



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 08:01 AM UTC

PDB ID : 2VLC / pdb_00002vlc
Title : Crystal structure of Natural Cinnamomin (Isoform III)
Authors : Azzi, A.; Wang, T.; Zhu, D.-W.; Zou, Y.-S.; Liu, W.-Y.; Lin, S.-X.
Deposited on : 2008-01-11
Resolution : 2.95 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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

A user guide is available at

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

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:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

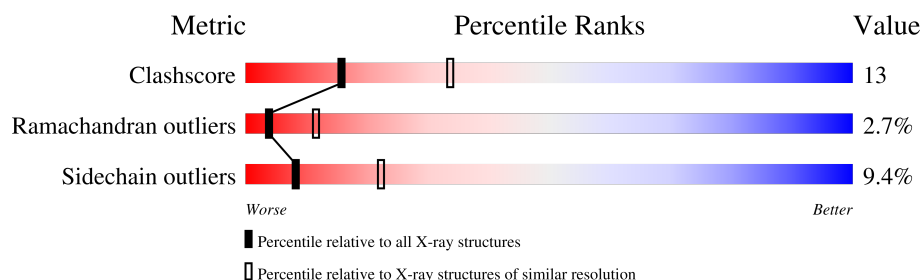
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	570	64% 22% . 9%
1	B	570	59% 25% 5% . 9%
2	C	2	50% 50%
2	D	2	50% 50%
2	E	2	50% 50%
2	F	2	100%
2	G	2	100%
2	H	2	100%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	C	1	X	-	-	-
2	NAG	C	2	X	-	-	-
2	NAG	D	1	X	-	-	-
2	NAG	D	2	X	-	-	-
2	NAG	E	1	X	-	-	-
2	NAG	E	2	X	-	-	-
2	NAG	F	1	X	-	-	-
2	NAG	F	2	X	-	-	-
2	NAG	G	1	X	-	-	-
2	NAG	G	2	X	-	-	-
2	NAG	H	1	X	-	-	-
2	NAG	H	2	X	-	-	-
4	XYP	B	601	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8406 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Ribosome-inactivating protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	518	Total	C	N	O	S	4	0	0
			4036	2529	713	773	21			
1	B	518	Total	C	N	O	S	5	0	0
			4038	2529	713	775	21			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	183	SER	THR	conflict	UNP Q94BW3
A	307	ARG	ASP	conflict	UNP Q94BW3
B	183	SER	THR	conflict	UNP Q94BW3
B	307	ARG	ASP	conflict	UNP Q94BW3

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



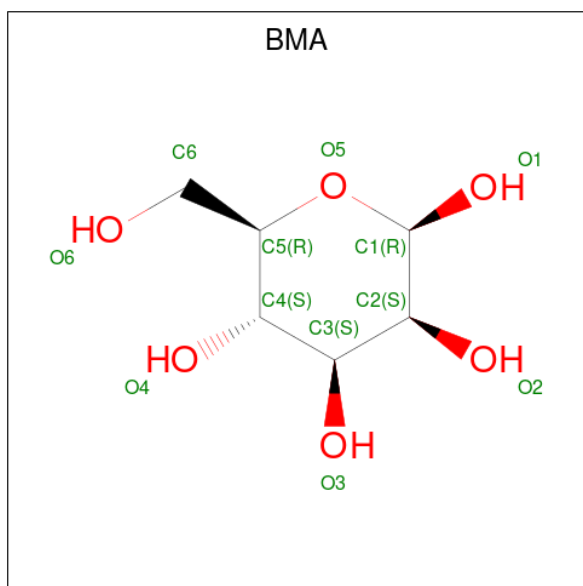
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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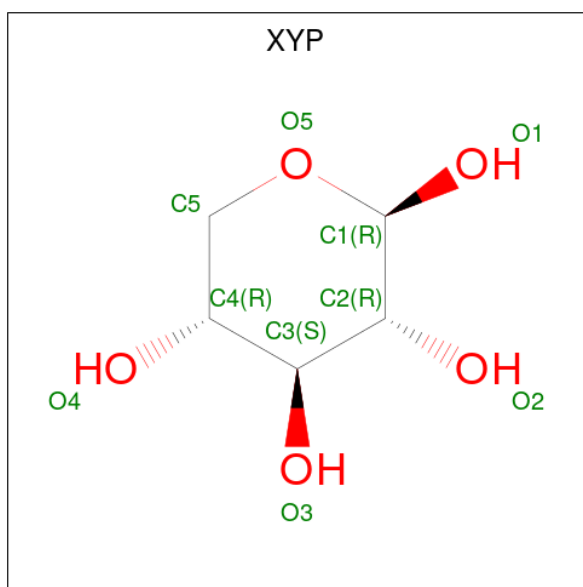
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is beta-D-mannopyranose (CCD ID: BMA) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 4 is beta-D-xylopyranose (CCD ID: XYP) (formula: $C_5H_{10}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	5	5		
4	B	1	Total	C	O	0	0
			10	5	5		

- Molecule 5 is water.

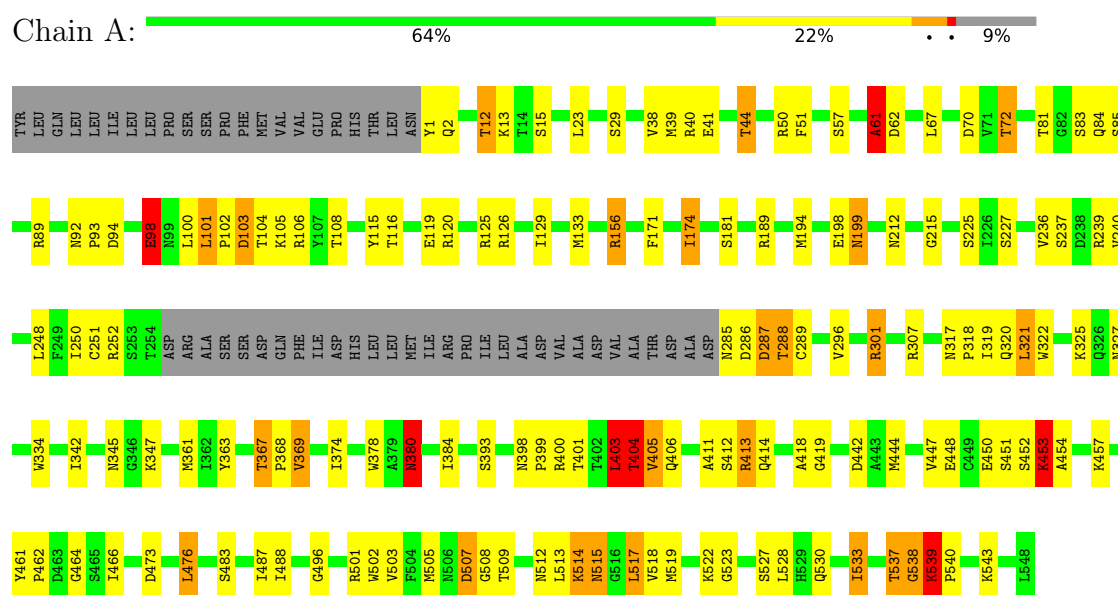
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	70	Total	O	0	0
			70	70		
5	B	50	Total	O	0	0
			50	50		

3 Residue-property plots

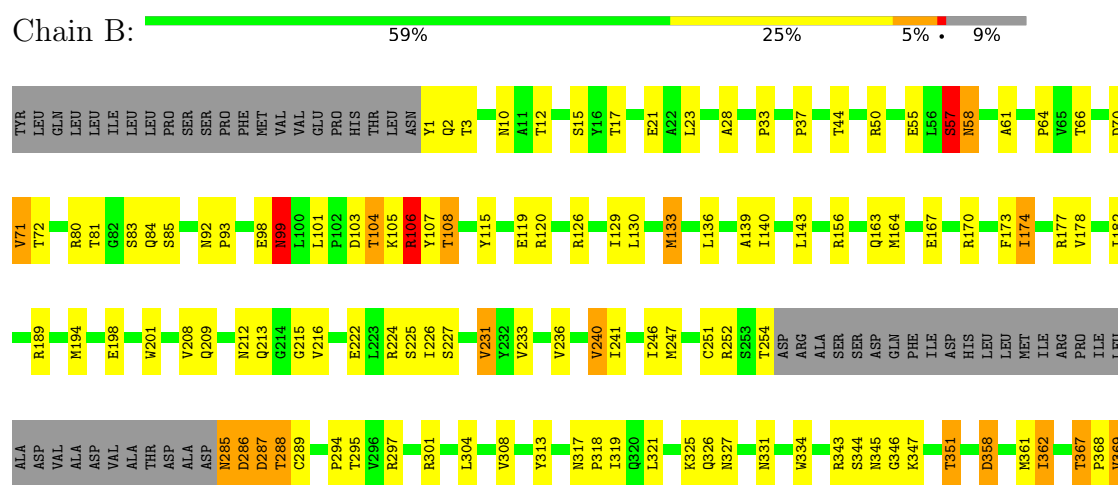
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

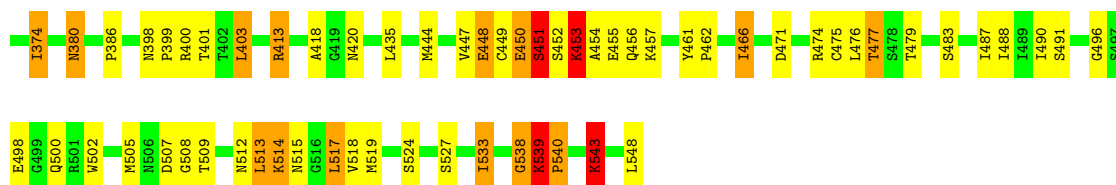
Note EDS was not executed.

• Molecule 1: Ribosome-inactivating protein



• Molecule 1: Ribosome-inactivating protein





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

Chain C: 50% 50%



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

Chain D: 50% 50%



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

Chain E: 50% 50%



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

Chain F: 100%



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

Chain G: 100%



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

Chain H: 100%

MAG1
MAG2

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.39Å 126.33Å 161.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.00 – 2.95	Depositor
% Data completeness (in resolution range)	98.9 (17.00-2.95)	Depositor
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.238 , 0.309	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8406	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.21	8/4122 (0.2%)	1.22	30/5610 (0.5%)
1	B	1.13	9/4124 (0.2%)	1.37	26/5612 (0.5%)
All	All	1.17	17/8246 (0.2%)	1.30	56/11222 (0.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	9
1	B	0	6
All	All	0	15

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	368	PRO	N-CD	-41.17	0.90	1.47
1	B	358	ASP	CG-OD2	28.71	1.79	1.25
1	A	462	PRO	N-CD	-16.56	1.24	1.47
1	B	462	PRO	CB-CG	-15.85	0.70	1.49
1	A	462	PRO	CB-CG	-14.89	0.75	1.49

The worst 5 of 56 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	358	ASP	CB-CG-OD2	-40.74	24.69	118.40
1	B	462	PRO	CA-N-CD	-15.91	89.73	112.00
1	B	451	SER	N-CA-C	-14.57	93.31	112.68
1	B	368	PRO	CA-N-CD	-14.17	92.16	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	368	PRO	N-CD-CG	11.04	119.77	103.20

There are no chirality outliers.

5 of 15 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	367	THR	Peptide
1	A	403	LEU	Peptide
1	A	57	SER	Peptide
1	A	61	ALA	Peptide
1	A	98	GLU	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4036	0	3954	100	0
1	B	4038	0	3954	117	0
2	C	28	0	17	0	0
2	D	28	0	25	0	0
2	E	28	0	24	4	0
2	F	28	0	21	4	0
2	G	28	0	24	2	0
2	H	28	0	25	2	0
3	A	12	0	9	0	0
3	B	12	0	11	0	0
4	A	10	0	0	0	0
4	B	10	0	0	0	0
5	A	70	0	0	2	0
5	B	50	0	0	4	0
All	All	8406	0	8064	217	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 13.

The worst 5 of 217 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:743:HOH:O	2:H:2:NAG:O7	1.56	1.21
1:A:380:ASN:ND2	1:A:513:LEU:HD11	1.55	1.18
1:A:61:ALA:HB1	1:A:62:ASP:HA	1.44	1.00
1:B:251:CYS:HB2	1:B:288:THR:O	1.61	0.98
1:A:380:ASN:CG	1:A:380:ASN:O	2.08	0.96

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	514/570 (90%)	467 (91%)	33 (6%)	14 (3%)	4	11
1	B	514/570 (90%)	469 (91%)	31 (6%)	14 (3%)	4	11
All	All	1028/1140 (90%)	936 (91%)	64 (6%)	28 (3%)	4	11

5 of 28 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	287	ASP
1	A	288	THR
1	A	404	THR
1	A	514	LYS
1	A	538	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	446/496 (90%)	407 (91%)	39 (9%)	9	25
1	B	446/496 (90%)	401 (90%)	45 (10%)	7	20
All	All	892/992 (90%)	808 (91%)	84 (9%)	8	22

5 of 84 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	213	GLN
1	B	451	SER
1	B	231	VAL
1	B	367	THR
1	B	487	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	326	GLN
1	B	385	ASN
1	B	530	GLN
1	B	529	HIS
1	B	213	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

12 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.80	0	17,19,21	2.24	7 (41%)
2	NAG	C	2	2	14,14,15	2.47	6 (42%)	17,19,21	7.14	11 (64%)
2	NAG	D	1	2	14,14,15	0.66	0	17,19,21	1.88	2 (11%)
2	NAG	D	2	2	14,14,15	1.29	1 (7%)	17,19,21	2.83	5 (29%)
2	NAG	E	1	1,2	14,14,15	0.83	0	17,19,21	3.19	5 (29%)
2	NAG	E	2	2	14,14,15	2.48	2 (14%)	17,19,21	2.81	7 (41%)
2	NAG	F	1	1,2	14,14,15	0.50	0	17,19,21	2.37	7 (41%)
2	NAG	F	2	2	14,14,15	2.56	1 (7%)	17,19,21	3.02	5 (29%)
2	NAG	G	1	2	14,14,15	0.68	0	17,19,21	1.30	2 (11%)
2	NAG	G	2	2	14,14,15	0.81	1 (7%)	17,19,21	10.80	6 (35%)
2	NAG	H	1	1,2	14,14,15	1.03	1 (7%)	17,19,21	2.17	4 (23%)
2	NAG	H	2	2	14,14,15	1.24	2 (14%)	17,19,21	1.95	4 (23%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	1,2	1/1/5/7	2/6/23/26	0/1/1/1
2	NAG	C	2	2	3/3/5/7	5/6/23/26	0/1/1/1
2	NAG	D	1	2	2/2/5/7	3/6/23/26	0/1/1/1
2	NAG	D	2	2	1/1/5/7	4/6/23/26	0/1/1/1
2	NAG	E	1	1,2	1/1/5/7	5/6/23/26	1/1/1/1
2	NAG	E	2	2	1/1/5/7	5/6/23/26	0/1/1/1
2	NAG	F	1	1,2	1/1/5/7	3/6/23/26	0/1/1/1
2	NAG	F	2	2	4/4/5/7	2/6/23/26	0/1/1/1
2	NAG	G	1	2	2/2/5/7	5/6/23/26	0/1/1/1
2	NAG	G	2	2	4/4/5/7	5/6/23/26	0/1/1/1
2	NAG	H	1	1,2	2/2/5/7	3/6/23/26	0/1/1/1
2	NAG	H	2	2	4/4/5/7	4/6/23/26	0/1/1/1

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	NAG	O3-C3	-9.16	1.20	1.43
2	E	2	NAG	O3-C3	-8.29	1.22	1.43
2	C	2	NAG	O6-C6	4.90	1.63	1.42
2	C	2	NAG	O3-C3	-4.05	1.32	1.43
2	D	2	NAG	O3-C3	4.01	1.52	1.43

The worst 5 of 65 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	NAG	C8-C7-N2	26.21	159.59	116.12
2	C	2	NAG	O3-C3-C2	25.01	161.35	109.40
2	G	2	NAG	O7-C7-N2	-24.18	79.26	121.98
2	G	2	NAG	O7-C7-C8	-23.36	80.46	122.05
2	G	2	NAG	O3-C3-C2	11.41	133.09	109.40

5 of 26 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	NAG	C2
2	C	2	NAG	C5
2	C	2	NAG	C3
2	C	2	NAG	C2
2	D	1	NAG	C3

5 of 46 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	C	2	NAG	C3-C2-N2-C7
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C1-C2-C3-C4-C5-O5

6 monomers are involved in 12 short contacts:

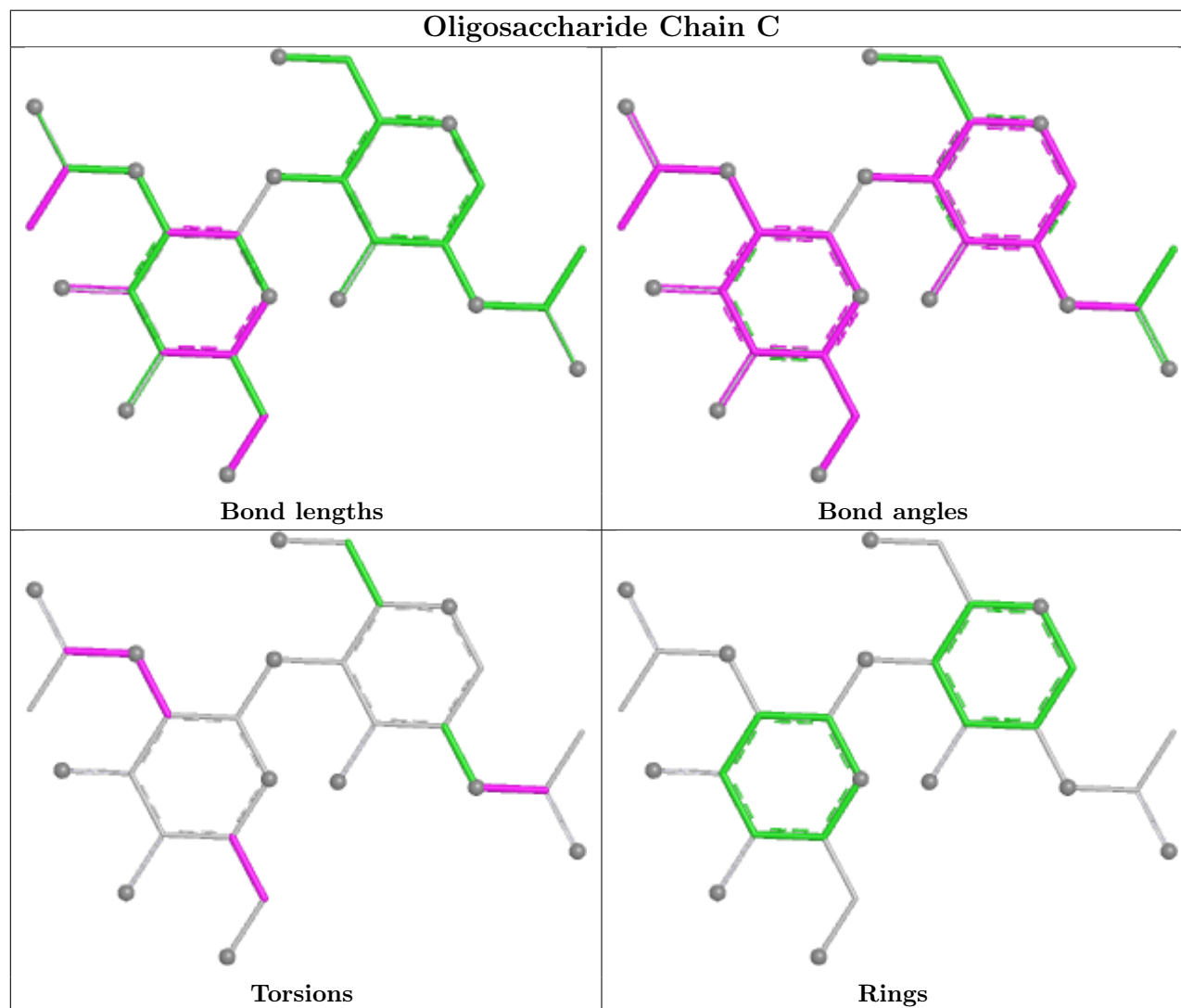
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	NAG	1	0
2	H	1	NAG	1	0

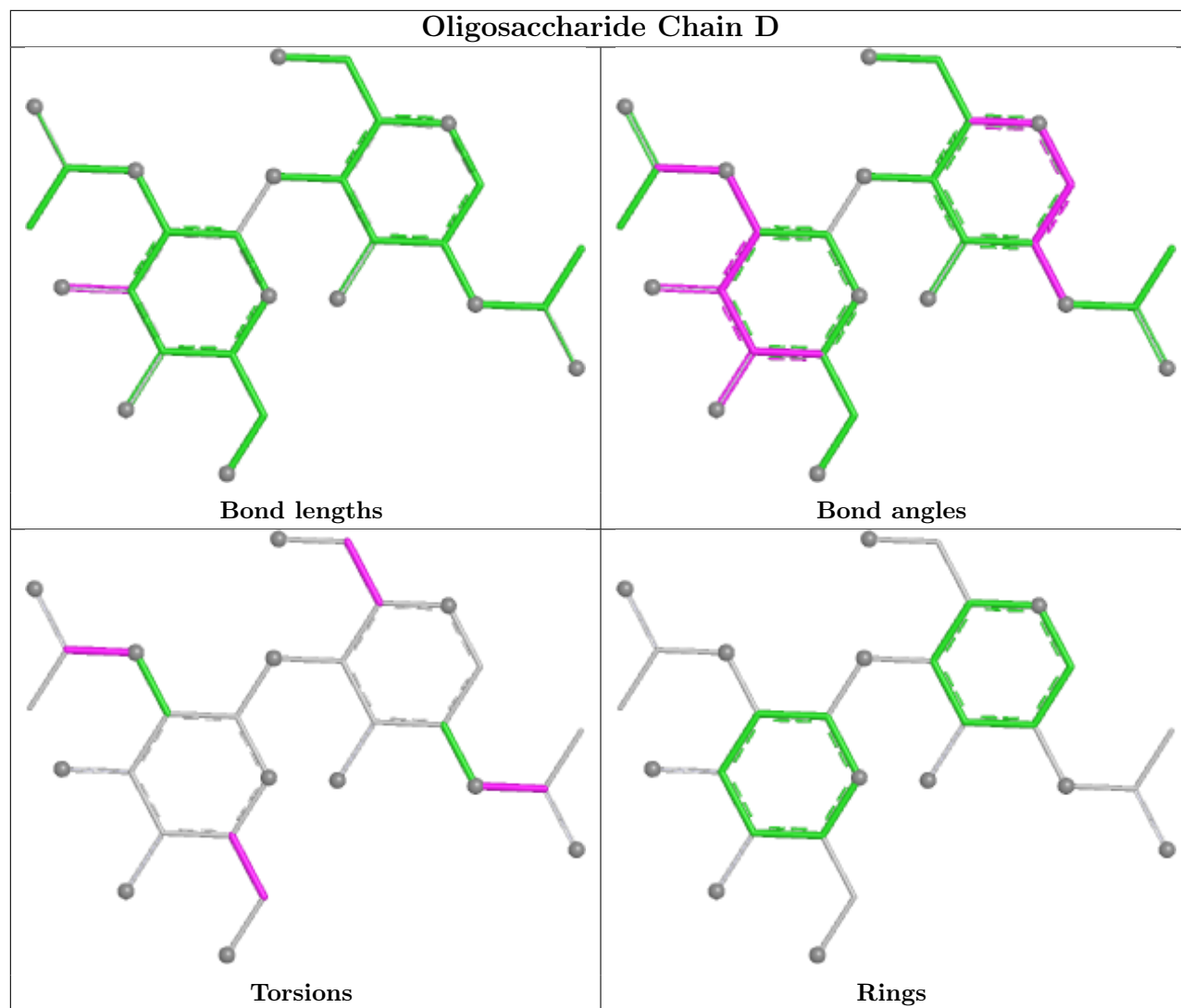
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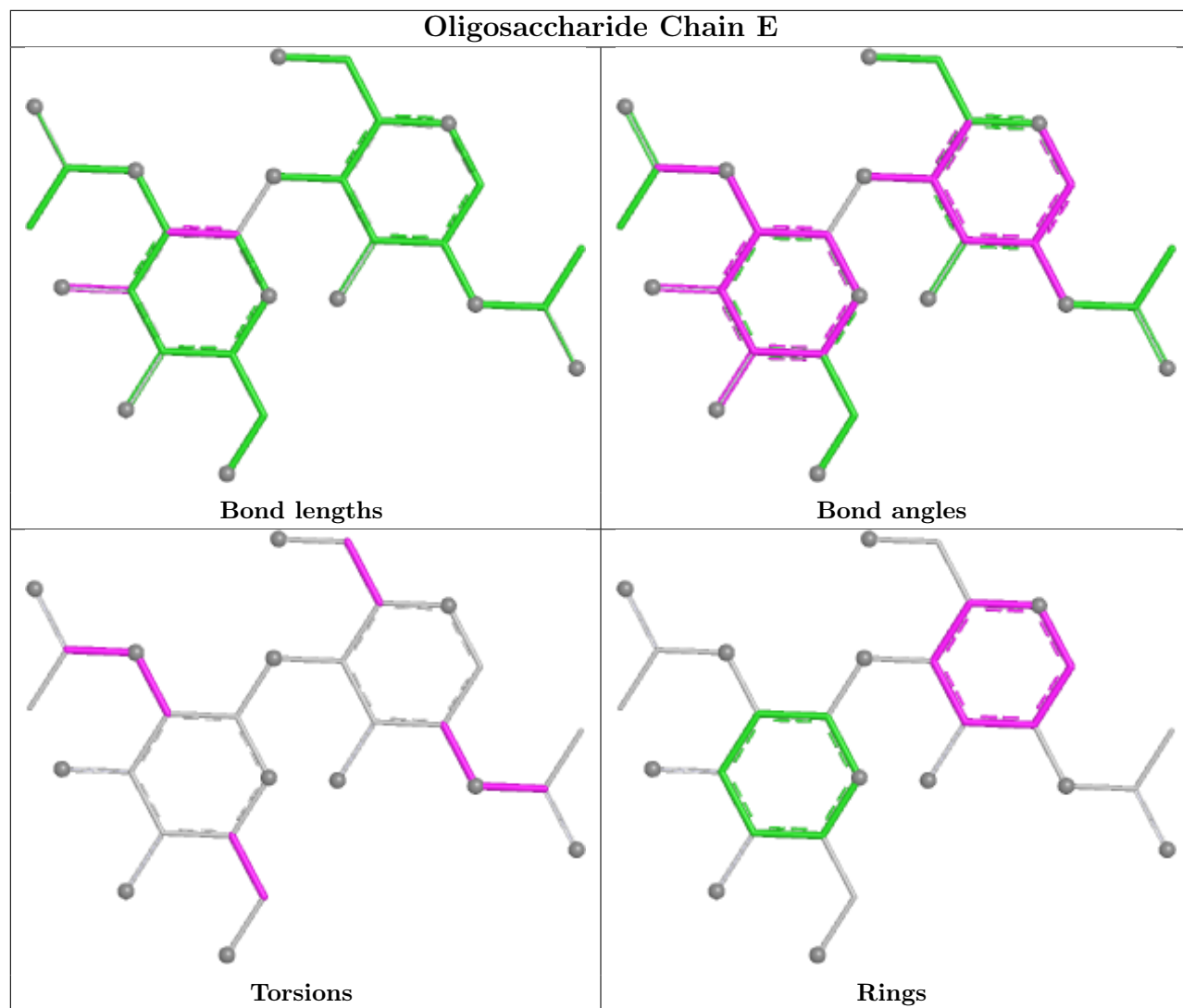
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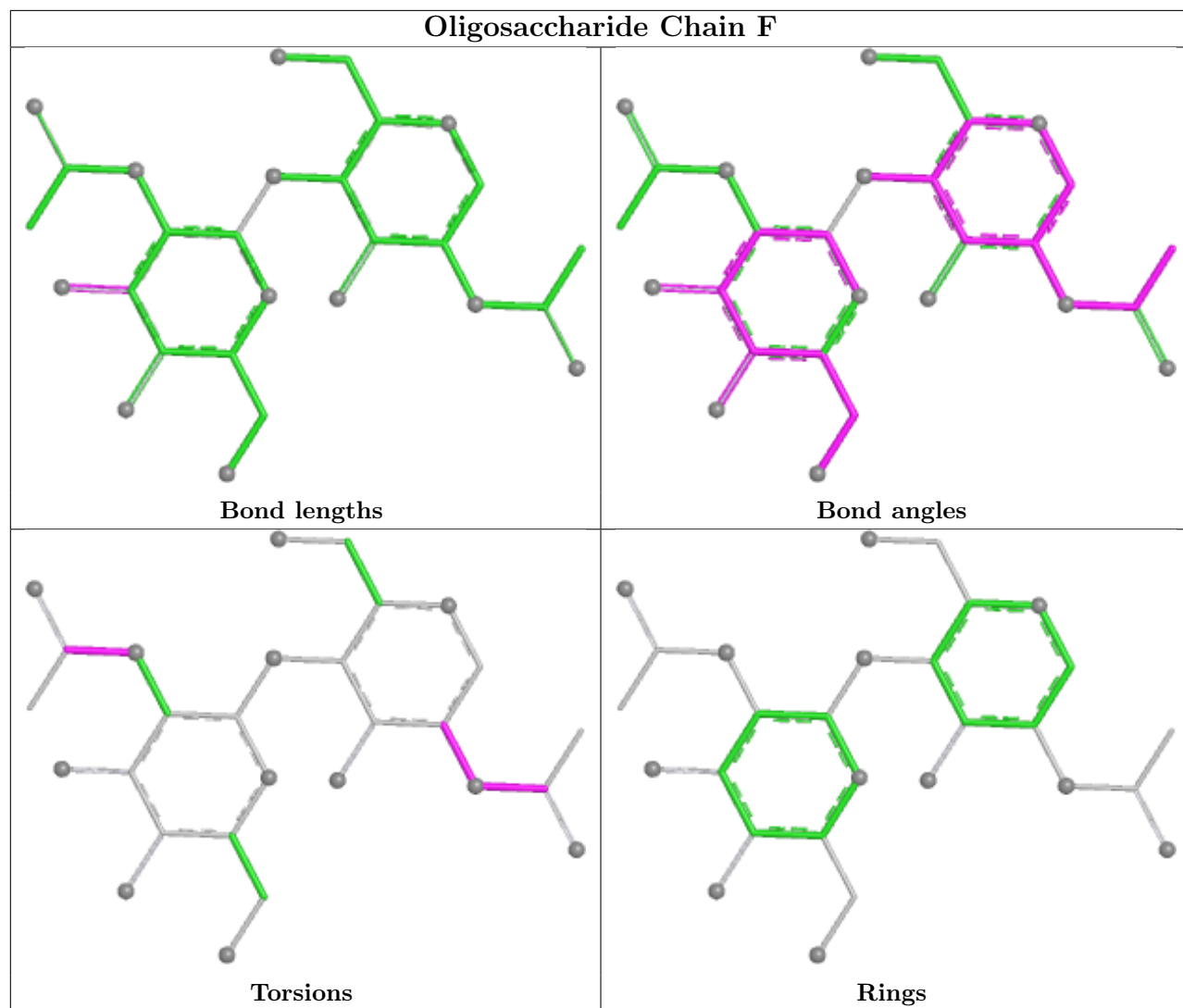
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1	NAG	4	0
2	G	2	NAG	1	0
2	G	1	NAG	1	0
2	E	1	NAG	4	0

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

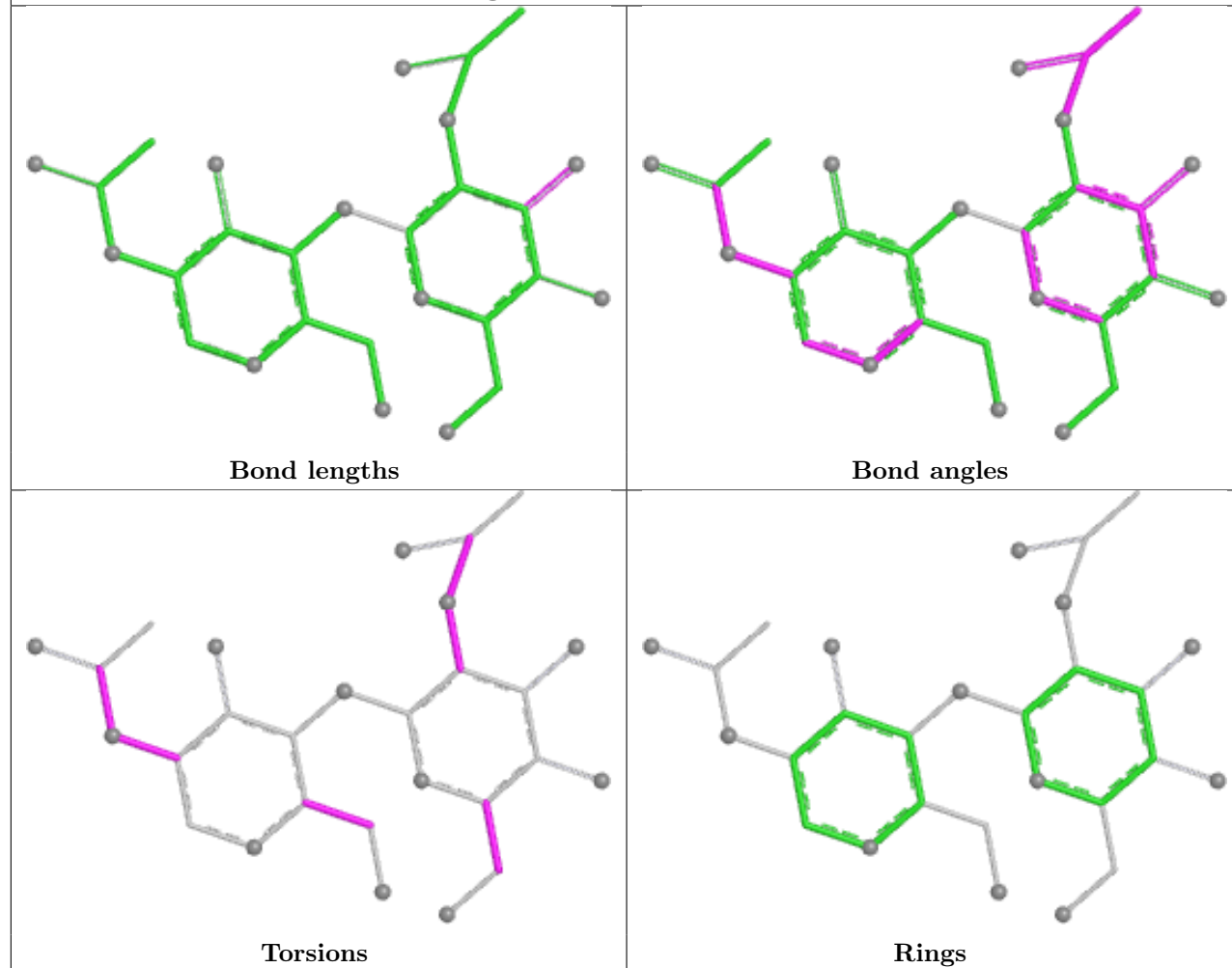


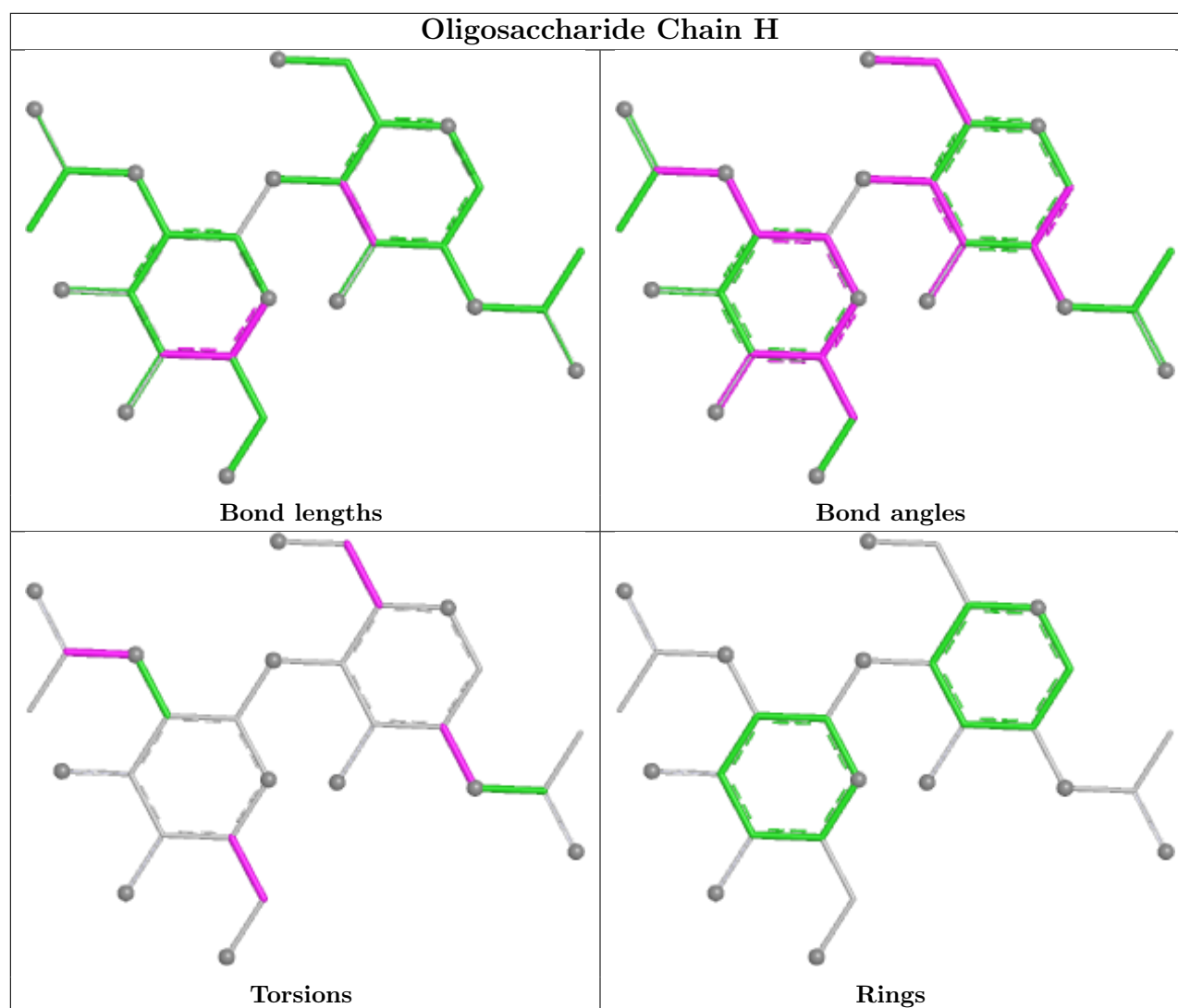






Oligosaccharide Chain G





5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
4	XYP	B	601	-	10,10,10	1.23	1 (10%)	14,14,14	1.29	2 (14%)
4	XYP	A	602	-	10,10,10	1.24	1 (10%)	14,14,14	1.28	3 (21%)
3	BMA	B	602	-	12,12,12	0.79	0	17,17,17	1.88	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BMA	A	601	-	12,12,12	0.75	0	17,17,17	2.36	7 (41%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	XYP	B	601	-	1/1/4/4	-	0/1/1/1
4	XYP	A	602	-	-	-	0/1/1/1
3	BMA	B	602	-	-	2/2/22/22	0/1/1/1
3	BMA	A	601	-	-	2/2/22/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	602	XYP	O5-C1	-3.48	1.37	1.43
4	B	601	XYP	O5-C1	-3.24	1.37	1.43

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	BMA	O6-C6-C5	-4.55	95.82	111.33
3	A	601	BMA	C4-C3-C2	-4.42	103.08	110.83
3	B	602	BMA	O6-C6-C5	-4.06	97.51	111.33
3	A	601	BMA	O4-C4-C5	3.94	119.04	109.32
3	A	601	BMA	O5-C5-C4	-3.27	103.81	109.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	601	XYP	C1

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	602	BMA	O5-C5-C6-O6
3	A	601	BMA	C4-C5-C6-O6
3	B	602	BMA	C4-C5-C6-O6
3	A	601	BMA	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.