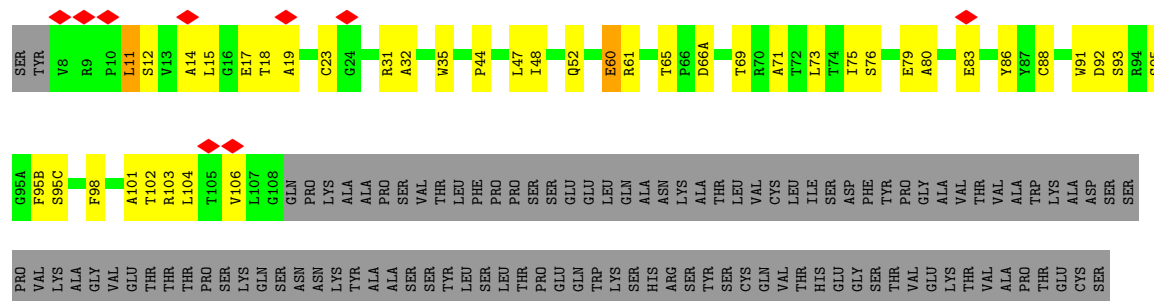
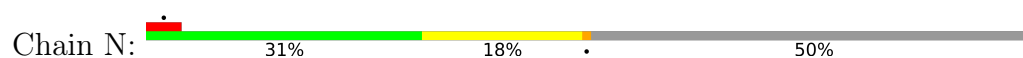
Q1 V2 Q3 L4 Q5 E6 S7 G8 P9 G10 L11 V12 K13 P14 S15 E16 T17 L18 S19 V20 V21 T21 C22 V23 V24 D27 S28 M29 N30 Y32 Y33 V34 T35 W36 I37 R38 Q39 S40 P41 G42 K43 E46 W47 I48 G49 Y50 I51 S52 D53 R54 E55 S56 A57 N60 S65 R66



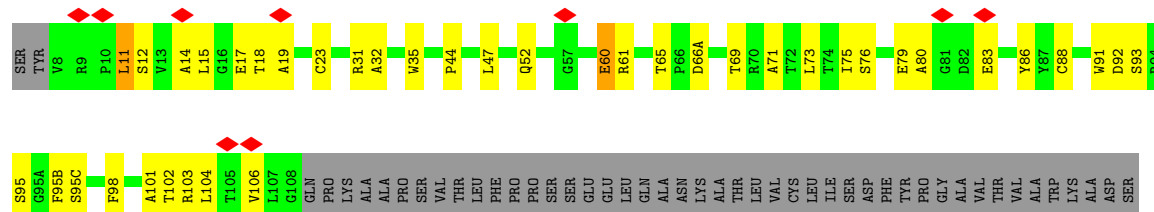
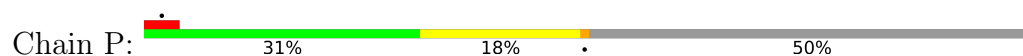
• Molecule 5: Heavy chain of 10-1074



• Molecule 6: Light chain of 10-1074



• Molecule 6: Light chain of 10-1074



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

• Molecule 6: Light chain of 10-1074

Chain R: 

SER TRP R8 R9 P10 L11 S12 V13 A14 L15 G16 E17 T18 A19 C23 A27 R31 A32 W35 P44 L47 Q52 D53 G57 E60 E61 T65 P66 D66A T69 E70 A71 T72 L73 S76 E79 A80 G81 D82 E83 C88 W91 D92 S93 R94

S95 G95A F95B S95C F98 A101 T102 R103 L104 T105 V106 L107 G108 GLN PRO LYS ALA ALA PRO SER VAL LEU PHE PRO SER GLU GLU LEU GLN ALA ASN LYS ALA THR LEU VAL CYS ILE SER ASP PHE TYR PRO GLY VAL LYS THR VAL ALA THR TRP LYS ASP SER

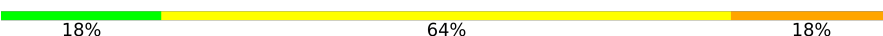
SER PRO VAL LYS ALA GLY VAL GLU THR THR PRO SER LYS GLN SER ASN ASN LYS TYR ALA ALA SER SER TYR LEU LEU SER PRO GLU GLN TRP LYS SER HIS ARG SER TYR CYS GLN VAL THR HIS GLY SER THR VAL GLU LYS THR VAL ALA PRO THR CYS SER

• Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S: 

MAN1 MAN2 MAN3 MAN4 MAN5 MAN6 MAN7 MAN8 MAN9 MAN10 MAN11

• Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a: 

MAN1 MAN2 MAN3 MAN4 MAN5 MAN6 MAN7 MAN8 MAN9 MAN10 MAN11

• Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain j: 

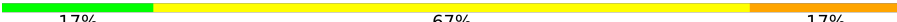
MAN1 MAN2 MAN3 MAN4 MAN5 MAN6 MAN7 MAN8 MAN9 MAN10 MAN11

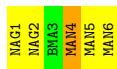
• Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  50% 50%



• Molecule 8: 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 Y:  17% 67% 17%

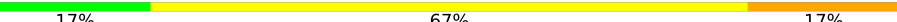


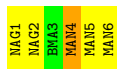
• Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:  50% 50%



• Molecule 8: 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 h:  17% 67% 17%

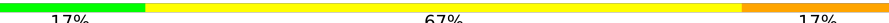


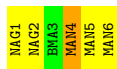
• Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain k:  50% 50%



• Molecule 8: 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 r:  17% 67% 17%

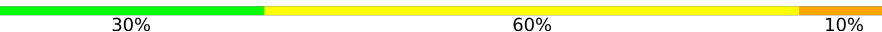


● Molecule 9: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain U: 

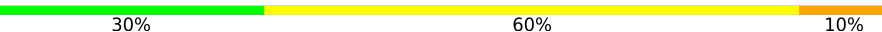


● Molecule 9: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain d: 

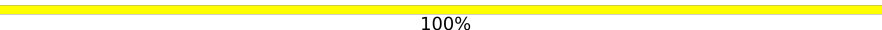


● Molecule 9: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain n: 

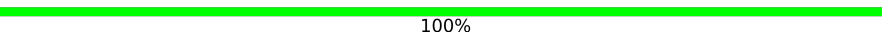


● Molecule 10: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain V: 

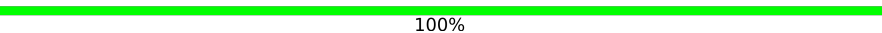


● Molecule 10: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain X: 



● Molecule 10: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain c: 



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

Chain f:  100%



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

Chain e:  100%



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

Chain g:  100%



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

Chain i:  100%



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

Chain m:  50% 50%



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

Chain o:  100%



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

Chain p:  100%



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

Chain q:  100%



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

Chain t:  100%



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

Chain u:  100%



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

Chain v:  50% 50%



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

Chain w:  50% 50%



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

Chain x:  50% 50%

NAG1
NAG2

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

Chain y:  50% 50%

NAG1
NAG2

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

Chain z:  50% 50%

NAG1
NAG2

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

Chain 0:  50% 50%

NAG1
NAG2

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

Chain 1:  50% 50%

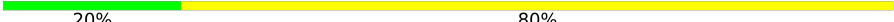
NAG1
NAG2

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

Chain 2:  50% 50%

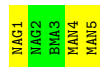
NAG1
NAG2

- 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

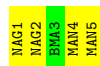
Chain Z:  20% 80%



- 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



- 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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	549061	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 ($e^-/\text{\AA}^2$)	54	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2700	Depositor
Magnification	60241	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	25.188	Depositor
Minimum map value	-13.205	Depositor
Average map value	-0.014	Depositor
Map value standard deviation	0.602	Depositor
Recommended contour level	1.4	Depositor
Map size (Å)	403.38, 403.38, 403.38	wwPDB
Map dimensions	486, 486, 486	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.18	0/3588	0.48	3/4867 (0.1%)
1	E	0.18	0/3588	0.48	3/4867 (0.1%)
1	I	0.18	0/3588	0.48	3/4867 (0.1%)
2	B	0.34	0/983	0.63	3/1333 (0.2%)
2	F	0.34	0/983	0.63	3/1333 (0.2%)
2	J	0.34	0/983	0.63	3/1333 (0.2%)
3	C	0.18	0/1030	0.32	0/1403
3	G	0.18	0/1030	0.32	0/1403
3	K	0.18	0/1030	0.32	0/1403
4	D	0.12	0/832	0.32	0/1130
4	H	0.12	0/832	0.32	0/1130
4	L	0.12	0/832	0.32	0/1130
5	M	0.23	0/1055	0.61	3/1436 (0.2%)
5	O	0.23	0/1055	0.61	3/1436 (0.2%)
5	Q	0.23	0/1055	0.61	3/1436 (0.2%)
6	N	0.19	0/845	0.62	2/1148 (0.2%)
6	P	0.19	0/845	0.62	2/1148 (0.2%)
6	R	0.19	0/845	0.62	2/1148 (0.2%)
All	All	0.21	0/24999	0.50	33/33951 (0.1%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	O	41	PRO	CB-CA-C	7.94	124.66	111.56
5	Q	41	PRO	CB-CA-C	7.93	124.65	111.56
5	M	41	PRO	CB-CA-C	7.93	124.64	111.56
2	F	542	ARG	N-CA-C	6.77	119.28	107.49
2	B	542	ARG	N-CA-C	6.77	119.27	107.49
2	J	542	ARG	N-CA-C	6.75	119.23	107.49

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Q	41	PRO	N-CA-CB	-6.36	96.57	103.25
5	M	41	PRO	N-CA-CB	-6.33	96.61	103.25
5	O	41	PRO	N-CA-CB	-6.31	96.62	103.25
1	A	385	CYS	CA-C-N	-6.00	110.01	122.07
1	A	385	CYS	C-N-CA	-6.00	110.01	122.07
1	A	61	TYR	N-CA-C	-6.00	106.21	112.93
1	E	385	CYS	CA-C-N	-5.99	110.04	122.07
1	E	385	CYS	C-N-CA	-5.99	110.04	122.07
1	I	385	CYS	CA-C-N	-5.98	110.05	122.07
1	I	385	CYS	C-N-CA	-5.98	110.05	122.07
5	Q	41	PRO	CA-N-CD	-5.96	103.66	112.00
5	M	41	PRO	CA-N-CD	-5.95	103.67	112.00
1	I	61	TYR	N-CA-C	-5.95	106.27	112.93
5	O	41	PRO	CA-N-CD	-5.95	103.68	112.00
1	E	61	TYR	N-CA-C	-5.94	106.28	112.93
2	B	540	GLN	N-CA-C	-5.91	101.89	111.04
2	F	540	GLN	N-CA-C	-5.90	101.89	111.04
2	J	540	GLN	N-CA-C	-5.90	101.89	111.04
6	N	17	GLU	N-CA-C	-5.79	100.56	109.76
6	R	17	GLU	N-CA-C	-5.77	100.58	109.76
2	F	620	GLU	CA-CB-CG	5.76	125.62	114.10
2	J	620	GLU	CA-CB-CG	5.76	125.61	114.10
2	B	620	GLU	CA-CB-CG	5.75	125.60	114.10
6	P	17	GLU	N-CA-C	-5.74	100.64	109.76
6	R	61	ARG	N-CA-C	-5.24	105.61	113.89
6	N	61	ARG	N-CA-C	-5.24	105.61	113.89
6	P	61	ARG	N-CA-C	-5.22	105.65	113.89

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3510	0	3435	100	0
1	E	3510	0	3435	100	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	I	3510	0	3435	102	0
2	B	964	0	944	75	0
2	F	964	0	944	76	0
2	J	964	0	944	72	0
3	C	1003	0	974	21	0
3	G	1003	0	974	25	0
3	K	1003	0	974	23	0
4	D	814	0	787	14	0
4	H	814	0	787	14	0
4	L	814	0	787	16	0
5	M	1030	0	995	43	0
5	O	1030	0	995	44	0
5	Q	1030	0	995	43	0
6	N	824	0	788	24	0
6	P	824	0	788	24	0
6	R	824	0	788	22	0
7	S	127	0	106	1	0
7	a	127	0	106	2	0
7	j	127	0	106	1	0
8	T	72	0	61	0	0
8	Y	72	0	61	3	0
8	b	72	0	61	0	0
8	h	72	0	61	3	0
8	k	72	0	61	0	0
8	r	72	0	61	3	0
9	U	116	0	97	4	0
9	d	116	0	97	4	0
9	n	116	0	97	4	0
10	0	28	0	25	0	0
10	1	28	0	25	0	0
10	2	28	0	25	0	0
10	V	28	0	25	1	0
10	X	28	0	25	0	0
10	c	28	0	25	0	0
10	e	28	0	25	1	0
10	f	28	0	25	0	0
10	g	28	0	25	0	0
10	l	28	0	25	0	0
10	m	28	0	25	0	0
10	o	28	0	25	1	0
10	p	28	0	25	0	0
10	q	28	0	25	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	t	28	0	25	0	0
10	u	28	0	25	0	0
10	v	28	0	25	1	0
10	w	28	0	25	1	0
10	x	28	0	25	1	0
10	y	28	0	25	1	0
10	z	28	0	25	1	0
11	Z	61	0	52	2	0
11	i	61	0	52	2	0
11	s	61	0	52	3	0
12	A	126	0	117	0	0
12	B	42	0	39	1	0
12	E	126	0	117	0	0
12	F	42	0	39	1	0
12	I	126	0	117	0	0
12	J	42	0	39	1	0
All	All	26871	0	25893	759	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 14.

All (759) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:TRP:CZ3	1:A:475:MET:SD	2.25	1.29
1:I:427:TRP:CE3	1:I:475:MET:HE1	1.67	1.29
1:E:427:TRP:CZ3	1:E:475:MET:SD	2.25	1.28
1:A:427:TRP:CE3	1:A:475:MET:HE1	1.68	1.27
1:I:427:TRP:CZ3	1:I:475:MET:SD	2.25	1.27
1:E:427:TRP:CE3	1:E:475:MET:HE1	1.67	1.26
1:I:427:TRP:CE3	1:I:475:MET:CE	2.21	1.23
1:E:427:TRP:CE3	1:E:475:MET:CE	2.21	1.23
1:A:427:TRP:CE3	1:A:475:MET:CE	2.22	1.23
1:E:427:TRP:CZ3	1:E:475:MET:HE1	1.76	1.20
1:I:427:TRP:CZ3	1:I:475:MET:HE1	1.76	1.20
1:A:427:TRP:CZ3	1:A:475:MET:HE1	1.77	1.20
2:B:651:THR:OG1	2:F:538:THR:HG21	1.40	1.18
1:E:427:TRP:CE3	1:E:475:MET:SD	2.39	1.16
1:I:427:TRP:CE3	1:I:475:MET:SD	2.39	1.15
1:I:427:TRP:CZ3	1:I:475:MET:CE	2.30	1.14
1:A:427:TRP:CZ3	1:A:475:MET:CE	2.31	1.14

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:TRP:CE3	1:A:475:MET:SD	2.39	1.14
1:E:427:TRP:CZ3	1:E:475:MET:CE	2.30	1.13
2:F:651:THR:OG1	2:J:538:THR:HG21	1.53	1.06
2:B:538:THR:HG21	2:J:651:THR:OG1	1.63	0.97
1:A:427:TRP:HZ3	1:A:475:MET:SD	1.84	0.95
1:E:427:TRP:HE3	1:E:475:MET:CE	1.71	0.95
1:E:427:TRP:HZ3	1:E:475:MET:SD	1.84	0.95
1:I:427:TRP:HE3	1:I:475:MET:CE	1.71	0.94
2:F:661:LEU:HD11	1:I:503:ARG:HG2	1.53	0.91
1:A:427:TRP:HE3	1:A:475:MET:CE	1.72	0.89
2:F:620:GLU:OE1	2:F:620:GLU:N	2.06	0.89
1:I:427:TRP:HZ3	1:I:475:MET:SD	1.84	0.89
2:B:620:GLU:OE1	2:B:620:GLU:N	2.06	0.88
2:J:620:GLU:OE1	2:J:620:GLU:N	2.06	0.88
2:B:588:GLN:OE1	2:F:544:LEU:HD11	1.73	0.87
2:B:543:GLN:O	2:B:544:LEU:O	1.94	0.85
2:J:543:GLN:O	2:J:544:LEU:O	1.95	0.84
2:J:542:ARG:HH11	2:J:542:ARG:HB2	1.44	0.83
2:B:544:LEU:HD11	2:J:588:GLN:OE1	1.79	0.83
2:B:542:ARG:HH11	2:B:542:ARG:HB2	1.43	0.83
2:F:542:ARG:HH11	2:F:542:ARG:HB2	1.44	0.83
2:F:543:GLN:O	2:F:544:LEU:O	1.94	0.83
1:A:427:TRP:HZ3	1:A:475:MET:CE	1.89	0.82
5:Q:40:SER:HB3	5:Q:41:PRO:HD2	1.62	0.81
5:O:40:SER:HB3	5:O:41:PRO:HD2	1.62	0.81
1:E:448:ASN:OD1	1:E:448:ASN:N	2.14	0.81
5:M:40:SER:HB3	5:M:41:PRO:HD2	1.62	0.80
1:E:427:TRP:HZ3	1:E:475:MET:CE	1.88	0.79
5:O:38:ARG:HH21	5:O:46:GLU:HG2	1.48	0.79
1:A:448:ASN:N	1:A:448:ASN:OD1	2.14	0.78
5:M:38:ARG:HH21	5:M:46:GLU:HG2	1.48	0.78
2:B:588:GLN:CD	2:F:544:LEU:HD13	2.07	0.78
1:A:59:LYS:O	1:A:60:ALA:CB	2.32	0.78
1:E:59:LYS:O	1:E:60:ALA:CB	2.32	0.78
5:Q:38:ARG:HH21	5:Q:46:GLU:HG2	1.48	0.77
1:I:59:LYS:O	1:I:60:ALA:CB	2.32	0.77
1:I:448:ASN:N	1:I:448:ASN:OD1	2.14	0.77
6:R:52:GLN:HG3	9:n:6:MAN:H2	1.68	0.76
2:B:544:LEU:HD13	2:J:588:GLN:CD	2.11	0.76
2:B:588:GLN:OE1	2:F:544:LEU:CD1	2.33	0.75
6:P:52:GLN:HG3	9:d:6:MAN:H2	1.69	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:6:GLN:HE21	3:C:107:THR:HG23	1.52	0.74
3:G:12:LYS:HD2	3:G:18:VAL:HG22	1.70	0.74
3:K:6:GLN:HE21	3:K:107:THR:HG23	1.52	0.74
1:A:59:LYS:O	1:A:60:ALA:HB3	1.87	0.73
6:N:52:GLN:HG3	9:U:6:MAN:H2	1.69	0.73
2:F:661:LEU:HD21	1:I:503:ARG:HB2	1.70	0.73
3:K:12:LYS:HD2	3:K:18:VAL:HG22	1.70	0.73
2:B:592:PHE:HE1	2:F:544:LEU:HD13	1.53	0.72
3:C:12:LYS:HD2	3:C:18:VAL:HG22	1.70	0.72
1:E:59:LYS:O	1:E:60:ALA:HB3	1.87	0.72
1:I:59:LYS:O	1:I:60:ALA:HB3	1.87	0.72
5:Q:100(A):ILE:HG22	5:Q:100(J):PHE:HB3	1.70	0.72
6:R:11:LEU:HD21	6:R:102:THR:HA	1.71	0.72
2:F:592:PHE:HE1	2:J:544:LEU:HD13	1.53	0.72
3:G:6:GLN:HE21	3:G:107:THR:HG23	1.52	0.72
6:P:11:LEU:HD21	6:P:102:THR:HA	1.72	0.72
5:O:100(A):ILE:HG22	5:O:100(J):PHE:HB3	1.70	0.72
6:N:11:LEU:HD21	6:N:102:THR:HA	1.72	0.72
2:B:544:LEU:CD1	2:J:588:GLN:OE1	2.38	0.71
5:M:100(A):ILE:HG22	5:M:100(J):PHE:HB3	1.70	0.71
1:E:104:MET:O	1:E:108:ILE:HG12	1.90	0.71
1:A:104:MET:O	1:A:108:ILE:HG12	1.90	0.71
5:Q:47:TRP:HB3	5:Q:60:ASN:HD22	1.56	0.71
1:I:104:MET:O	1:I:108:ILE:HG12	1.90	0.70
3:G:6:GLN:H	3:G:105:GLN:HE22	1.39	0.70
2:B:606:THR:HG21	2:B:646:LEU:HD23	1.73	0.70
3:C:6:GLN:H	3:C:105:GLN:HE22	1.39	0.70
5:M:47:TRP:HB3	5:M:60:ASN:HD22	1.56	0.70
5:O:47:TRP:HB3	5:O:60:ASN:HD22	1.56	0.70
3:K:6:GLN:H	3:K:105:GLN:HE22	1.39	0.70
2:F:606:THR:HG21	2:F:646:LEU:HD23	1.73	0.70
2:J:606:THR:HG21	2:J:646:LEU:HD23	1.73	0.70
1:I:203:GLN:O	1:I:204:ALA:O	2.10	0.69
1:E:253:PRO:HA	1:E:479:TRP:HE1	1.57	0.69
1:A:253:PRO:HA	1:A:479:TRP:HE1	1.58	0.69
1:A:503:ARG:HG2	2:J:661:LEU:HD11	1.75	0.69
1:I:253:PRO:HA	1:I:479:TRP:HE1	1.57	0.69
1:A:203:GLN:O	1:A:204:ALA:O	2.10	0.69
1:E:203:GLN:O	1:E:204:ALA:O	2.10	0.69
2:F:522:PHE:HE1	2:F:541:ALA:O	1.76	0.69
3:G:66:ARG:HG2	3:G:66:ARG:HH11	1.58	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:427:TRP:HZ3	1:I:475:MET:CE	1.88	0.68
3:C:66:ARG:HG2	3:C:66:ARG:HH11	1.59	0.68
3:K:66:ARG:HH11	3:K:66:ARG:HG2	1.59	0.68
1:E:58:ALA:HB2	1:E:77:THR:CG2	2.24	0.68
1:I:58:ALA:HB2	1:I:77:THR:CG2	2.24	0.68
2:B:651:THR:OG1	2:F:538:THR:CG2	2.31	0.68
1:A:58:ALA:HB2	1:A:77:THR:CG2	2.24	0.68
2:F:592:PHE:CE1	2:J:544:LEU:CD1	2.78	0.67
1:I:55:ALA:HB3	1:I:216:HIS:HB2	1.76	0.67
2:B:522:PHE:HE1	2:B:541:ALA:O	1.77	0.67
2:F:533:ALA:O	2:F:537:LEU:N	2.25	0.67
2:J:522:PHE:HE1	2:J:541:ALA:O	1.76	0.66
1:A:55:ALA:HB3	1:A:216:HIS:HB2	1.76	0.66
1:E:55:ALA:HB3	1:E:216:HIS:HB2	1.76	0.66
2:F:644:ASP:OD1	2:J:543:GLN:OE1	2.13	0.66
5:Q:39:GLN:NE2	5:Q:43:LYS:O	2.30	0.65
2:F:543:GLN:C	2:F:544:LEU:O	2.40	0.65
2:B:588:GLN:CD	2:F:544:LEU:CD1	2.70	0.65
5:O:47:TRP:HE1	5:O:50:TYR:HD1	1.45	0.65
5:O:39:GLN:NE2	5:O:43:LYS:O	2.30	0.64
5:M:39:GLN:NE2	5:M:43:LYS:O	2.30	0.64
2:J:533:ALA:O	2:J:537:LEU:N	2.25	0.64
1:A:474:ASP:OD1	1:A:474:ASP:N	2.31	0.64
2:J:543:GLN:C	2:J:544:LEU:O	2.40	0.64
2:F:522:PHE:CE1	2:F:541:ALA:O	2.51	0.64
2:J:522:PHE:CE1	2:J:541:ALA:O	2.51	0.64
2:J:543:GLN:O	2:J:544:LEU:C	2.42	0.63
1:I:220:PRO:HB3	2:J:578:ALA:HB1	1.80	0.63
2:B:543:GLN:C	2:B:544:LEU:O	2.40	0.63
2:F:592:PHE:CE1	2:J:544:LEU:HD13	2.33	0.63
5:Q:47:TRP:HE1	5:Q:50:TYR:HD1	1.45	0.63
4:H:81:GLU:OE1	4:H:81:GLU:N	2.27	0.63
2:B:543:GLN:O	2:B:544:LEU:C	2.41	0.63
2:B:533:ALA:O	2:B:537:LEU:N	2.27	0.63
5:M:47:TRP:HE1	5:M:50:TYR:HD1	1.45	0.63
1:E:128:THR:OG1	1:I:167:ASP:OD2	2.15	0.62
2:J:611:ASN:HB3	12:J:700:NAG:HN2	1.64	0.62
2:F:543:GLN:O	2:F:544:LEU:C	2.41	0.62
1:A:220:PRO:HB3	2:B:578:ALA:HB1	1.80	0.62
1:A:261:LEU:HD13	8:Y:1:NAG:H82	1.81	0.62
2:B:522:PHE:CE1	2:B:541:ALA:O	2.52	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:220:PRO:HB3	2:F:578:ALA:HB1	1.80	0.62
1:E:261:LEU:HD13	8:h:1:NAG:H82	1.82	0.61
2:B:611:ASN:HB3	12:B:700:NAG:HN2	1.64	0.61
2:B:544:LEU:HD13	2:J:588:GLN:NE2	2.15	0.61
1:E:270:ILE:HD13	1:E:289:ASN:H	1.66	0.61
1:A:270:ILE:HD13	1:A:289:ASN:H	1.66	0.61
2:B:543:GLN:OE1	2:B:543:GLN:HA	2.01	0.61
2:F:592:PHE:HE1	2:J:544:LEU:CD1	2.14	0.61
2:F:611:ASN:HB3	12:F:700:NAG:HN2	1.64	0.61
4:L:37:GLN:HB2	4:L:47:LEU:HD11	1.83	0.61
1:A:421:LYS:NZ	1:A:423:ILE:O	2.34	0.61
5:Q:53:ASP:O	5:Q:71:ARG:NH2	2.34	0.61
4:D:37:GLN:HB2	4:D:47:LEU:HD11	1.83	0.61
1:I:421:LYS:NZ	1:I:423:ILE:O	2.34	0.60
2:J:543:GLN:OE1	2:J:543:GLN:HA	2.01	0.60
2:B:544:LEU:CD1	2:J:588:GLN:CD	2.74	0.60
5:M:53:ASP:O	5:M:71:ARG:NH2	2.34	0.60
1:E:421:LYS:NZ	1:E:423:ILE:O	2.34	0.60
4:H:37:GLN:HB2	4:H:47:LEU:HD11	1.83	0.60
2:B:592:PHE:CE1	2:F:544:LEU:CD1	2.85	0.60
1:I:270:ILE:HD13	1:I:289:ASN:H	1.66	0.60
1:I:261:LEU:HD13	8:r:1:NAG:H82	1.83	0.60
4:H:6:GLN:HB3	4:H:23:CYS:HB3	1.84	0.60
5:O:53:ASP:O	5:O:71:ARG:NH2	2.34	0.60
2:F:543:GLN:OE1	2:F:543:GLN:HA	2.01	0.59
5:M:7:SER:OG	5:M:21:THR:OG1	2.19	0.59
4:D:81:GLU:OE1	4:D:81:GLU:N	2.27	0.59
1:I:68:VAL:HG11	1:I:112:TRP:CE3	2.38	0.59
1:A:503:ARG:HB2	2:J:661:LEU:HD21	1.83	0.59
2:F:592:PHE:CE1	2:J:544:LEU:HD12	2.38	0.59
1:I:278:THR:O	1:I:456:ARG:NH2	2.36	0.59
3:K:66:ARG:HH11	3:K:66:ARG:CG	2.16	0.59
1:A:68:VAL:HG11	1:A:112:TRP:CE3	2.38	0.59
1:A:278:THR:O	1:A:456:ARG:NH2	2.36	0.59
2:B:580:VAL:HG21	2:F:576:LEU:HD11	1.84	0.59
1:E:68:VAL:HG11	1:E:112:TRP:CE3	2.37	0.59
4:L:6:GLN:HB3	4:L:23:CYS:HB3	1.84	0.59
5:Q:37:ILE:HG22	5:Q:91:TYR:HB2	1.85	0.59
2:B:592:PHE:HE1	2:F:544:LEU:CD1	2.16	0.59
3:C:66:ARG:HH11	3:C:66:ARG:CG	2.16	0.59
1:E:503:ARG:HB3	1:E:504:ARG:HG2	1.85	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:87:THR:HG23	5:M:110:THR:HA	1.85	0.58
1:E:58:ALA:HB2	1:E:77:THR:HG21	1.85	0.58
1:I:58:ALA:HB2	1:I:77:THR:HG21	1.85	0.58
5:Q:6:GLU:HB3	5:Q:107:THR:HG23	1.84	0.58
1:A:192:ARG:NH1	1:A:193:LEU:O	2.37	0.58
5:M:6:GLU:HB3	5:M:107:THR:HG23	1.84	0.58
3:G:66:ARG:HH11	3:G:66:ARG:CG	2.16	0.58
5:Q:100:ARG:HH11	9:n:5:MAN:H62	1.68	0.58
2:B:644:ASP:O	2:B:648:ILE:HG12	2.04	0.58
1:E:278:THR:O	1:E:456:ARG:NH2	2.36	0.58
2:F:651:THR:OG1	2:J:538:THR:CG2	2.43	0.58
5:O:6:GLU:HB3	5:O:107:THR:HG23	1.84	0.58
5:O:37:ILE:HG22	5:O:91:TYR:HB2	1.85	0.58
2:J:538:THR:HG22	2:J:538:THR:O	2.04	0.58
4:D:6:GLN:HB3	4:D:23:CYS:HB3	1.84	0.58
1:E:192:ARG:NH1	1:E:193:LEU:O	2.37	0.58
2:J:535:ILE:O	2:J:539:VAL:HG23	2.03	0.58
2:F:661:LEU:HD21	1:I:503:ARG:CB	2.34	0.58
3:G:83:THR:OG1	3:G:85:ASP:OD1	2.20	0.58
5:M:37:ILE:HG22	5:M:91:TYR:HB2	1.85	0.58
1:I:192:ARG:NH1	1:I:193:LEU:O	2.37	0.58
5:O:87:THR:HG23	5:O:110:THR:HA	1.85	0.57
6:P:18:THR:HG21	6:P:76:SER:HA	1.86	0.57
3:K:40:ALA:HB3	3:K:43:GLN:HB3	1.86	0.57
2:B:538:THR:HG22	2:B:538:THR:O	2.04	0.57
5:Q:87:THR:HG23	5:Q:110:THR:HA	1.86	0.57
1:A:503:ARG:HB3	1:A:504:ARG:HG2	1.85	0.57
2:F:535:ILE:O	2:F:539:VAL:HG23	2.03	0.57
4:L:81:GLU:OE1	4:L:81:GLU:N	2.27	0.57
2:F:644:ASP:O	2:F:648:ILE:HG12	2.04	0.57
2:B:544:LEU:HD13	2:J:592:PHE:HE1	1.70	0.57
1:I:474:ASP:OD1	1:I:474:ASP:N	2.31	0.57
6:P:83:GLU:HA	6:P:104:LEU:HD22	1.87	0.57
1:A:58:ALA:HB2	1:A:77:THR:HG21	1.85	0.57
2:B:535:ILE:O	2:B:539:VAL:HG23	2.05	0.57
1:E:297:THR:HG23	1:E:299:PRO:HD3	1.87	0.57
1:E:381:GLU:HB3	1:E:420:ILE:HD11	1.87	0.57
2:F:538:THR:HG22	2:F:538:THR:O	2.04	0.57
2:J:644:ASP:O	2:J:648:ILE:HG12	2.04	0.57
6:N:18:THR:HG21	6:N:76:SER:HA	1.86	0.56
6:N:83:GLU:HA	6:N:104:LEU:HD22	1.87	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:363:PRO:O	1:E:469:ARG:NH1	2.38	0.56
1:I:503:ARG:HB3	1:I:504:ARG:HG2	1.85	0.56
6:R:18:THR:HG21	6:R:76:SER:HA	1.86	0.56
6:R:83:GLU:HA	6:R:104:LEU:HD22	1.87	0.56
1:E:496:ILE:O	2:F:631:TRP:NE1	2.35	0.56
3:C:40:ALA:HB3	3:C:43:GLN:HB3	1.86	0.56
5:O:100:ARG:HH11	9:d:5:MAN:H62	1.69	0.56
1:I:95:MET:HE1	1:I:273:ARG:HD3	1.88	0.56
3:K:74:LEU:HD23	3:K:77(C):PRO:HD3	1.87	0.56
5:O:82:LEU:HG	5:O:82(C):VAL:HG12	1.88	0.56
5:Q:24:VAL:O	5:Q:76:ASN:ND2	2.39	0.56
2:B:536:THR:C	2:B:538:THR:H	2.13	0.56
5:M:100:ARG:HH11	9:U:5:MAN:H62	1.69	0.56
1:E:95:MET:HE1	1:E:273:ARG:HD3	1.88	0.56
2:J:536:THR:C	2:J:538:THR:H	2.14	0.56
1:A:227:ARG:HG2	1:A:229:ASN:ND2	2.21	0.56
1:A:363:PRO:O	1:A:469:ARG:NH1	2.38	0.56
1:I:297:THR:HG23	1:I:299:PRO:HD3	1.87	0.56
3:C:83:THR:OG1	3:C:85:ASP:OD1	2.20	0.56
3:G:40:ALA:HB3	3:G:43:GLN:HB3	1.86	0.56
1:I:381:GLU:HB3	1:I:420:ILE:HD11	1.87	0.56
5:Q:65:SER:O	5:Q:65:SER:OG	2.23	0.56
1:A:297:THR:HG23	1:A:299:PRO:HD3	1.87	0.56
1:A:95:MET:HE1	1:A:273:ARG:HD3	1.88	0.56
1:I:227:ARG:HG2	1:I:229:ASN:ND2	2.21	0.56
1:I:363:PRO:O	1:I:469:ARG:NH1	2.38	0.56
3:K:83:THR:OG1	3:K:85:ASP:OD1	2.20	0.56
2:B:620:GLU:H	2:B:620:GLU:CD	2.09	0.55
1:A:381:GLU:HB3	1:A:420:ILE:HD11	1.87	0.55
3:C:74:LEU:HD23	3:C:77(C):PRO:HD3	1.87	0.55
1:E:227:ARG:HG2	1:E:229:ASN:ND2	2.21	0.55
2:F:536:THR:C	2:F:538:THR:H	2.14	0.55
3:G:74:LEU:HD23	3:G:77(C):PRO:HD3	1.87	0.55
5:O:24:VAL:O	5:O:76:ASN:ND2	2.39	0.55
5:O:13:LYS:HG3	5:O:16:GLU:HG3	1.89	0.55
4:D:32:TRP:HB2	4:D:92:ASP:HB3	1.88	0.55
5:O:7:SER:OG	5:O:21:THR:OG1	2.19	0.55
1:E:168:LYS:HA	1:E:168:LYS:HE3	1.89	0.55
4:H:32:TRP:HB2	4:H:92:ASP:HB3	1.88	0.55
1:I:66:HIS:CE1	1:I:72:HIS:CE1	2.95	0.55
1:I:496:ILE:O	2:J:631:TRP:NE1	2.35	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:24:VAL:O	5:M:76:ASN:ND2	2.39	0.55
5:M:82:LEU:HG	5:M:82(C):VAL:HG12	1.87	0.55
1:E:66:HIS:CE1	1:E:72:HIS:CE1	2.95	0.55
2:F:588:GLN:OE1	2:J:544:LEU:HD11	2.07	0.54
2:B:588:GLN:HE22	2:F:544:LEU:HD22	1.71	0.54
2:B:601:LYS:HB2	2:B:601:LYS:NZ	2.23	0.54
3:G:100:ARG:HB3	3:G:100:ARG:NH1	2.22	0.54
1:I:65:VAL:HG21	8:r:4:MAN:H4	1.88	0.54
2:F:601:LYS:HB2	2:F:601:LYS:NZ	2.23	0.54
5:Q:82:LEU:HG	5:Q:82(C):VAL:HG12	1.88	0.54
1:A:66:HIS:CE1	1:A:72:HIS:CE1	2.95	0.54
4:L:32:TRP:HB2	4:L:92:ASP:HB3	1.89	0.54
1:A:128:THR:OG1	1:E:167:ASP:OD2	2.22	0.54
2:B:588:GLN:NE2	2:F:544:LEU:HD13	2.22	0.54
1:I:260:LEU:HD12	1:I:451:GLY:HA3	1.90	0.54
1:E:265:LEU:HD11	1:E:291:SER:HB3	1.90	0.54
6:P:32:ALA:HB3	6:P:91:TRP:HB2	1.90	0.54
2:J:601:LYS:HB2	2:J:601:LYS:NZ	2.22	0.54
1:A:76:PRO:HD3	2:B:575:GLN:NE2	2.23	0.54
1:I:168:LYS:HE3	1:I:168:LYS:HA	1.89	0.54
1:A:168:LYS:HE3	1:A:168:LYS:HA	1.89	0.54
1:A:265:LEU:HD11	1:A:291:SER:HB3	1.90	0.54
1:E:65:VAL:HG21	8:h:4:MAN:H4	1.88	0.54
1:A:496:ILE:O	2:B:631:TRP:NE1	2.35	0.54
6:R:32:ALA:HB3	6:R:91:TRP:HB2	1.90	0.54
1:E:427:TRP:HD1	1:E:431:GLY:HA3	1.72	0.54
1:A:386:ASN:HB3	1:A:388:THR:HG23	1.90	0.53
5:M:13:LYS:HG3	5:M:16:GLU:HG3	1.89	0.53
1:E:260:LEU:HD12	1:E:451:GLY:HA3	1.90	0.53
1:I:265:LEU:HD11	1:I:291:SER:HB3	1.90	0.53
3:K:100:ARG:NH1	3:K:100:ARG:HB3	2.22	0.53
1:I:386:ASN:HB3	1:I:388:THR:HG23	1.90	0.53
5:Q:13:LYS:HG3	5:Q:16:GLU:HG3	1.89	0.53
5:M:65:SER:O	5:M:65:SER:OG	2.23	0.53
5:O:47:TRP:HZ3	5:O:100(P):MET:HE3	1.74	0.53
1:A:37:THR:OG1	1:A:499:THR:OG1	2.20	0.53
1:A:446:LEU:HD12	8:Y:2:NAG:H81	1.90	0.53
5:M:51:ILE:HD11	5:M:55:GLU:HA	1.90	0.53
3:C:100:ARG:HB3	3:C:100:ARG:NH1	2.22	0.53
1:I:76:PRO:HD3	2:J:575:GLN:NE2	2.23	0.53
1:A:102:GLU:O	1:A:106:GLU:HG2	2.09	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:N:32:ALA:HB3	6:N:91:TRP:HB2	1.90	0.53
1:E:76:PRO:HD3	2:F:575:GLN:NE2	2.23	0.53
1:I:102:GLU:O	1:I:106:GLU:HG2	2.09	0.53
1:E:386:ASN:HB3	1:E:388:THR:HG23	1.90	0.53
5:O:51:ILE:HD11	5:O:55:GLU:HA	1.90	0.53
1:A:260:LEU:HD12	1:A:451:GLY:HA3	1.90	0.52
1:E:446:LEU:HD12	8:h:2:NAG:H81	1.91	0.52
1:E:474:ASP:OD1	1:E:474:ASP:N	2.31	0.52
2:F:587:LEU:O	2:F:591:LYS:HG3	2.10	0.52
5:Q:51:ILE:HD11	5:Q:55:GLU:HA	1.90	0.52
1:A:427:TRP:HD1	1:A:431:GLY:HA3	1.72	0.52
1:E:102:GLU:O	1:E:106:GLU:HG2	2.09	0.52
5:M:29:MET:HE1	5:M:73:THR:HG21	1.92	0.52
1:I:427:TRP:HD1	1:I:431:GLY:HA3	1.72	0.52
5:O:29:MET:HE1	5:O:73:THR:HG21	1.92	0.52
5:Q:47:TRP:HZ3	5:Q:100(P):MET:HE3	1.74	0.52
1:E:128:THR:N	1:I:167:ASP:OD2	2.34	0.52
5:O:100(M):TYR:HB2	6:P:91:TRP:CD1	2.45	0.52
5:O:18:LEU:HB3	5:O:82:LEU:HB3	1.92	0.52
1:A:379:ARG:HB3	1:A:443:ILE:HD12	1.92	0.52
5:M:18:LEU:HB3	5:M:82:LEU:HB3	1.92	0.52
2:F:620:GLU:H	2:F:620:GLU:CD	2.09	0.52
2:B:587:LEU:O	2:B:591:LYS:HG3	2.10	0.52
1:I:379:ARG:HB3	1:I:443:ILE:HD12	1.92	0.52
5:Q:18:LEU:HB3	5:Q:82:LEU:HB3	1.92	0.52
5:Q:100(M):TYR:HB2	6:R:91:TRP:CD1	2.45	0.52
5:M:100(M):TYR:HB2	6:N:91:TRP:CD1	2.45	0.51
2:J:587:LEU:O	2:J:591:LYS:HG3	2.10	0.51
1:A:65:VAL:HG21	8:Y:4:MAN:H4	1.92	0.51
5:M:47:TRP:HZ3	5:M:100(P):MET:HE3	1.74	0.51
1:E:216:HIS:HB3	1:E:247:CYS:HB3	1.92	0.51
3:G:9:THR:OG1	3:G:108:LEU:O	2.26	0.51
5:Q:77:GLN:O	5:Q:78:LEU:HD23	2.10	0.51
2:B:651:THR:CB	2:F:538:THR:HG21	2.37	0.51
1:E:379:ARG:HB3	1:E:443:ILE:HD12	1.92	0.51
5:Q:67:VAL:HB	5:Q:82:LEU:HD13	1.93	0.51
1:A:216:HIS:HB3	1:A:247:CYS:HB3	1.92	0.51
2:B:637:ASN:HB3	11:Z:1:NAG:O5	2.11	0.51
5:O:20:VAL:HB	5:O:80:LEU:HD12	1.93	0.51
2:B:592:PHE:CE1	2:F:544:LEU:HD12	2.46	0.51
5:M:77:GLN:O	5:M:78:LEU:HD23	2.11	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:O:67:VAL:HB	5:O:82:LEU:HD13	1.93	0.51
5:O:77:GLN:O	5:O:78:LEU:HD23	2.10	0.51
1:I:216:HIS:HB3	1:I:247:CYS:HB3	1.92	0.51
5:Q:20:VAL:HB	5:Q:80:LEU:HD12	1.93	0.51
1:A:158:SER:HB2	1:A:171:LYS:HE2	1.93	0.51
5:Q:100:ARG:NH1	9:n:5:MAN:H62	2.26	0.51
4:D:79:GLN:HG3	4:D:80:ALA:H	1.75	0.50
5:Q:29:MET:HE1	5:Q:73:THR:HG21	1.92	0.50
1:A:503:ARG:CB	2:J:661:LEU:HD21	2.42	0.50
5:M:20:VAL:HB	5:M:80:LEU:HD12	1.93	0.50
4:H:79:GLN:HG3	4:H:80:ALA:H	1.75	0.50
4:L:79:GLN:HG3	4:L:80:ALA:H	1.75	0.50
1:E:205:CYS:SG	1:E:206:PRO:HD2	2.52	0.50
2:F:637:ASN:HB3	11:i:1:NAG:O5	2.12	0.50
6:R:31:ARG:NH2	6:R:69:THR:OG1	2.45	0.50
1:A:177:TYR:HE2	1:A:422:GLN:HE21	1.60	0.50
5:O:100:ARG:NH1	9:d:5:MAN:H62	2.27	0.50
5:M:67:VAL:HB	5:M:82:LEU:HD13	1.93	0.50
1:I:78:ASP:OD1	1:I:78:ASP:N	2.45	0.50
1:A:60:ALA:O	1:A:63:THR:O	2.30	0.49
1:A:382:PHE:O	1:A:420:ILE:HG13	2.12	0.49
1:E:158:SER:HB2	1:E:171:LYS:HE2	1.93	0.49
2:J:637:ASN:HB3	11:s:1:NAG:O5	2.12	0.49
6:P:31:ARG:NH2	6:P:69:THR:OG1	2.45	0.49
1:I:158:SER:HB2	1:I:171:LYS:HE2	1.93	0.49
1:I:177:TYR:HE2	1:I:422:GLN:HE21	1.60	0.49
1:I:382:PHE:O	1:I:420:ILE:HG13	2.12	0.49
2:B:618:THR:O	2:B:621:GLU:HG2	2.12	0.49
1:I:60:ALA:O	1:I:63:THR:O	2.30	0.49
1:A:205:CYS:SG	1:A:206:PRO:HD2	2.52	0.49
3:C:81:GLU:OE1	7:S:6:MAN:O3	2.27	0.49
2:J:618:THR:O	2:J:621:GLU:HG2	2.12	0.49
1:A:426:MET:HE3	1:A:427:TRP:CZ2	2.47	0.49
1:E:385:CYS:HA	1:E:418:CYS:HA	1.95	0.49
1:I:205:CYS:SG	1:I:206:PRO:HD2	2.52	0.49
2:J:522:PHE:CZ	2:J:542:ARG:NH1	2.81	0.49
1:E:37:THR:OG1	1:E:499:THR:OG1	2.20	0.49
1:E:55:ALA:O	1:E:216:HIS:HB2	2.13	0.49
1:E:382:PHE:O	1:E:420:ILE:HG13	2.12	0.49
1:A:55:ALA:O	1:A:216:HIS:HB2	2.13	0.49
1:E:78:ASP:N	1:E:78:ASP:OD1	2.45	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:177:TYR:HE2	1:E:422:GLN:HE21	1.60	0.49
1:I:426:MET:HE3	1:I:427:TRP:CZ2	2.47	0.49
2:J:537:LEU:O	2:J:537:LEU:HG	2.13	0.49
2:B:540:GLN:O	2:B:541:ALA:HB3	2.13	0.49
5:M:47:TRP:HE3	6:N:98:PHE:HZ	1.61	0.49
5:Q:47:TRP:HE3	6:R:98:PHE:HZ	1.61	0.49
6:R:66(A):ASP:HB3	9:n:5:MAN:H3	1.95	0.49
6:N:31:ARG:NH2	6:N:69:THR:OG1	2.45	0.49
1:E:60:ALA:O	1:E:63:THR:O	2.30	0.49
1:E:276:ASN:O	1:E:279:ASN:N	2.46	0.49
2:F:522:PHE:CZ	2:F:542:ARG:NH1	2.81	0.49
2:F:618:THR:O	2:F:621:GLU:HG2	2.12	0.49
4:H:65:SER:OG	4:H:72:THR:OG1	2.26	0.49
2:B:537:LEU:HG	2:B:537:LEU:O	2.13	0.48
1:A:78:ASP:OD1	1:A:78:ASP:N	2.45	0.48
10:o:1:NAG:O3	10:o:2:NAG:O5	2.29	0.48
1:E:426:MET:HE3	1:E:427:TRP:CZ2	2.47	0.48
1:A:276:ASN:O	1:A:279:ASN:N	2.46	0.48
5:O:47:TRP:HE3	6:P:98:PHE:HZ	1.61	0.48
1:I:229:ASN:HB3	10:z:1:NAG:H61	1.95	0.48
10:e:1:NAG:O3	10:e:2:NAG:O5	2.29	0.48
1:I:55:ALA:O	1:I:216:HIS:HB2	2.13	0.48
5:Q:4:LEU:HD12	5:Q:92:CYS:HB3	1.96	0.48
4:D:39:ARG:HE	4:D:40:PRO:HD2	1.78	0.48
1:E:42:VAL:HA	2:F:537:LEU:HD12	1.96	0.48
2:F:540:GLN:O	2:F:541:ALA:HB3	2.13	0.48
1:I:444:LYS:HE3	1:I:444:LYS:HB2	1.54	0.48
1:A:229:ASN:HB3	10:x:1:NAG:H61	1.94	0.48
5:M:47:TRP:HB3	5:M:60:ASN:ND2	2.27	0.48
6:N:35:TRP:HB2	6:N:47:LEU:HB2	1.96	0.48
5:M:4:LEU:HD12	5:M:92:CYS:HB3	1.96	0.48
5:M:100:ARG:NH1	9:U:5:MAN:H62	2.29	0.48
2:F:537:LEU:HG	2:F:537:LEU:O	2.13	0.48
4:H:38:GLN:HB2	4:H:44:PRO:HG3	1.96	0.48
4:H:39:ARG:HE	4:H:40:PRO:HD2	1.78	0.48
5:O:4:LEU:HD12	5:O:92:CYS:HB3	1.96	0.48
6:P:66(A):ASP:HB3	9:d:5:MAN:H3	1.96	0.48
1:A:385:CYS:HA	1:A:418:CYS:HA	1.95	0.47
3:C:100(E):HIS:HB3	3:C:100(J):MET:HE2	1.96	0.47
2:B:522:PHE:CZ	2:B:542:ARG:NH1	2.82	0.47
4:D:38:GLN:HB2	4:D:44:PRO:HG3	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:446:LEU:HD12	8:r:2:NAG:H81	1.96	0.47
5:M:103:TRP:CG	6:N:44:PRO:HD2	2.49	0.47
1:I:276:ASN:O	1:I:279:ASN:N	2.46	0.47
2:J:540:GLN:O	2:J:541:ALA:HB3	2.13	0.47
1:A:42:VAL:HA	2:B:537:LEU:HD12	1.96	0.47
3:C:35(A):ASN:O	3:C:93:THR:OG1	2.32	0.47
1:E:427:TRP:CD1	1:E:431:GLY:HA3	2.49	0.47
1:I:358:ILE:HD12	1:I:468:PHE:HE2	1.80	0.47
1:I:427:TRP:CD1	1:I:431:GLY:HA3	2.49	0.47
3:K:100:ARG:HB3	3:K:100:ARG:HH11	1.79	0.47
2:B:654:GLU:O	2:B:657:GLU:HG3	2.14	0.47
2:F:654:GLU:O	2:F:657:GLU:HG3	2.14	0.47
3:G:100:ARG:HB3	3:G:100:ARG:HH11	1.79	0.47
3:G:100(E):HIS:HB3	3:G:100(J):MET:HE2	1.96	0.47
3:K:100(E):HIS:HB3	3:K:100(J):MET:HE2	1.96	0.47
4:L:39:ARG:HE	4:L:40:PRO:HD2	1.78	0.47
1:A:121:LYS:HB2	1:A:121:LYS:HE2	1.73	0.47
3:C:100:ARG:HB3	3:C:100:ARG:HH11	1.79	0.47
1:I:385:CYS:HA	1:I:418:CYS:HA	1.95	0.47
2:J:654:GLU:O	2:J:657:GLU:HG3	2.14	0.47
5:Q:103:TRP:CG	6:R:44:PRO:HD2	2.49	0.47
1:A:358:ILE:HD12	1:A:468:PHE:HE2	1.80	0.47
6:N:15:LEU:HD12	6:N:106:VAL:HG23	1.97	0.47
1:I:350:LYS:HB3	1:I:355:ASN:OD1	2.15	0.47
5:Q:51:ILE:HG21	5:Q:70:SER:O	2.14	0.47
1:A:427:TRP:HE3	1:A:475:MET:SD	2.13	0.47
1:A:427:TRP:CD1	1:A:431:GLY:HA3	2.49	0.47
1:A:444:LYS:HB2	1:A:444:LYS:HE3	1.54	0.47
5:M:51:ILE:HG21	5:M:70:SER:O	2.14	0.47
1:E:66:HIS:HE1	1:E:72:HIS:CE1	2.33	0.47
3:G:25:TYR:CE1	3:G:77(B):PRO:HG3	2.50	0.47
3:K:25:TYR:CE1	3:K:77(B):PRO:HG3	2.50	0.47
4:L:38:GLN:HB2	4:L:44:PRO:HG3	1.96	0.47
5:M:21:THR:HA	5:M:78:LEU:O	2.15	0.47
1:E:229:ASN:HB3	10:y:1:NAG:H61	1.96	0.47
5:O:21:THR:HA	5:O:78:LEU:O	2.15	0.47
5:O:51:ILE:HG21	5:O:70:SER:O	2.14	0.47
5:O:103:TRP:CG	6:P:44:PRO:HD2	2.49	0.47
1:I:503:ARG:HE	1:I:504:ARG:H	1.63	0.47
5:Q:21:THR:HA	5:Q:78:LEU:O	2.15	0.47
6:R:15:LEU:HD12	6:R:106:VAL:HG23	1.97	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:N:66(A):ASP:HB3	9:U:5:MAN:H3	1.97	0.46
1:I:42:VAL:HA	2:J:537:LEU:HD12	1.96	0.46
6:R:35:TRP:HB2	6:R:47:LEU:HB2	1.96	0.46
5:Q:47:TRP:HB3	5:Q:60:ASN:ND2	2.27	0.46
2:B:536:THR:C	2:B:538:THR:N	2.73	0.46
1:I:37:THR:OG1	1:I:499:THR:OG1	2.20	0.46
10:V:1:NAG:O3	10:V:2:NAG:O5	2.29	0.46
1:E:350:LYS:HB3	1:E:355:ASN:OD1	2.15	0.46
1:A:503:ARG:HE	1:A:504:ARG:H	1.63	0.46
4:D:52:ALA:O	11:Z:2:NAG:O6	2.33	0.46
6:P:35:TRP:HB2	6:P:47:LEU:HB2	1.96	0.46
2:B:544:LEU:CD1	2:J:592:PHE:HE1	2.28	0.46
3:C:25:TYR:CE1	3:C:77(B):PRO:HG3	2.50	0.46
3:C:93:THR:HB	3:C:100(L):PHE:HB3	1.98	0.46
5:Q:76:ASN:O	5:Q:77:GLN:HG3	2.16	0.46
2:B:544:LEU:CD1	2:J:592:PHE:CE1	2.99	0.46
1:E:358:ILE:HD12	1:E:468:PHE:HE2	1.80	0.46
1:E:444:LYS:HB2	1:E:444:LYS:HE3	1.54	0.46
6:R:60:GLU:CD	6:R:60:GLU:H	2.24	0.46
1:A:66:HIS:HE1	1:A:72:HIS:CE1	2.33	0.46
5:M:76:ASN:O	5:M:77:GLN:HG3	2.16	0.46
1:E:333:ILE:HG13	1:E:390:LEU:HD21	1.97	0.46
3:G:93:THR:HB	3:G:100(L):PHE:HB3	1.98	0.46
5:Q:12:VAL:HG11	5:Q:18:LEU:HD13	1.98	0.46
1:E:321(A):ASP:N	1:E:321(A):ASP:OD1	2.48	0.46
2:F:544:LEU:HD23	2:F:544:LEU:HA	1.74	0.46
1:I:321(A):ASP:OD1	1:I:321(A):ASP:N	2.48	0.46
1:A:333:ILE:HG13	1:A:390:LEU:HD21	1.97	0.45
2:B:542:ARG:HH11	2:B:542:ARG:CB	2.22	0.45
5:M:12:VAL:HG11	5:M:18:LEU:HD13	1.98	0.45
1:E:503:ARG:HE	1:E:504:ARG:H	1.63	0.45
2:F:587:LEU:HD11	2:J:583:VAL:HG13	1.97	0.45
2:F:658:LYS:HD2	2:J:603:ILE:HG21	1.98	0.45
3:G:35(A):ASN:O	3:G:93:THR:OG1	2.32	0.45
1:I:82:GLN:OE1	1:I:82:GLN:HA	2.16	0.45
1:A:350:LYS:HB3	1:A:355:ASN:OD1	2.15	0.45
3:K:35(A):ASN:O	3:K:93:THR:OG1	2.32	0.45
3:K:100:ARG:HH11	3:K:100:ARG:CB	2.29	0.45
5:O:47:TRP:HB3	5:O:60:ASN:ND2	2.27	0.45
3:C:100:ARG:HH11	3:C:100:ARG:CB	2.29	0.45
1:I:309:ILE:HD13	1:I:317:PHE:HB2	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:536:THR:C	2:J:538:THR:N	2.74	0.45
3:K:93:THR:HB	3:K:100(L):PHE:HB3	1.98	0.45
6:R:23:CYS:HB3	6:R:88:CYS:HB2	1.82	0.45
1:A:34:LEU:HB3	1:A:498:PRO:HB2	1.99	0.45
3:G:73:ASP:CG	3:G:74:LEU:H	2.25	0.45
5:O:76:ASN:O	5:O:77:GLN:HG3	2.16	0.45
1:I:333:ILE:HG13	1:I:390:LEU:HD21	1.97	0.45
1:A:368:ASP:O	1:A:372:THR:HG22	2.17	0.45
1:A:408:THR:O	1:A:410:ASP:N	2.47	0.45
2:B:532:ALA:HA	2:B:535:ILE:HG12	1.98	0.45
3:G:28:ASN:OD1	3:G:29:THR:N	2.46	0.45
1:A:321(A):ASP:OD1	1:A:321(A):ASP:N	2.48	0.45
1:E:34:LEU:HB3	1:E:498:PRO:HB2	1.99	0.45
2:F:584:GLU:OE1	2:J:579:ARG:HD2	2.16	0.45
6:P:15:LEU:HD12	6:P:106:VAL:HG23	1.97	0.45
4:L:4:MET:HE1	4:L:90:GLN:HB3	1.99	0.45
1:A:411:ASN:OD1	1:A:411:ASN:N	2.50	0.45
3:C:9:THR:OG1	3:C:108:LEU:O	2.26	0.45
1:E:368:ASP:O	1:E:372:THR:HG22	2.17	0.45
3:G:100:ARG:HH11	3:G:100:ARG:CB	2.29	0.45
4:H:81:GLU:H	4:H:81:GLU:CD	2.22	0.45
1:I:66:HIS:HE1	1:I:72:HIS:CE1	2.33	0.45
1:I:203:GLN:C	1:I:204:ALA:O	2.60	0.45
1:I:411:ASN:OD1	1:I:411:ASN:N	2.50	0.45
4:L:51:GLY:HA3	11:s:1:NAG:H83	1.98	0.45
1:A:293:LYS:HA	1:A:448:ASN:HA	1.99	0.45
1:A:309:ILE:HD13	1:A:317:PHE:HB2	1.98	0.45
2:B:592:PHE:CE1	2:F:544:LEU:HD13	2.40	0.45
3:C:28:ASN:OD1	3:C:29:THR:N	2.46	0.45
6:N:60:GLU:CD	6:N:60:GLU:H	2.24	0.45
1:E:121:LYS:HB2	1:E:121:LYS:HE2	1.73	0.45
1:E:293:LYS:HA	1:E:448:ASN:HA	1.99	0.45
5:O:12:VAL:HG11	5:O:18:LEU:HD13	1.98	0.45
6:P:60:GLU:CD	6:P:60:GLU:H	2.24	0.45
4:L:50:ARG:HD2	4:L:50:ARG:HA	1.82	0.45
1:E:491:ILE:HG21	2:F:523:LEU:HD21	2.00	0.44
2:F:532:ALA:HA	2:F:535:ILE:HG12	1.98	0.44
4:D:4:MET:HE1	4:D:90:GLN:HB3	1.99	0.44
1:E:203:GLN:C	1:E:204:ALA:O	2.60	0.44
5:O:2:VAL:HG13	5:O:24:VAL:HG23	1.99	0.44
2:J:542:ARG:HH11	2:J:542:ARG:CB	2.22	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:494:LEU:HD23	1:E:494:LEU:HA	1.89	0.44
2:F:536:THR:C	2:F:538:THR:N	2.74	0.44
5:O:27:ASP:CG	5:O:31:ASN:HD21	2.25	0.44
1:I:293:LYS:HA	1:I:448:ASN:HA	1.99	0.44
3:K:28:ASN:OD1	3:K:29:THR:N	2.46	0.44
1:E:82:GLN:OE1	1:E:82:GLN:HA	2.16	0.44
1:E:240:LYS:HD2	1:E:240:LYS:HA	1.71	0.44
4:L:52:ALA:O	11:s:2:NAG:O6	2.34	0.44
1:A:82:GLN:HA	1:A:82:GLN:OE1	2.16	0.44
3:C:73:ASP:CG	3:C:74:LEU:H	2.25	0.44
5:M:27:ASP:OD1	5:M:31:ASN:ND2	2.40	0.44
4:H:4:MET:HE1	4:H:90:GLN:HB3	1.99	0.44
1:I:287:HIS:CD2	1:I:288:PHE:H	2.36	0.44
4:L:36:TYR:HB2	4:L:87:TYR:HB2	1.99	0.44
1:A:287:HIS:CD2	1:A:288:PHE:H	2.36	0.44
1:A:491:ILE:HG21	2:B:523:LEU:HD21	2.00	0.44
4:D:36:TYR:HB2	4:D:87:TYR:HB2	1.99	0.44
4:H:35:TRP:CZ3	4:H:88:CYS:HB3	2.53	0.44
5:O:22:CYS:SG	5:O:78:LEU:HD12	2.58	0.44
1:I:368:ASP:O	1:I:372:THR:HG22	2.17	0.44
1:E:253:PRO:HA	1:E:479:TRP:NE1	2.30	0.44
3:K:73:ASP:CG	3:K:74:LEU:H	2.25	0.44
5:Q:2:VAL:HG13	5:Q:24:VAL:HG23	1.99	0.44
6:R:92:ASP:OD1	6:R:93:SER:N	2.51	0.44
6:N:23:CYS:HB3	6:N:88:CYS:HB2	1.82	0.44
1:E:58:ALA:HB2	1:E:77:THR:HG23	1.98	0.44
1:E:287:HIS:CD2	1:E:288:PHE:H	2.36	0.44
6:P:92:ASP:OD1	6:P:93:SER:N	2.51	0.44
3:K:6:GLN:H	3:K:105:GLN:NE2	2.11	0.44
4:L:35:TRP:CZ3	4:L:88:CYS:HB3	2.53	0.44
2:F:591:LYS:HG2	2:J:586:TYR:CZ	2.52	0.43
4:H:36:TYR:HB2	4:H:87:TYR:HB2	1.99	0.43
4:D:35:TRP:CZ3	4:D:88:CYS:HB3	2.53	0.43
5:M:22:CYS:SG	5:M:78:LEU:HD12	2.58	0.43
1:E:503:ARG:NE	1:E:503:ARG:HA	2.34	0.43
1:I:34:LEU:HB3	1:I:498:PRO:HB2	1.99	0.43
2:J:532:ALA:HA	2:J:535:ILE:HG12	1.98	0.43
3:C:100:ARG:HA	3:C:100(E):HIS:CE1	2.53	0.43
5:M:31:ASN:OD1	5:M:32:TYR:N	2.51	0.43
1:E:426:MET:HE3	1:E:427:TRP:CE2	2.54	0.43
5:O:10:GLY:HA2	5:O:109:VAL:HA	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:491:ILE:HG21	2:J:523:LEU:HD21	2.00	0.43
3:K:100:ARG:HA	3:K:100(E):HIS:CE1	2.53	0.43
6:N:18:THR:HG22	6:N:19:ALA:N	2.33	0.43
6:N:65:THR:HG21	6:N:71:ALA:HA	2.00	0.43
4:H:79:GLN:HG3	4:H:80:ALA:N	2.34	0.43
1:E:309:ILE:HD13	1:E:317:PHE:HB2	1.98	0.43
5:M:27:ASP:CG	5:M:31:ASN:HD21	2.25	0.43
1:E:408:THR:O	1:E:410:ASP:N	2.47	0.43
2:F:619:PHE:HB3	2:F:620:GLU:OE1	2.19	0.43
1:I:240:LYS:HD2	1:I:240:LYS:HA	1.71	0.43
5:Q:10:GLY:HA2	5:Q:109:VAL:HA	2.00	0.43
1:A:426:MET:HE3	1:A:427:TRP:CE2	2.54	0.43
3:K:51:ILE:O	3:K:51:ILE:HG23	2.19	0.43
1:A:240:LYS:HA	1:A:240:LYS:HD2	1.71	0.43
1:A:253:PRO:HA	1:A:479:TRP:NE1	2.30	0.43
5:M:2:VAL:HG13	5:M:24:VAL:HG23	1.99	0.43
2:F:651:THR:CB	2:J:538:THR:HG21	2.46	0.43
1:I:328:LYS:HB2	1:I:328:LYS:HE3	1.81	0.43
1:I:426:MET:HE3	1:I:427:TRP:CE2	2.54	0.43
1:I:503:ARG:HA	1:I:503:ARG:NE	2.34	0.43
5:Q:22:CYS:SG	5:Q:78:LEU:HD12	2.58	0.43
5:Q:27:ASP:CG	5:Q:31:ASN:HD21	2.25	0.43
6:N:92:ASP:OD1	6:N:93:SER:N	2.51	0.43
2:J:597:GLY:CA	2:J:650:GLN:HE21	2.32	0.43
5:Q:33:TYR:N	5:Q:95:ALA:O	2.52	0.43
6:R:18:THR:HG22	6:R:19:ALA:N	2.33	0.43
6:R:65:THR:HG21	6:R:71:ALA:HA	2.00	0.43
1:A:287:HIS:HD2	1:A:288:PHE:H	1.66	0.42
2:B:544:LEU:HD22	2:J:588:GLN:HE22	1.84	0.42
6:N:14:ALA:HB1	6:N:106:VAL:HG23	2.00	0.42
1:E:261:LEU:HD23	1:E:261:LEU:HA	1.83	0.42
3:G:51:ILE:HG23	3:G:51:ILE:O	2.19	0.42
1:I:111:LEU:HD11	2:J:571:TRP:CZ2	2.54	0.42
2:J:619:PHE:HB3	2:J:620:GLU:OE1	2.19	0.42
1:A:203:GLN:C	1:A:204:ALA:O	2.60	0.42
6:P:101:ALA:O	6:P:103:ARG:NH1	2.53	0.42
1:I:494:LEU:HD23	1:I:494:LEU:HA	1.89	0.42
4:L:65:SER:OG	4:L:72:THR:OG1	2.26	0.42
1:A:407:GLU:O	1:A:408:THR:OG1	2.33	0.42
5:M:35:THR:HB	5:M:95:ALA:HB2	2.02	0.42
1:E:153:ASP:N	1:E:153:ASP:OD1	2.52	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:99:ASP:O	3:G:100(E):HIS:NE2	2.47	0.42
1:I:287:HIS:HD2	1:I:288:PHE:H	1.66	0.42
2:J:573:ILE:O	2:J:577:GLN:HG3	2.19	0.42
5:M:33:TYR:N	5:M:95:ALA:O	2.52	0.42
1:I:253:PRO:HA	1:I:479:TRP:NE1	2.30	0.42
6:R:14:ALA:HB1	6:R:106:VAL:HG23	2.00	0.42
2:B:584:GLU:OE1	2:F:579:ARG:HD2	2.19	0.42
2:B:619:PHE:HB3	2:B:620:GLU:OE1	2.19	0.42
1:E:111:LEU:HD11	2:F:571:TRP:CZ2	2.54	0.42
1:E:411:ASN:OD1	1:E:411:ASN:N	2.50	0.42
6:P:18:THR:HG22	6:P:19:ALA:N	2.34	0.42
1:I:475:MET:HE3	1:I:475:MET:HB3	1.95	0.42
4:L:94:TYR:HA	4:L:95:PRO:HA	1.84	0.42
1:A:58:ALA:HB2	1:A:77:THR:HG23	1.99	0.42
2:B:573:ILE:O	2:B:577:GLN:HG3	2.19	0.42
4:D:94:TYR:HA	4:D:95:PRO:HA	1.84	0.42
1:E:218:CYS:HA	1:E:246:GLN:O	2.20	0.42
1:E:287:HIS:HD2	1:E:288:PHE:H	1.66	0.42
2:F:641:GLN:O	2:F:645:ILE:HG12	2.20	0.42
10:v:1:NAG:H62	10:v:2:NAG:H82	2.00	0.42
1:A:100:MET:HE2	1:A:483:LEU:HD22	2.01	0.42
2:B:538:THR:HG21	2:J:651:THR:CB	2.49	0.42
2:B:545:LEU:H	2:B:545:LEU:HG	1.67	0.42
1:I:100:MET:HE2	1:I:483:LEU:HD22	2.01	0.42
3:K:73:ASP:OD1	7:j:1:NAG:H83	2.20	0.42
4:L:79:GLN:HG3	4:L:80:ALA:N	2.34	0.42
2:F:597:GLY:CA	2:F:650:GLN:HE21	2.32	0.42
3:G:6:GLN:H	3:G:105:GLN:NE2	2.11	0.42
5:O:35:THR:HB	5:O:95:ALA:HB2	2.02	0.42
6:P:14:ALA:HB1	6:P:106:VAL:HG23	2.00	0.42
6:P:35:TRP:CG	6:P:73:LEU:HD22	2.55	0.42
1:I:218:CYS:HA	1:I:246:GLN:O	2.20	0.42
5:Q:31:ASN:OD1	5:Q:32:TYR:N	2.51	0.42
6:R:80:ALA:HA	6:R:83:GLU:HB2	2.02	0.42
1:A:111:LEU:HD11	2:B:571:TRP:CZ2	2.54	0.42
1:A:218:CYS:HA	1:A:246:GLN:O	2.20	0.42
1:A:503:ARG:NE	1:A:503:ARG:HA	2.33	0.42
2:B:597:GLY:CA	2:B:650:GLN:HE21	2.32	0.42
3:C:51:ILE:HG23	3:C:51:ILE:O	2.19	0.42
4:D:79:GLN:HG3	4:D:80:ALA:N	2.34	0.42
5:M:10:GLY:HA2	5:M:109:VAL:HA	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:100:ARG:HA	3:G:100(E):HIS:CE1	2.53	0.42
5:O:65:SER:O	5:O:65:SER:OG	2.23	0.42
1:I:261:LEU:HD23	1:I:261:LEU:HA	1.83	0.42
1:I:276:ASN:HB3	1:I:282:LYS:HG3	2.02	0.42
2:J:641:GLN:O	2:J:645:ILE:HG12	2.20	0.42
5:O:31:ASN:OD1	5:O:32:TYR:N	2.51	0.42
6:P:65:THR:HG21	6:P:71:ALA:HA	2.00	0.42
6:P:83:GLU:O	6:P:104:LEU:HB2	2.20	0.42
1:I:121:LYS:HB2	1:I:121:LYS:HE2	1.73	0.42
6:R:35:TRP:CG	6:R:73:LEU:HD22	2.55	0.42
10:w:1:NAG:H62	10:w:2:NAG:H82	2.01	0.42
6:N:83:GLU:O	6:N:104:LEU:HB2	2.20	0.41
2:F:542:ARG:HH11	2:F:542:ARG:CB	2.22	0.41
2:F:573:ILE:O	2:F:577:GLN:HG3	2.19	0.41
2:F:601:LYS:HB2	2:F:601:LYS:HZ2	1.85	0.41
4:H:51:GLY:HA3	11:i:1:NAG:H83	2.01	0.41
5:Q:18:LEU:HD12	5:Q:19:SER:H	1.85	0.41
6:R:83:GLU:O	6:R:104:LEU:HB2	2.20	0.41
6:N:35:TRP:CG	6:N:73:LEU:HD22	2.55	0.41
6:P:80:ALA:HA	6:P:83:GLU:HB2	2.02	0.41
1:I:58:ALA:HB2	1:I:77:THR:HG23	1.98	0.41
1:A:193:LEU:HB2	1:A:196:CYS:SG	2.61	0.41
5:O:18:LEU:HD12	5:O:19:SER:H	1.85	0.41
5:O:33:TYR:N	5:O:95:ALA:O	2.52	0.41
2:J:545:LEU:H	2:J:545:LEU:HG	1.67	0.41
5:Q:35:THR:HB	5:Q:95:ALA:HB2	2.02	0.41
5:O:30:ASN:OD1	5:O:30:ASN:N	2.53	0.41
5:Q:48:ILE:HG22	5:Q:49:GLY:N	2.35	0.41
6:R:101:ALA:O	6:R:103:ARG:NH1	2.53	0.41
2:B:641:GLN:O	2:B:645:ILE:HG12	2.20	0.41
5:M:18:LEU:HD12	5:M:19:SER:H	1.85	0.41
6:N:101:ALA:O	6:N:103:ARG:NH1	2.53	0.41
1:I:100:MET:HE2	1:I:100:MET:HB3	1.91	0.41
5:Q:27:ASP:OD1	5:Q:31:ASN:ND2	2.40	0.41
1:A:276:ASN:HB3	1:A:282:LYS:HG3	2.02	0.41
1:E:100:MET:HE2	1:E:483:LEU:HD22	2.01	0.41
5:O:48:ILE:HG22	5:O:49:GLY:N	2.35	0.41
1:I:75:VAL:HB	2:J:575:GLN:HE22	1.85	0.41
2:J:538:THR:CG2	2:J:538:THR:O	2.69	0.41
1:A:153:ASP:OD1	1:A:153:ASP:N	2.52	0.41
1:A:232:LYS:HB2	1:A:232:LYS:HE2	1.86	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:544:LEU:HD12	2:J:592:PHE:CE1	2.56	0.41
4:D:50:ARG:HA	4:D:50:ARG:HD2	1.82	0.41
5:M:48:ILE:HG22	5:M:49:GLY:N	2.35	0.41
6:N:75:ILE:HD13	6:N:86:TYR:HE2	1.86	0.41
1:E:183:PRO:HA	1:E:191:TYR:CD1	2.56	0.41
3:G:81:GLU:OE1	7:a:6:MAN:O3	2.34	0.41
1:I:427:TRP:HE3	1:I:475:MET:SD	2.13	0.41
6:N:80:ALA:HA	6:N:83:GLU:HB2	2.02	0.41
1:A:75:VAL:HB	2:B:575:GLN:HE22	1.85	0.41
1:A:183:PRO:HA	1:A:191:TYR:CD1	2.56	0.41
2:B:601:LYS:HB2	2:B:601:LYS:HZ2	1.84	0.41
2:B:651:THR:CB	2:F:538:THR:CG2	2.98	0.41
1:E:193:LEU:HB2	1:E:196:CYS:SG	2.61	0.41
1:E:276:ASN:HB3	1:E:282:LYS:HG3	2.02	0.41
1:E:475:MET:HE3	1:E:475:MET:HB3	1.94	0.41
2:F:538:THR:CG2	2:F:538:THR:O	2.69	0.41
2:F:545:LEU:H	2:F:545:LEU:HG	1.66	0.41
1:I:35:TRP:N	1:I:499:THR:O	2.39	0.41
1:I:153:ASP:OD1	1:I:153:ASP:N	2.52	0.41
1:I:437:PRO:HA	1:I:438:PRO:HD3	1.95	0.41
1:E:249:HIS:O	1:E:251:ILE:HG23	2.21	0.41
5:Q:33:TYR:CD1	5:Q:52:SER:HA	2.56	0.41
2:B:543:GLN:OE1	2:J:644:ASP:OD1	2.38	0.40
3:G:73:ASP:OD1	7:a:1:NAG:H83	2.21	0.40
2:J:544:LEU:HD23	2:J:544:LEU:HA	1.74	0.40
2:J:580:VAL:O	2:J:584:GLU:HG3	2.21	0.40
3:K:99:ASP:O	3:K:100(E):HIS:NE2	2.47	0.40
2:B:544:LEU:HD23	2:B:544:LEU:HA	1.74	0.40
2:B:580:VAL:O	2:B:584:GLU:HG3	2.21	0.40
2:B:627:THR:OG1	2:B:630:GLU:HG3	2.21	0.40
2:B:661:LEU:HD11	1:E:503:ARG:HG2	2.03	0.40
1:E:368:ASP:HB3	1:E:371:ILE:HD13	2.03	0.40
1:I:232:LYS:HB2	1:I:232:LYS:HE2	1.86	0.40
6:N:48:ILE:HD12	6:N:48:ILE:HA	1.98	0.40
1:E:407:GLU:O	1:E:408:THR:OG1	2.33	0.40
2:J:618:THR:O	2:J:620:GLU:N	2.55	0.40
1:A:42:VAL:HA	2:B:537:LEU:CD1	2.52	0.40
1:A:72:HIS:C	1:A:74:CYS:H	2.30	0.40
3:G:4:LEU:HD12	3:G:22:CYS:SG	2.62	0.40
6:P:23:CYS:HB3	6:P:88:CYS:HB2	1.82	0.40
6:P:75:ILE:HD13	6:P:86:TYR:HE2	1.86	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:597:GLY:HA3	2:F:650:GLN:HE21	1.87	0.40
5:O:33:TYR:CD1	5:O:52:SER:HA	2.56	0.40
6:P:102:THR:O	6:P:102:THR:OG1	2.39	0.40
3:K:4:LEU:HD12	3:K:22:CYS:SG	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	435/485 (90%)	390 (90%)	40 (9%)	5 (1%)	11	39
1	E	435/485 (90%)	390 (90%)	40 (9%)	5 (1%)	11	39
1	I	435/485 (90%)	390 (90%)	40 (9%)	5 (1%)	11	39
2	B	116/155 (75%)	103 (89%)	10 (9%)	3 (3%)	4	21
2	F	116/155 (75%)	103 (89%)	10 (9%)	3 (3%)	4	21
2	J	116/155 (75%)	103 (89%)	10 (9%)	3 (3%)	4	21
3	C	128/238 (54%)	124 (97%)	4 (3%)	0	100	100
3	G	128/238 (54%)	124 (97%)	4 (3%)	0	100	100
3	K	128/238 (54%)	124 (97%)	4 (3%)	0	100	100
4	D	105/215 (49%)	101 (96%)	4 (4%)	0	100	100
4	H	105/215 (49%)	101 (96%)	4 (4%)	0	100	100
4	L	105/215 (49%)	101 (96%)	4 (4%)	0	100	100
5	M	129/238 (54%)	116 (90%)	12 (9%)	1 (1%)	16	46
5	O	129/238 (54%)	116 (90%)	12 (9%)	1 (1%)	16	46
5	Q	129/238 (54%)	116 (90%)	12 (9%)	1 (1%)	16	46
6	N	105/214 (49%)	97 (92%)	7 (7%)	1 (1%)	12	41

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	P	105/214 (49%)	97 (92%)	7 (7%)	1 (1%)	12	41
6	R	105/214 (49%)	97 (92%)	7 (7%)	1 (1%)	12	41
All	All	3054/4635 (66%)	2793 (92%)	231 (8%)	30 (1%)	15	41

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	VAL
1	A	204	ALA
1	A	408	THR
2	B	544	LEU
2	B	619	PHE
5	M	41	PRO
6	N	12	SER
1	E	154	VAL
1	E	204	ALA
1	E	408	THR
2	F	544	LEU
2	F	619	PHE
5	O	41	PRO
6	P	12	SER
1	I	154	VAL
1	I	204	ALA
1	I	408	THR
2	J	544	LEU
2	J	619	PHE
5	Q	41	PRO
6	R	12	SER
1	A	60	ALA
1	E	60	ALA
1	I	60	ALA
2	B	541	ALA
2	F	541	ALA
2	J	541	ALA
1	A	65	VAL
1	E	65	VAL
1	I	65	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	397/433 (92%)	388 (98%)	9 (2%)	44	66
1	E	397/433 (92%)	388 (98%)	9 (2%)	44	66
1	I	397/433 (92%)	388 (98%)	9 (2%)	44	66
2	B	104/130 (80%)	95 (91%)	9 (9%)	9	32
2	F	104/130 (80%)	95 (91%)	9 (9%)	9	32
2	J	104/130 (80%)	95 (91%)	9 (9%)	9	32
3	C	111/204 (54%)	106 (96%)	5 (4%)	24	53
3	G	111/204 (54%)	106 (96%)	5 (4%)	24	53
3	K	111/204 (54%)	106 (96%)	5 (4%)	24	53
4	D	85/182 (47%)	82 (96%)	3 (4%)	32	59
4	H	85/182 (47%)	82 (96%)	3 (4%)	32	59
4	L	85/182 (47%)	82 (96%)	3 (4%)	32	59
5	M	115/208 (55%)	113 (98%)	2 (2%)	53	70
5	O	115/208 (55%)	113 (98%)	2 (2%)	53	70
5	Q	115/208 (55%)	113 (98%)	2 (2%)	53	70
6	N	85/178 (48%)	79 (93%)	6 (7%)	13	40
6	P	85/178 (48%)	79 (93%)	6 (7%)	13	40
6	R	85/178 (48%)	79 (93%)	6 (7%)	13	40
All	All	2691/4005 (67%)	2589 (96%)	102 (4%)	30	58

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	GLN
1	A	128	THR
1	A	190	GLU
1	A	315	GLN
1	A	444	LYS
1	A	448	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	474	ASP
1	A	475	MET
1	A	503	ARG
2	B	537	LEU
2	B	542	ARG
2	B	543	GLN
2	B	544	LEU
2	B	545	LEU
2	B	601	LYS
2	B	620	GLU
2	B	626	MET
2	B	650	GLN
3	C	13	LYS
3	C	66	ARG
3	C	78	SER
3	C	84	SER
3	C	101	SER
4	D	14	SER
4	D	42	LYS
4	D	65	SER
5	M	41	PRO
5	M	65	SER
6	N	11	LEU
6	N	60	GLU
6	N	79	GLU
6	N	95	SER
6	N	95(B)	PHE
6	N	95(C)	SER
1	E	103	GLN
1	E	128	THR
1	E	190	GLU
1	E	315	GLN
1	E	444	LYS
1	E	448	ASN
1	E	474	ASP
1	E	475	MET
1	E	503	ARG
2	F	537	LEU
2	F	542	ARG
2	F	543	GLN
2	F	544	LEU
2	F	545	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	F	601	LYS
2	F	620	GLU
2	F	626	MET
2	F	650	GLN
3	G	13	LYS
3	G	66	ARG
3	G	78	SER
3	G	84	SER
3	G	101	SER
4	H	14	SER
4	H	42	LYS
4	H	65	SER
5	O	41	PRO
5	O	65	SER
6	P	11	LEU
6	P	60	GLU
6	P	79	GLU
6	P	95	SER
6	P	95(B)	PHE
6	P	95(C)	SER
1	I	103	GLN
1	I	128	THR
1	I	190	GLU
1	I	315	GLN
1	I	444	LYS
1	I	448	ASN
1	I	474	ASP
1	I	475	MET
1	I	503	ARG
2	J	537	LEU
2	J	542	ARG
2	J	543	GLN
2	J	544	LEU
2	J	545	LEU
2	J	601	LYS
2	J	620	GLU
2	J	626	MET
2	J	650	GLN
3	K	13	LYS
3	K	66	ARG
3	K	78	SER
3	K	84	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	K	101	SER
4	L	14	SER
4	L	42	LYS
4	L	65	SER
5	Q	41	PRO
5	Q	65	SER
6	R	11	LEU
6	R	60	GLU
6	R	79	GLU
6	R	95	SER
6	R	95(B)	PHE
6	R	95(C)	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	72	HIS
1	A	80	ASN
1	A	103	GLN
1	A	114	GLN
1	A	422	GLN
1	A	425	ASN
2	B	540	GLN
2	B	575	GLN
2	B	650	GLN
3	C	6	GLN
3	C	39	GLN
4	D	6	GLN
4	D	38	GLN
5	M	3	GLN
5	M	99	GLN
1	E	66	HIS
1	E	72	HIS
1	E	80	ASN
1	E	103	GLN
1	E	114	GLN
1	E	460	ASN
2	F	540	GLN
2	F	575	GLN
2	F	653	GLN
3	G	6	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	G	39	GLN
3	G	62	HIS
4	H	6	GLN
4	H	38	GLN
5	O	3	GLN
5	O	99	GLN
6	P	89	HIS
1	I	66	HIS
1	I	72	HIS
1	I	80	ASN
1	I	103	GLN
1	I	114	GLN
1	I	315	GLN
1	I	422	GLN
1	I	425	ASN
1	I	460	ASN
2	J	540	GLN
2	J	575	GLN
2	J	650	GLN
3	K	6	GLN
3	K	39	GLN
4	L	6	GLN
4	L	38	GLN
5	Q	3	GLN
5	Q	99	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

156 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	$\# Z > 2$
10	NAG	0	1	10,1	14,14,15	0.71	1 (7%)	17,19,21	0.71	0
10	NAG	0	2	10	14,14,15	0.31	0	17,19,21	0.46	0
10	NAG	1	1	10,1	14,14,15	0.71	1 (7%)	17,19,21	0.72	0
10	NAG	1	2	10	14,14,15	0.30	0	17,19,21	0.46	0
10	NAG	2	1	10,1	14,14,15	0.72	1 (7%)	17,19,21	0.72	0
10	NAG	2	2	10	14,14,15	0.29	0	17,19,21	0.46	0
7	NAG	S	1	7,1	14,14,15	0.38	0	17,19,21	0.74	1 (5%)
7	MAN	S	10	7	11,11,12	0.56	0	15,15,17	1.10	2 (13%)
7	MAN	S	11	7	11,11,12	0.58	0	15,15,17	0.96	2 (13%)
7	NAG	S	2	7	14,14,15	0.34	0	17,19,21	0.45	0
7	BMA	S	3	7	11,11,12	0.57	0	15,15,17	0.78	0
7	MAN	S	4	7	11,11,12	0.59	0	15,15,17	0.93	2 (13%)
7	MAN	S	5	7	11,11,12	0.56	0	15,15,17	1.00	2 (13%)
7	MAN	S	6	7	11,11,12	0.59	0	15,15,17	1.05	2 (13%)
7	MAN	S	7	7	11,11,12	0.61	0	15,15,17	1.04	2 (13%)
7	MAN	S	8	7	11,11,12	0.62	0	15,15,17	0.93	2 (13%)
7	MAN	S	9	7	11,11,12	0.64	0	15,15,17	0.96	2 (13%)
8	NAG	T	1	8,1	14,14,15	0.29	0	17,19,21	0.41	0
8	NAG	T	2	8	14,14,15	0.31	0	17,19,21	0.48	0
8	BMA	T	3	8	11,11,12	0.49	0	15,15,17	0.66	0
8	MAN	T	4	8	11,11,12	0.79	0	15,15,17	1.00	2 (13%)
8	MAN	T	5	8	11,11,12	0.79	0	15,15,17	1.23	2 (13%)
8	MAN	T	6	8	11,11,12	0.61	0	15,15,17	0.94	2 (13%)
9	NAG	U	1	9,1	14,14,15	0.26	0	17,19,21	0.46	0
9	MAN	U	10	9	11,11,12	0.62	0	15,15,17	0.90	2 (13%)
9	NAG	U	2	9	14,14,15	0.21	0	17,19,21	0.48	0
9	BMA	U	3	9	11,11,12	0.92	1 (9%)	15,15,17	1.07	1 (6%)
9	MAN	U	4	9	11,11,12	0.69	0	15,15,17	0.86	0
9	MAN	U	5	9	11,11,12	0.82	0	15,15,17	0.95	0
9	MAN	U	6	9	11,11,12	0.63	0	15,15,17	1.18	2 (13%)
9	MAN	U	7	9	11,11,12	0.60	0	15,15,17	1.17	2 (13%)
9	MAN	U	8	9	11,11,12	0.74	0	15,15,17	0.79	1 (6%)
9	MAN	U	9	9	11,11,12	0.59	0	15,15,17	0.97	2 (13%)
10	NAG	V	1	10,1	14,14,15	0.43	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	V	2	10	14,14,15	0.45	0	17,19,21	0.45	0
10	NAG	X	1	10,1	14,14,15	0.22	0	17,19,21	0.43	0
10	NAG	X	2	10	14,14,15	0.23	0	17,19,21	0.43	0
8	NAG	Y	1	8,1	14,14,15	0.23	0	17,19,21	0.57	0
8	NAG	Y	2	8	14,14,15	0.21	0	17,19,21	0.41	0
8	BMA	Y	3	8	11,11,12	0.44	0	15,15,17	0.68	0
8	MAN	Y	4	8	11,11,12	0.69	0	15,15,17	0.98	2 (13%)
8	MAN	Y	5	8	11,11,12	0.79	0	15,15,17	1.25	2 (13%)
8	MAN	Y	6	8	11,11,12	0.61	0	15,15,17	0.95	2 (13%)
11	NAG	Z	1	11,2	14,14,15	0.18	0	17,19,21	0.49	0
11	NAG	Z	2	11	14,14,15	0.25	0	17,19,21	0.46	0
11	BMA	Z	3	11	11,11,12	0.58	0	15,15,17	0.74	0
11	MAN	Z	4	11	11,11,12	0.78	0	15,15,17	1.01	2 (13%)
11	MAN	Z	5	11	11,11,12	0.60	0	15,15,17	0.98	2 (13%)
7	NAG	a	1	7,1	14,14,15	0.39	0	17,19,21	0.75	1 (5%)
7	MAN	a	10	7	11,11,12	0.56	0	15,15,17	1.08	2 (13%)
7	MAN	a	11	7	11,11,12	0.58	0	15,15,17	0.96	2 (13%)
7	NAG	a	2	7	14,14,15	0.32	0	17,19,21	0.45	0
7	BMA	a	3	7	11,11,12	0.57	0	15,15,17	0.77	0
7	MAN	a	4	7	11,11,12	0.60	0	15,15,17	0.92	2 (13%)
7	MAN	a	5	7	11,11,12	0.57	0	15,15,17	1.00	2 (13%)
7	MAN	a	6	7	11,11,12	0.61	0	15,15,17	1.04	2 (13%)
7	MAN	a	7	7	11,11,12	0.61	0	15,15,17	1.05	2 (13%)
7	MAN	a	8	7	11,11,12	0.62	0	15,15,17	0.93	2 (13%)
7	MAN	a	9	7	11,11,12	0.63	0	15,15,17	0.97	2 (13%)
8	NAG	b	1	8,1	14,14,15	0.28	0	17,19,21	0.40	0
8	NAG	b	2	8	14,14,15	0.30	0	17,19,21	0.46	0
8	BMA	b	3	8	11,11,12	0.49	0	15,15,17	0.67	0
8	MAN	b	4	8	11,11,12	0.78	0	15,15,17	1.01	2 (13%)
8	MAN	b	5	8	11,11,12	0.80	0	15,15,17	1.22	2 (13%)
8	MAN	b	6	8	11,11,12	0.62	0	15,15,17	0.95	2 (13%)
10	NAG	c	1	10,1	14,14,15	0.20	0	17,19,21	0.46	0
10	NAG	c	2	10	14,14,15	0.23	0	17,19,21	0.44	0
9	NAG	d	1	9,1	14,14,15	0.27	0	17,19,21	0.46	0
9	MAN	d	10	9	11,11,12	0.64	0	15,15,17	0.90	2 (13%)
9	NAG	d	2	9	14,14,15	0.20	0	17,19,21	0.48	0
9	BMA	d	3	9	11,11,12	0.90	1 (9%)	15,15,17	1.08	1 (6%)
9	MAN	d	4	9	11,11,12	0.67	0	15,15,17	0.87	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	MAN	d	5	9	11,11,12	0.78	0	15,15,17	0.95	0
9	MAN	d	6	9	11,11,12	0.63	0	15,15,17	1.17	2 (13%)
9	MAN	d	7	9	11,11,12	0.63	0	15,15,17	1.17	2 (13%)
9	MAN	d	8	9	11,11,12	0.76	1 (9%)	15,15,17	0.81	1 (6%)
9	MAN	d	9	9	11,11,12	0.60	0	15,15,17	0.96	2 (13%)
10	NAG	e	1	10,1	14,14,15	0.42	0	17,19,21	0.54	0
10	NAG	e	2	10	14,14,15	0.47	0	17,19,21	0.45	0
10	NAG	f	1	10,1	14,14,15	0.21	0	17,19,21	0.46	0
10	NAG	f	2	10	14,14,15	0.23	0	17,19,21	0.43	0
10	NAG	g	1	10,1	14,14,15	0.23	0	17,19,21	0.44	0
10	NAG	g	2	10	14,14,15	0.23	0	17,19,21	0.43	0
8	NAG	h	1	8,1	14,14,15	0.23	0	17,19,21	0.55	0
8	NAG	h	2	8	14,14,15	0.22	0	17,19,21	0.40	0
8	BMA	h	3	8	11,11,12	0.42	0	15,15,17	0.68	0
8	MAN	h	4	8	11,11,12	0.64	0	15,15,17	0.97	2 (13%)
8	MAN	h	5	8	11,11,12	0.78	0	15,15,17	1.24	2 (13%)
8	MAN	h	6	8	11,11,12	0.63	0	15,15,17	0.95	2 (13%)
11	NAG	i	1	11,2	14,14,15	0.18	0	17,19,21	0.47	0
11	NAG	i	2	11	14,14,15	0.26	0	17,19,21	0.44	0
11	BMA	i	3	11	11,11,12	0.57	0	15,15,17	0.75	0
11	MAN	i	4	11	11,11,12	0.77	0	15,15,17	1.01	2 (13%)
11	MAN	i	5	11	11,11,12	0.61	0	15,15,17	0.97	2 (13%)
7	NAG	j	1	7,1	14,14,15	0.38	0	17,19,21	0.74	1 (5%)
7	MAN	j	10	7	11,11,12	0.56	0	15,15,17	1.07	2 (13%)
7	MAN	j	11	7	11,11,12	0.59	0	15,15,17	0.95	2 (13%)
7	NAG	j	2	7	14,14,15	0.33	0	17,19,21	0.45	0
7	BMA	j	3	7	11,11,12	0.56	0	15,15,17	0.76	0
7	MAN	j	4	7	11,11,12	0.60	0	15,15,17	0.92	2 (13%)
7	MAN	j	5	7	11,11,12	0.56	0	15,15,17	1.01	2 (13%)
7	MAN	j	6	7	11,11,12	0.60	0	15,15,17	1.04	2 (13%)
7	MAN	j	7	7	11,11,12	0.60	0	15,15,17	1.04	2 (13%)
7	MAN	j	8	7	11,11,12	0.61	0	15,15,17	0.92	2 (13%)
7	MAN	j	9	7	11,11,12	0.63	0	15,15,17	0.97	2 (13%)
8	NAG	k	1	8,1	14,14,15	0.28	0	17,19,21	0.40	0
8	NAG	k	2	8	14,14,15	0.31	0	17,19,21	0.45	0
8	BMA	k	3	8	11,11,12	0.50	0	15,15,17	0.66	0
8	MAN	k	4	8	11,11,12	0.79	0	15,15,17	1.01	2 (13%)
8	MAN	k	5	8	11,11,12	0.79	0	15,15,17	1.22	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	MAN	k	6	8	11,11,12	0.61	0	15,15,17	0.94	2 (13%)
10	NAG	l	1	10,1	14,14,15	0.19	0	17,19,21	0.46	0
10	NAG	l	2	10	14,14,15	0.23	0	17,19,21	0.43	0
10	NAG	m	1	10,1	14,14,15	0.63	1 (7%)	17,19,21	0.58	0
10	NAG	m	2	10	14,14,15	0.25	0	17,19,21	0.46	0
9	NAG	n	1	9,1	14,14,15	0.25	0	17,19,21	0.47	0
9	MAN	n	10	9	11,11,12	0.63	0	15,15,17	0.91	2 (13%)
9	NAG	n	2	9	14,14,15	0.20	0	17,19,21	0.48	0
9	BMA	n	3	9	11,11,12	0.90	1 (9%)	15,15,17	1.09	1 (6%)
9	MAN	n	4	9	11,11,12	0.68	0	15,15,17	0.86	0
9	MAN	n	5	9	11,11,12	0.77	0	15,15,17	0.96	0
9	MAN	n	6	9	11,11,12	0.62	0	15,15,17	1.16	2 (13%)
9	MAN	n	7	9	11,11,12	0.61	0	15,15,17	1.18	2 (13%)
9	MAN	n	8	9	11,11,12	0.76	1 (9%)	15,15,17	0.80	1 (6%)
9	MAN	n	9	9	11,11,12	0.60	0	15,15,17	0.96	2 (13%)
10	NAG	o	1	10,1	14,14,15	0.43	0	17,19,21	0.53	0
10	NAG	o	2	10	14,14,15	0.46	0	17,19,21	0.45	0
10	NAG	p	1	10,1	14,14,15	0.23	0	17,19,21	0.44	0
10	NAG	p	2	10	14,14,15	0.22	0	17,19,21	0.44	0
10	NAG	q	1	10,1	14,14,15	0.23	0	17,19,21	0.45	0
10	NAG	q	2	10	14,14,15	0.24	0	17,19,21	0.43	0
8	NAG	r	1	8,1	14,14,15	0.23	0	17,19,21	0.54	0
8	NAG	r	2	8	14,14,15	0.21	0	17,19,21	0.41	0
8	BMA	r	3	8	11,11,12	0.42	0	15,15,17	0.69	0
8	MAN	r	4	8	11,11,12	0.63	0	15,15,17	0.96	2 (13%)
8	MAN	r	5	8	11,11,12	0.79	0	15,15,17	1.24	2 (13%)
8	MAN	r	6	8	11,11,12	0.61	0	15,15,17	0.97	2 (13%)
11	NAG	s	1	11,2	14,14,15	0.18	0	17,19,21	0.46	0
11	NAG	s	2	11	14,14,15	0.27	0	17,19,21	0.42	0
11	BMA	s	3	11	11,11,12	0.56	0	15,15,17	0.76	0
11	MAN	s	4	11	11,11,12	0.78	0	15,15,17	1.00	2 (13%)
11	MAN	s	5	11	11,11,12	0.61	0	15,15,17	0.97	2 (13%)
10	NAG	t	1	10,1	14,14,15	0.20	0	17,19,21	0.46	0
10	NAG	t	2	10	14,14,15	0.24	0	17,19,21	0.44	0
10	NAG	u	1	10,1	14,14,15	0.20	0	17,19,21	0.45	0
10	NAG	u	2	10	14,14,15	0.23	0	17,19,21	0.43	0
10	NAG	v	1	10,1	14,14,15	0.63	1 (7%)	17,19,21	0.57	0
10	NAG	v	2	10	14,14,15	0.25	0	17,19,21	0.46	0
10	NAG	w	1	10,1	14,14,15	0.62	1 (7%)	17,19,21	0.56	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	NAG	w	2	10	14,14,15	0.24	0	17,19,21	0.46	0
10	NAG	x	1	10,1	14,14,15	0.51	0	17,19,21	0.64	0
10	NAG	x	2	10	14,14,15	0.25	0	17,19,21	0.49	0
10	NAG	y	1	10,1	14,14,15	0.52	0	17,19,21	0.64	0
10	NAG	y	2	10	14,14,15	0.24	0	17,19,21	0.48	0
10	NAG	z	1	10,1	14,14,15	0.51	0	17,19,21	0.63	0
10	NAG	z	2	10	14,14,15	0.23	0	17,19,21	0.49	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
10	NAG	0	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	0	2	10	-	2/6/23/26	0/1/1/1
10	NAG	1	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	1	2	10	-	2/6/23/26	0/1/1/1
10	NAG	2	1	10,1	-	1/6/23/26	0/1/1/1
10	NAG	2	2	10	-	2/6/23/26	0/1/1/1
7	NAG	S	1	7,1	-	2/6/23/26	0/1/1/1
7	MAN	S	10	7	-	2/2/19/22	0/1/1/1
7	MAN	S	11	7	-	0/2/19/22	0/1/1/1
7	NAG	S	2	7	-	1/6/23/26	0/1/1/1
7	BMA	S	3	7	-	0/2/19/22	0/1/1/1
7	MAN	S	4	7	-	0/2/19/22	0/1/1/1
7	MAN	S	5	7	-	2/2/19/22	0/1/1/1
7	MAN	S	6	7	-	1/2/19/22	0/1/1/1
7	MAN	S	7	7	-	0/2/19/22	0/1/1/1
7	MAN	S	8	7	-	0/2/19/22	0/1/1/1
7	MAN	S	9	7	-	1/2/19/22	0/1/1/1
8	NAG	T	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	T	2	8	-	2/6/23/26	0/1/1/1
8	BMA	T	3	8	-	0/2/19/22	0/1/1/1
8	MAN	T	4	8	-	0/2/19/22	0/1/1/1
8	MAN	T	5	8	-	0/2/19/22	0/1/1/1
8	MAN	T	6	8	-	0/2/19/22	0/1/1/1
9	NAG	U	1	9,1	-	2/6/23/26	0/1/1/1
9	MAN	U	10	9	-	0/2/19/22	0/1/1/1
9	NAG	U	2	9	-	0/6/23/26	0/1/1/1
9	BMA	U	3	9	-	0/2/19/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	U	4	9	-	0/2/19/22	0/1/1/1
9	MAN	U	5	9	-	2/2/19/22	0/1/1/1
9	MAN	U	6	9	-	2/2/19/22	0/1/1/1
9	MAN	U	7	9	-	0/2/19/22	0/1/1/1
9	MAN	U	8	9	-	0/2/19/22	0/1/1/1
9	MAN	U	9	9	-	0/2/19/22	0/1/1/1
10	NAG	V	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	V	2	10	-	2/6/23/26	0/1/1/1
10	NAG	X	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	X	2	10	-	2/6/23/26	0/1/1/1
8	NAG	Y	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	Y	2	8	-	0/6/23/26	0/1/1/1
8	BMA	Y	3	8	-	0/2/19/22	0/1/1/1
8	MAN	Y	4	8	-	0/2/19/22	0/1/1/1
8	MAN	Y	5	8	-	0/2/19/22	0/1/1/1
8	MAN	Y	6	8	-	0/2/19/22	0/1/1/1
11	NAG	Z	1	11,2	-	1/6/23/26	0/1/1/1
11	NAG	Z	2	11	-	1/6/23/26	0/1/1/1
11	BMA	Z	3	11	-	2/2/19/22	0/1/1/1
11	MAN	Z	4	11	-	0/2/19/22	0/1/1/1
11	MAN	Z	5	11	-	0/2/19/22	0/1/1/1
7	NAG	a	1	7,1	-	2/6/23/26	0/1/1/1
7	MAN	a	10	7	-	2/2/19/22	0/1/1/1
7	MAN	a	11	7	-	0/2/19/22	0/1/1/1
7	NAG	a	2	7	-	1/6/23/26	0/1/1/1
7	BMA	a	3	7	-	0/2/19/22	0/1/1/1
7	MAN	a	4	7	-	0/2/19/22	0/1/1/1
7	MAN	a	5	7	-	2/2/19/22	0/1/1/1
7	MAN	a	6	7	-	1/2/19/22	0/1/1/1
7	MAN	a	7	7	-	0/2/19/22	0/1/1/1
7	MAN	a	8	7	-	0/2/19/22	0/1/1/1
7	MAN	a	9	7	-	1/2/19/22	0/1/1/1
8	NAG	b	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	b	2	8	-	2/6/23/26	0/1/1/1
8	BMA	b	3	8	-	0/2/19/22	0/1/1/1
8	MAN	b	4	8	-	0/2/19/22	0/1/1/1
8	MAN	b	5	8	-	0/2/19/22	0/1/1/1
8	MAN	b	6	8	-	0/2/19/22	0/1/1/1
10	NAG	c	1	10,1	-	2/6/23/26	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	c	2	10	-	3/6/23/26	0/1/1/1
9	NAG	d	1	9,1	-	2/6/23/26	0/1/1/1
9	MAN	d	10	9	-	0/2/19/22	0/1/1/1
9	NAG	d	2	9	-	0/6/23/26	0/1/1/1
9	BMA	d	3	9	-	0/2/19/22	0/1/1/1
9	MAN	d	4	9	-	0/2/19/22	0/1/1/1
9	MAN	d	5	9	-	2/2/19/22	0/1/1/1
9	MAN	d	6	9	-	2/2/19/22	0/1/1/1
9	MAN	d	7	9	-	0/2/19/22	0/1/1/1
9	MAN	d	8	9	-	0/2/19/22	0/1/1/1
9	MAN	d	9	9	-	0/2/19/22	0/1/1/1
10	NAG	e	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	e	2	10	-	2/6/23/26	0/1/1/1
10	NAG	f	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	f	2	10	-	2/6/23/26	0/1/1/1
10	NAG	g	1	10,1	-	1/6/23/26	0/1/1/1
10	NAG	g	2	10	-	2/6/23/26	0/1/1/1
8	NAG	h	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	h	2	8	-	0/6/23/26	0/1/1/1
8	BMA	h	3	8	-	0/2/19/22	0/1/1/1
8	MAN	h	4	8	-	0/2/19/22	0/1/1/1
8	MAN	h	5	8	-	0/2/19/22	0/1/1/1
8	MAN	h	6	8	-	0/2/19/22	0/1/1/1
11	NAG	i	1	11,2	-	1/6/23/26	0/1/1/1
11	NAG	i	2	11	-	0/6/23/26	0/1/1/1
11	BMA	i	3	11	-	2/2/19/22	0/1/1/1
11	MAN	i	4	11	-	0/2/19/22	0/1/1/1
11	MAN	i	5	11	-	0/2/19/22	0/1/1/1
7	NAG	j	1	7,1	-	2/6/23/26	0/1/1/1
7	MAN	j	10	7	-	0/2/19/22	0/1/1/1
7	MAN	j	11	7	-	0/2/19/22	0/1/1/1
7	NAG	j	2	7	-	0/6/23/26	0/1/1/1
7	BMA	j	3	7	-	0/2/19/22	0/1/1/1
7	MAN	j	4	7	-	0/2/19/22	0/1/1/1
7	MAN	j	5	7	-	2/2/19/22	0/1/1/1
7	MAN	j	6	7	-	1/2/19/22	0/1/1/1
7	MAN	j	7	7	-	0/2/19/22	0/1/1/1
7	MAN	j	8	7	-	0/2/19/22	0/1/1/1
7	MAN	j	9	7	-	1/2/19/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	k	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	k	2	8	-	2/6/23/26	0/1/1/1
8	BMA	k	3	8	-	0/2/19/22	0/1/1/1
8	MAN	k	4	8	-	0/2/19/22	0/1/1/1
8	MAN	k	5	8	-	0/2/19/22	0/1/1/1
8	MAN	k	6	8	-	0/2/19/22	0/1/1/1
10	NAG	l	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	l	2	10	-	2/6/23/26	0/1/1/1
10	NAG	m	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	m	2	10	-	2/6/23/26	0/1/1/1
9	NAG	n	1	9,1	-	2/6/23/26	0/1/1/1
9	MAN	n	10	9	-	0/2/19/22	0/1/1/1
9	NAG	n	2	9	-	0/6/23/26	0/1/1/1
9	BMA	n	3	9	-	0/2/19/22	0/1/1/1
9	MAN	n	4	9	-	0/2/19/22	0/1/1/1
9	MAN	n	5	9	-	2/2/19/22	0/1/1/1
9	MAN	n	6	9	-	2/2/19/22	0/1/1/1
9	MAN	n	7	9	-	0/2/19/22	0/1/1/1
9	MAN	n	8	9	-	0/2/19/22	0/1/1/1
9	MAN	n	9	9	-	0/2/19/22	0/1/1/1
10	NAG	o	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	o	2	10	-	2/6/23/26	0/1/1/1
10	NAG	p	1	10,1	-	0/6/23/26	0/1/1/1
10	NAG	p	2	10	-	2/6/23/26	0/1/1/1
10	NAG	q	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	q	2	10	-	3/6/23/26	0/1/1/1
8	NAG	r	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	r	2	8	-	0/6/23/26	0/1/1/1
8	BMA	r	3	8	-	0/2/19/22	0/1/1/1
8	MAN	r	4	8	-	0/2/19/22	0/1/1/1
8	MAN	r	5	8	-	0/2/19/22	0/1/1/1
8	MAN	r	6	8	-	0/2/19/22	0/1/1/1
11	NAG	s	1	11,2	-	1/6/23/26	0/1/1/1
11	NAG	s	2	11	-	0/6/23/26	0/1/1/1
11	BMA	s	3	11	-	2/2/19/22	0/1/1/1
11	MAN	s	4	11	-	0/2/19/22	0/1/1/1
11	MAN	s	5	11	-	0/2/19/22	0/1/1/1
10	NAG	t	1	10,1	-	2/6/23/26	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	t	2	10	-	3/6/23/26	0/1/1/1
10	NAG	u	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	u	2	10	-	2/6/23/26	0/1/1/1
10	NAG	v	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	v	2	10	-	2/6/23/26	0/1/1/1
10	NAG	w	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	w	2	10	-	2/6/23/26	0/1/1/1
10	NAG	x	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	x	2	10	-	2/6/23/26	0/1/1/1
10	NAG	y	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	y	2	10	-	2/6/23/26	0/1/1/1
10	NAG	z	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	z	2	10	-	2/6/23/26	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	2	1	NAG	C1-C2	-2.34	1.49	1.52
10	1	1	NAG	C1-C2	-2.33	1.49	1.52
10	0	1	NAG	C1-C2	-2.32	1.49	1.52
10	m	1	NAG	O5-C1	-2.20	1.40	1.43
9	U	3	BMA	O5-C1	-2.20	1.40	1.43
10	v	1	NAG	O5-C1	-2.20	1.40	1.43
10	w	1	NAG	O5-C1	-2.19	1.40	1.43
9	d	3	BMA	O5-C1	-2.18	1.40	1.43
9	n	3	BMA	O5-C1	-2.12	1.40	1.43
9	d	8	MAN	O5-C1	-2.01	1.40	1.43
9	n	8	MAN	O5-C1	-2.01	1.40	1.43

All (129) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	Y	5	MAN	C1-O5-C5	3.51	116.90	112.19
9	U	6	MAN	C1-O5-C5	3.48	116.85	112.19
8	h	5	MAN	C1-O5-C5	3.47	116.83	112.19
8	T	5	MAN	C1-O5-C5	3.47	116.83	112.19
8	r	5	MAN	C1-O5-C5	3.46	116.83	112.19
8	b	5	MAN	C1-O5-C5	3.44	116.80	112.19
8	k	5	MAN	C1-O5-C5	3.44	116.80	112.19

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	d	6	MAN	C1-O5-C5	3.38	116.71	112.19
9	n	6	MAN	C1-O5-C5	3.34	116.66	112.19
9	d	7	MAN	C1-O5-C5	3.30	116.61	112.19
9	n	7	MAN	C1-O5-C5	3.29	116.60	112.19
9	U	7	MAN	C1-O5-C5	3.26	116.56	112.19
7	S	10	MAN	C1-O5-C5	3.01	116.22	112.19
7	a	10	MAN	C1-O5-C5	2.94	116.12	112.19
7	S	6	MAN	C1-O5-C5	2.93	116.11	112.19
7	j	10	MAN	C1-O5-C5	2.88	116.05	112.19
7	a	6	MAN	C1-O5-C5	2.87	116.03	112.19
7	j	6	MAN	C1-O5-C5	2.82	115.97	112.19
7	j	7	MAN	C1-O5-C5	2.69	115.79	112.19
7	j	5	MAN	C1-O5-C5	2.69	115.79	112.19
7	a	7	MAN	C1-O5-C5	2.68	115.78	112.19
7	S	7	MAN	C1-O5-C5	2.66	115.76	112.19
7	a	5	MAN	C1-O5-C5	2.64	115.72	112.19
7	S	5	MAN	C1-O5-C5	2.63	115.71	112.19
7	a	1	NAG	C1-O5-C5	2.48	115.51	112.19
7	j	1	NAG	C1-O5-C5	2.47	115.50	112.19
7	S	1	NAG	C1-O5-C5	2.45	115.47	112.19
9	U	9	MAN	C1-O5-C5	2.41	115.41	112.19
11	Z	5	MAN	C1-O5-C5	2.39	115.39	112.19
9	n	9	MAN	C1-O5-C5	2.37	115.36	112.19
11	i	5	MAN	C1-O5-C5	2.36	115.35	112.19
7	j	9	MAN	C1-O5-C5	2.35	115.34	112.19
9	d	9	MAN	C1-O5-C5	2.35	115.33	112.19
8	r	6	MAN	C1-O5-C5	2.34	115.32	112.19
11	s	5	MAN	C1-O5-C5	2.33	115.31	112.19
7	a	9	MAN	C1-O5-C5	2.32	115.30	112.19
7	S	9	MAN	C1-O5-C5	2.31	115.28	112.19
7	S	4	MAN	O2-C2-C3	-2.28	105.42	110.15
7	a	10	MAN	O2-C2-C3	-2.28	105.42	110.15
7	a	11	MAN	C1-O5-C5	2.28	115.24	112.19
7	a	4	MAN	O2-C2-C3	-2.28	105.43	110.15
7	a	7	MAN	O2-C2-C3	-2.28	105.44	110.15
7	j	10	MAN	O2-C2-C3	-2.28	105.44	110.15
7	S	10	MAN	O2-C2-C3	-2.27	105.45	110.15
7	S	7	MAN	O2-C2-C3	-2.26	105.47	110.15
7	S	11	MAN	C1-O5-C5	2.26	115.21	112.19
8	Y	6	MAN	C1-O5-C5	2.26	115.21	112.19
7	j	7	MAN	O2-C2-C3	-2.26	105.47	110.15
9	U	6	MAN	O2-C2-C3	-2.26	105.47	110.15

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	n	6	MAN	O2-C2-C3	-2.26	105.47	110.15
9	d	6	MAN	O2-C2-C3	-2.26	105.48	110.15
7	j	4	MAN	O2-C2-C3	-2.25	105.50	110.15
7	j	11	MAN	O2-C2-C3	-2.24	105.50	110.15
7	a	8	MAN	O2-C2-C3	-2.24	105.51	110.15
7	a	11	MAN	O2-C2-C3	-2.24	105.51	110.15
8	h	6	MAN	C1-O5-C5	2.24	115.19	112.19
8	T	6	MAN	C1-O5-C5	2.23	115.18	112.19
8	b	6	MAN	C1-O5-C5	2.23	115.17	112.19
7	S	11	MAN	O2-C2-C3	-2.23	105.53	110.15
8	k	6	MAN	C1-O5-C5	2.22	115.16	112.19
7	S	8	MAN	O2-C2-C3	-2.22	105.56	110.15
7	j	11	MAN	C1-O5-C5	2.22	115.16	112.19
11	s	5	MAN	O2-C2-C3	-2.21	105.57	110.15
11	Z	5	MAN	O2-C2-C3	-2.20	105.59	110.15
11	Z	4	MAN	O2-C2-C3	-2.19	105.62	110.15
7	j	8	MAN	O2-C2-C3	-2.18	105.63	110.15
8	b	6	MAN	O2-C2-C3	-2.18	105.63	110.15
7	a	9	MAN	O2-C2-C3	-2.18	105.63	110.15
9	n	7	MAN	O2-C2-C3	-2.18	105.64	110.15
11	i	5	MAN	O2-C2-C3	-2.18	105.64	110.15
7	j	9	MAN	O2-C2-C3	-2.18	105.64	110.15
9	U	7	MAN	O2-C2-C3	-2.17	105.65	110.15
7	S	9	MAN	O2-C2-C3	-2.17	105.65	110.15
8	Y	4	MAN	O2-C2-C3	-2.17	105.65	110.15
7	S	6	MAN	O2-C2-C3	-2.17	105.65	110.15
8	r	4	MAN	O2-C2-C3	-2.17	105.65	110.15
7	j	6	MAN	O2-C2-C3	-2.17	105.66	110.15
9	d	9	MAN	O2-C2-C3	-2.16	105.67	110.15
8	h	4	MAN	O2-C2-C3	-2.16	105.67	110.15
8	r	6	MAN	O2-C2-C3	-2.15	105.69	110.15
8	r	5	MAN	O2-C2-C3	-2.15	105.69	110.15
8	Y	5	MAN	O2-C2-C3	-2.15	105.70	110.15
8	k	6	MAN	O2-C2-C3	-2.15	105.70	110.15
7	j	5	MAN	O2-C2-C3	-2.15	105.70	110.15
8	T	6	MAN	O2-C2-C3	-2.15	105.70	110.15
9	d	7	MAN	O2-C2-C3	-2.15	105.70	110.15
7	a	6	MAN	O2-C2-C3	-2.15	105.71	110.15
8	h	6	MAN	O2-C2-C3	-2.15	105.71	110.15
8	Y	6	MAN	O2-C2-C3	-2.14	105.71	110.15
7	j	8	MAN	C1-O5-C5	2.14	115.06	112.19
7	S	8	MAN	C1-O5-C5	2.14	115.05	112.19

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	a	8	MAN	C1-O5-C5	2.14	115.05	112.19
9	n	10	MAN	O2-C2-C3	-2.14	105.72	110.15
9	U	10	MAN	O2-C2-C3	-2.14	105.72	110.15
9	U	9	MAN	O2-C2-C3	-2.14	105.72	110.15
9	d	10	MAN	O2-C2-C3	-2.14	105.73	110.15
7	a	5	MAN	O2-C2-C3	-2.13	105.73	110.15
8	h	4	MAN	C1-O5-C5	2.13	115.05	112.19
8	h	5	MAN	O2-C2-C3	-2.13	105.74	110.15
8	b	5	MAN	O2-C2-C3	-2.13	105.74	110.15
8	Y	4	MAN	C1-O5-C5	2.13	115.04	112.19
8	k	4	MAN	C1-O5-C5	2.13	115.04	112.19
7	S	4	MAN	C1-O5-C5	2.13	115.03	112.19
7	j	4	MAN	C1-O5-C5	2.12	115.03	112.19
8	T	5	MAN	O2-C2-C3	-2.12	105.75	110.15
8	k	5	MAN	O2-C2-C3	-2.12	105.77	110.15
9	n	9	MAN	O2-C2-C3	-2.11	105.77	110.15
7	S	5	MAN	O2-C2-C3	-2.11	105.78	110.15
9	n	3	BMA	C1-O5-C5	2.11	115.01	112.19
8	r	4	MAN	C1-O5-C5	2.10	115.01	112.19
8	b	4	MAN	C1-O5-C5	2.10	115.00	112.19
11	i	4	MAN	C1-O5-C5	2.10	115.00	112.19
11	i	4	MAN	O2-C2-C3	-2.10	105.81	110.15
11	s	4	MAN	O2-C2-C3	-2.09	105.81	110.15
8	b	4	MAN	O2-C2-C3	-2.09	105.82	110.15
11	s	4	MAN	C1-O5-C5	2.09	114.98	112.19
8	k	4	MAN	O2-C2-C3	-2.08	105.83	110.15
7	a	4	MAN	C1-O5-C5	2.08	114.97	112.19
8	T	4	MAN	O2-C2-C3	-2.07	105.86	110.15
9	d	3	BMA	C1-O5-C5	2.07	114.96	112.19
8	T	4	MAN	C1-O5-C5	2.07	114.95	112.19
9	n	8	MAN	O2-C2-C3	-2.05	105.90	110.15
9	n	10	MAN	C1-O5-C5	2.05	114.93	112.19
9	d	8	MAN	O2-C2-C3	-2.05	105.91	110.15
9	U	10	MAN	C1-O5-C5	2.04	114.92	112.19
11	Z	4	MAN	C1-O5-C5	2.04	114.91	112.19
9	U	8	MAN	O2-C2-C3	-2.02	105.97	110.15
9	d	10	MAN	C1-O5-C5	2.02	114.89	112.19
9	U	3	BMA	C1-O5-C5	2.00	114.87	112.19

There are no chirality outliers.

All (147) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	V	1	NAG	C4-C5-C6-O6
10	e	1	NAG	C4-C5-C6-O6
10	o	1	NAG	C4-C5-C6-O6
10	x	2	NAG	O5-C5-C6-O6
10	y	2	NAG	O5-C5-C6-O6
9	U	6	MAN	C4-C5-C6-O6
10	o	2	NAG	O5-C5-C6-O6
10	z	2	NAG	O5-C5-C6-O6
9	d	6	MAN	C4-C5-C6-O6
9	n	6	MAN	C4-C5-C6-O6
10	e	2	NAG	O5-C5-C6-O6
10	z	1	NAG	O5-C5-C6-O6
9	n	1	NAG	O5-C5-C6-O6
10	V	2	NAG	O5-C5-C6-O6
10	m	1	NAG	O5-C5-C6-O6
10	v	1	NAG	O5-C5-C6-O6
10	v	2	NAG	O5-C5-C6-O6
10	w	1	NAG	O5-C5-C6-O6
10	0	2	NAG	O5-C5-C6-O6
10	1	2	NAG	O5-C5-C6-O6
10	2	2	NAG	O5-C5-C6-O6
7	S	5	MAN	C4-C5-C6-O6
10	y	2	NAG	C4-C5-C6-O6
7	j	5	MAN	O5-C5-C6-O6
9	d	1	NAG	O5-C5-C6-O6
10	m	2	NAG	O5-C5-C6-O6
10	p	2	NAG	O5-C5-C6-O6
10	t	1	NAG	O5-C5-C6-O6
10	w	2	NAG	O5-C5-C6-O6
10	x	1	NAG	C4-C5-C6-O6
10	x	2	NAG	C4-C5-C6-O6
10	y	1	NAG	C4-C5-C6-O6
10	z	1	NAG	C4-C5-C6-O6
10	2	2	NAG	C4-C5-C6-O6
10	g	2	NAG	O5-C5-C6-O6
10	q	1	NAG	O5-C5-C6-O6
10	u	1	NAG	O5-C5-C6-O6
10	z	2	NAG	C4-C5-C6-O6
10	x	1	NAG	O5-C5-C6-O6
10	y	1	NAG	O5-C5-C6-O6
10	0	2	NAG	C4-C5-C6-O6
10	1	2	NAG	C4-C5-C6-O6
7	a	5	MAN	O5-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
8	T	2	NAG	O5-C5-C6-O6
8	b	2	NAG	O5-C5-C6-O6
9	U	1	NAG	O5-C5-C6-O6
10	X	2	NAG	O5-C5-C6-O6
10	e	1	NAG	O5-C5-C6-O6
10	o	2	NAG	C4-C5-C6-O6
9	d	6	MAN	O5-C5-C6-O6
10	V	1	NAG	O5-C5-C6-O6
7	a	5	MAN	C4-C5-C6-O6
8	k	2	NAG	O5-C5-C6-O6
9	U	6	MAN	O5-C5-C6-O6
10	f	2	NAG	O5-C5-C6-O6
10	q	2	NAG	O5-C5-C6-O6
10	m	2	NAG	C4-C5-C6-O6
7	S	5	MAN	O5-C5-C6-O6
9	n	6	MAN	O5-C5-C6-O6
10	o	1	NAG	O5-C5-C6-O6
9	n	1	NAG	C4-C5-C6-O6
10	v	2	NAG	C4-C5-C6-O6
10	w	2	NAG	C4-C5-C6-O6
9	d	1	NAG	C4-C5-C6-O6
10	e	2	NAG	C4-C5-C6-O6
10	m	1	NAG	C4-C5-C6-O6
10	c	2	NAG	O5-C5-C6-O6
10	t	2	NAG	O5-C5-C6-O6
8	k	2	NAG	C4-C5-C6-O6
10	g	2	NAG	C4-C5-C6-O6
10	q	1	NAG	C4-C5-C6-O6
10	t	1	NAG	C4-C5-C6-O6
10	u	2	NAG	O5-C5-C6-O6
10	v	1	NAG	C4-C5-C6-O6
9	U	1	NAG	C4-C5-C6-O6
10	p	2	NAG	C4-C5-C6-O6
10	u	1	NAG	C4-C5-C6-O6
10	w	1	NAG	C4-C5-C6-O6
10	l	2	NAG	O5-C5-C6-O6
7	j	5	MAN	C4-C5-C6-O6
8	b	2	NAG	C4-C5-C6-O6
10	V	2	NAG	C4-C5-C6-O6
8	T	2	NAG	C4-C5-C6-O6
8	b	1	NAG	O5-C5-C6-O6
8	k	1	NAG	O5-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
8	k	1	NAG	C4-C5-C6-O6
10	X	2	NAG	C4-C5-C6-O6
8	b	1	NAG	C4-C5-C6-O6
8	r	1	NAG	O5-C5-C6-O6
8	Y	1	NAG	O5-C5-C6-O6
8	h	1	NAG	O5-C5-C6-O6
10	f	2	NAG	C4-C5-C6-O6
10	q	2	NAG	C4-C5-C6-O6
8	T	1	NAG	O5-C5-C6-O6
8	T	1	NAG	C4-C5-C6-O6
7	S	1	NAG	O5-C5-C6-O6
10	t	2	NAG	C4-C5-C6-O6
10	c	2	NAG	C4-C5-C6-O6
10	u	2	NAG	C4-C5-C6-O6
7	j	9	MAN	O5-C5-C6-O6
7	a	1	NAG	O5-C5-C6-O6
7	a	9	MAN	O5-C5-C6-O6
7	S	1	NAG	C4-C5-C6-O6
10	l	2	NAG	C4-C5-C6-O6
11	Z	3	BMA	C4-C5-C6-O6
7	S	9	MAN	O5-C5-C6-O6
11	i	3	BMA	C4-C5-C6-O6
7	a	1	NAG	C4-C5-C6-O6
11	s	3	BMA	C4-C5-C6-O6
7	j	1	NAG	O5-C5-C6-O6
7	j	1	NAG	C4-C5-C6-O6
7	a	6	MAN	O5-C5-C6-O6
7	j	6	MAN	O5-C5-C6-O6
11	Z	1	NAG	O5-C5-C6-O6
7	S	6	MAN	O5-C5-C6-O6
11	i	1	NAG	O5-C5-C6-O6
11	s	1	NAG	O5-C5-C6-O6
9	n	5	MAN	O5-C5-C6-O6
11	Z	3	BMA	O5-C5-C6-O6
10	X	1	NAG	C4-C5-C6-O6
9	d	5	MAN	O5-C5-C6-O6
9	U	5	MAN	O5-C5-C6-O6
11	i	3	BMA	O5-C5-C6-O6
11	s	3	BMA	O5-C5-C6-O6
10	X	1	NAG	O5-C5-C6-O6
9	n	5	MAN	C4-C5-C6-O6
10	l	1	NAG	C4-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
9	d	5	MAN	C4-C5-C6-O6
9	U	5	MAN	C4-C5-C6-O6
7	S	2	NAG	C4-C5-C6-O6
8	r	1	NAG	C4-C5-C6-O6
8	h	1	NAG	C4-C5-C6-O6
10	l	1	NAG	O5-C5-C6-O6
8	Y	1	NAG	C4-C5-C6-O6
7	S	10	MAN	O5-C5-C6-O6
7	S	10	MAN	C4-C5-C6-O6
7	a	2	NAG	C4-C5-C6-O6
10	c	1	NAG	C4-C5-C6-O6
10	c	2	NAG	C1-C2-N2-C7
10	q	2	NAG	C1-C2-N2-C7
10	t	2	NAG	C1-C2-N2-C7
10	2	1	NAG	C4-C5-C6-O6
10	c	1	NAG	O5-C5-C6-O6
7	a	10	MAN	O5-C5-C6-O6
10	g	1	NAG	C4-C5-C6-O6
7	a	10	MAN	C4-C5-C6-O6
11	Z	2	NAG	C4-C5-C6-O6

There are no ring outliers.

37 monomers are involved in 40 short contacts:

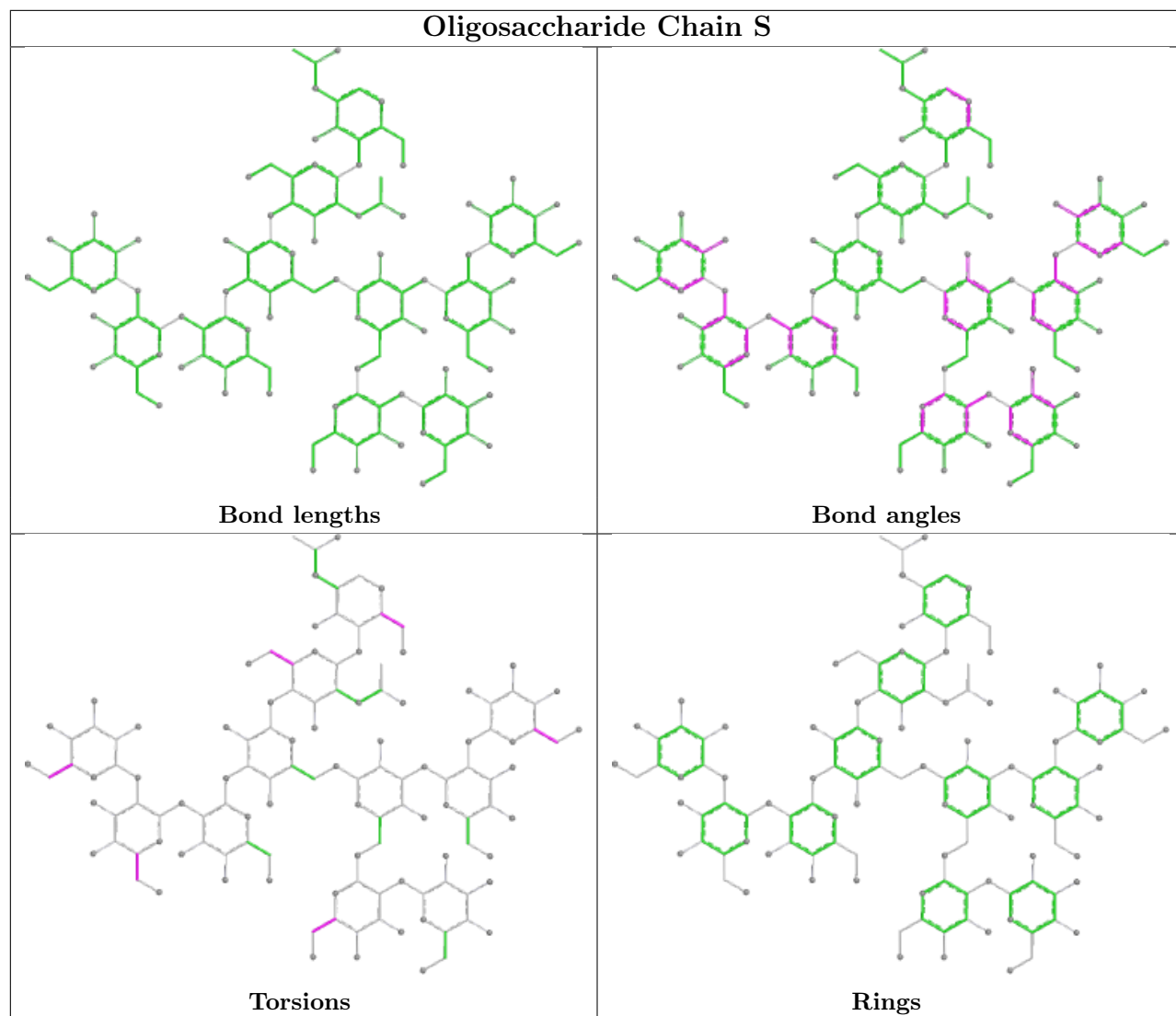
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	S	6	MAN	1	0
10	e	2	NAG	1	0
8	r	1	NAG	1	0
9	d	5	MAN	3	0
9	d	6	MAN	1	0
11	s	2	NAG	1	0
10	e	1	NAG	1	0
8	r	4	MAN	1	0
8	r	2	NAG	1	0
9	U	5	MAN	3	0
11	Z	2	NAG	1	0
10	o	2	NAG	1	0
10	V	1	NAG	1	0
10	x	1	NAG	1	0
9	n	6	MAN	1	0
10	w	1	NAG	1	0
7	a	1	NAG	1	0

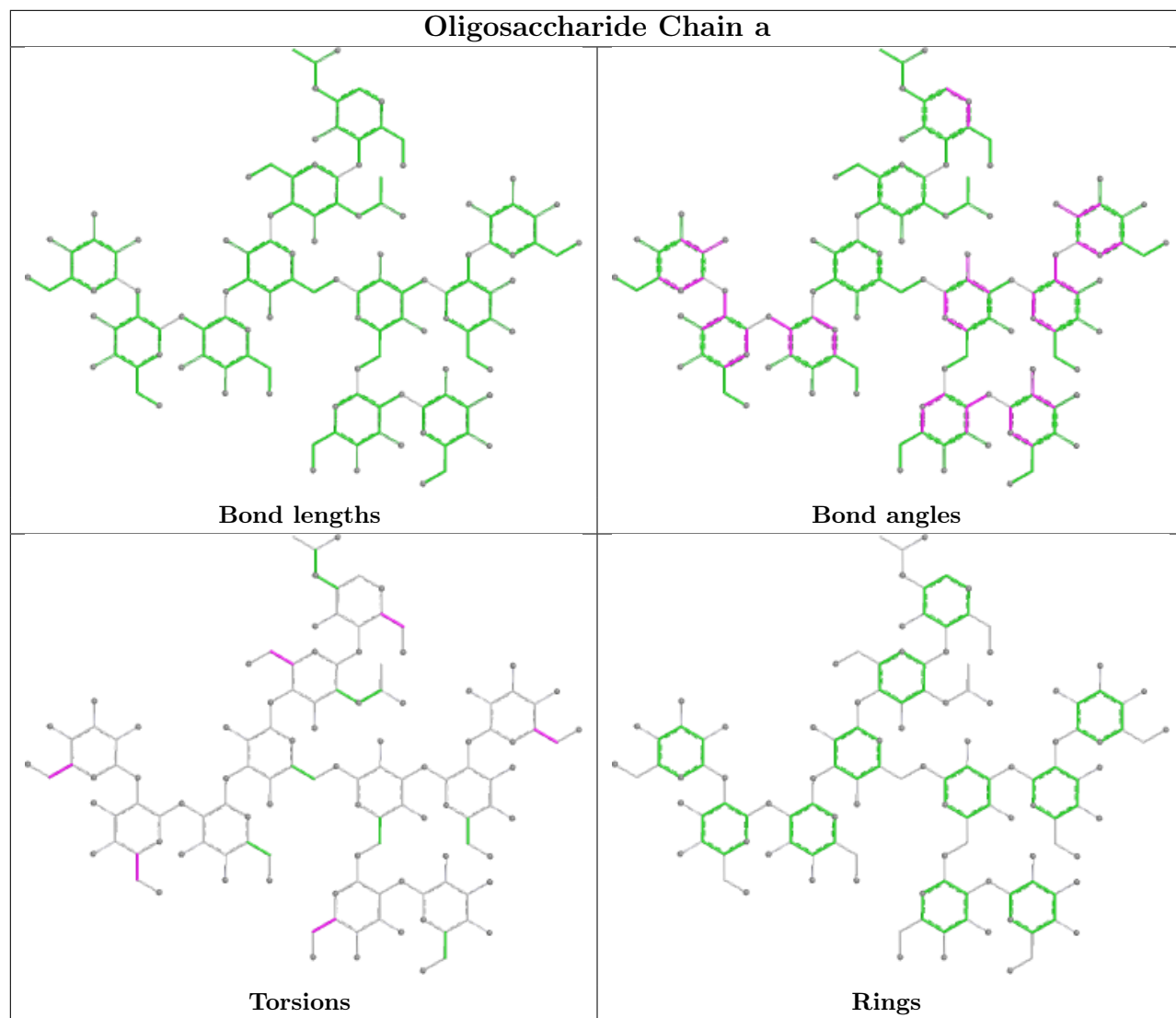
Continued on next page...

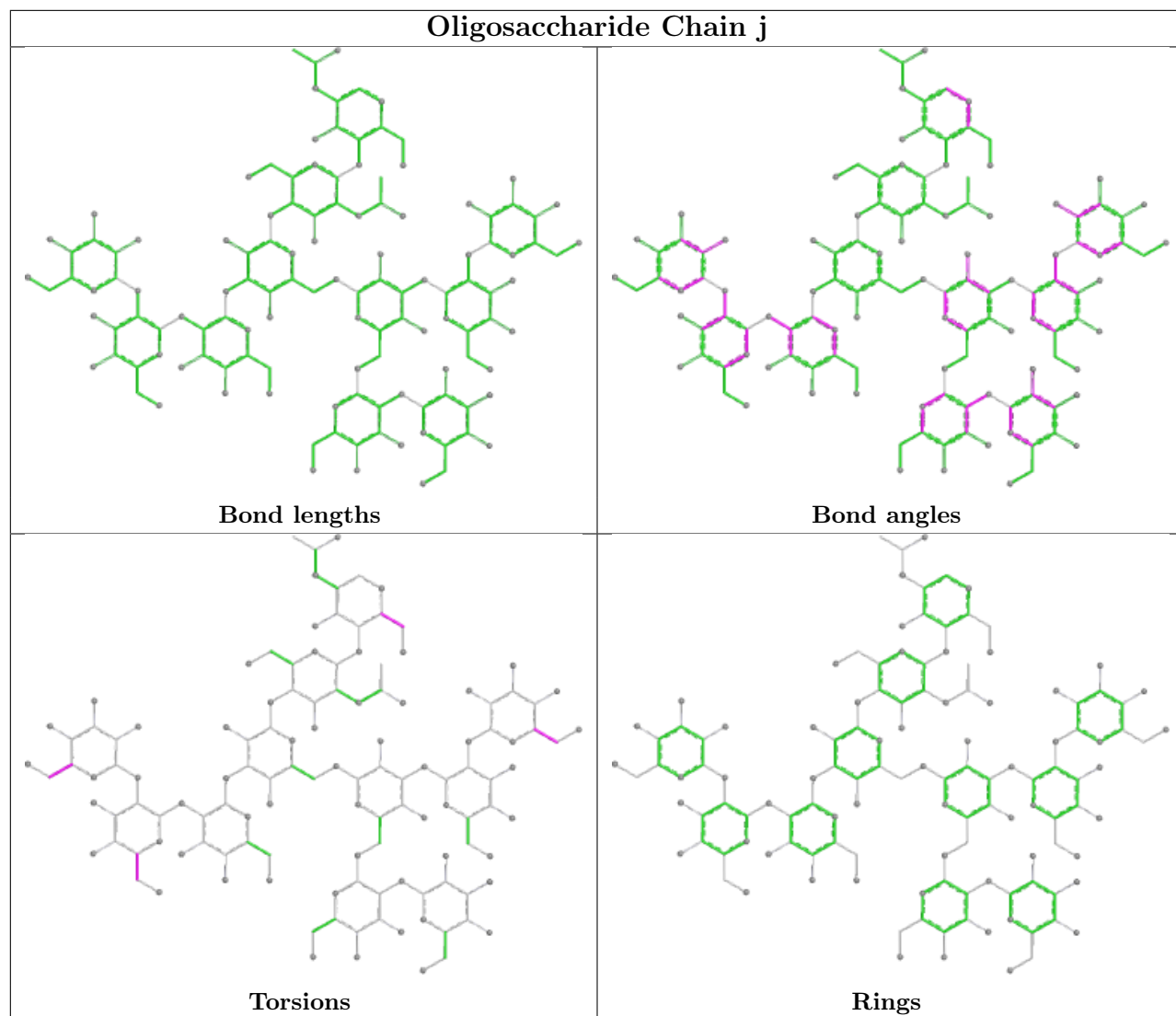
Continued from previous page...

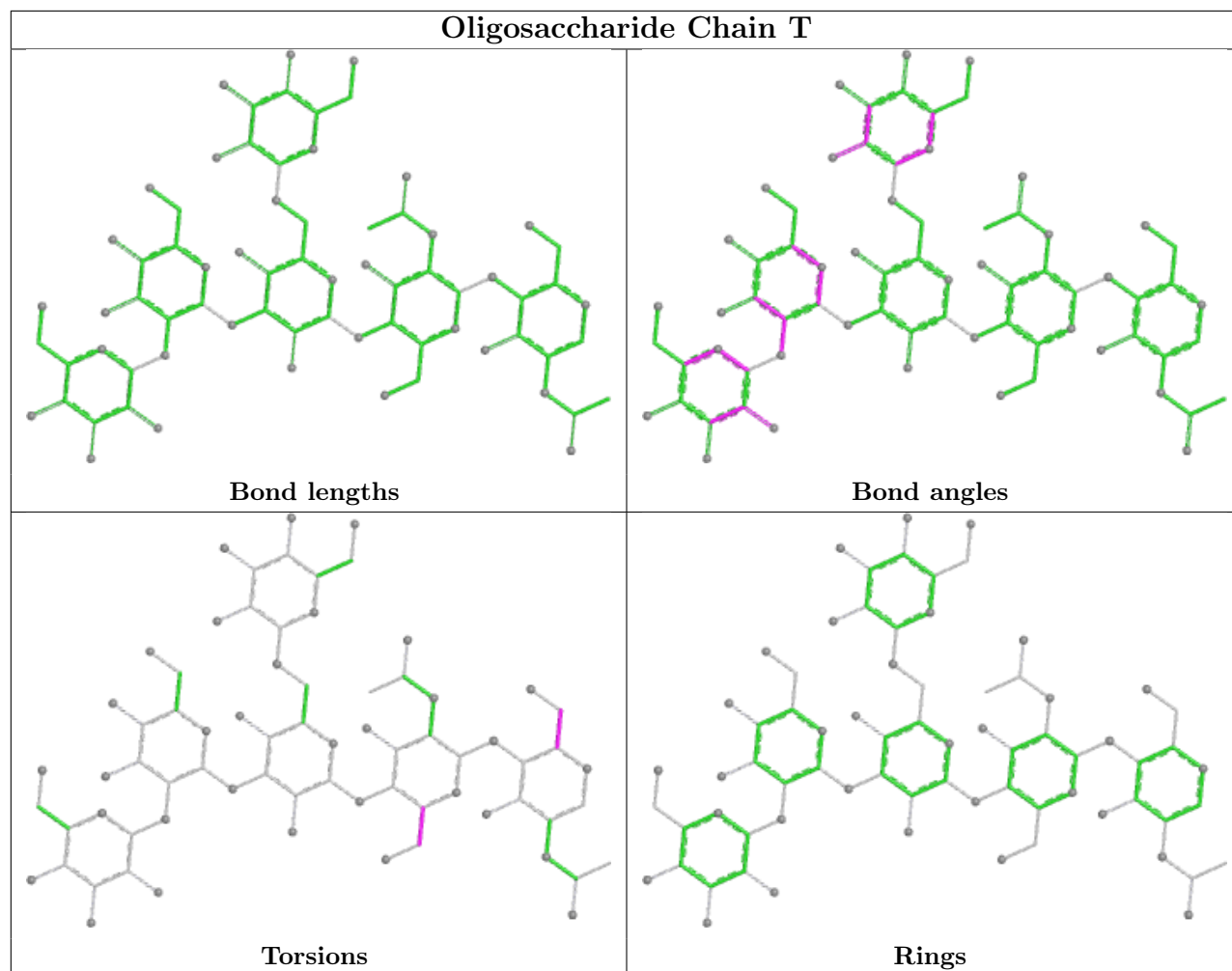
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	h	1	NAG	1	0
11	s	1	NAG	2	0
9	U	6	MAN	1	0
10	z	1	NAG	1	0
10	y	1	NAG	1	0
10	o	1	NAG	1	0
10	w	2	NAG	1	0
8	Y	1	NAG	1	0
10	v	2	NAG	1	0
8	h	4	MAN	1	0
7	a	6	MAN	1	0
8	Y	2	NAG	1	0
11	i	1	NAG	2	0
7	j	1	NAG	1	0
11	Z	1	NAG	1	0
10	V	2	NAG	1	0
9	n	5	MAN	3	0
8	Y	4	MAN	1	0
10	v	1	NAG	1	0
8	h	2	NAG	1	0

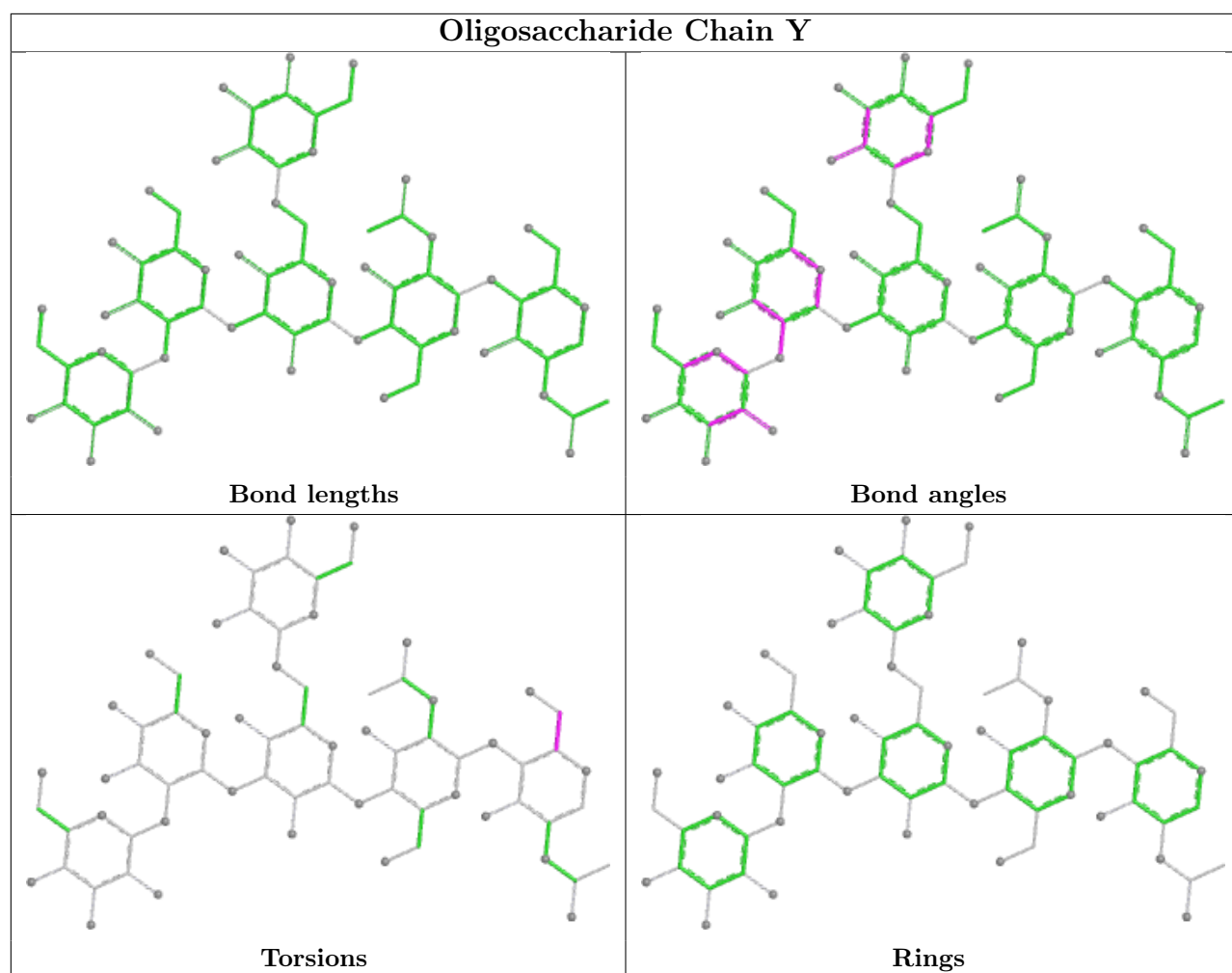
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

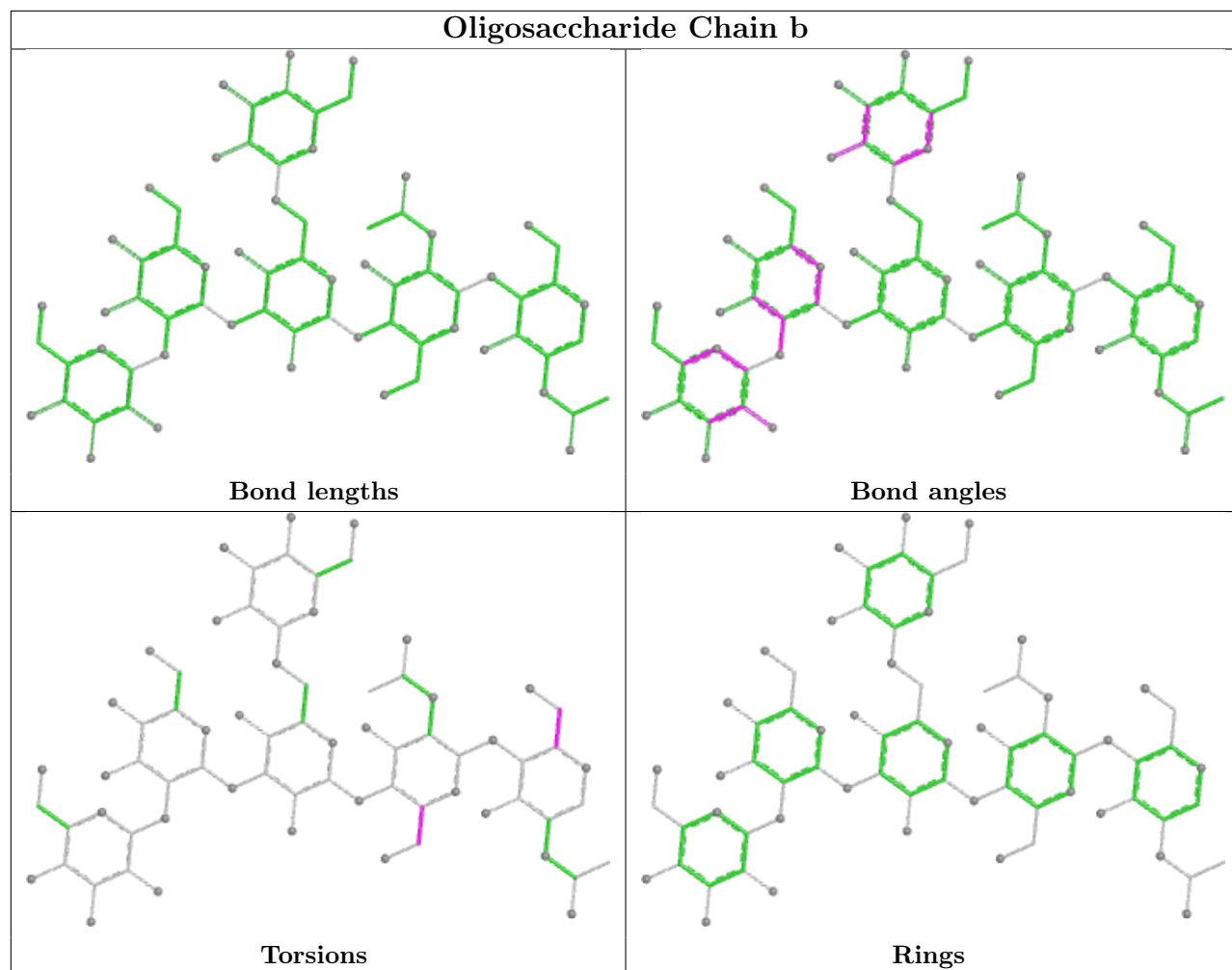


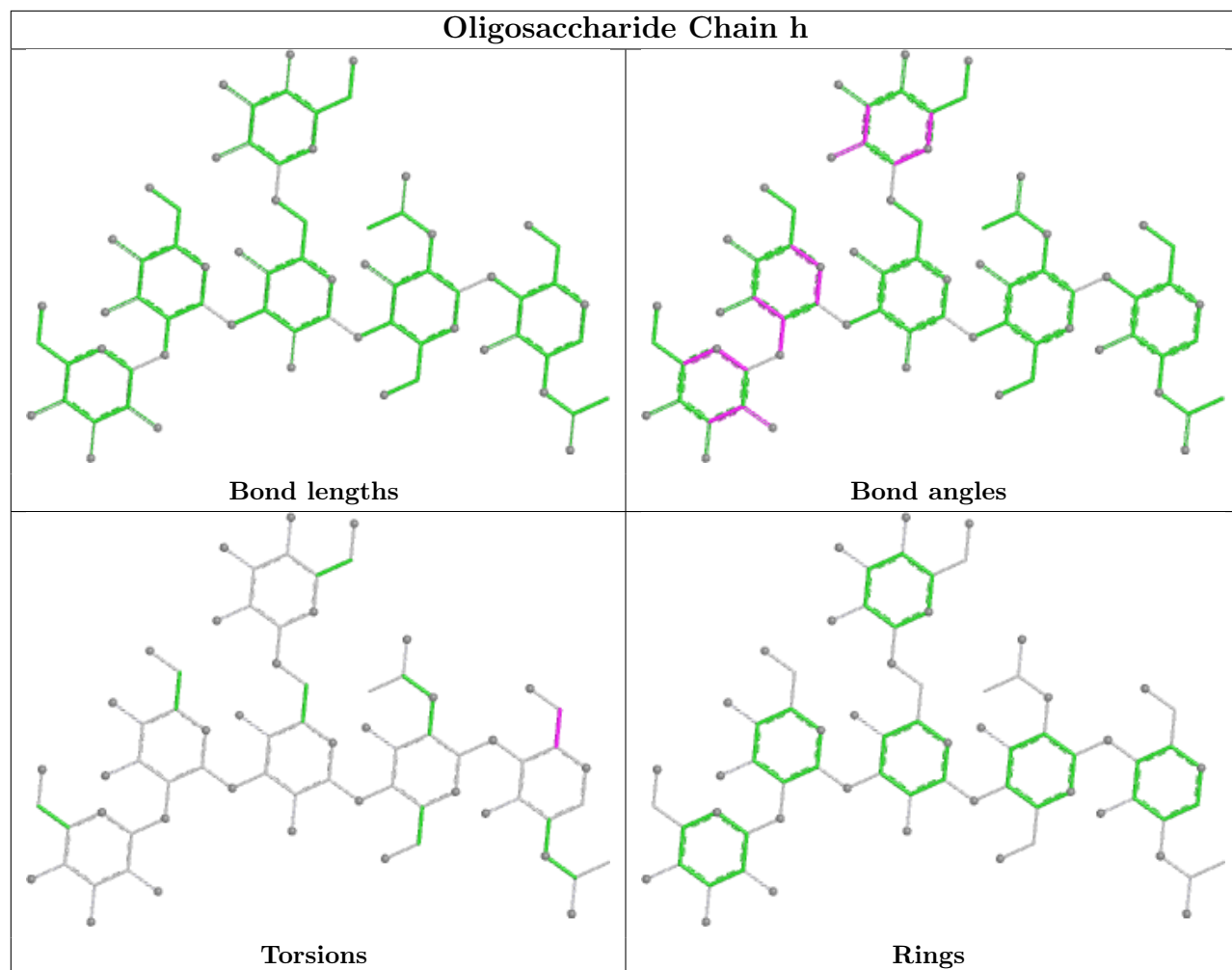


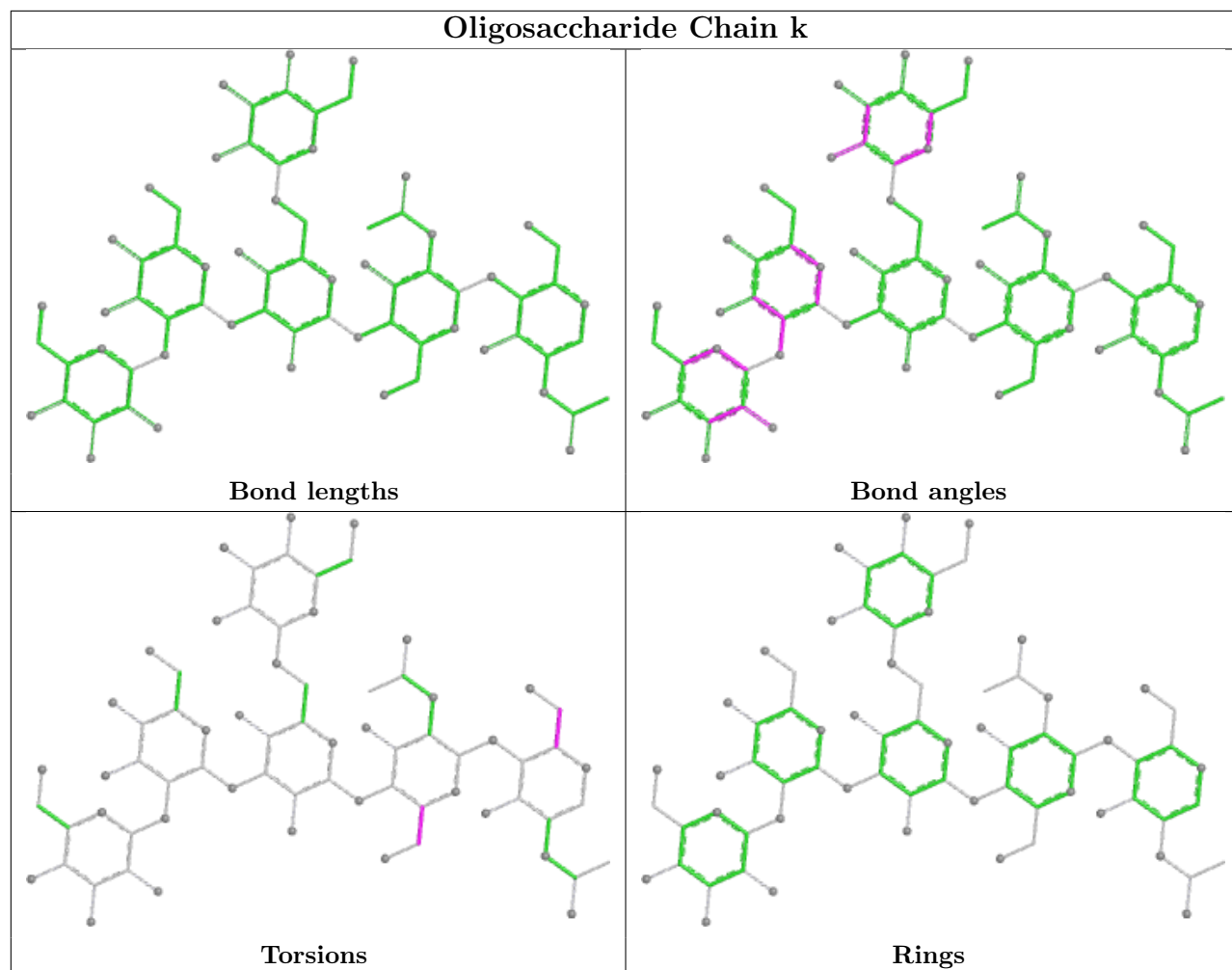


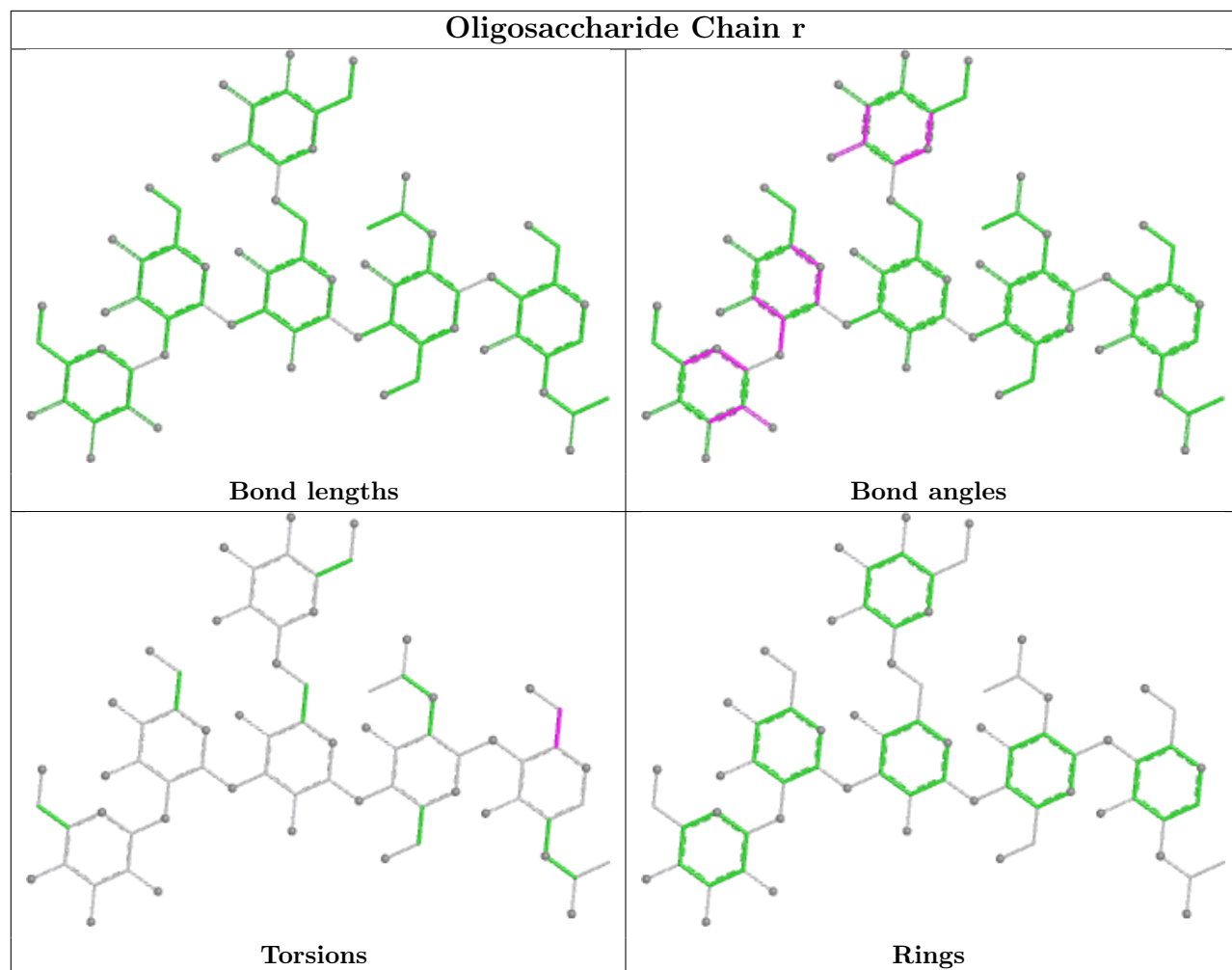


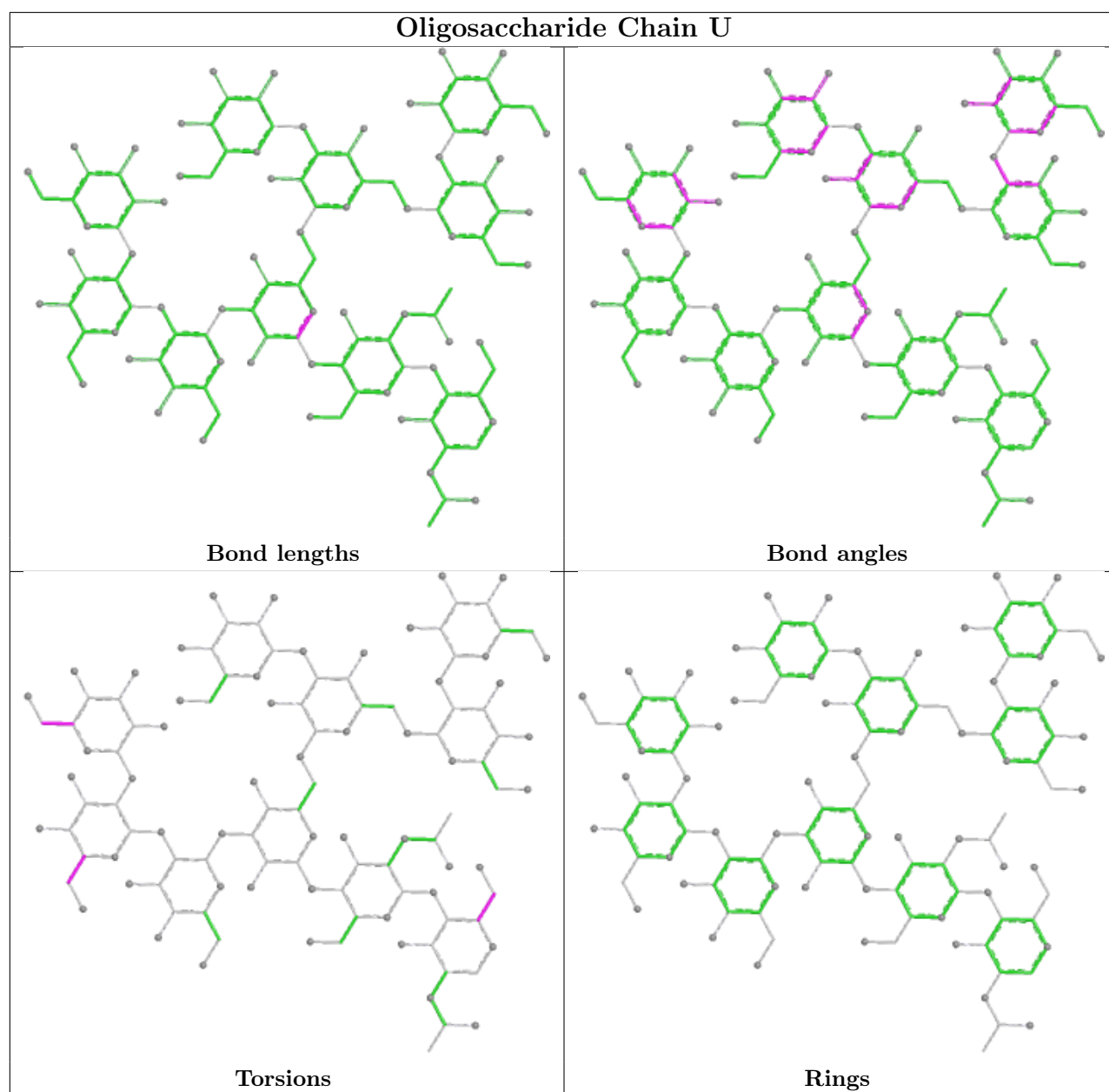


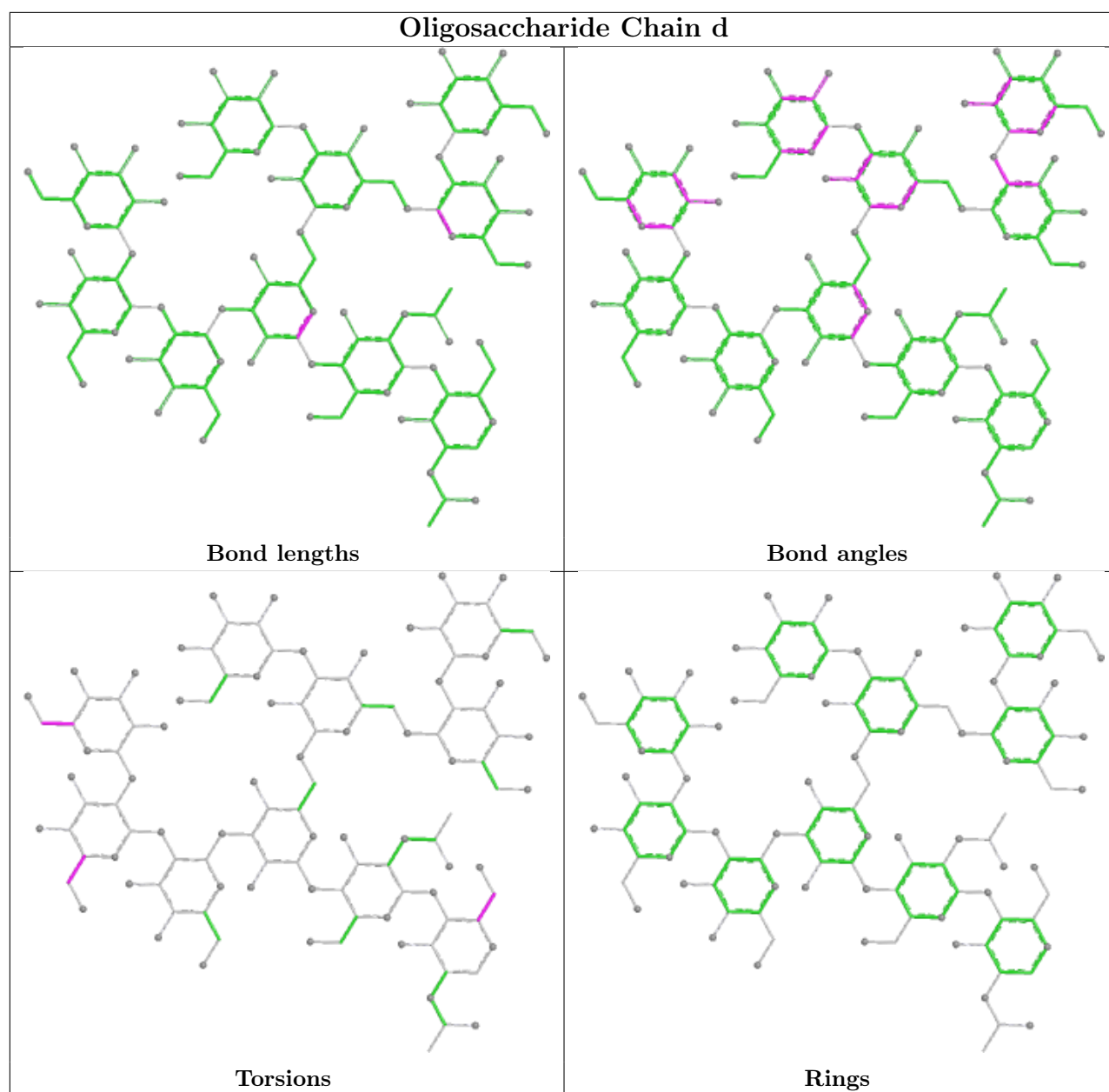


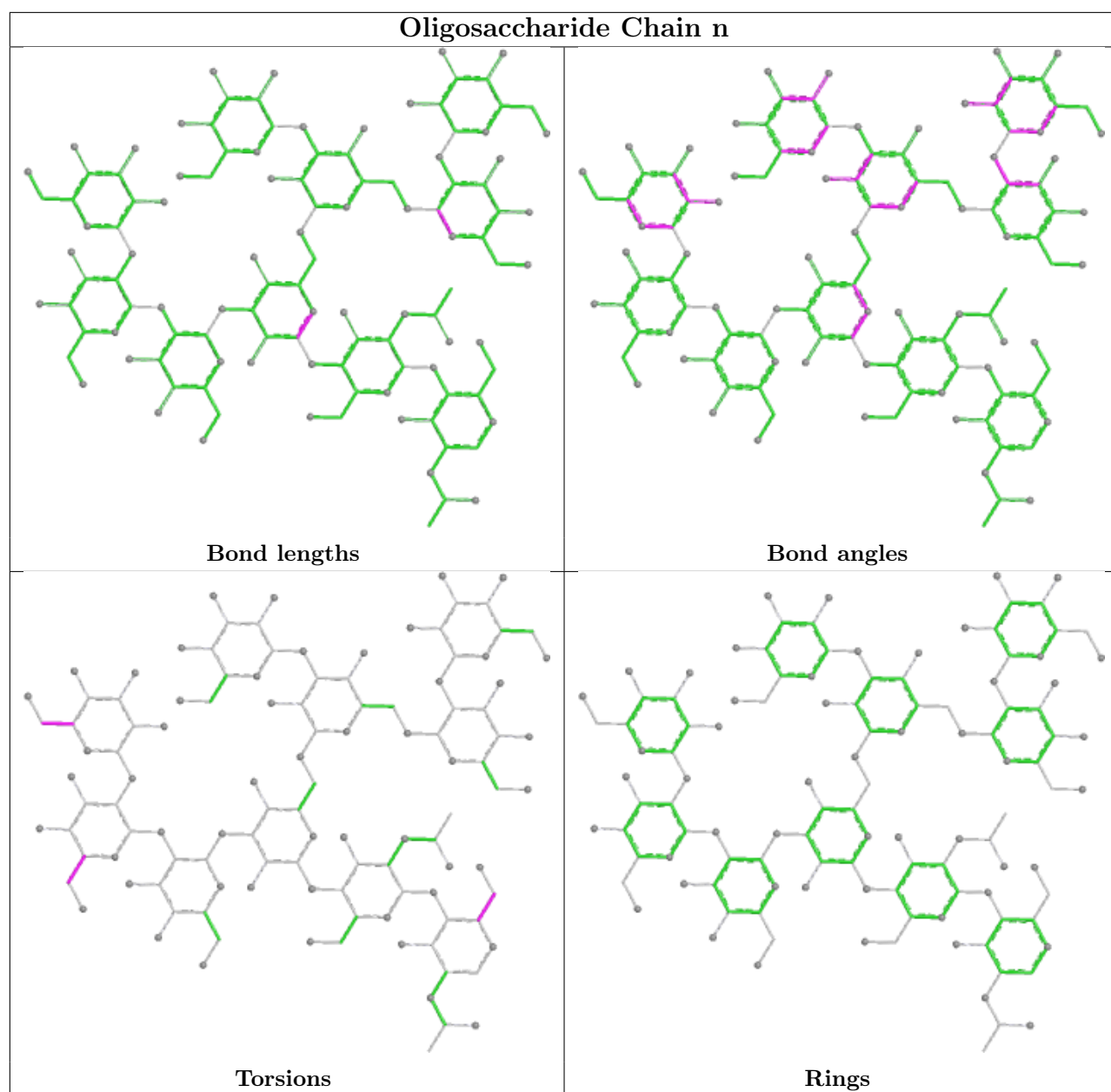


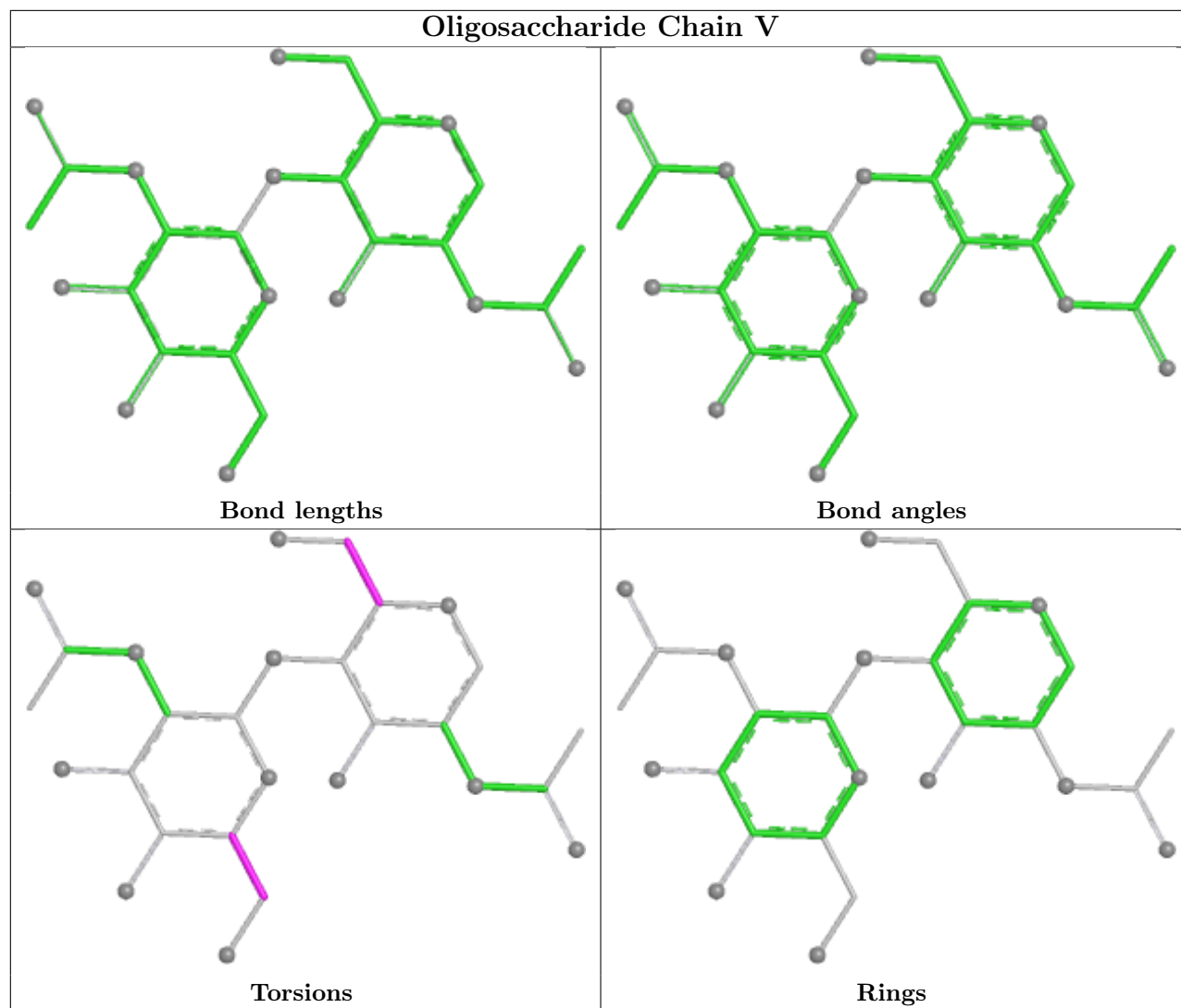


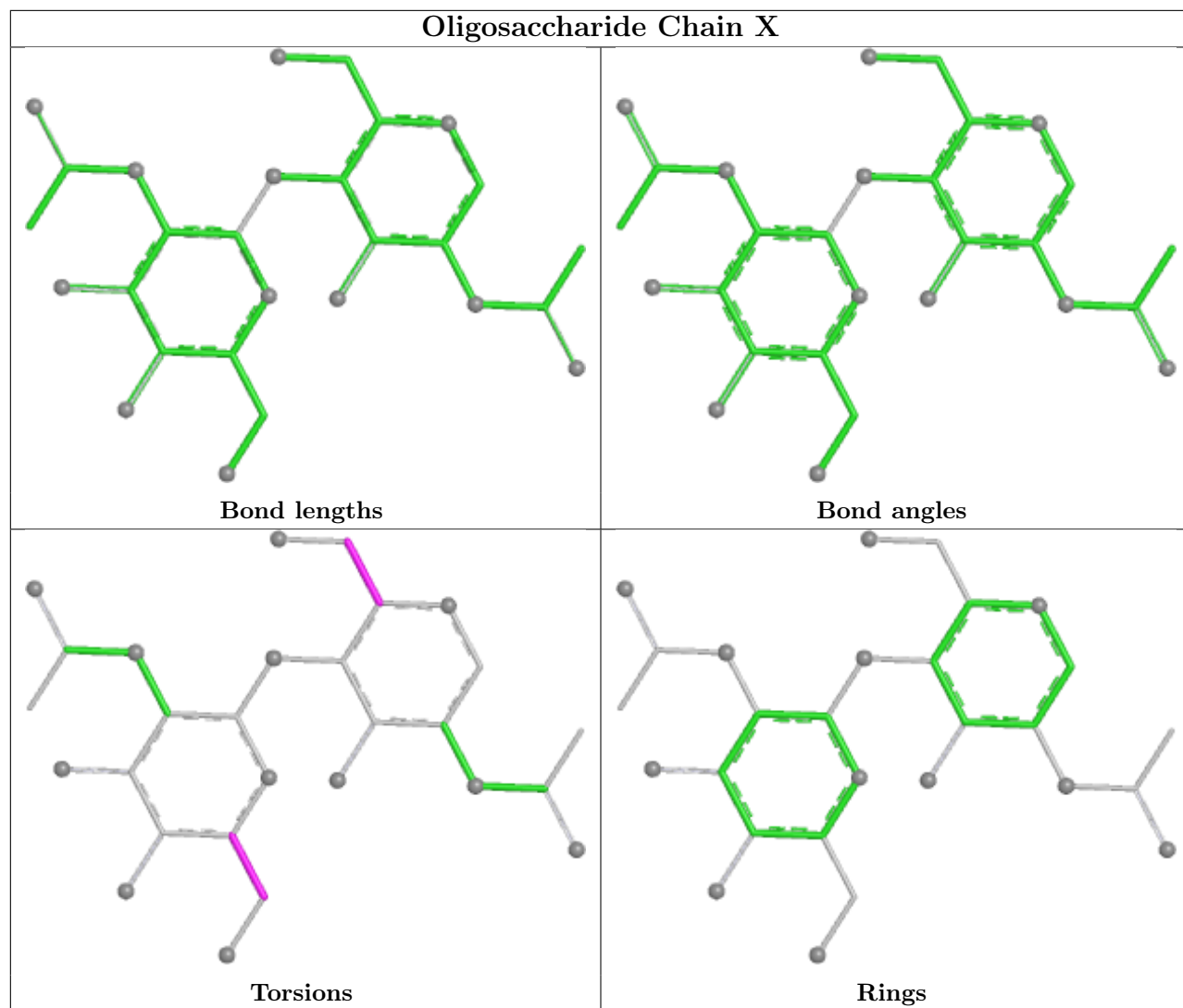


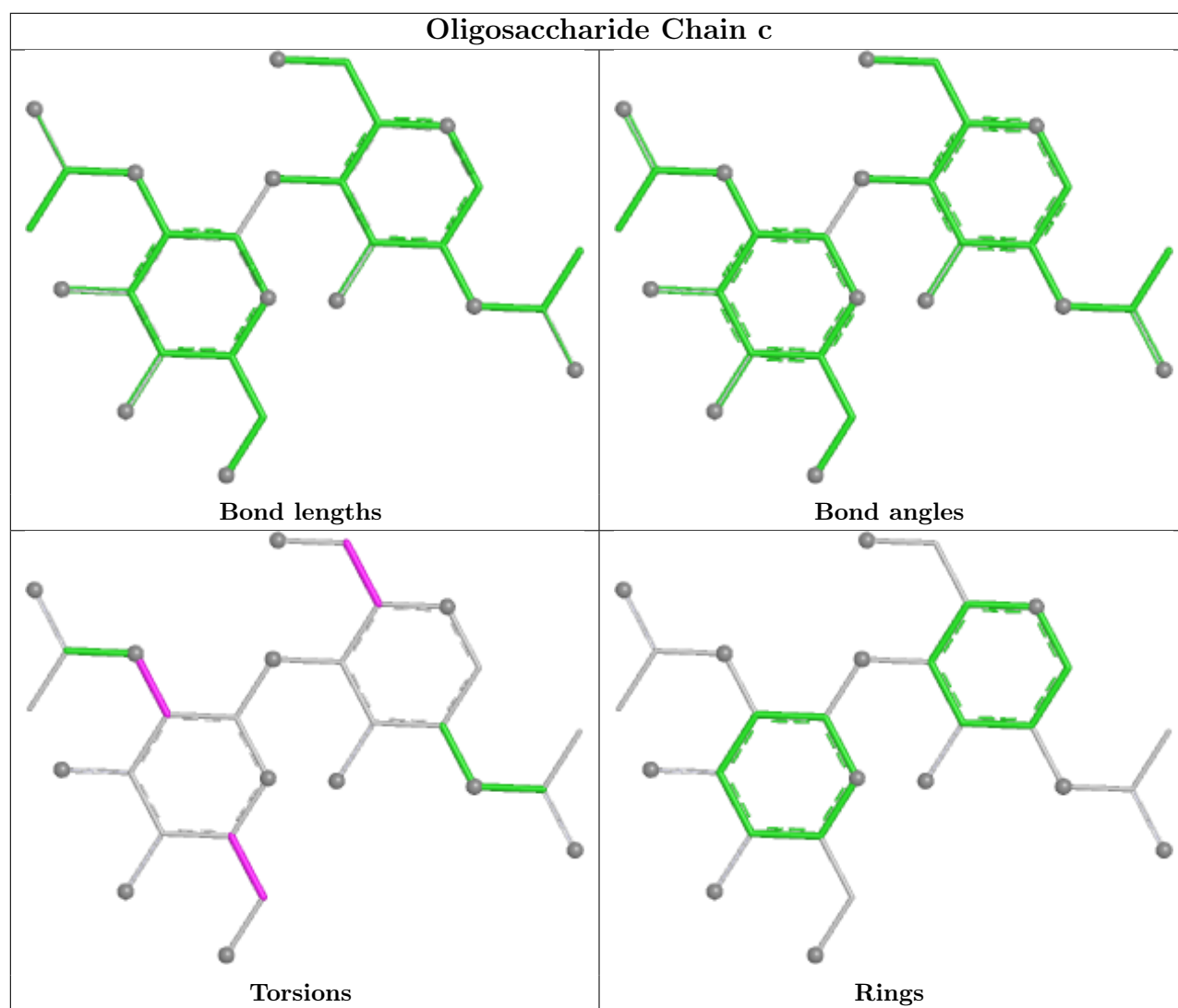


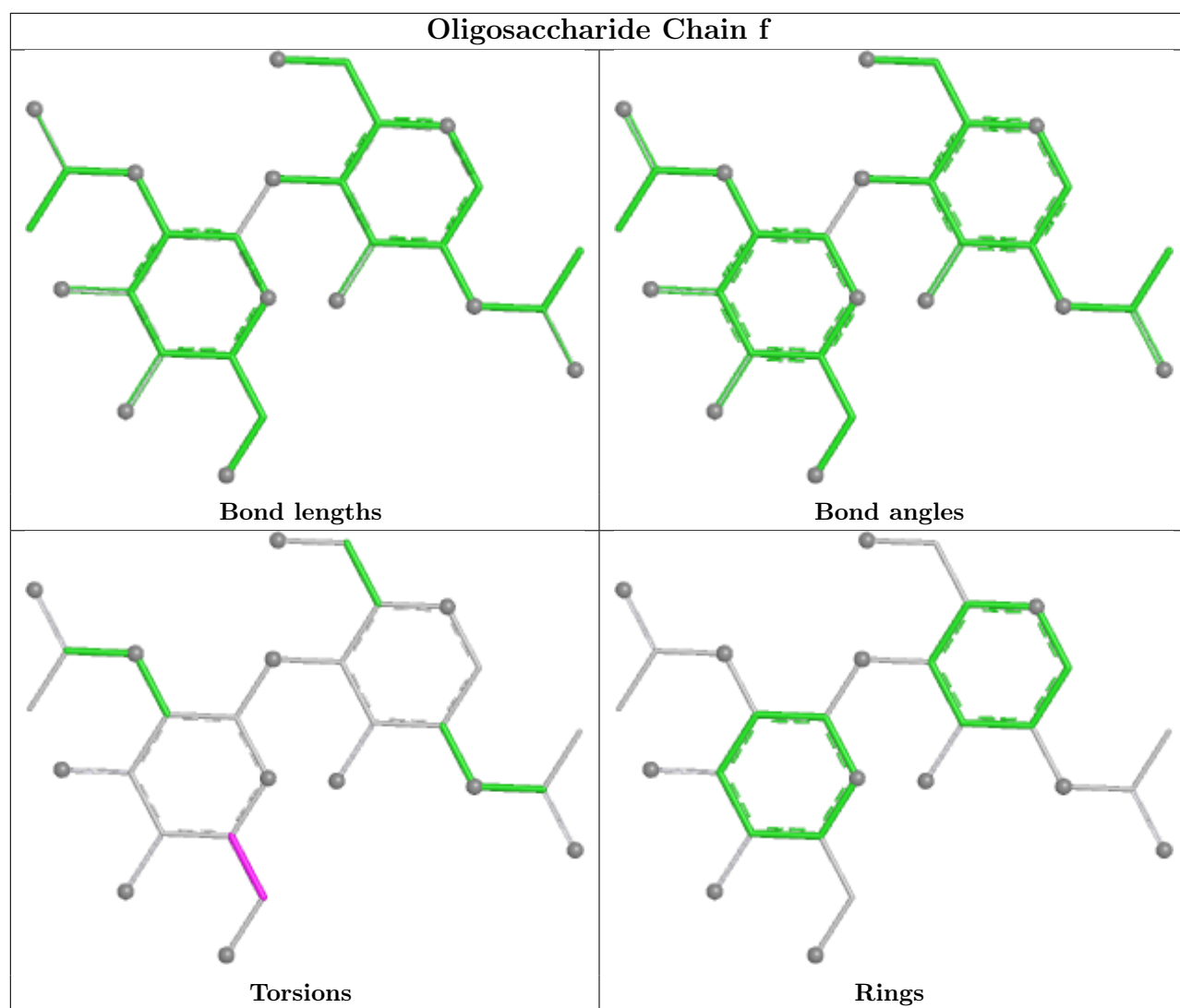


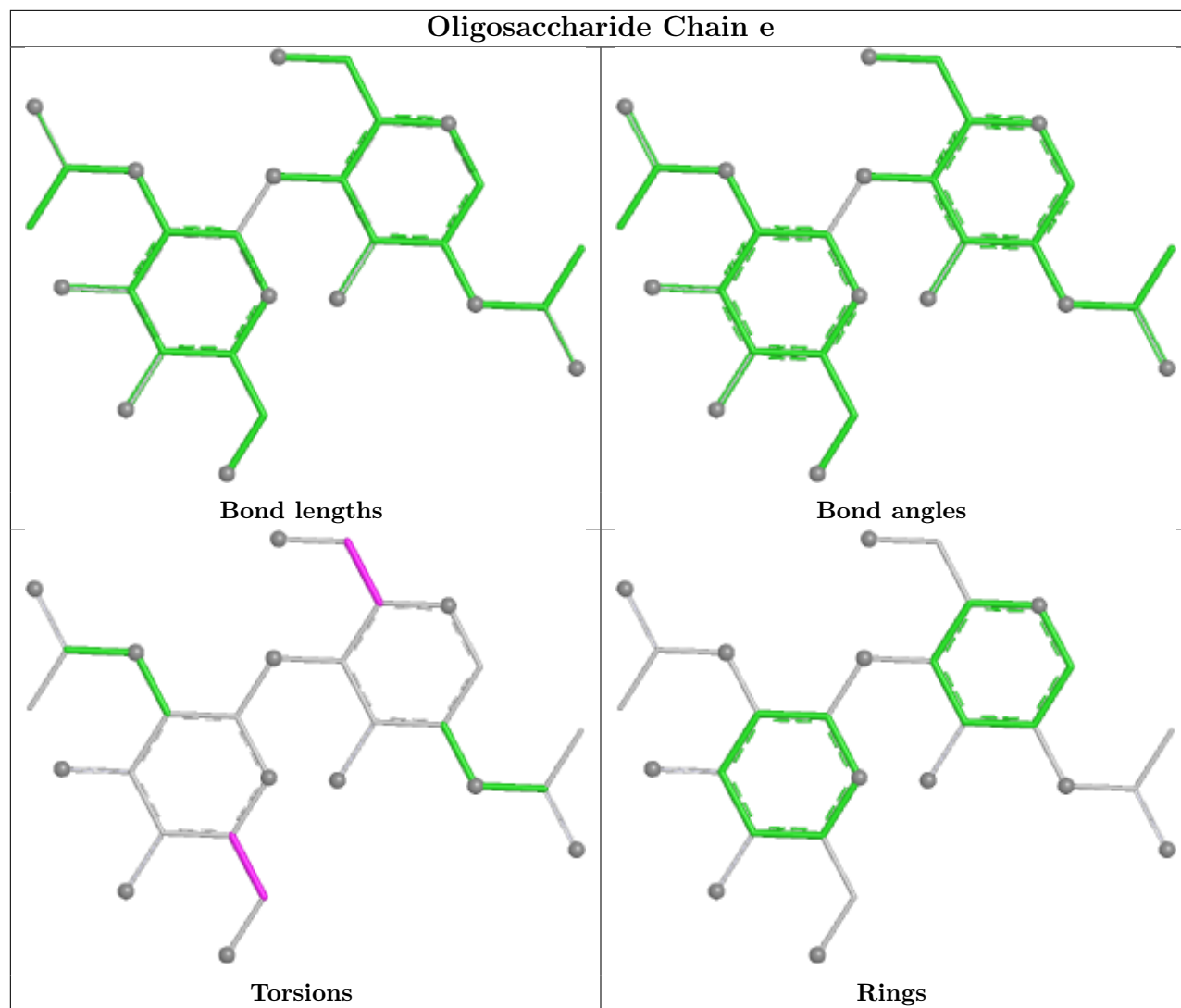


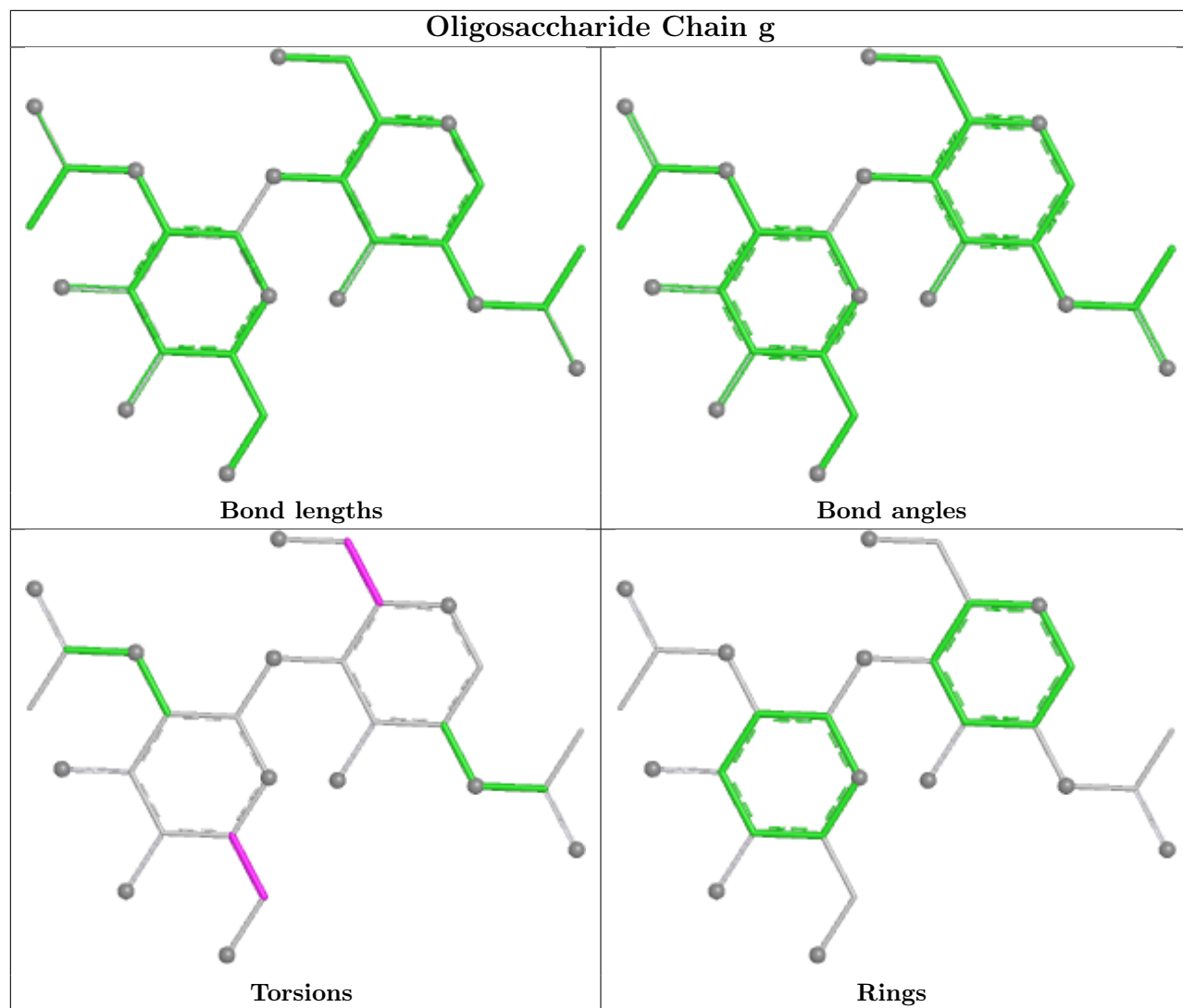


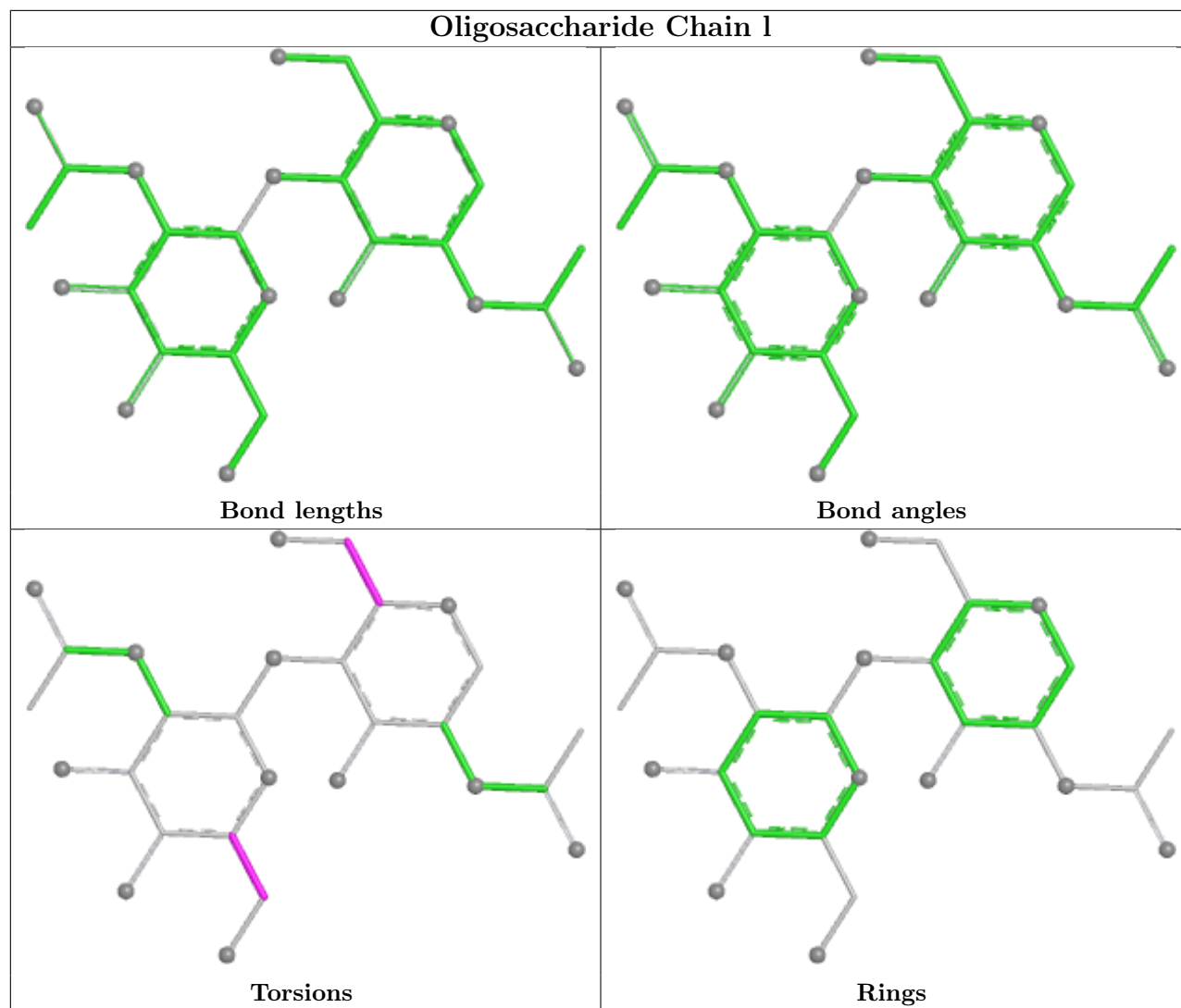


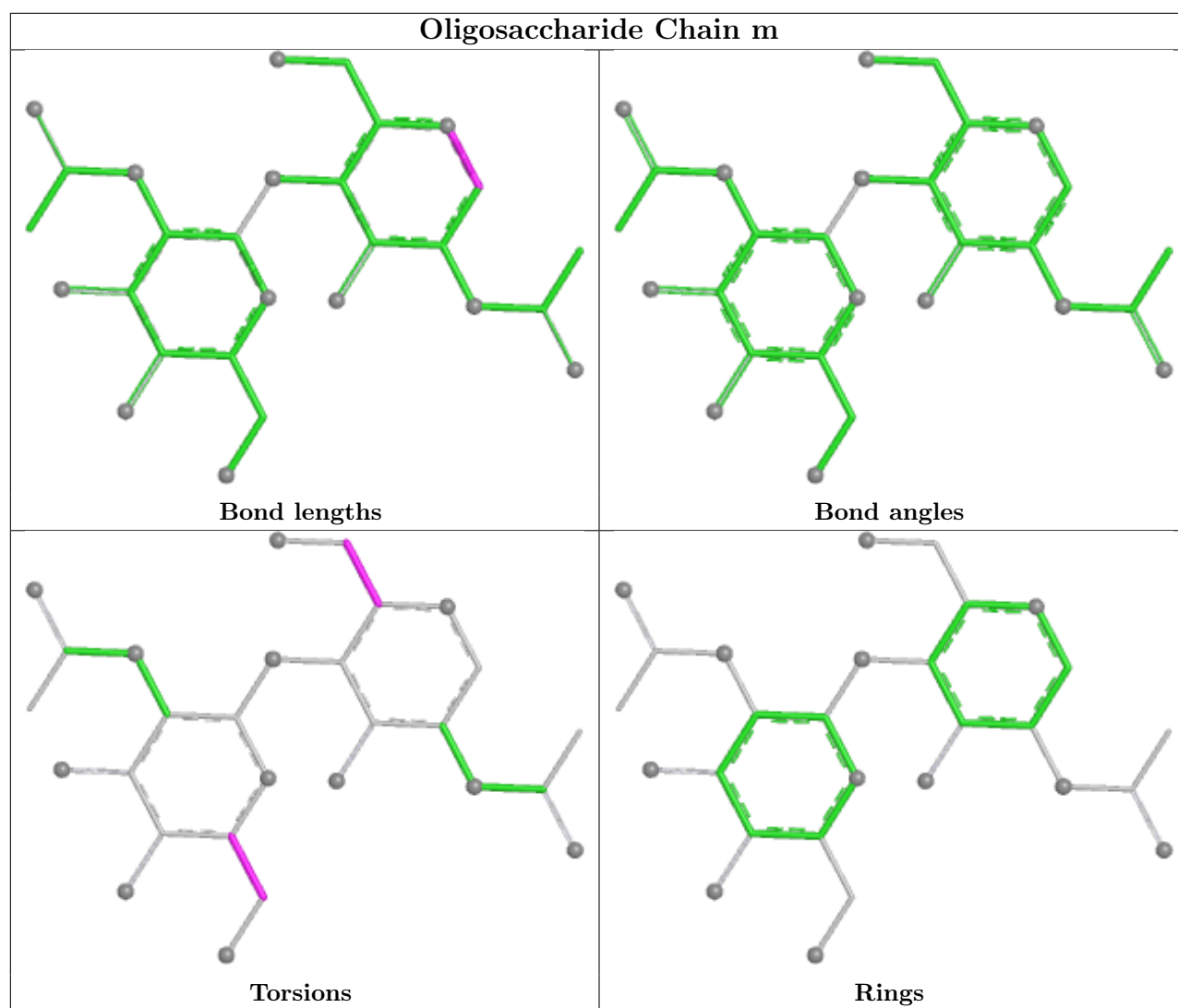


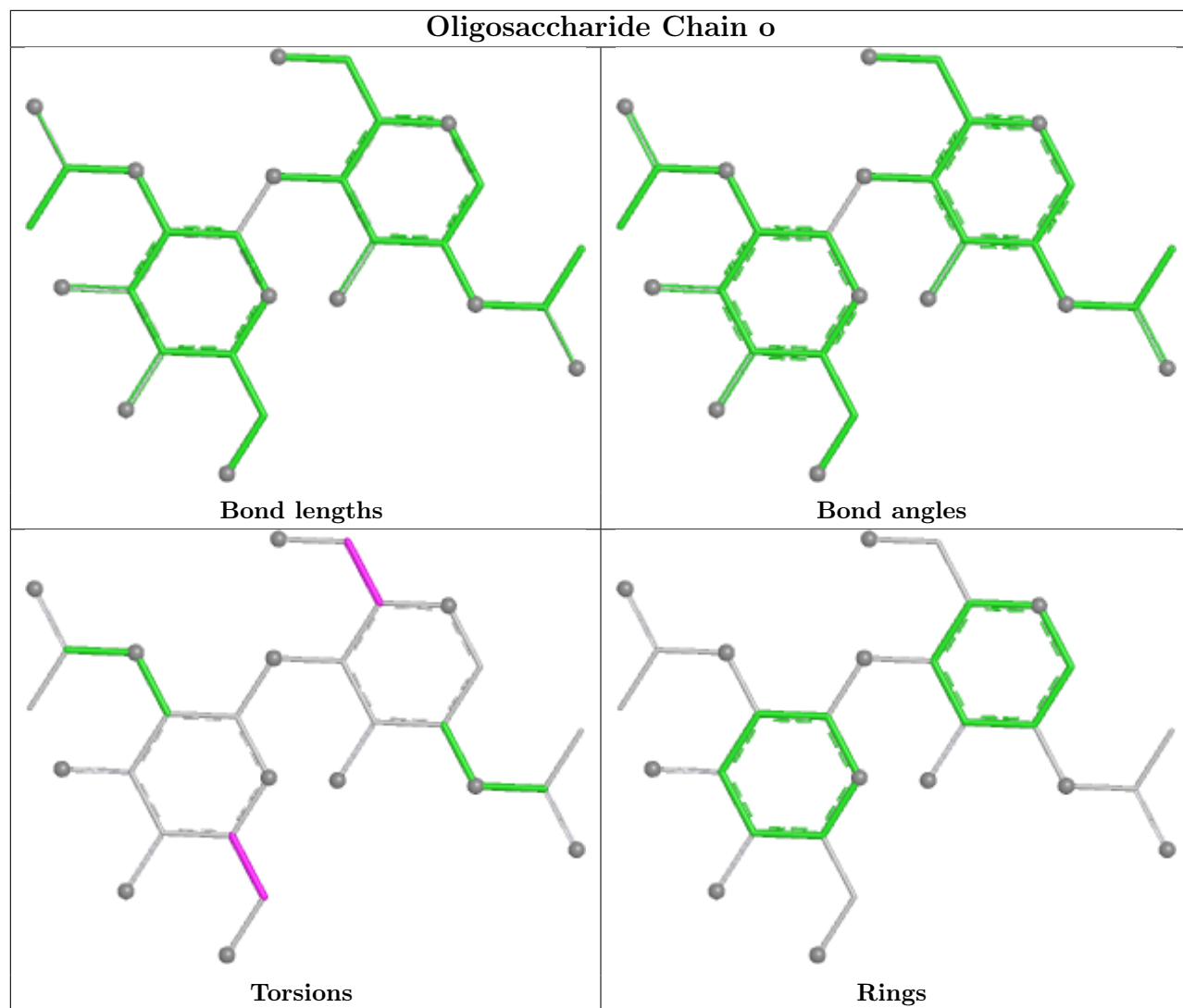


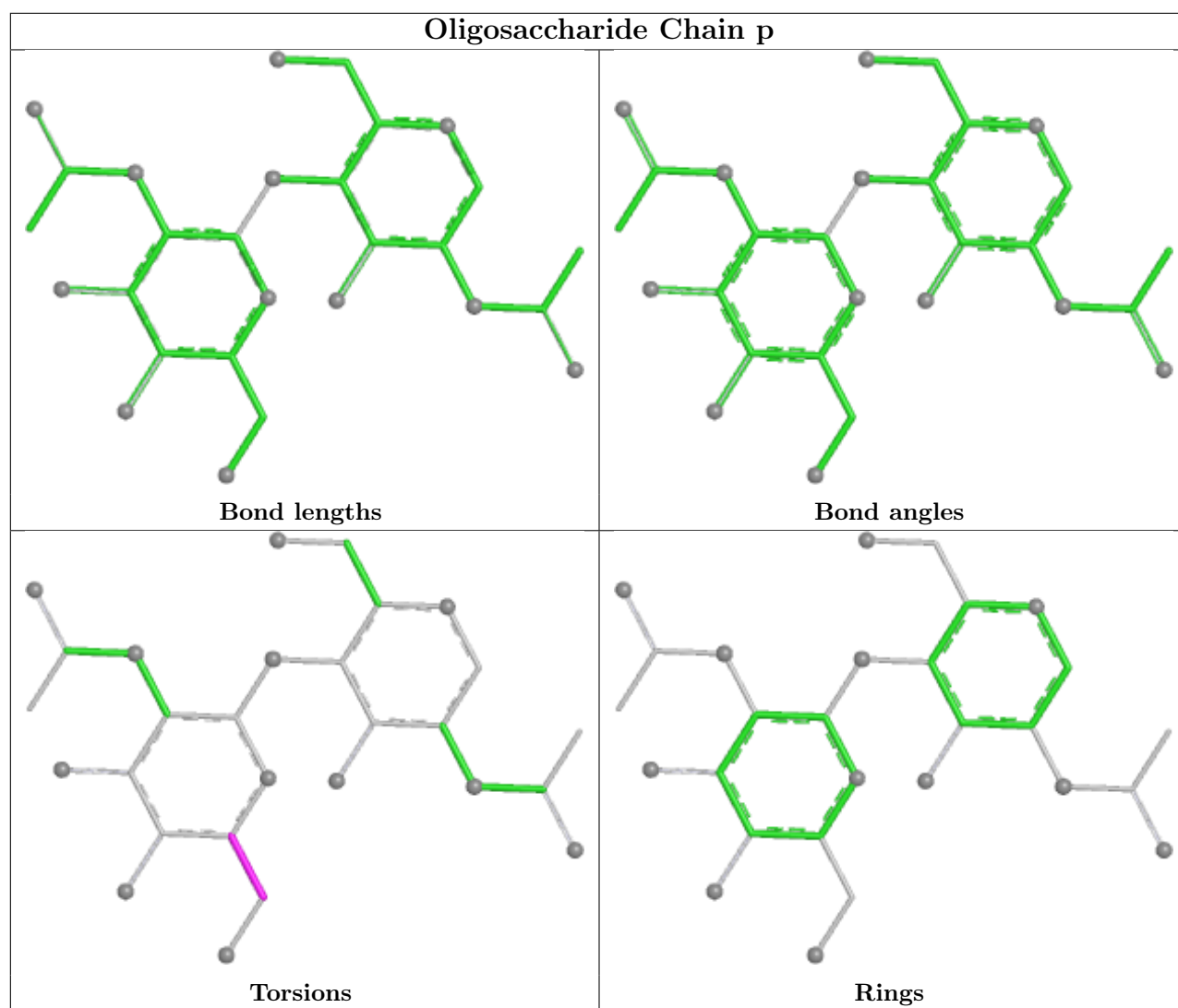


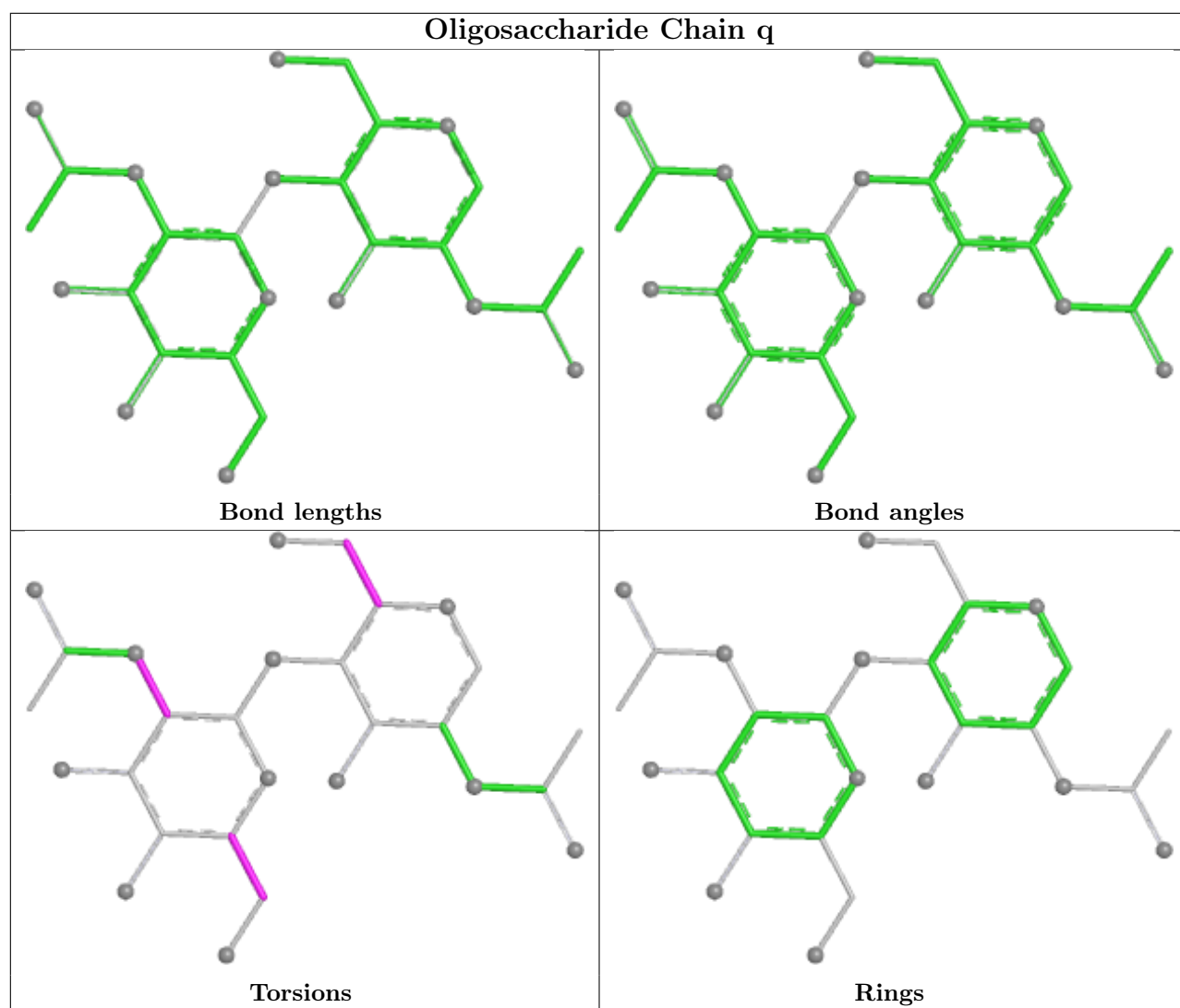


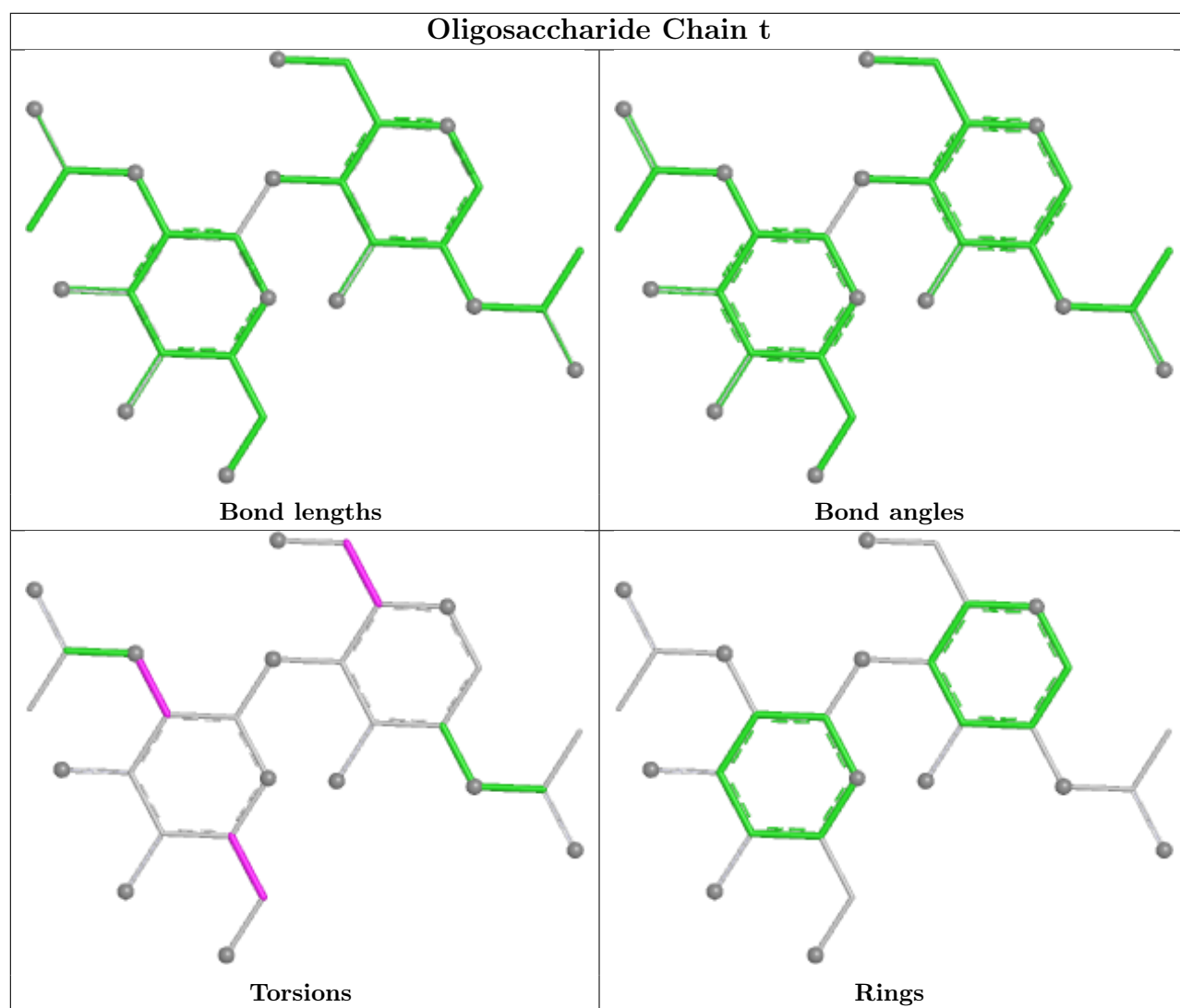


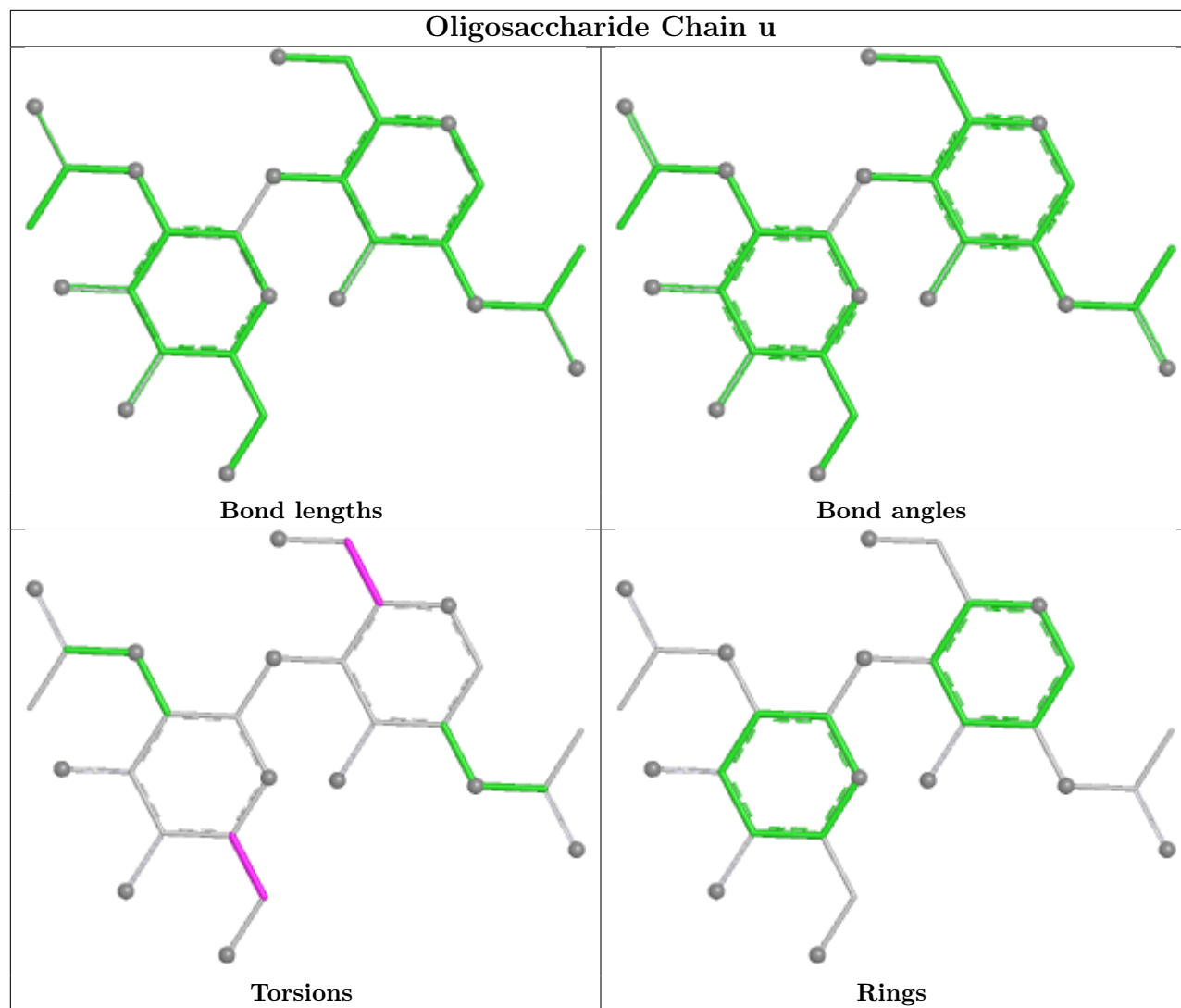


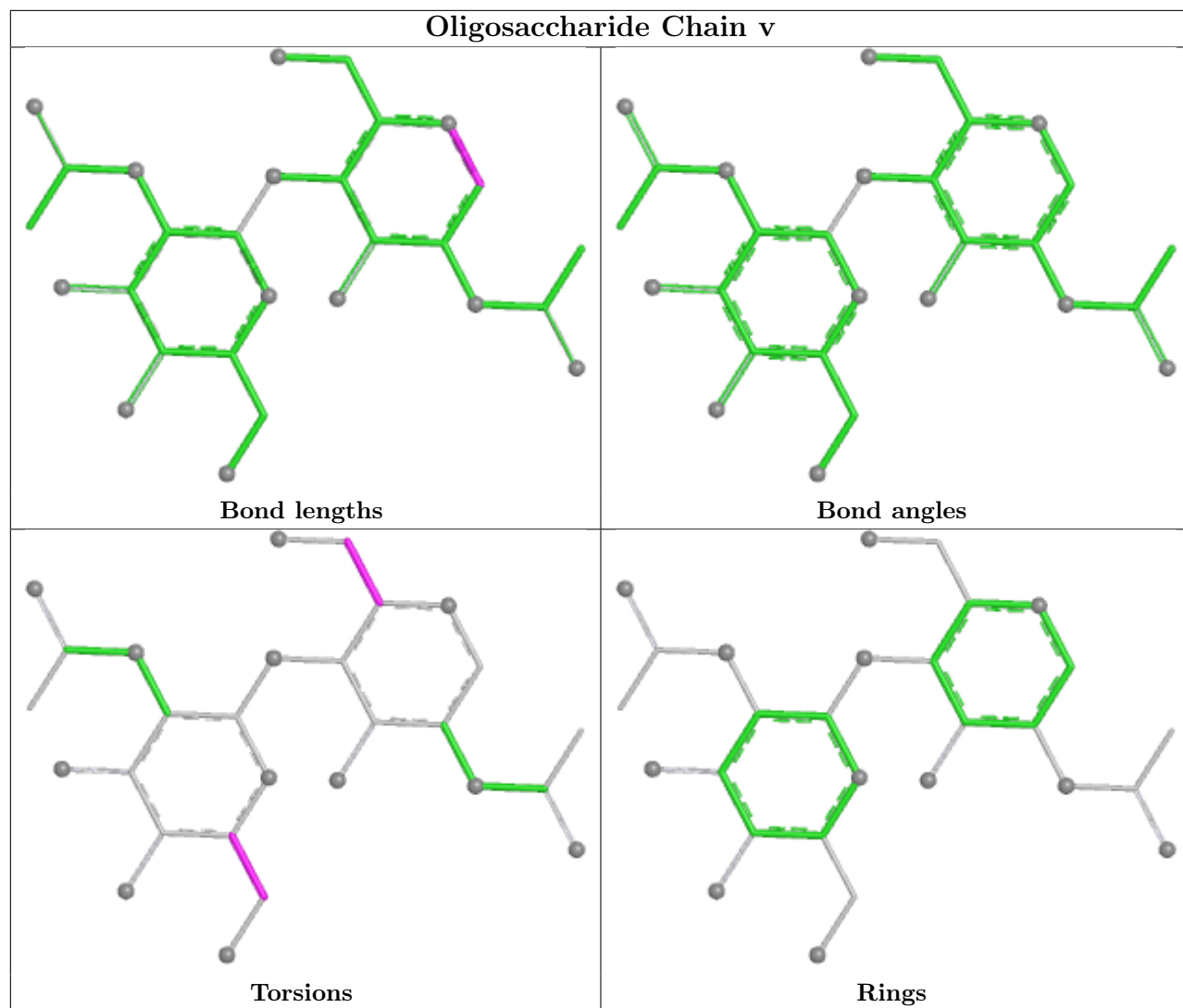


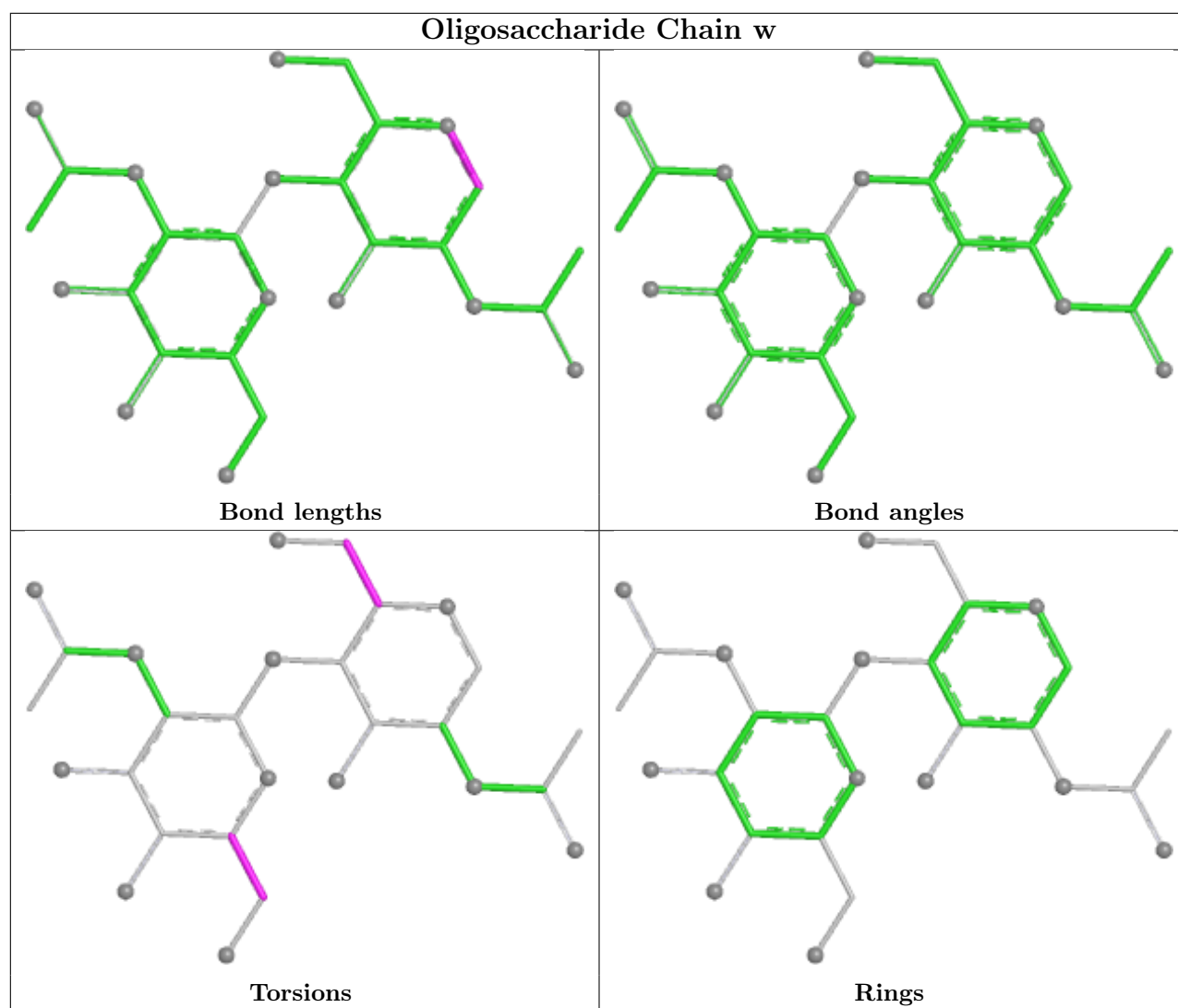


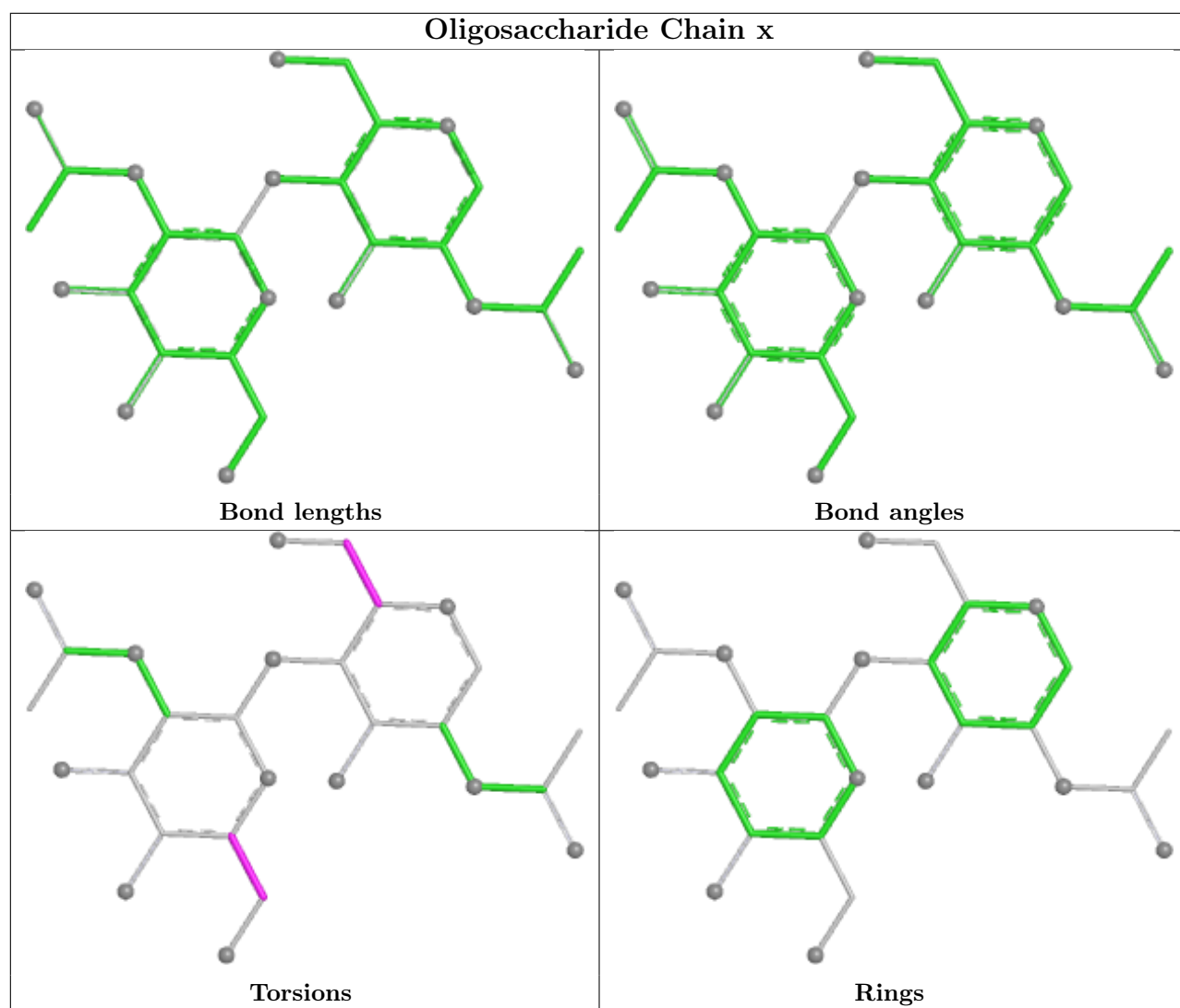


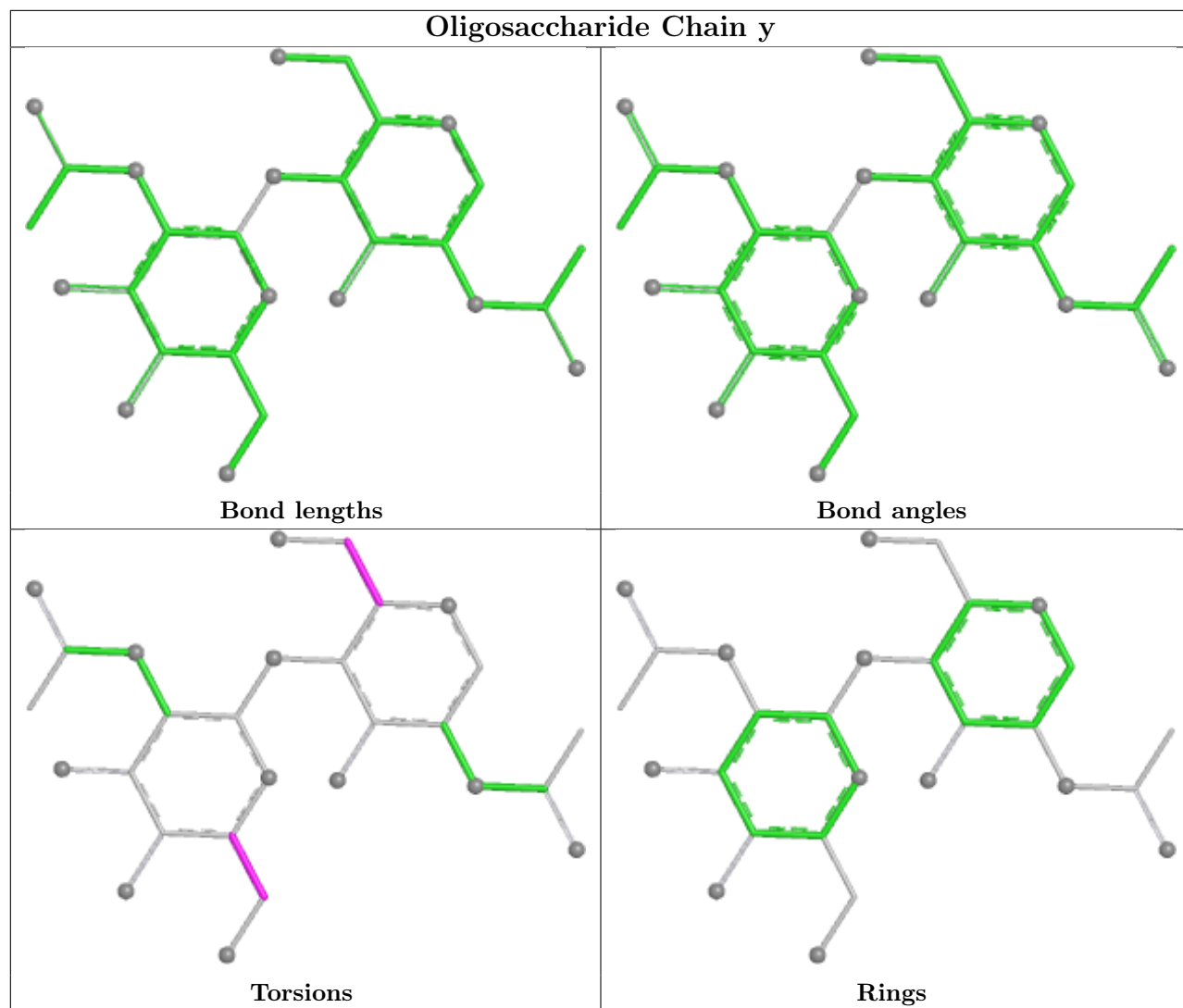


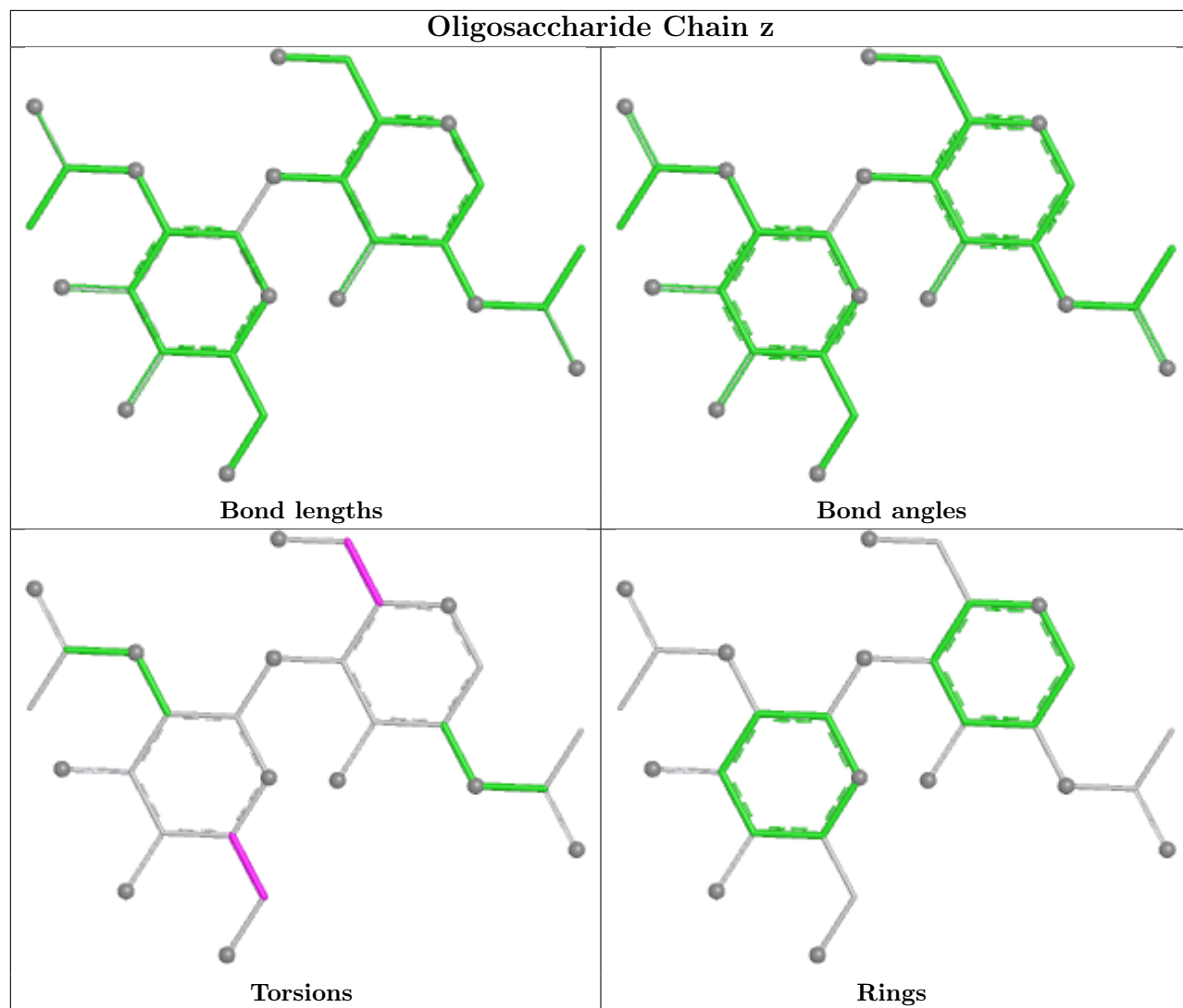


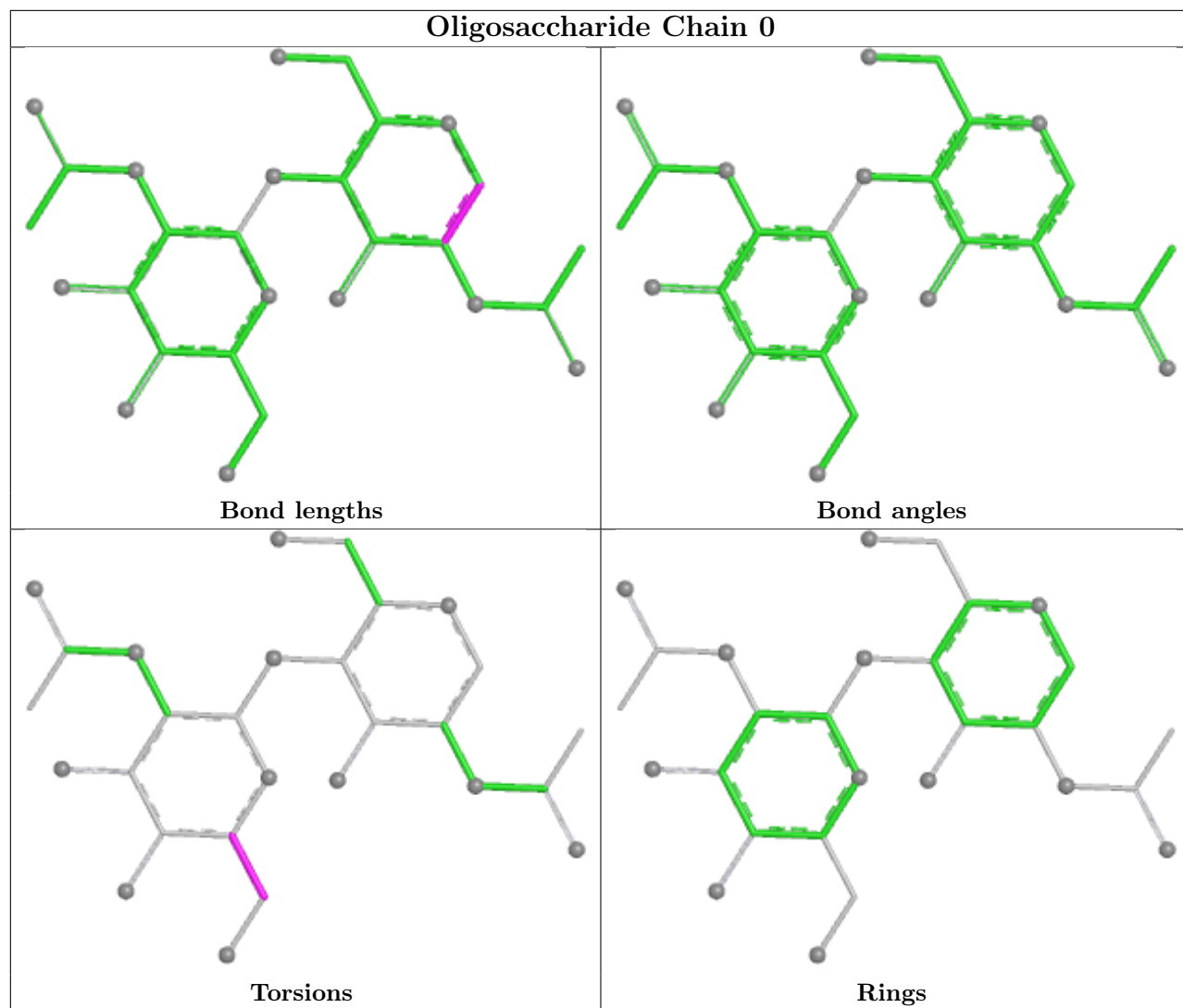


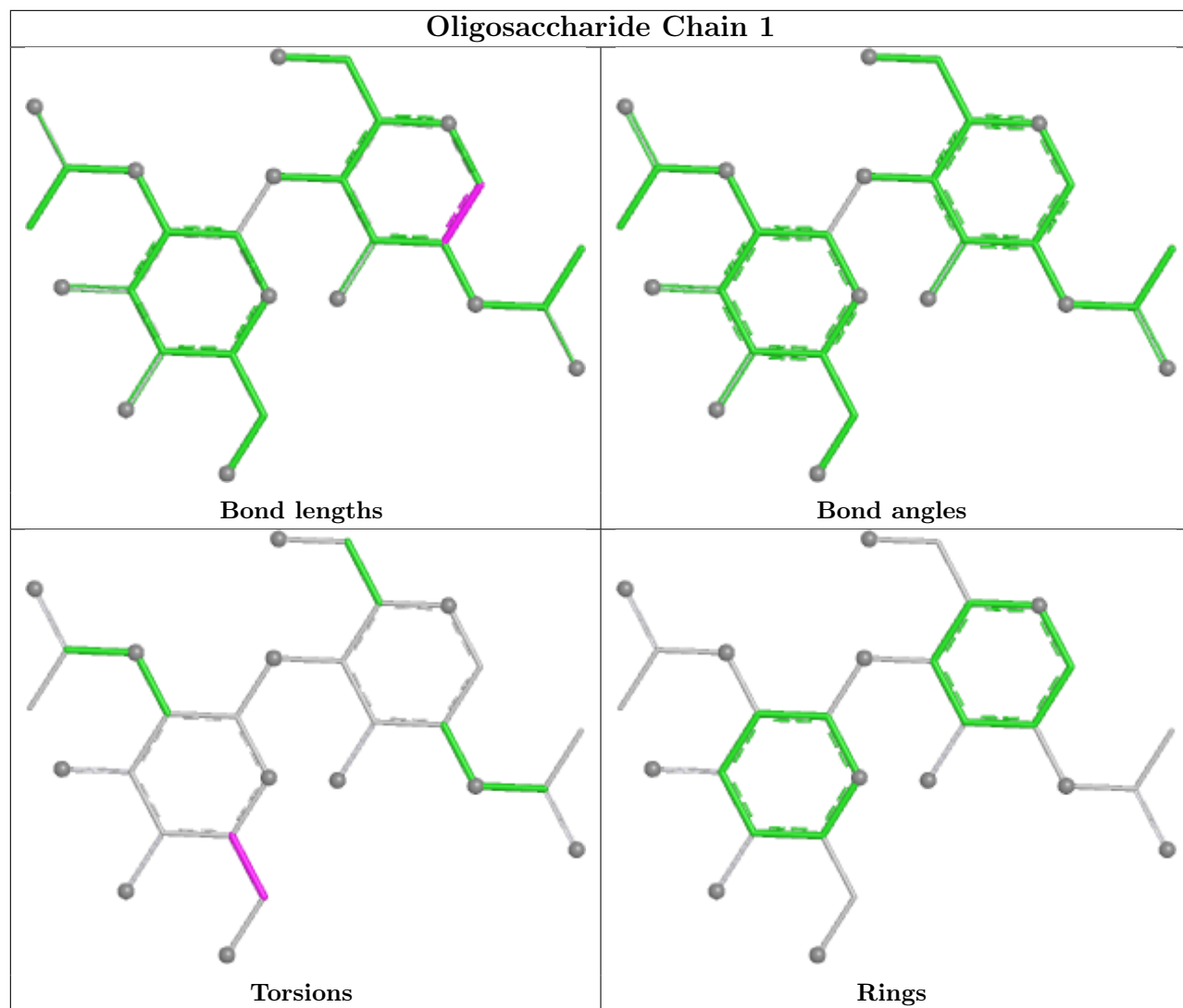


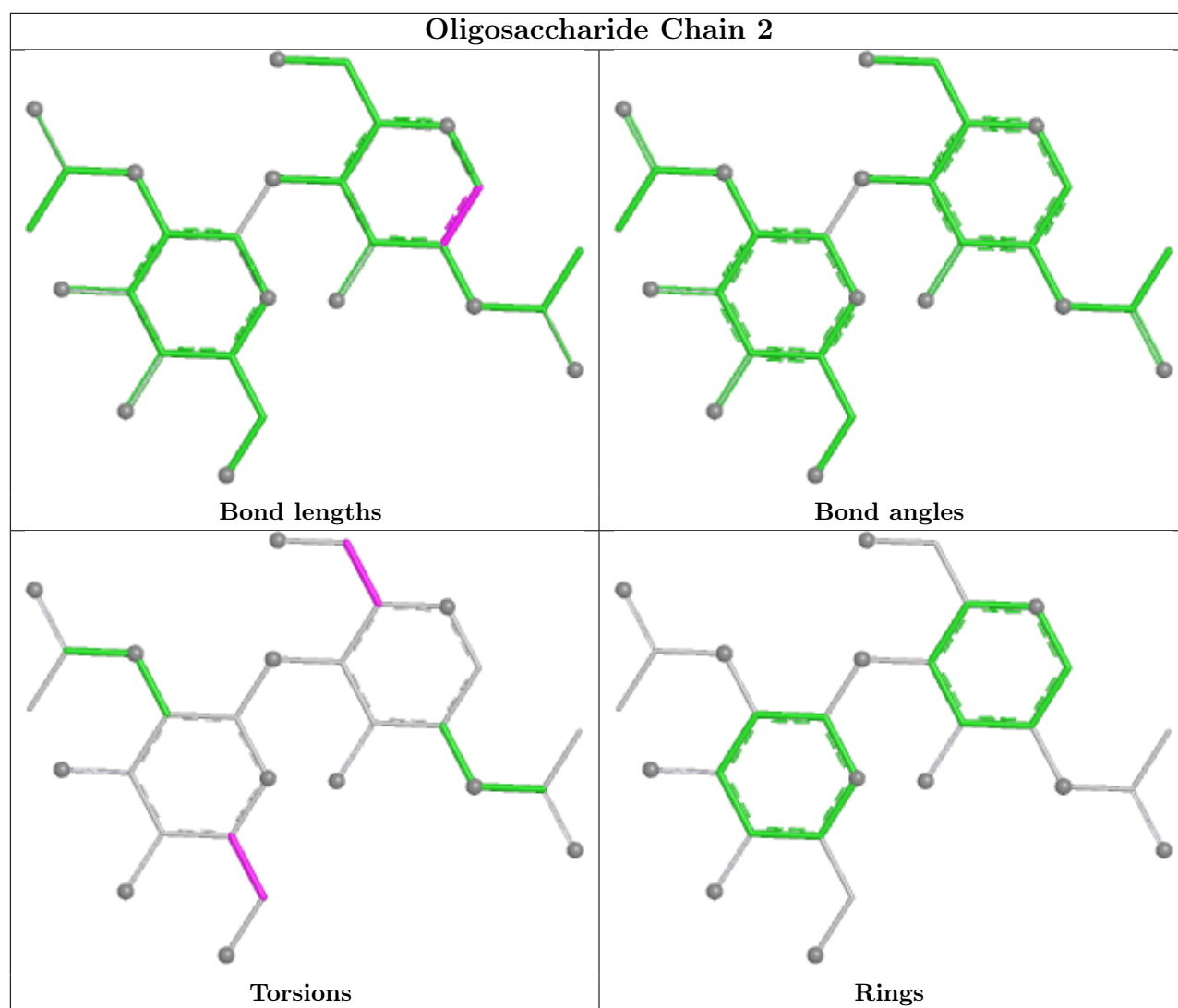


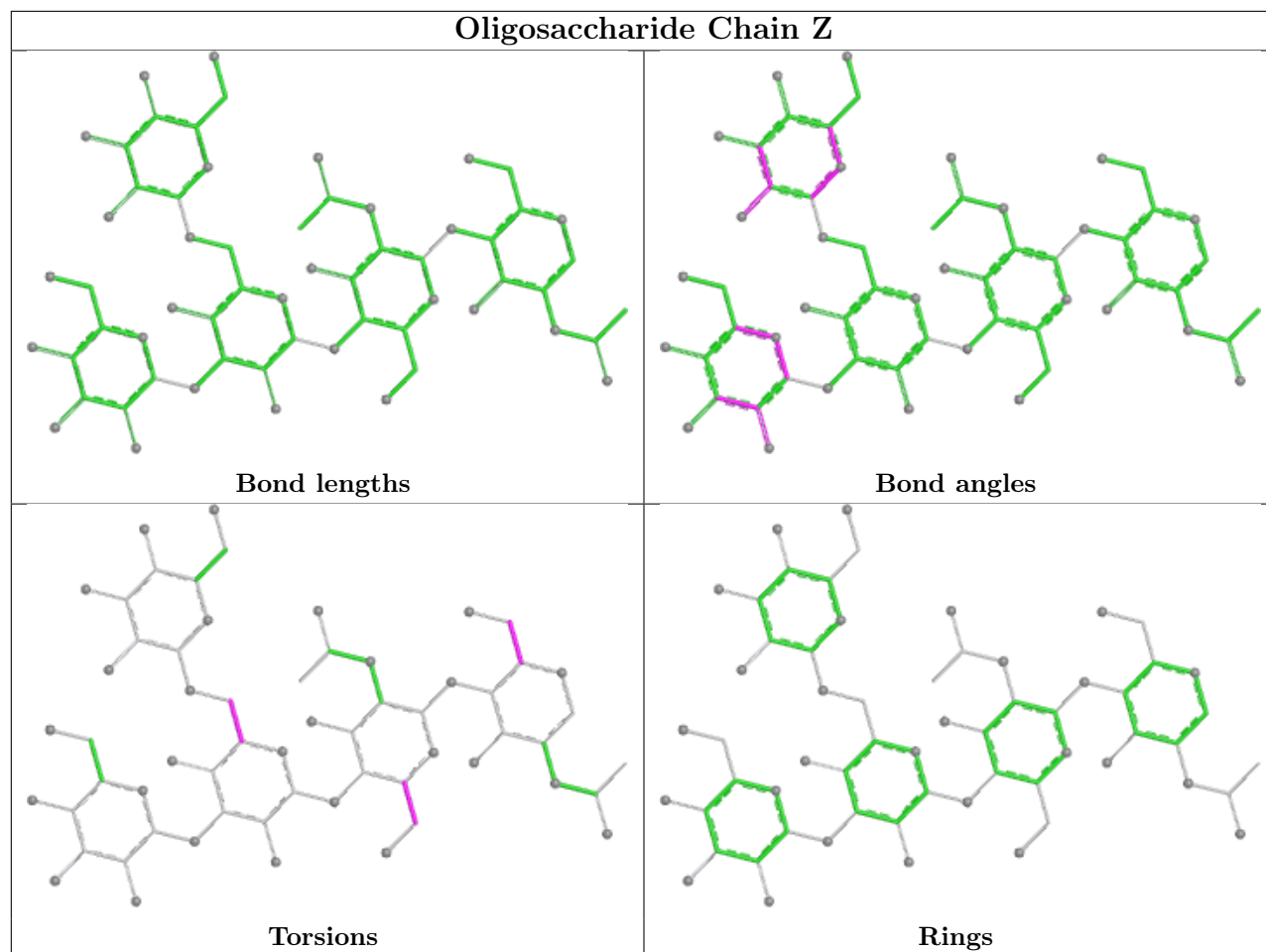


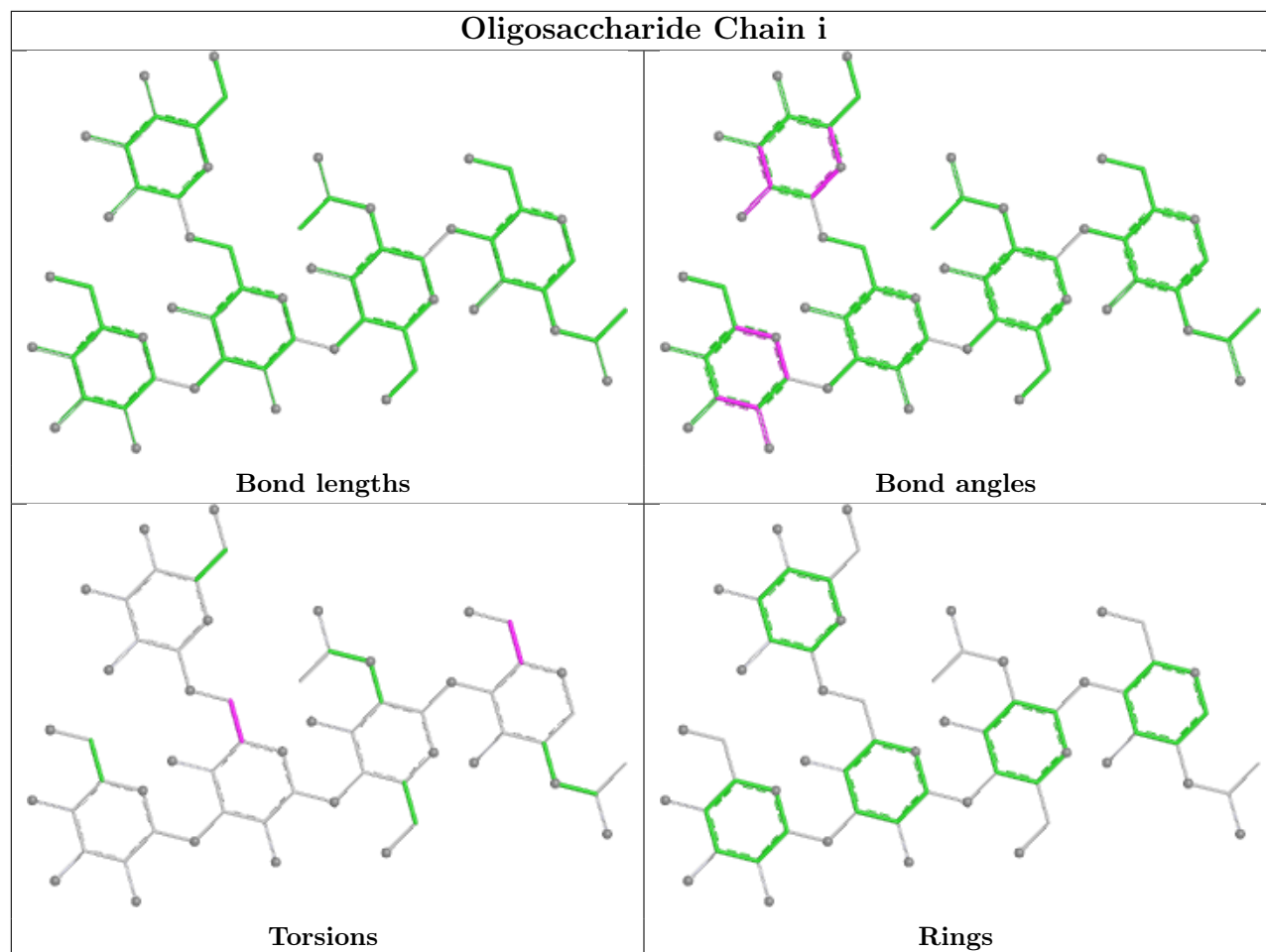


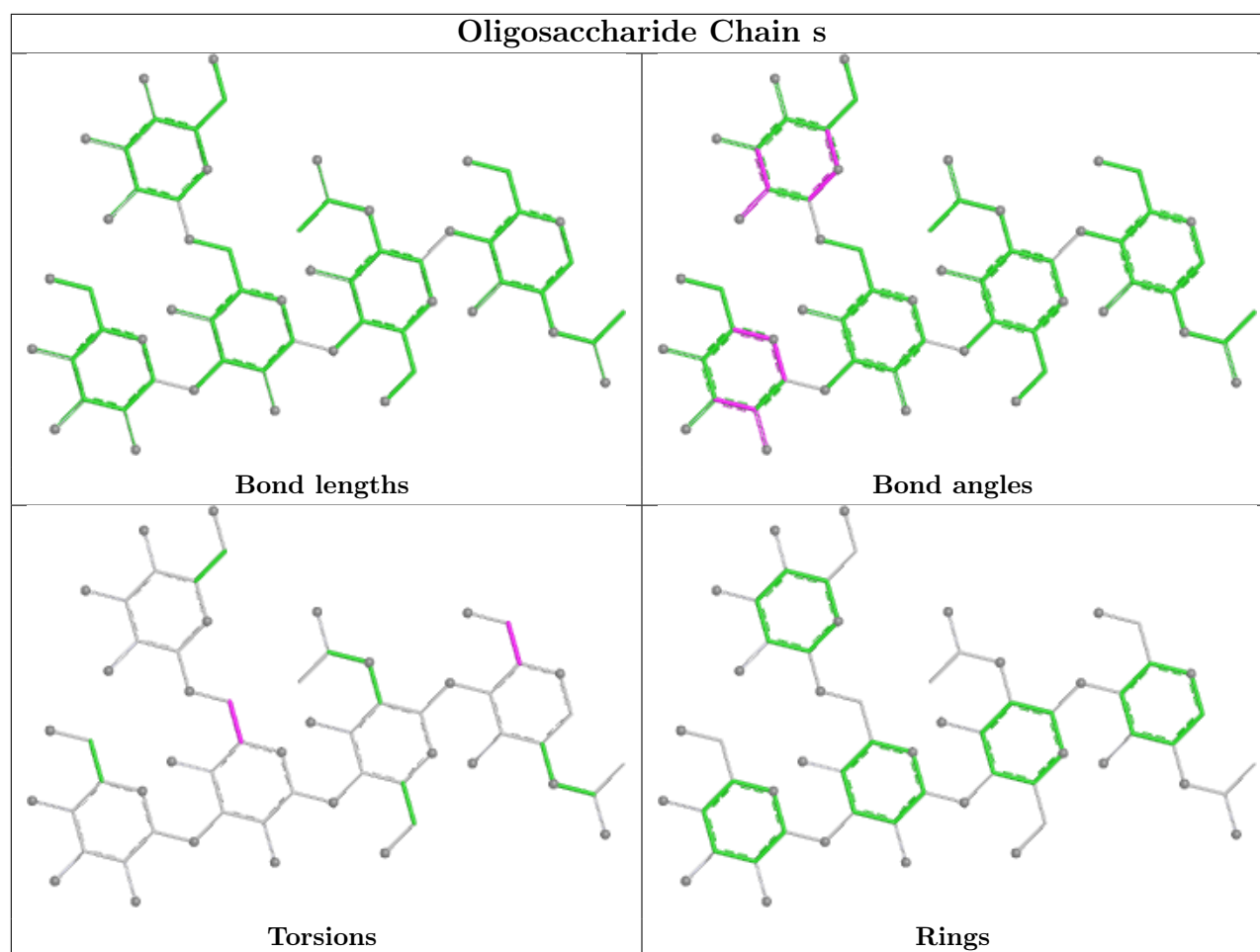












5.6 Ligand geometry [i](#)

36 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
12	NAG	I	601	1	14,14,15	0.25	0	17,19,21	0.43	0
12	NAG	B	702	2	14,14,15	0.20	0	17,19,21	0.37	0
12	NAG	E	605	1	14,14,15	0.25	0	17,19,21	0.45	0
12	NAG	A	602	1	14,14,15	0.24	0	17,19,21	0.42	0
12	NAG	I	604	1	14,14,15	0.19	0	17,19,21	0.40	0
12	NAG	A	608	1	14,14,15	0.40	0	17,19,21	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
12	NAG	F	700	2	14,14,15	0.46	0	17,19,21	0.58	0
12	NAG	F	702	2	14,14,15	0.20	0	17,19,21	0.37	0
12	NAG	A	601	1	14,14,15	0.25	0	17,19,21	0.42	0
12	NAG	I	608	1	14,14,15	0.40	0	17,19,21	0.58	0
12	NAG	B	701	2	14,14,15	0.22	0	17,19,21	0.43	0
12	NAG	J	700	2	14,14,15	0.47	0	17,19,21	0.57	0
12	NAG	J	701	2	14,14,15	0.21	0	17,19,21	0.44	0
12	NAG	B	700	2	14,14,15	0.47	0	17,19,21	0.58	0
12	NAG	I	602	1	14,14,15	0.24	0	17,19,21	0.42	0
12	NAG	A	607	1	14,14,15	0.25	0	17,19,21	0.45	0
12	NAG	A	606	1	14,14,15	0.32	0	17,19,21	0.69	1 (5%)
12	NAG	I	607	1	14,14,15	0.24	0	17,19,21	0.46	0
12	NAG	I	603	1	14,14,15	0.88	1 (7%)	17,19,21	1.08	1 (5%)
12	NAG	F	701	2	14,14,15	0.23	0	17,19,21	0.44	0
12	NAG	E	604	1	14,14,15	0.20	0	17,19,21	0.40	0
12	NAG	E	607	1	14,14,15	0.25	0	17,19,21	0.45	0
12	NAG	E	609	1	14,14,15	0.22	0	17,19,21	0.43	0
12	NAG	E	603	1	14,14,15	0.88	1 (7%)	17,19,21	1.10	1 (5%)
12	NAG	I	605	1	14,14,15	0.24	0	17,19,21	0.46	0
12	NAG	E	601	1	14,14,15	0.25	0	17,19,21	0.43	0
12	NAG	A	609	1	14,14,15	0.21	0	17,19,21	0.43	0
12	NAG	A	605	1	14,14,15	0.24	0	17,19,21	0.45	0
12	NAG	A	603	1	14,14,15	0.86	1 (7%)	17,19,21	1.09	1 (5%)
12	NAG	I	606	1	14,14,15	0.30	0	17,19,21	0.69	1 (5%)
12	NAG	J	702	2	14,14,15	0.20	0	17,19,21	0.37	0
12	NAG	A	604	1	14,14,15	0.20	0	17,19,21	0.40	0
12	NAG	E	606	1	14,14,15	0.30	0	17,19,21	0.69	1 (5%)
12	NAG	I	609	1	14,14,15	0.22	0	17,19,21	0.42	0
12	NAG	E	608	1	14,14,15	0.39	0	17,19,21	0.58	0
12	NAG	E	602	1	14,14,15	0.24	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
12	NAG	I	601	1	-	0/6/23/26	0/1/1/1
12	NAG	B	702	2	-	0/6/23/26	0/1/1/1
12	NAG	E	605	1	-	2/6/23/26	0/1/1/1
12	NAG	A	602	1	-	2/6/23/26	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	I	604	1	-	2/6/23/26	0/1/1/1
12	NAG	A	608	1	-	4/6/23/26	0/1/1/1
12	NAG	F	700	2	-	3/6/23/26	0/1/1/1
12	NAG	F	702	2	-	0/6/23/26	0/1/1/1
12	NAG	A	601	1	-	0/6/23/26	0/1/1/1
12	NAG	I	608	1	-	4/6/23/26	0/1/1/1
12	NAG	B	701	2	-	2/6/23/26	0/1/1/1
12	NAG	J	700	2	-	3/6/23/26	0/1/1/1
12	NAG	J	701	2	-	2/6/23/26	0/1/1/1
12	NAG	B	700	2	-	3/6/23/26	0/1/1/1
12	NAG	I	602	1	-	2/6/23/26	0/1/1/1
12	NAG	A	607	1	-	2/6/23/26	0/1/1/1
12	NAG	A	606	1	-	0/6/23/26	0/1/1/1
12	NAG	I	607	1	-	2/6/23/26	0/1/1/1
12	NAG	I	603	1	-	4/6/23/26	0/1/1/1
12	NAG	F	701	2	-	2/6/23/26	0/1/1/1
12	NAG	E	604	1	-	2/6/23/26	0/1/1/1
12	NAG	E	607	1	-	2/6/23/26	0/1/1/1
12	NAG	E	609	1	-	2/6/23/26	0/1/1/1
12	NAG	E	603	1	-	4/6/23/26	0/1/1/1
12	NAG	I	605	1	-	2/6/23/26	0/1/1/1
12	NAG	E	601	1	-	0/6/23/26	0/1/1/1
12	NAG	A	609	1	-	2/6/23/26	0/1/1/1
12	NAG	A	605	1	-	2/6/23/26	0/1/1/1
12	NAG	A	603	1	-	4/6/23/26	0/1/1/1
12	NAG	I	606	1	-	0/6/23/26	0/1/1/1
12	NAG	J	702	2	-	0/6/23/26	0/1/1/1
12	NAG	A	604	1	-	2/6/23/26	0/1/1/1
12	NAG	E	606	1	-	0/6/23/26	0/1/1/1
12	NAG	I	609	1	-	2/6/23/26	0/1/1/1
12	NAG	E	608	1	-	4/6/23/26	0/1/1/1
12	NAG	E	602	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)
12	E	603	NAG	O5-C1	2.86	1.48	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	I	603	NAG	O5-C1	2.84	1.48	1.43
12	A	603	NAG	O5-C1	2.80	1.48	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	E	603	NAG	C1-O5-C5	4.07	117.64	112.19
12	A	603	NAG	C1-O5-C5	4.06	117.63	112.19
12	I	603	NAG	C1-O5-C5	4.03	117.59	112.19
12	I	606	NAG	C1-O5-C5	2.44	115.46	112.19
12	A	606	NAG	C1-O5-C5	2.43	115.44	112.19
12	E	606	NAG	C1-O5-C5	2.43	115.44	112.19

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	A	604	NAG	O5-C5-C6-O6
12	E	604	NAG	O5-C5-C6-O6
12	I	604	NAG	O5-C5-C6-O6
12	A	609	NAG	O5-C5-C6-O6
12	E	609	NAG	O5-C5-C6-O6
12	I	609	NAG	O5-C5-C6-O6
12	B	700	NAG	C4-C5-C6-O6
12	F	700	NAG	C4-C5-C6-O6
12	J	700	NAG	C4-C5-C6-O6
12	B	701	NAG	O5-C5-C6-O6
12	F	701	NAG	O5-C5-C6-O6
12	J	701	NAG	O5-C5-C6-O6
12	A	607	NAG	O5-C5-C6-O6
12	E	607	NAG	O5-C5-C6-O6
12	I	607	NAG	O5-C5-C6-O6
12	A	609	NAG	C4-C5-C6-O6
12	B	701	NAG	C4-C5-C6-O6
12	E	609	NAG	C4-C5-C6-O6
12	F	701	NAG	C4-C5-C6-O6
12	I	609	NAG	C4-C5-C6-O6
12	J	701	NAG	C4-C5-C6-O6
12	A	604	NAG	C4-C5-C6-O6
12	A	607	NAG	C4-C5-C6-O6
12	E	604	NAG	C4-C5-C6-O6
12	E	607	NAG	C4-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
12	I	604	NAG	C4-C5-C6-O6
12	I	607	NAG	C4-C5-C6-O6
12	B	700	NAG	O5-C5-C6-O6
12	F	700	NAG	O5-C5-C6-O6
12	J	700	NAG	O5-C5-C6-O6
12	A	605	NAG	C4-C5-C6-O6
12	E	605	NAG	C4-C5-C6-O6
12	I	605	NAG	C4-C5-C6-O6
12	A	602	NAG	O5-C5-C6-O6
12	E	602	NAG	O5-C5-C6-O6
12	I	602	NAG	O5-C5-C6-O6
12	A	602	NAG	C4-C5-C6-O6
12	E	602	NAG	C4-C5-C6-O6
12	I	602	NAG	C4-C5-C6-O6
12	A	608	NAG	O5-C5-C6-O6
12	E	608	NAG	O5-C5-C6-O6
12	I	608	NAG	O5-C5-C6-O6
12	A	605	NAG	O5-C5-C6-O6
12	E	605	NAG	O5-C5-C6-O6
12	I	605	NAG	O5-C5-C6-O6
12	A	608	NAG	C4-C5-C6-O6
12	E	608	NAG	C4-C5-C6-O6
12	I	608	NAG	C4-C5-C6-O6
12	A	603	NAG	C1-C2-N2-C7
12	E	603	NAG	C1-C2-N2-C7
12	I	603	NAG	C1-C2-N2-C7
12	A	603	NAG	C4-C5-C6-O6
12	E	603	NAG	C4-C5-C6-O6
12	I	603	NAG	C4-C5-C6-O6
12	A	603	NAG	C3-C2-N2-C7
12	A	608	NAG	C3-C2-N2-C7
12	E	603	NAG	C3-C2-N2-C7
12	E	608	NAG	C3-C2-N2-C7
12	I	603	NAG	C3-C2-N2-C7
12	I	608	NAG	C3-C2-N2-C7
12	E	603	NAG	O5-C5-C6-O6
12	I	603	NAG	O5-C5-C6-O6
12	A	603	NAG	O5-C5-C6-O6
12	A	608	NAG	C1-C2-N2-C7
12	E	608	NAG	C1-C2-N2-C7
12	I	608	NAG	C1-C2-N2-C7
12	B	700	NAG	C3-C2-N2-C7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
12	F	700	NAG	C3-C2-N2-C7
12	J	700	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	F	700	NAG	1	0
12	J	700	NAG	1	0
12	B	700	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

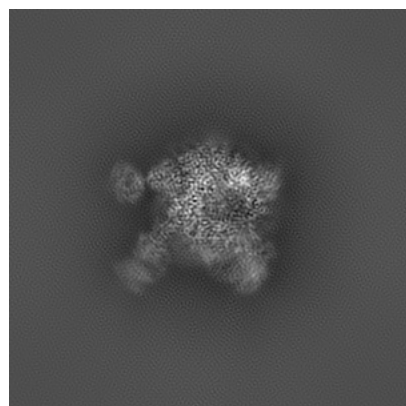
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29783. These allow visual inspection of the internal detail of the map and identification of artifacts.

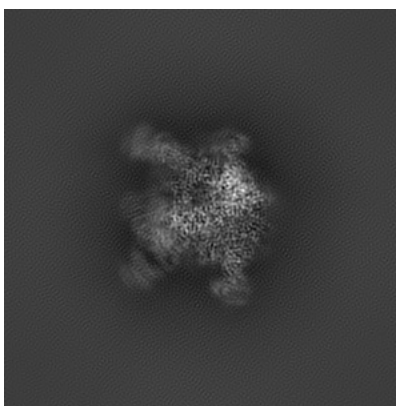
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

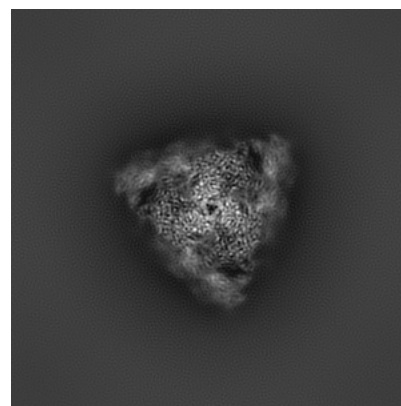
6.1.1 Primary map



X

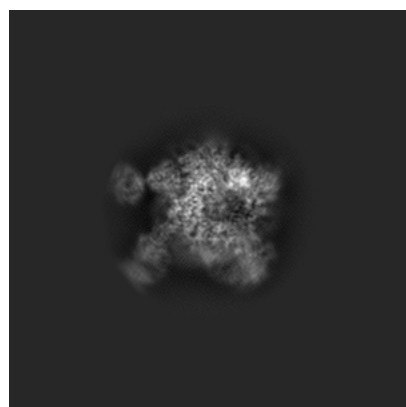


Y

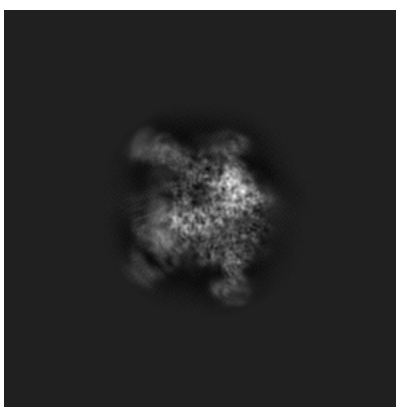


Z

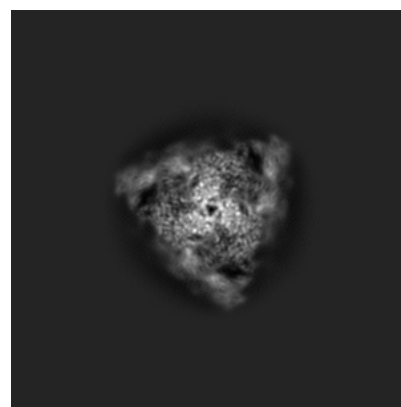
6.1.2 Raw map



X



Y

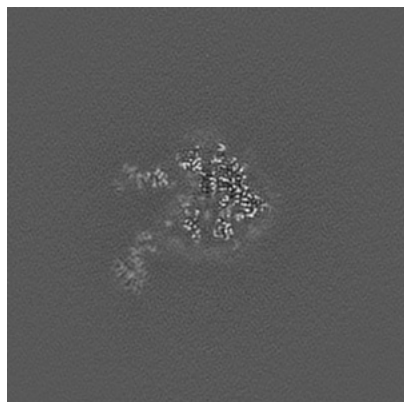


Z

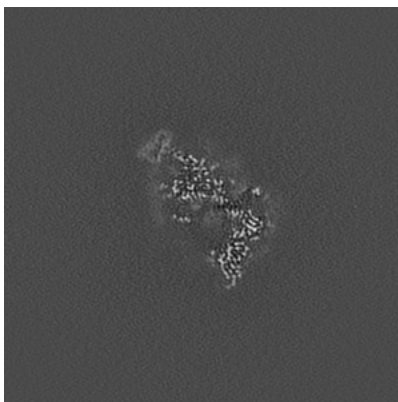
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

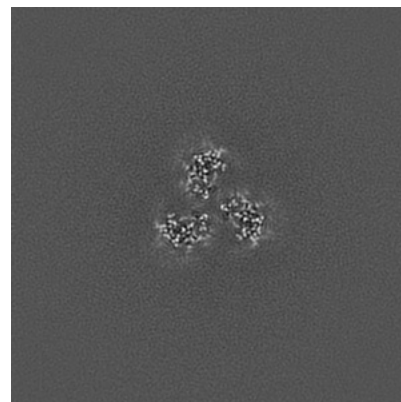
6.2.1 Primary map



X Index: 243

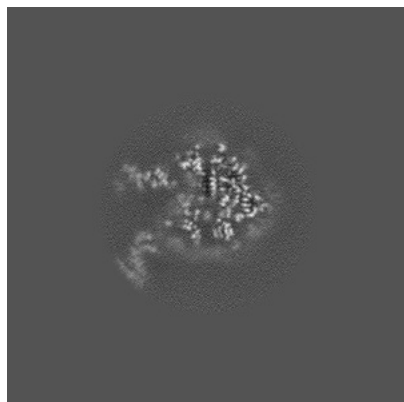


Y Index: 243

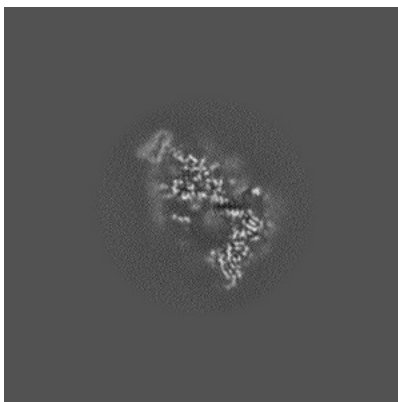


Z Index: 243

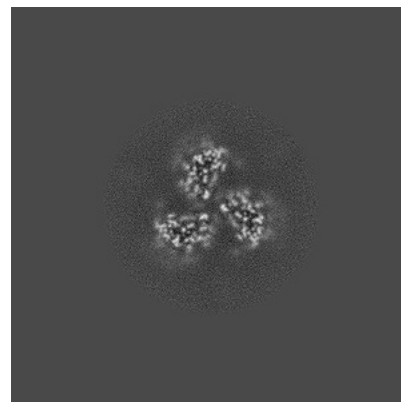
6.2.2 Raw map



X Index: 243



Y Index: 243

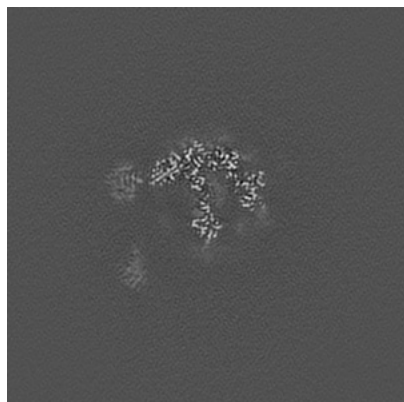


Z Index: 243

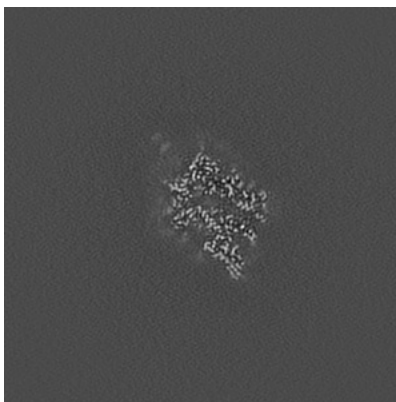
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

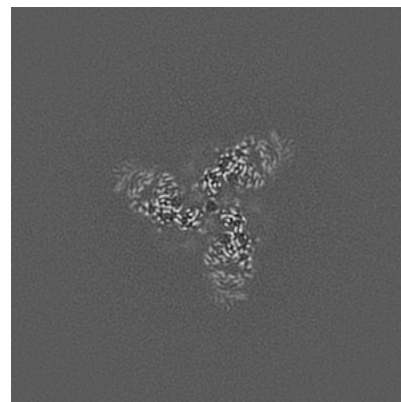
6.3.1 Primary map



X Index: 257

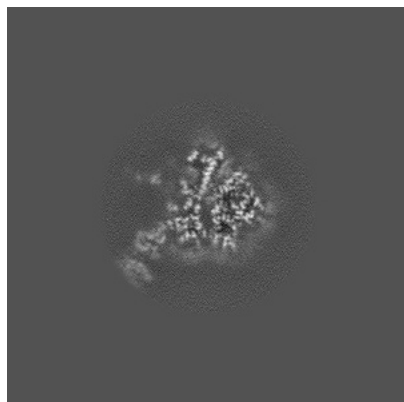


Y Index: 232

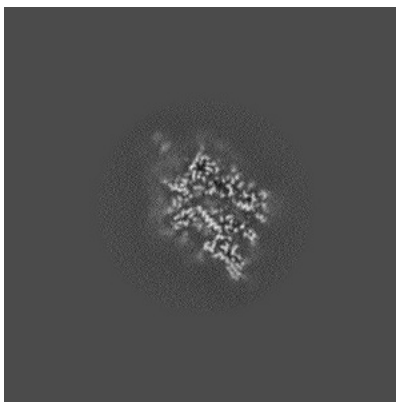


Z Index: 274

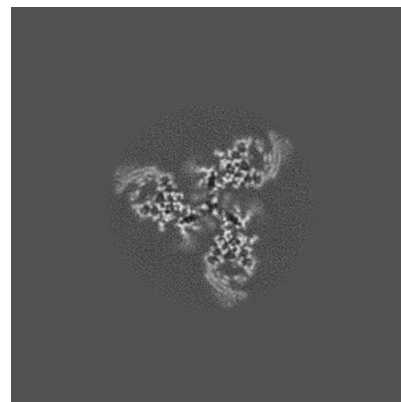
6.3.2 Raw map



X Index: 235



Y Index: 232

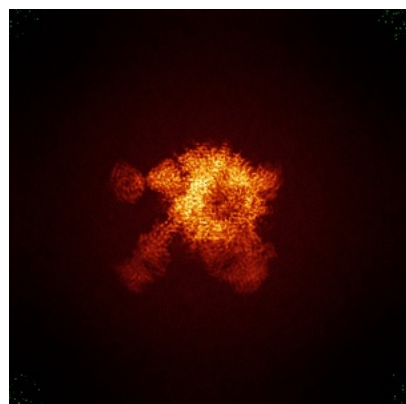


Z Index: 279

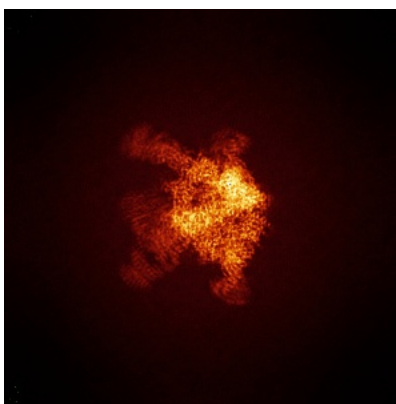
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

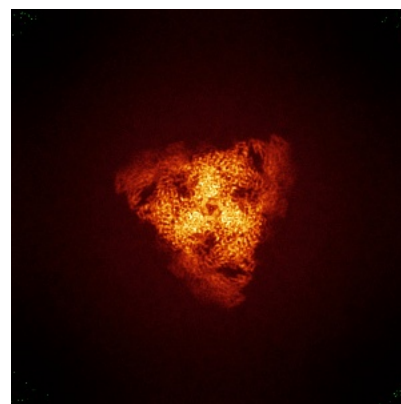
6.4.1 Primary map



X

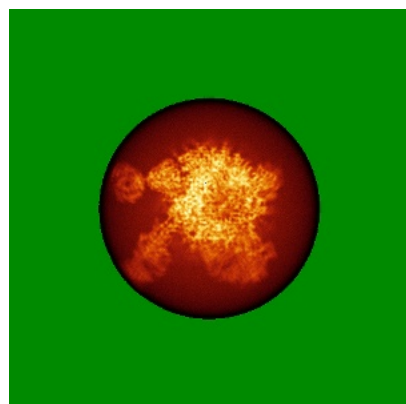


Y

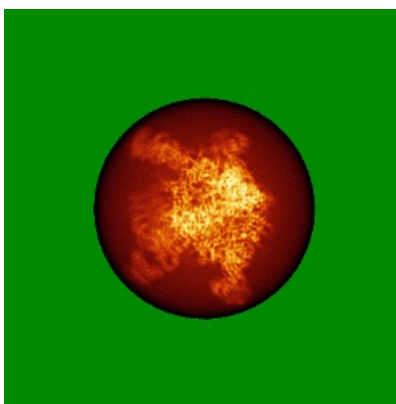


Z

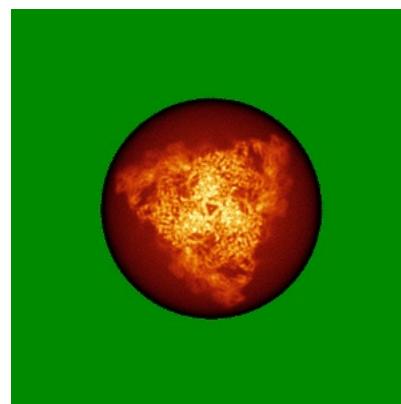
6.4.2 Raw map



X



Y

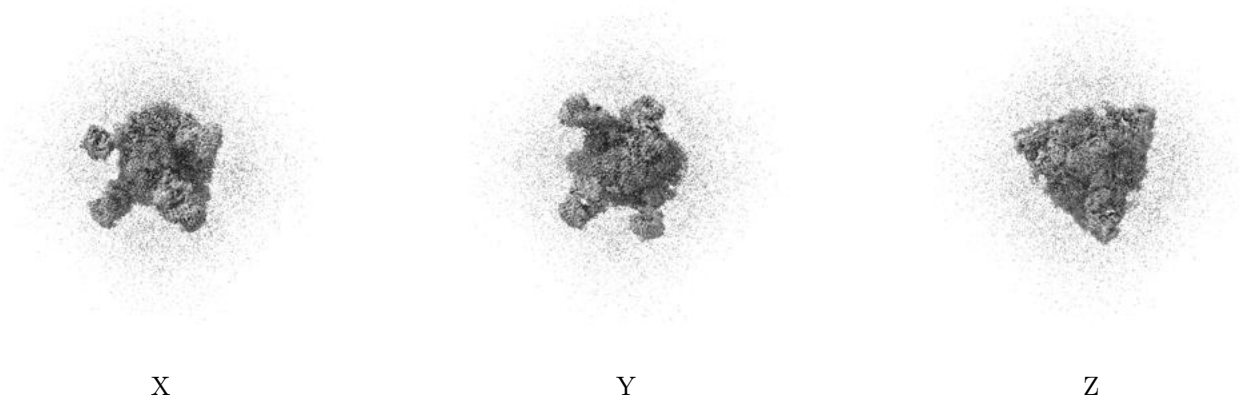


Z

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.

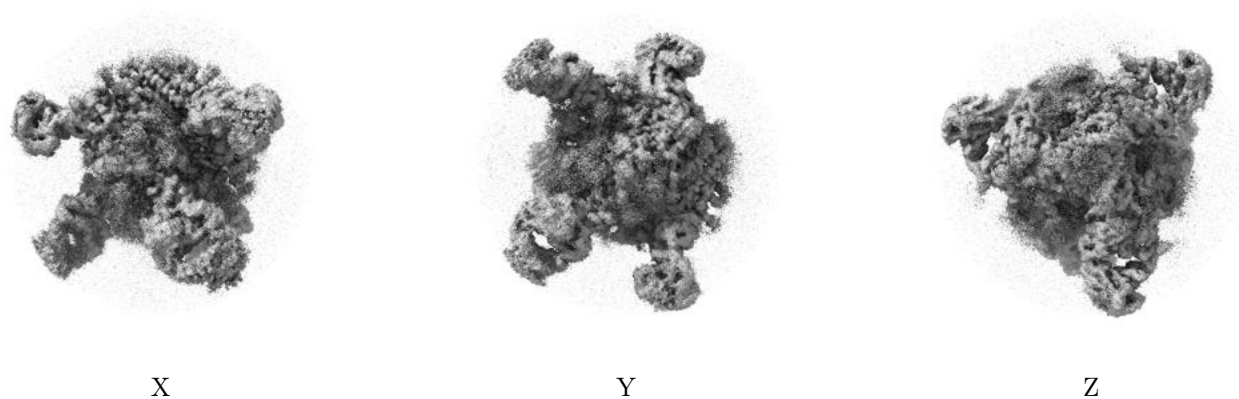
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 1.4. 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.

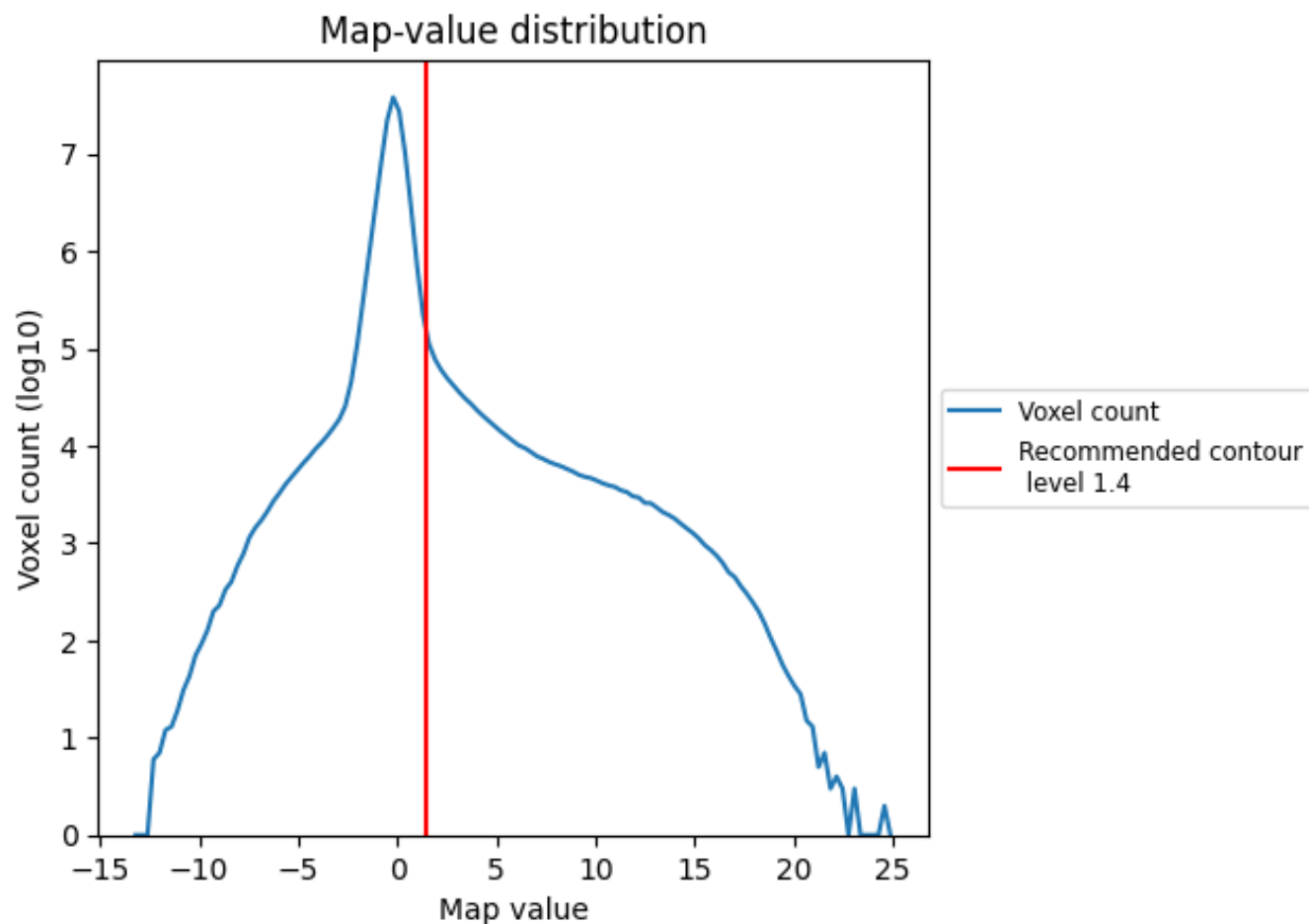
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

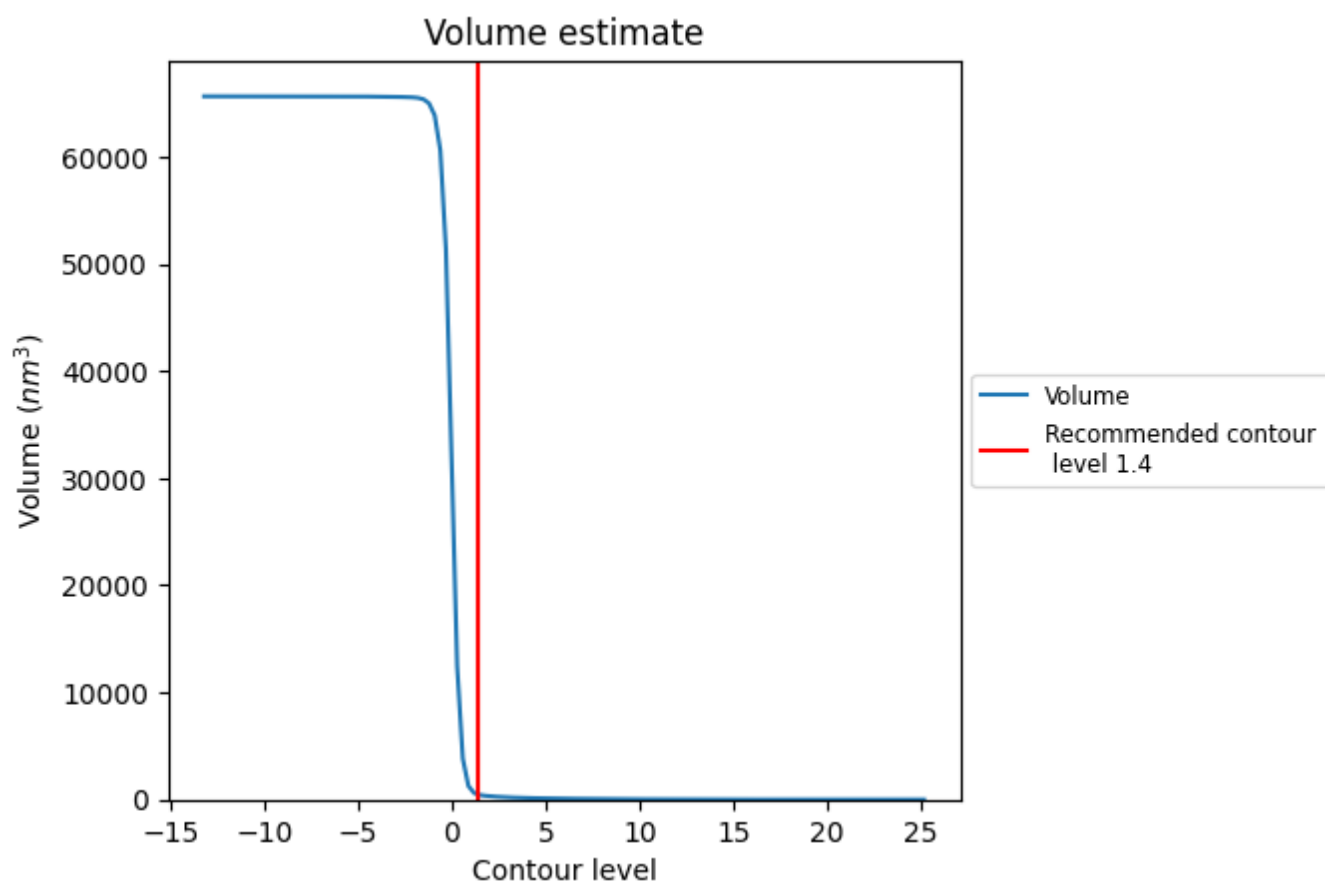
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

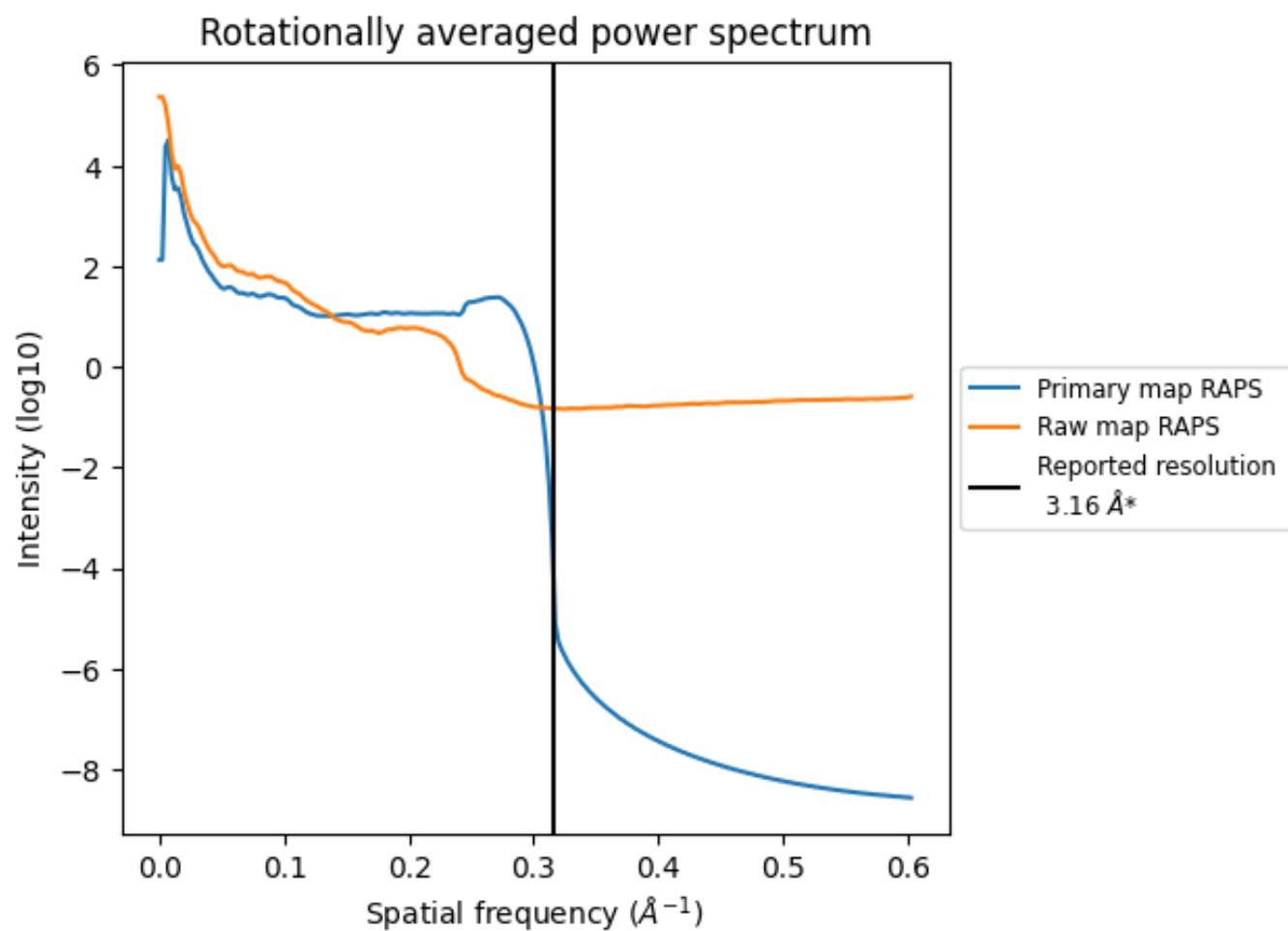
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 487 nm³; this corresponds to an approximate mass of 440 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 ⓘ

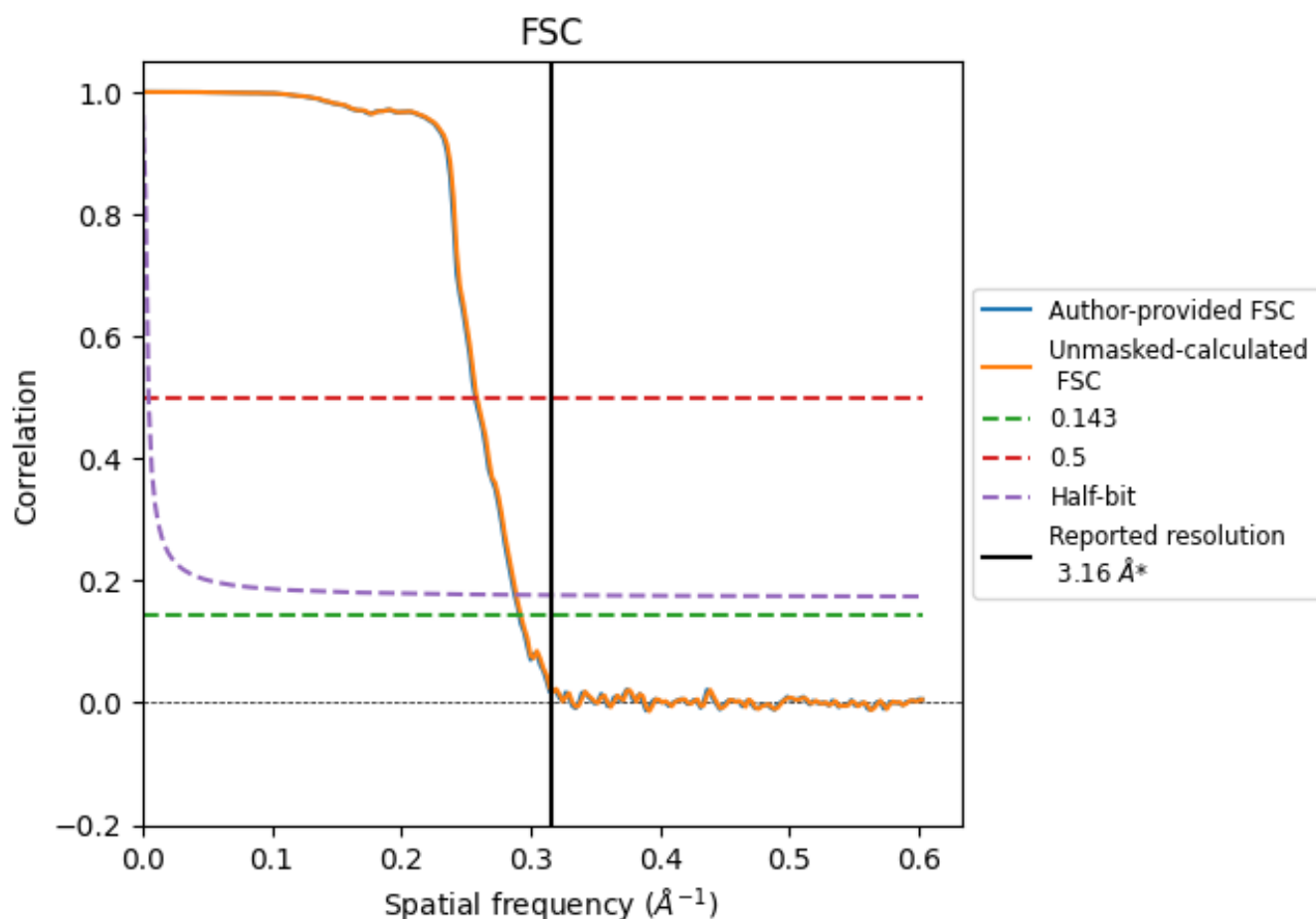


*Reported resolution corresponds to spatial frequency of 0.316 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.316 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.16	-	-
Author-provided FSC curve	3.43	3.89	3.48
Unmasked-calculated*	3.42	3.87	3.46

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

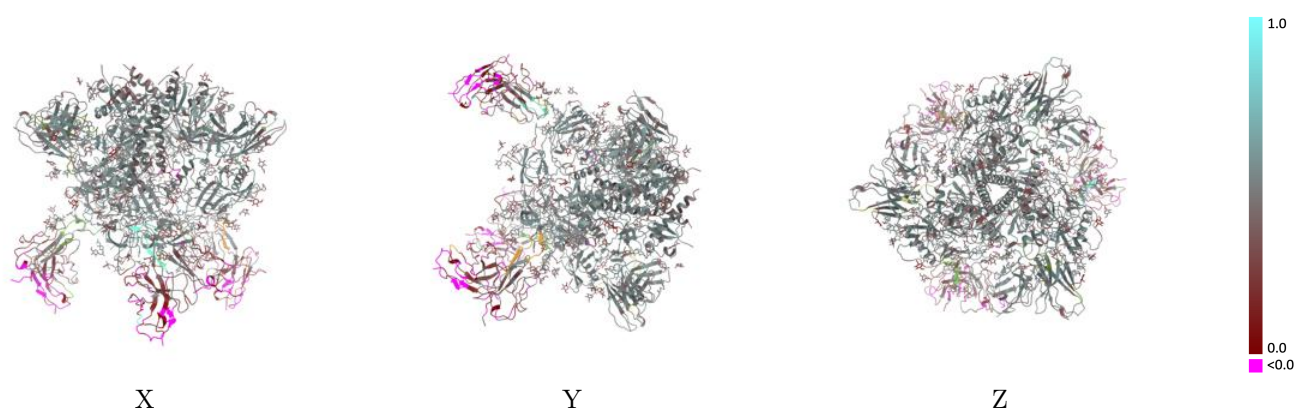
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29783 and PDB model 8G6U. Per-residue inclusion information can be found in section 3 on page 14.

9.1 Map-model overlay [i](#)

This section was not generated.

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

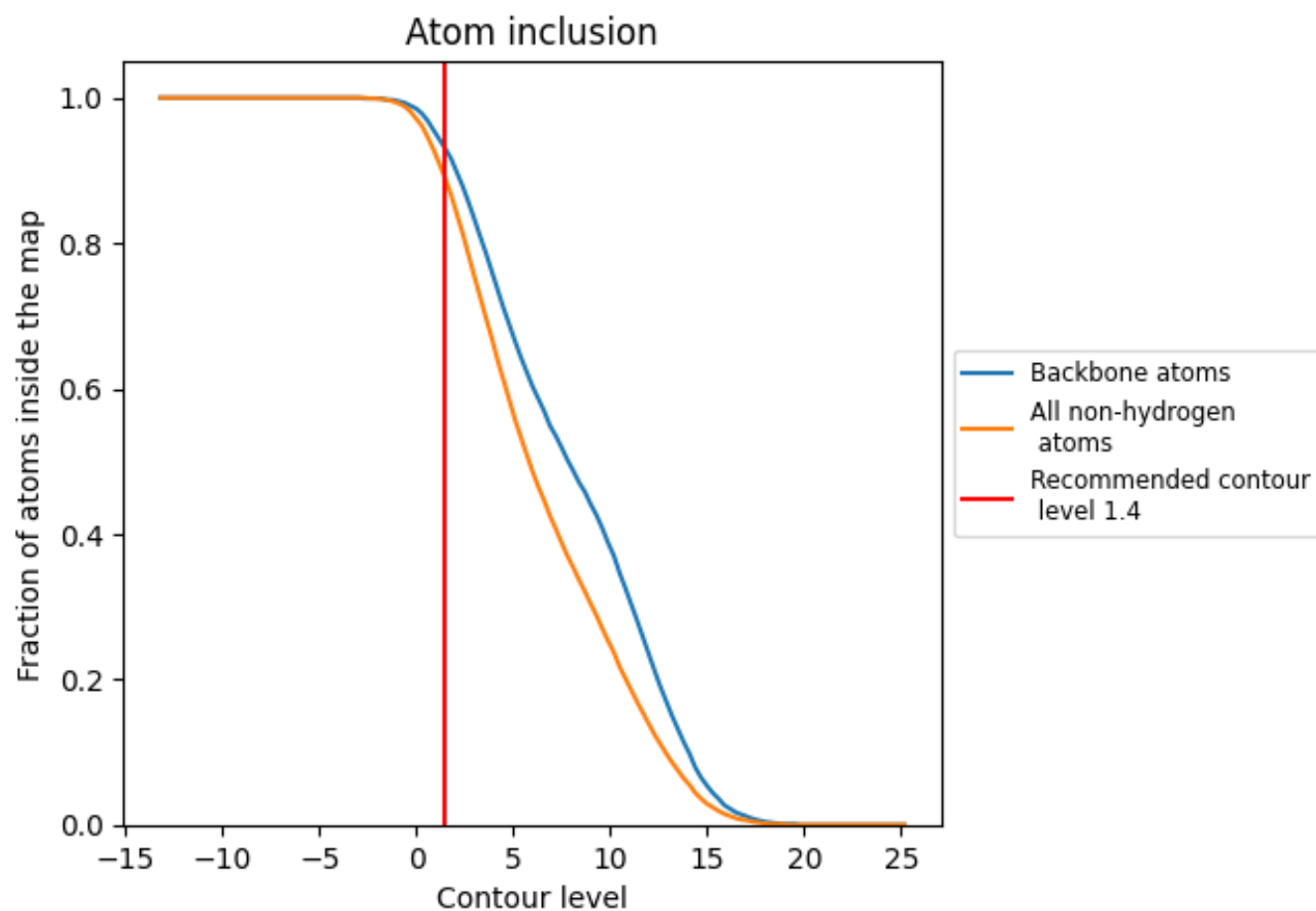


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% 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 (1.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8950	 0.4330
0	 0.8210	 0.3450
1	 0.7860	 0.3280
2	 0.8210	 0.3170
A	 0.9370	 0.4990
B	 0.9220	 0.4660
C	 0.9400	 0.5010
D	 0.9400	 0.4890
E	 0.9370	 0.4970
F	 0.9230	 0.4710
G	 0.9500	 0.5070
H	 0.9470	 0.5010
I	 0.9280	 0.4910
J	 0.9170	 0.4680
K	 0.9480	 0.5030
L	 0.9430	 0.4920
M	 0.7470	 0.2110
N	 0.7990	 0.2960
O	 0.7360	 0.1970
P	 0.8070	 0.2890
Q	 0.7110	 0.1720
R	 0.8000	 0.2740
S	 0.8980	 0.4230
T	 0.8890	 0.4170
U	 0.8620	 0.4040
V	 0.8570	 0.4000
X	 0.8210	 0.3400
Y	 0.9170	 0.3930
Z	 0.8850	 0.4180
a	 0.8900	 0.4300
b	 0.9030	 0.4190
c	 0.8930	 0.3990
d	 0.8530	 0.3960
e	 0.8570	 0.3740
f	 0.8930	 0.4150



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
g	 0.8210	 0.3320
h	 0.9170	 0.3970
i	 0.8850	 0.4220
j	 0.9050	 0.4580
k	 0.8750	 0.4180
l	 0.8930	 0.4250
m	 0.6790	 0.2440
n	 0.8360	 0.4050
o	 0.8570	 0.4210
p	 0.8210	 0.3320
q	 0.9640	 0.4730
r	 0.8890	 0.4020
s	 0.9020	 0.4210
t	 0.9640	 0.4660
u	 1.0000	 0.4690
v	 0.6790	 0.2460
w	 0.6790	 0.2400
x	 0.8930	 0.4570
y	 0.8930	 0.4600
z	 0.8930	 0.4380