



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:21 AM UTC

PDB ID : 1LTT / pdb_00001ltt
Title : LACTOSE BINDING TO HEAT-LABILE ENTEROTOXIN REVEALED BY
X-RAY CRYSTALLOGRAPHY
Authors : Sixma, T.K.; Hol, W.G.J.
Deposited on : 1992-07-15
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure 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/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

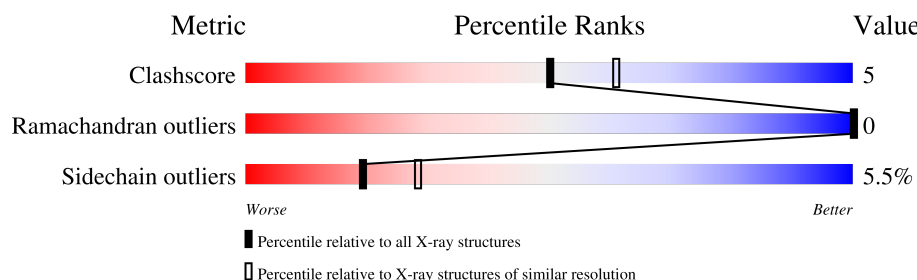
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

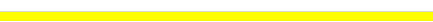
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	D	103	76% 22% .
1	E	103	80% 19% .
1	F	103	75% 17% 7% .
1	G	103	70% 27% ..
1	H	103	78% 19% .
2	A	185	65% 32% .
3	C	41	63% 32% 5%
4	B	2	50% 50%

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Mol	Chain	Length	Quality of chain
4	I	2	 100%
4	J	2	 100%
4	K	2	 100%
4	L	2	 50%  50%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6427 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 HEAT-LABILE ENTEROTOXIN, SUBUNIT B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	103	Total	C	N	O	S	0	0	0
			824	516	139	163	6			
1	E	103	Total	C	N	O	S	0	0	0
			824	516	139	163	6			
1	F	103	Total	C	N	O	S	0	0	0
			824	516	139	163	6			
1	G	103	Total	C	N	O	S	0	0	0
			824	516	139	163	6			
1	H	103	Total	C	N	O	S	0	0	0
			824	516	139	163	6			

- Molecule 2 is a protein called HEAT-LABILE ENTEROTOXIN, SUBUNIT A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	185	Total	C	N	O	S	0	0	0
			1511	953	276	278	4			

- Molecule 3 is a protein called HEAT-LABILE ENTEROTOXIN, SUBUNIT A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	41	Total	C	N	O	S	0	0	0
			347	214	59	73	1			

- Molecule 4 is an oligosaccharide called beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	B	2	Total	C	O	0	0	0
			23	12	11			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	I	2	Total	C	O	0	0	0
			23	12	11			
4	J	2	Total	C	O	0	0	0
			23	12	11			
4	K	2	Total	C	O	0	0	0
			23	12	11			
4	L	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	65	Total	O	0	0
			65	65		
5	E	56	Total	O	0	0
			56	56		
5	F	37	Total	O	0	0
			37	37		
5	G	53	Total	O	0	0
			53	53		
5	H	51	Total	O	0	0
			51	51		
5	A	43	Total	O	0	0
			43	43		
5	C	29	Total	O	0	0
			29	29		

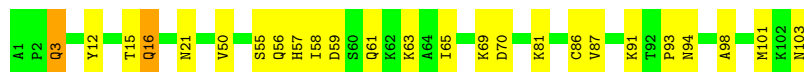
3 Residue-property plots [i](#)

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

Chain D: 



- Molecule 1: HEAT-LABILE ENTEROTOXIN, SUBUNIT B

Chain E: 



- Molecule 1: HEAT-LABILE ENTEROTOXIN, SUBUNIT B

Chain F: 



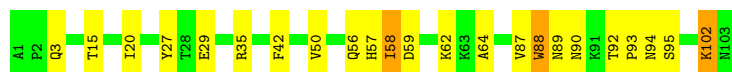
- Molecule 1: HEAT-LABILE ENTEROTOXIN, SUBUNIT B

Chain G: 



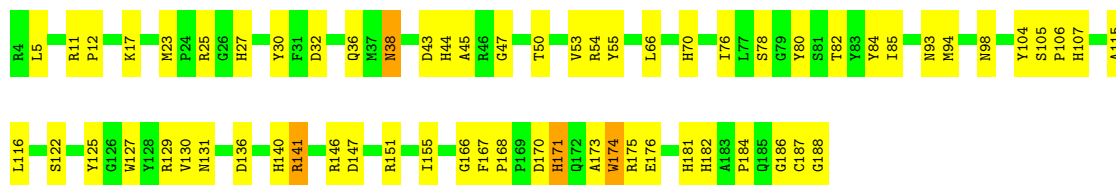
- Molecule 1: HEAT-LABILE ENTEROTOXIN, SUBUNIT B

Chain H: 



- Molecule 2: HEAT-LABILE ENTEROTOXIN, SUBUNIT A

Chain A:  65% 32%



- Molecule 3: HEAT-LABILE ENTEROTOXIN, SUBUNIT A

Chain C:  63% 32% 5%

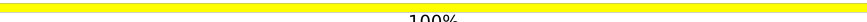


- Molecule 4: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain B:  50% 50%

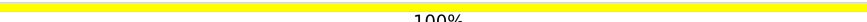


- Molecule 4: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain I:  100%

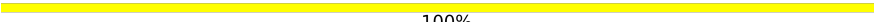


- Molecule 4: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain J:  100%



- Molecule 4: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain K:  100%



- Molecule 4: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain L:  50% 50%



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	119.80Å 101.20Å 64.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.182 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6427	wwPDB-VP
Average B, all atoms (Å ²)	23.0	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: GAL, BGC

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	D	0.98	1/835 (0.1%)	1.69	10/1124 (0.9%)
1	E	0.95	1/835 (0.1%)	1.71	14/1124 (1.2%)
1	F	1.04	1/835 (0.1%)	1.83	18/1124 (1.6%)
1	G	1.06	4/835 (0.5%)	1.83	21/1124 (1.9%)
1	H	0.98	2/835 (0.2%)	1.77	13/1124 (1.2%)
2	A	1.01	8/1559 (0.5%)	1.73	32/2120 (1.5%)
3	C	0.83	0/351	1.78	8/472 (1.7%)
All	All	1.00	17/6085 (0.3%)	1.76	116/8212 (1.4%)

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	E	0	1
1	G	0	1
All	All	0	2

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	57	HIS	CD2-NE2	-7.44	1.29	1.37
1	E	57	HIS	CD2-NE2	-6.95	1.30	1.37
2	A	140	HIS	CD2-NE2	-6.80	1.30	1.37
2	A	107	HIS	CD2-NE2	-6.59	1.30	1.37
2	A	181	HIS	CD2-NE2	-6.59	1.30	1.37
2	A	70	HIS	CD2-NE2	-6.53	1.30	1.37
1	D	57	HIS	CD2-NE2	-6.38	1.30	1.37
2	A	171	HIS	CD2-NE2	-6.26	1.30	1.37
1	F	57	HIS	CD2-NE2	-6.18	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	27	HIS	CD2-NE2	-6.15	1.31	1.37
2	A	182	HIS	CD2-NE2	-5.93	1.31	1.37
2	A	44	HIS	CD2-NE2	-5.76	1.31	1.37
1	H	57	HIS	CD2-NE2	-5.69	1.31	1.37
1	G	74	ILE	CA-CB	5.37	1.61	1.54
1	G	57	HIS	CG-ND1	-5.17	1.32	1.38
1	H	58	ILE	CA-CB	5.15	1.61	1.54
1	G	90	ASN	CG-OD1	5.03	1.33	1.23

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	50	VAL	N-CA-C	-8.70	95.29	107.99
1	F	103	ASN	CA-CB-CG	8.37	120.97	112.60
1	G	50	VAL	N-CA-C	-8.32	95.60	107.75
1	E	50	VAL	N-CA-C	-7.96	96.26	107.80
2	A	80	TYR	N-CA-C	-7.55	96.10	108.41
1	E	94	ASN	OD1-CG-ND2	-7.54	115.06	122.60
1	H	50	VAL	N-CA-C	-7.49	96.93	107.80
1	G	56	GLN	OE1-CD-NE2	-7.45	115.15	122.60
1	F	69	LYS	CG-CD-CE	7.26	127.99	111.30
1	D	103	ASN	OD1-CG-ND2	-7.22	115.38	122.60
1	E	59	ASP	N-CA-C	7.19	121.15	112.38
1	G	103	ASN	CA-CB-CG	-7.15	105.45	112.60
2	A	45	ALA	N-CA-C	7.08	119.85	111.71
1	G	99	ILE	CB-CA-C	-7.05	99.41	110.69
1	F	94	ASN	OD1-CG-ND2	-6.99	115.61	122.60
2	A	27	HIS	CB-CG-CD2	-6.95	122.16	131.20
2	A	32	ASP	CA-CB-CG	6.92	119.52	112.60
3	C	213	GLU	CA-CB-CG	-6.90	100.30	114.10
1	H	102	LYS	CA-C-N	6.83	133.99	121.70
1	H	102	LYS	C-N-CA	6.83	133.99	121.70
1	G	57	HIS	CB-CG-CD2	-6.82	122.33	131.20
1	E	3	GLN	N-CA-C	-6.74	105.61	114.31
1	F	87	VAL	O-C-N	-6.73	115.80	123.00
1	G	90	ASN	OD1-CG-ND2	-6.60	116.00	122.60
2	A	167	PHE	CA-C-N	6.58	124.46	119.66
2	A	167	PHE	C-N-CA	6.58	124.46	119.66
3	C	197	ASP	CA-C-O	-6.57	112.95	120.24
1	G	16	GLN	OE1-CD-NE2	-6.42	116.17	122.60
3	C	235	ARG	NE-CZ-NH2	-6.40	113.44	119.20
2	A	93	ASN	CA-CB-CG	6.38	118.98	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	171	HIS	CB-CG-CD2	-6.36	122.94	131.20
2	A	136	ASP	CA-CB-CG	6.34	118.94	112.60
1	F	94	ASN	CB-CG-ND2	6.32	125.88	116.40
2	A	107	HIS	CB-CG-CD2	-6.30	123.01	131.20
1	E	94	ASN	CB-CG-ND2	6.29	125.84	116.40
1	D	94	ASN	CB-CG-ND2	6.25	125.78	116.40
2	A	38	ASN	OD1-CG-ND2	-6.25	116.34	122.60
2	A	140	HIS	CA-CB-CG	-6.25	107.55	113.80
1	F	56	GLN	OE1-CD-NE2	-6.21	116.39	122.60
1	D	94	ASN	OD1-CG-ND2	-6.20	116.40	122.60
1	F	23	LYS	CA-CB-CG	6.16	126.41	114.10
1	F	59	ASP	N-CA-C	6.14	118.76	111.33
3	C	211	LEU	CA-C-O	-6.13	114.38	120.70
1	D	59	ASP	N-CA-C	6.12	118.75	111.71
1	F	3	GLN	CA-CB-CG	-6.07	101.96	114.10
1	F	91	LYS	CA-CB-CG	-6.07	101.97	114.10
1	G	94	ASN	OD1-CG-ND2	-6.07	116.53	122.60
1	D	57	HIS	CB-CG-CD2	-6.06	123.33	131.20
1	F	7	GLU	CB-CA-C	-6.05	100.57	110.85
1	G	102	LYS	CA-C-N	6.04	132.58	121.70
1	G	102	LYS	C-N-CA	6.04	132.58	121.70
2	A	182	HIS	CB-CG-CD2	-5.99	123.41	131.20
1	D	50	VAL	N-CA-C	-5.97	99.14	107.80
1	H	89	ASN	CA-CB-CG	5.97	118.57	112.60
1	D	21	ASN	OD1-CG-ND2	-5.95	116.65	122.60
2	A	98	ASN	CA-CB-CG	5.93	118.53	112.60
2	A	174	TRP	CG-CD2-CE3	5.88	139.78	133.90
2	A	181	HIS	CB-CG-CD2	-5.83	123.62	131.20
1	H	58	ILE	N-CA-C	-5.83	103.69	110.05
2	A	127	TRP	CG-CD2-CE3	5.82	139.72	133.90
2	A	170	ASP	CA-CB-CG	5.82	118.42	112.60
1	G	57	HIS	CB-CG-ND1	5.81	131.41	122.70
1	G	52	VAL	N-CA-C	-5.80	102.64	108.96
2	A	27	HIS	CB-CG-ND1	5.75	131.32	122.70
1	D	56	GLN	OE1-CD-NE2	-5.70	116.90	122.60
1	E	57	HIS	CB-CG-CD2	-5.69	123.80	131.20
1	H	88	TRP	CG-CD2-CE3	5.67	139.57	133.90
1	G	99	ILE	CB-CG1-CD1	-5.66	101.92	113.80
1	D	16	GLN	OE1-CD-NE2	-5.62	116.97	122.60
3	C	205	ASN	OD1-CG-ND2	-5.62	116.98	122.60
1	H	35	ARG	NE-CZ-NH2	-5.61	114.15	119.20
1	G	102	LYS	O-C-N	5.60	130.11	123.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	71	THR	CA-CB-OG1	-5.59	101.22	109.60
1	H	94	ASN	OD1-CG-ND2	-5.58	117.03	122.60
2	A	174	TRP	CB-CG-CD1	-5.54	118.59	126.90
3	C	236	ILE	CA-CB-CG1	-5.52	101.01	110.40
1	H	35	ARG	CG-CD-NE	-5.52	99.86	112.00
1	F	1	ALA	CA-C-N	5.52	125.46	119.78
1	F	1	ALA	C-N-CA	5.52	125.46	119.78
1	E	88	TRP	CG-CD2-CE3	5.45	139.35	133.90
1	E	49	GLN	OE1-CD-NE2	-5.37	117.23	122.60
1	H	64	ALA	CA-C-N	5.36	127.31	120.56
1	H	64	ALA	C-N-CA	5.36	127.31	120.56
1	G	42	PHE	CA-CB-CG	-5.35	108.45	113.80
1	D	70	ASP	CA-CB-CG	5.34	117.94	112.60
2	A	11	ARG	CA-C-N	5.33	125.87	120.38
2	A	11	ARG	C-N-CA	5.33	125.87	120.38
1	E	70	ASP	CA-CB-CG	5.32	117.92	112.60
2	A	43	ASP	CA-CB-CG	5.31	117.91	112.60
2	A	53	VAL	N-CA-C	-5.30	102.74	109.58
1	G	60	SER	CA-CB-OG	5.30	121.70	111.10
3	C	236	ILE	N-CA-CB	-5.29	102.51	111.50
1	F	90	ASN	N-CA-C	-5.28	107.02	112.93
1	F	88	TRP	CB-CG-CD1	-5.27	119.00	126.90
1	H	88	TRP	CB-CG-CD1	-5.23	119.05	126.90
1	G	13	ARG	NE-CZ-NH2	-5.23	114.50	119.20
2	A	106	PRO	O-C-N	-5.21	115.98	122.23
1	G	16	GLN	CG-CD-NE2	5.20	124.19	116.40
2	A	12	PRO	N-CA-C	5.17	117.01	110.70
1	E	21	ASN	OD1-CG-ND2	-5.16	117.44	122.60
1	E	88	TRP	CE2-CD2-CG	-5.15	101.02	107.20
1	E	101	MET	CG-SD-CE	-5.15	89.58	100.90
2	A	44	HIS	CA-C-N	5.12	127.66	120.28
2	A	44	HIS	C-N-CA	5.12	127.66	120.28
1	G	98	ALA	CA-C-O	-5.12	115.80	121.33
1	G	61	GLN	CG-CD-NE2	5.11	124.07	116.40
1	E	83	ASP	CA-CB-CG	5.10	117.70	112.60
2	A	174	TRP	CE2-CD2-CG	-5.10	101.08	107.20
1	H	102	LYS	O-C-N	5.08	129.56	123.36
1	F	69	LYS	O-C-N	-5.07	116.29	122.22
1	G	14	ASN	OD1-CG-ND2	-5.04	117.56	122.60
2	A	105	SER	CA-C-N	5.03	124.63	119.05
2	A	105	SER	C-N-CA	5.03	124.63	119.05
3	C	200	ASN	OD1-CG-ND2	-5.02	117.58	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	88	TRP	CB-CG-CD1	-5.01	119.38	126.90
2	A	140	HIS	CB-CG-CD2	-5.00	124.70	131.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	76	TYR	Sidechain
1	G	76	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	D	824	0	841	8	0
1	E	824	0	841	5	0
1	F	824	0	841	9	0
1	G	824	0	841	13	0
1	H	824	0	841	10	0
2	A	1511	0	1407	29	0
3	C	347	0	327	6	0
4	B	23	0	21	0	0
4	I	23	0	21	0	0
4	J	23	0	21	0	0
4	K	23	0	21	0	0
4	L	23	0	21	0	0
5	A	43	0	0	1	0
5	C	29	0	0	0	0
5	D	65	0	0	0	0
5	E	56	0	0	0	0
5	F	37	0	0	1	0
5	G	53	0	0	0	0
5	H	51	0	0	0	0
All	All	6427	0	6044	64	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 5.

All (64) 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:G:54:GLY:H	1:G:57:HIS:HD2	1.32	0.77
2:A:174:TRP:HB3	2:A:188:GLY:HA2	1.66	0.76
1:D:93:PRO:HG3	1:H:3:GLN:HG2	1.70	0.73
1:H:59:ASP:HA	1:H:62:LYS:HD3	1.75	0.69
3:C:198:THR:HA	3:C:201:GLU:HG2	1.75	0.68
1:G:1:ALA:HA	1:H:92:THR:O	1.93	0.68
1:G:3:GLN:HG2	1:H:93:PRO:HG3	1.76	0.68
1:F:82:ILE:HD12	1:F:99:ILE:HD11	1.78	0.66
1:F:61:GLN:O	1:F:65:ILE:HG13	1.98	0.64
2:A:84:TYR:CE2	2:A:129:ARG:HG2	2.33	0.63
1:G:53:PRO:HA	1:G:57:HIS:CD2	2.35	0.61
2:A:23:MET:HE1	2:A:30:TYR:HD1	1.68	0.58
1:E:37:MET:HE2	1:E:52:VAL:HG23	1.86	0.57
2:A:125:TYR:O	2:A:141:ARG:HD2	2.04	0.57
1:F:25:LEU:HB2	1:F:43:LYS:HD2	1.89	0.55
2:A:84:TYR:HE2	2:A:129:ARG:HG2	1.72	0.55
1:G:15:THR:HA	1:G:87:VAL:O	2.07	0.54
2:A:147:ASP:O	2:A:151:ARG:HB2	2.08	0.53
2:A:54:ARG:HG2	5:A:217:HOH:O	2.08	0.53
2:A:23:MET:HE1	2:A:30:TYR:CD1	2.44	0.52
2:A:82:THR:HA	2:A:130:VAL:O	2.10	0.51
2:A:122:SER:HB2	2:A:146:ARG:HG2	1.92	0.51
2:A:168:PRO:HG2	2:A:171:HIS:HB2	1.93	0.51
1:G:54:GLY:N	1:G:57:HIS:HD2	2.05	0.51
1:D:12:TYR:OH	1:E:35:ARG:HG3	2.12	0.50
1:G:99:ILE:HD12	1:H:29:GLU:HB3	1.94	0.49
2:A:94:MET:SD	2:A:115:ALA:HB2	2.53	0.49
2:A:125:TYR:CE1	2:A:141:ARG:HD3	2.48	0.49
2:A:129:ARG:NH2	2:A:131:ASN:HD21	2.12	0.48
1:H:15:THR:HA	1:H:87:VAL:O	2.14	0.48
2:A:104:TYR:CD1	2:A:171:HIS:HE1	2.32	0.47
1:F:103:ASN:HB2	2:A:151:ARG:HH12	1.78	0.47
1:D:3:GLN:HG3	1:E:47:THR:HG21	1.98	0.46
1:F:65:ILE:HG22	1:F:69:LYS:HE2	1.97	0.46
1:H:88:TRP:HB3	1:H:90:ASN:OD1	2.15	0.46
2:A:184:PRO:HB3	3:C:202:GLU:HB3	1.99	0.45
1:D:86:CYS:HB3	1:D:98:ALA:HB3	1.98	0.45
1:G:85:LEU:HD23	1:G:99:ILE:HG23	1.98	0.45
2:A:5:LEU:HD12	2:A:94:MET:HE3	1.98	0.45
1:G:54:GLY:H	1:G:57:HIS:CD2	2.23	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:25:ARG:HH11	2:A:25:ARG:HG2	1.82	0.45
3:C:217:LYS:HA	3:C:217:LYS:HD3	1.64	0.45
1:H:20:ILE:HG21	1:H:42:PHE:CE1	2.53	0.43
1:H:88:TRP:HE3	1:H:95:SER:HB2	1.83	0.43
2:A:25:ARG:HG2	2:A:25:ARG:NH1	2.33	0.43
2:A:38:ASN:HA	3:C:204:GLN:OE1	2.19	0.43
1:F:12:TYR:OH	1:G:35:ARG:HB2	2.19	0.43
2:A:174:TRP:HB3	2:A:188:GLY:CA	2.43	0.42
2:A:174:TRP:CD1	2:A:187:CYS:HB3	2.54	0.42
1:D:61:GLN:HG2	1:E:31:MET:O	2.19	0.42
1:F:23:LYS:HD3	5:F:113:HOH:O	2.20	0.42
1:D:15:THR:HA	1:D:87:VAL:O	2.20	0.42
1:G:85:LEU:CD2	1:G:99:ILE:HG23	2.50	0.41
2:A:47:GLY:O	2:A:54:ARG:NH1	2.49	0.41
2:A:186:GLY:O	3:C:199:CYS:HB2	2.20	0.41
2:A:173:ALA:HA	2:A:176:GLU:HG2	2.02	0.41
1:D:65:ILE:HG22	1:D:69:LYS:HE2	2.03	0.41
2:A:66:LEU:HD12	2:A:85:ILE:HG21	2.02	0.41
1:F:1:ALA:N	1:G:92:THR:O	2.52	0.40
1:G:101:MET:HE3	1:H:27:TYR:CD1	2.56	0.40
1:D:58:ILE:HD12	1:E:34:LYS:HE2	2.03	0.40
1:F:4:THR:HB	1:F:7:GLU:HG3	2.02	0.40
2:A:5:LEU:HD11	2:A:155:ILE:HD11	2.01	0.40
2:A:166:GLY:HA2	3:C:203:THR:HG21	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	D	101/103 (98%)	98 (97%)	3 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	101/103 (98%)	98 (97%)	3 (3%)	0	100	100
1	F	101/103 (98%)	97 (96%)	4 (4%)	0	100	100
1	G	101/103 (98%)	96 (95%)	5 (5%)	0	100	100
1	H	101/103 (98%)	98 (97%)	3 (3%)	0	100	100
2	A	183/185 (99%)	175 (96%)	8 (4%)	0	100	100
3	C	39/41 (95%)	39 (100%)	0	0	100	100
All	All	727/741 (98%)	701 (96%)	26 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	D	95/95 (100%)	88 (93%)	7 (7%)	13	17
1	E	95/95 (100%)	91 (96%)	4 (4%)	26	40
1	F	95/95 (100%)	89 (94%)	6 (6%)	16	23
1	G	95/95 (100%)	90 (95%)	5 (5%)	20	30
1	H	95/95 (100%)	92 (97%)	3 (3%)	34	51
2	A	155/155 (100%)	146 (94%)	9 (6%)	18	26
3	C	40/40 (100%)	37 (92%)	3 (8%)	12	17
All	All	670/670 (100%)	633 (94%)	37 (6%)	19	28

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	3	GLN
1	D	16	GLN
1	D	55	SER
1	D	63	LYS
1	D	81	LYS

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Mol	Chain	Res	Type
1	D	91	LYS
1	D	101	MET
1	E	23	LYS
1	E	35	ARG
1	E	95	SER
1	E	99	ILE
1	F	4	THR
1	F	43	LYS
1	F	60	SER
1	F	91	LYS
1	F	101	MET
1	F	103	ASN
1	G	17	ILE
1	G	35	ARG
1	G	59	ASP
1	G	91	LYS
1	G	99	ILE
1	H	56	GLN
1	H	58	ILE
1	H	102	LYS
2	A	17	LYS
2	A	36	GLN
2	A	50	THR
2	A	55	TYR
2	A	76	ILE
2	A	78	SER
2	A	116	LEU
2	A	141	ARG
2	A	175	ARG
3	C	200	ASN
3	C	234	ASN
3	C	236	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	3	GLN
1	D	16	GLN
1	D	21	ASN
1	F	3	GLN
1	F	14	ASN
1	F	49	GLN

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Mol	Chain	Res	Type
1	F	94	ASN
1	F	103	ASN
1	G	16	GLN
1	G	21	ASN
1	G	49	GLN
1	G	57	HIS
1	G	103	ASN
1	H	89	ASN
1	H	94	ASN
2	A	38	ASN
2	A	131	ASN
2	A	171	HIS

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 ⓘ

10 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	$\# Z > 2$
4	BGC	B	1	4	12,12,12	0.47	0	17,17,17	0.83	1 (5%)
4	GAL	B	2	4	11,11,12	0.82	0	15,15,17	0.76	0
4	BGC	I	1	4	12,12,12	0.72	0	17,17,17	1.33	2 (11%)
4	GAL	I	2	4	11,11,12	0.90	1 (9%)	15,15,17	1.37	2 (13%)
4	BGC	J	1	4	12,12,12	1.13	1 (8%)	17,17,17	1.46	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GAL	J	2	4	11,11,12	0.70	0	15,15,17	1.22	2 (13%)
4	BGC	K	1	4	12,12,12	0.99	1 (8%)	17,17,17	0.85	0
4	GAL	K	2	4	11,11,12	1.05	1 (9%)	15,15,17	0.83	0
4	BGC	L	1	4	12,12,12	0.91	1 (8%)	17,17,17	0.96	0
4	GAL	L	2	4	11,11,12	0.88	0	15,15,17	1.00	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BGC	B	1	4	-	0/2/22/22	0/1/1/1
4	GAL	B	2	4	-	0/2/19/22	0/1/1/1
4	BGC	I	1	4	-	1/2/22/22	0/1/1/1
4	GAL	I	2	4	-	0/2/19/22	0/1/1/1
4	BGC	J	1	4	-	2/2/22/22	0/1/1/1
4	GAL	J	2	4	-	0/2/19/22	0/1/1/1
4	BGC	K	1	4	-	0/2/22/22	0/1/1/1
4	GAL	K	2	4	-	0/2/19/22	0/1/1/1
4	BGC	L	1	4	-	2/2/22/22	0/1/1/1
4	GAL	L	2	4	-	0/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	K	2	GAL	C4-C5	2.57	1.58	1.53
4	I	2	GAL	C2-C3	2.48	1.56	1.52
4	K	1	BGC	C4-C5	2.23	1.57	1.53
4	J	1	BGC	C6-C5	2.09	1.58	1.51
4	L	1	BGC	C6-C5	2.07	1.58	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	1	BGC	C4-C3-C2	-3.94	103.91	110.83
4	I	1	BGC	C3-C4-C5	3.82	117.16	110.23
4	J	2	GAL	C1-O5-C5	3.57	116.98	112.19
4	I	2	GAL	C1-C2-C3	-3.42	104.67	109.64
4	B	1	BGC	C3-C4-C5	2.39	114.56	110.23
4	J	1	BGC	O4-C4-C3	2.37	115.96	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	1	BGC	O4-C4-C5	-2.33	103.58	109.32
4	J	2	GAL	O3-C3-C2	-2.32	105.32	110.05
4	J	1	BGC	O3-C3-C4	2.07	115.27	110.38
4	I	2	GAL	O2-C2-C3	2.06	114.41	110.15

There are no chirality outliers.

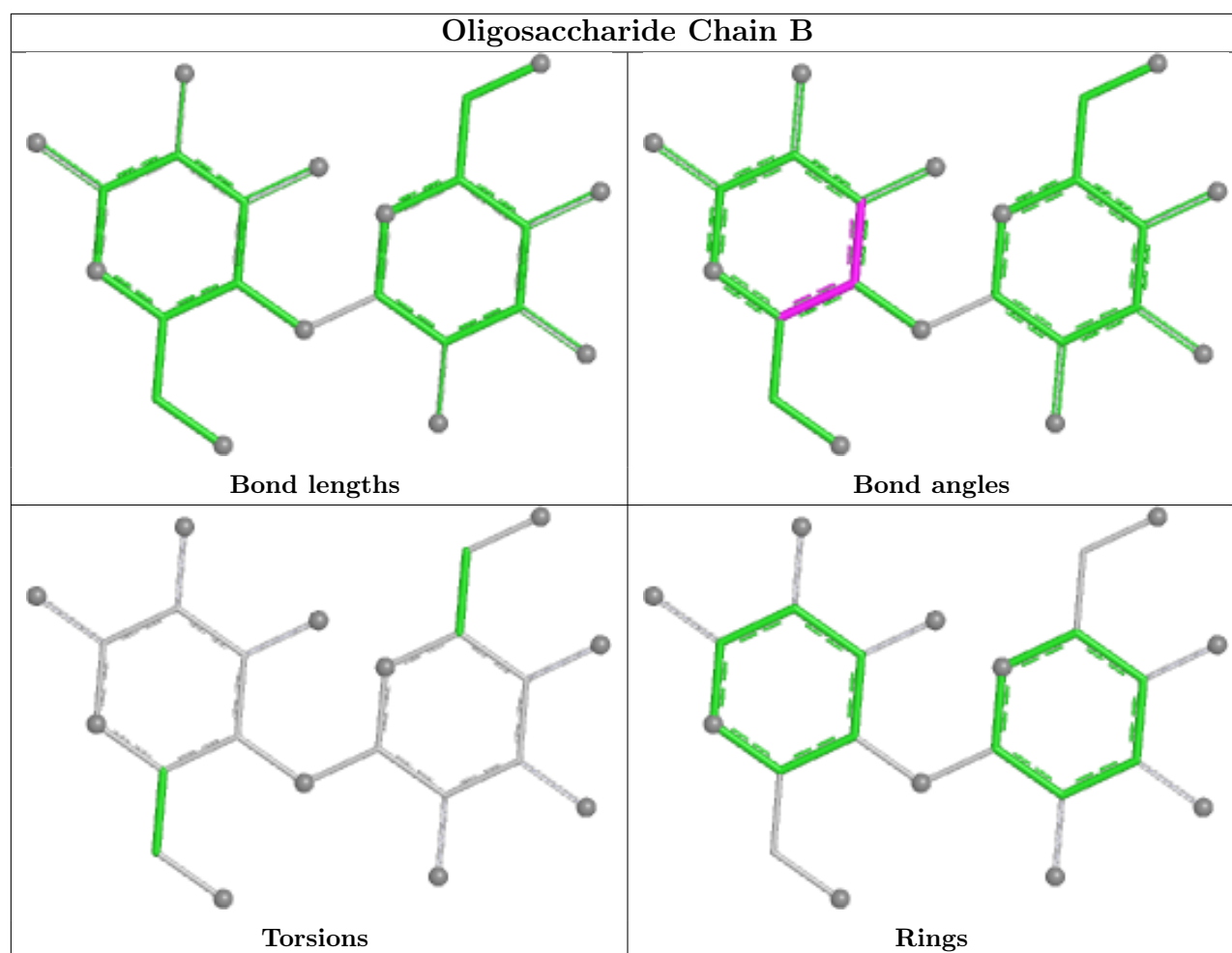
All (5) torsion outliers are listed below:

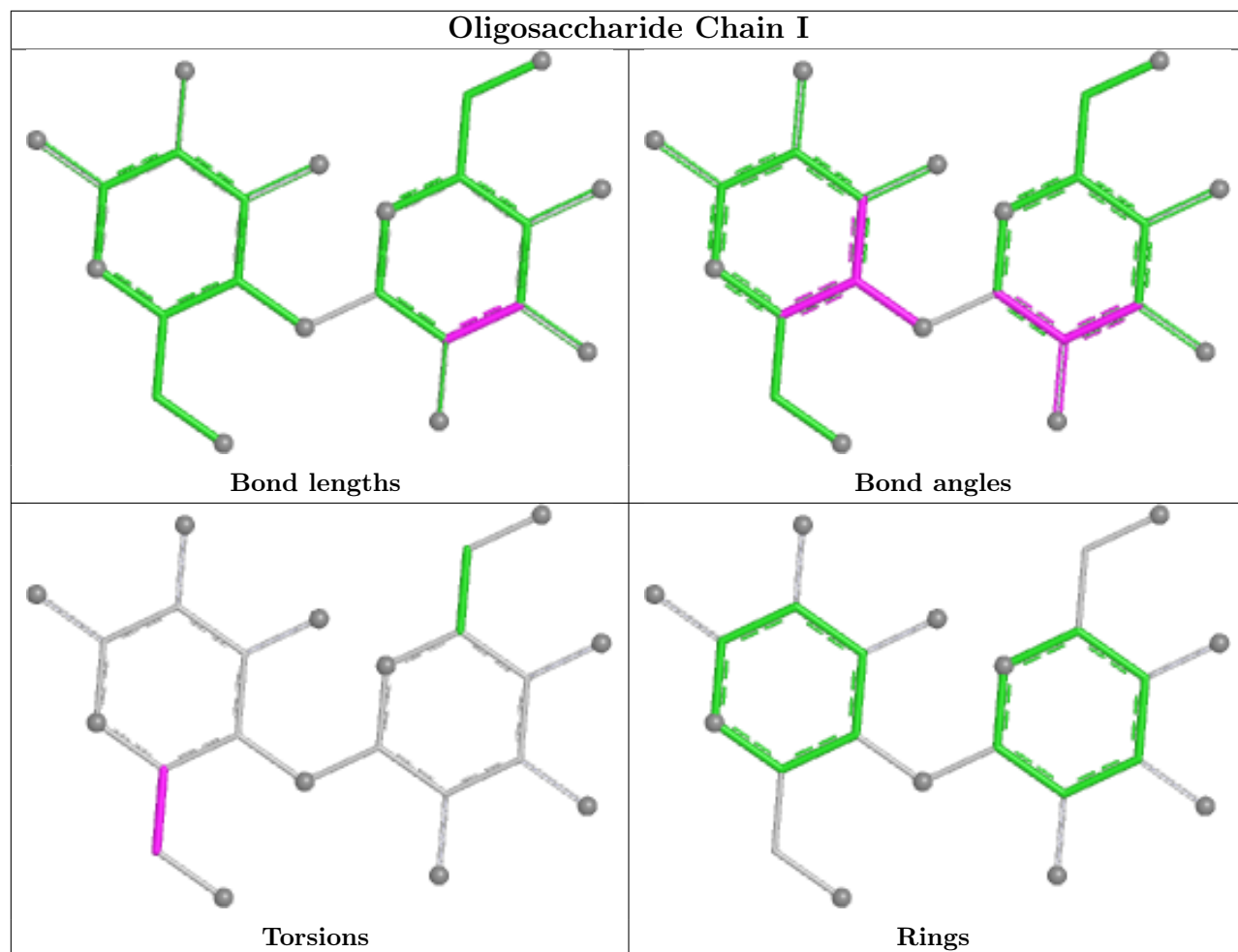
Mol	Chain	Res	Type	Atoms
4	J	1	BGC	O5-C5-C6-O6
4	I	1	BGC	O5-C5-C6-O6
4	J	1	BGC	C4-C5-C6-O6
4	L	1	BGC	C4-C5-C6-O6
4	L	1	BGC	O5-C5-C6-O6

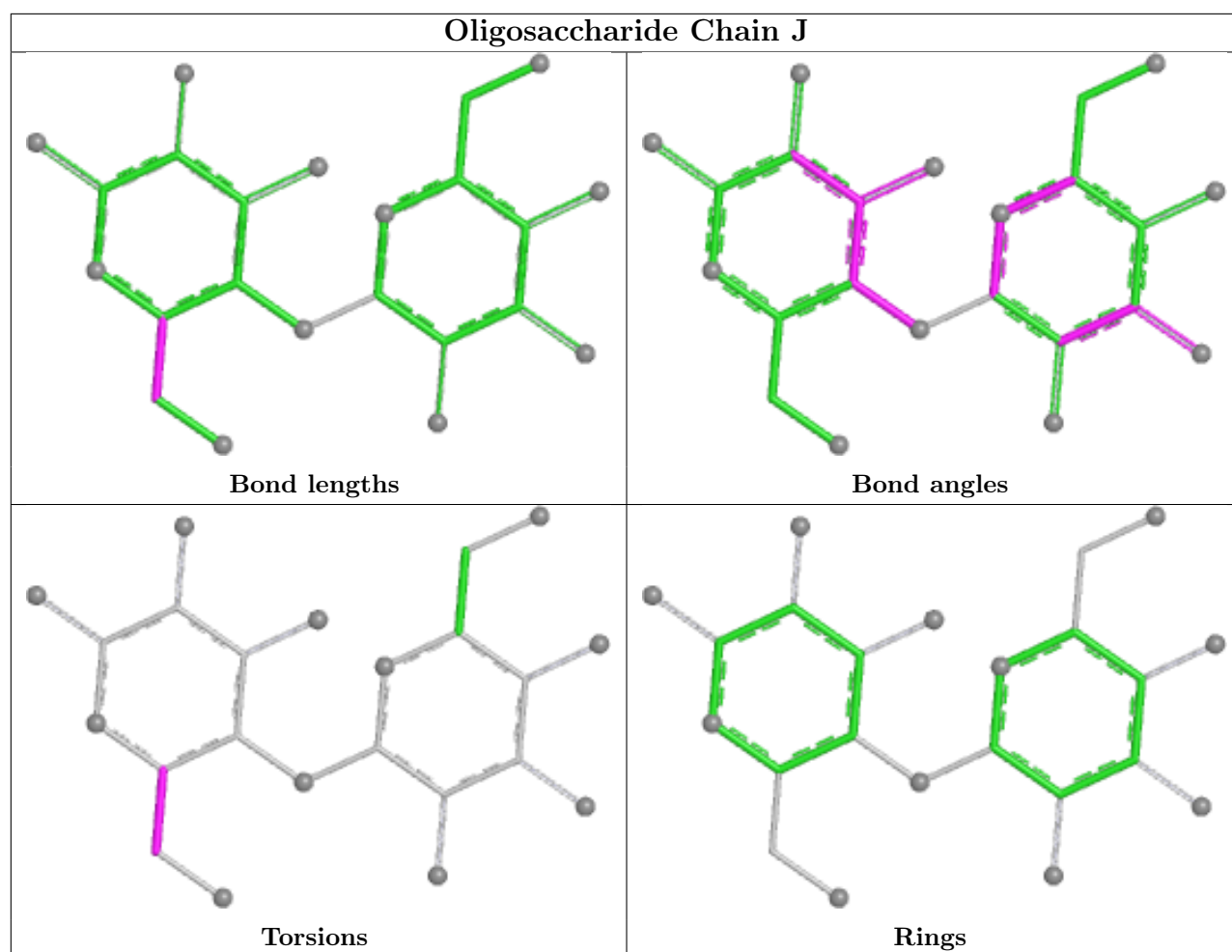
There are no ring outliers.

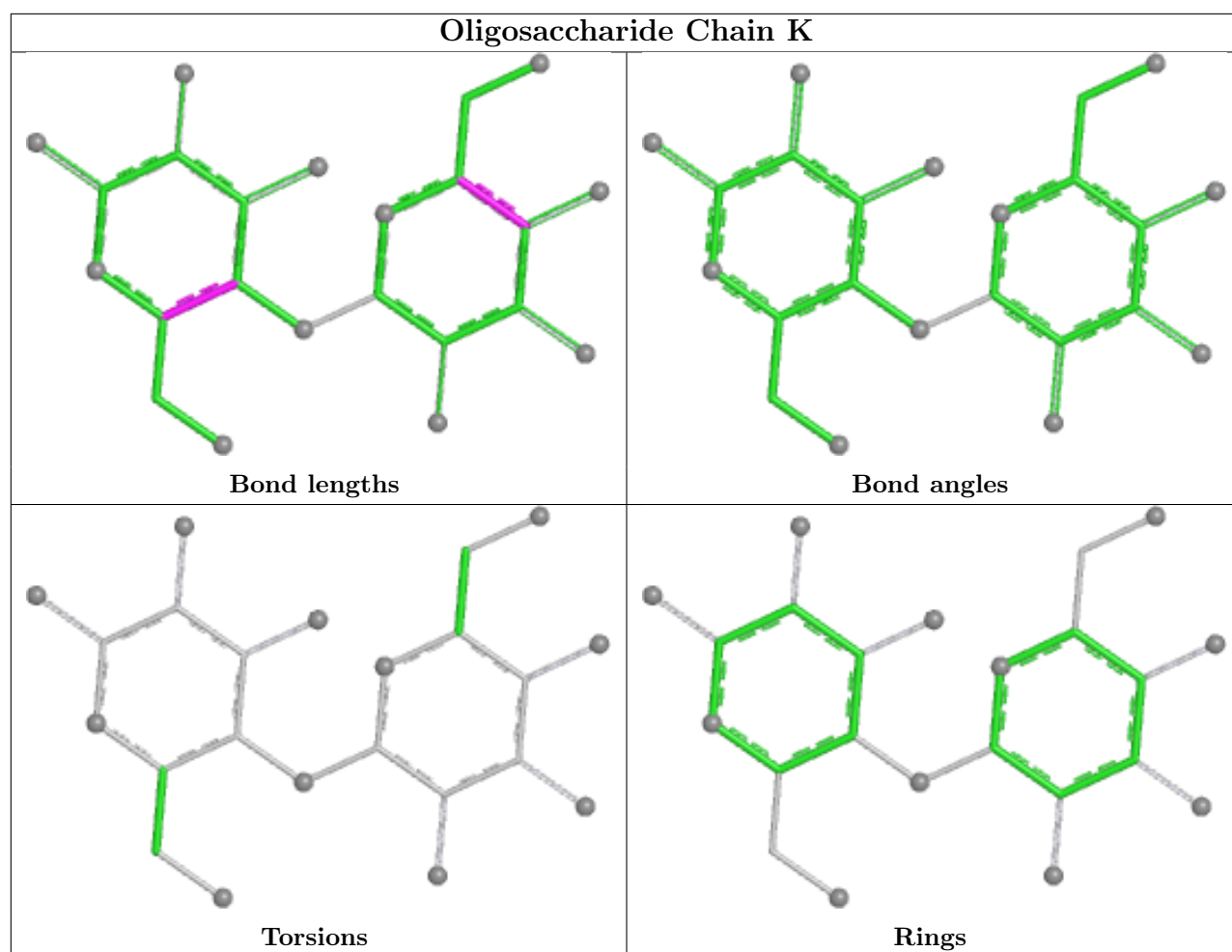
No monomer is involved in short contacts.

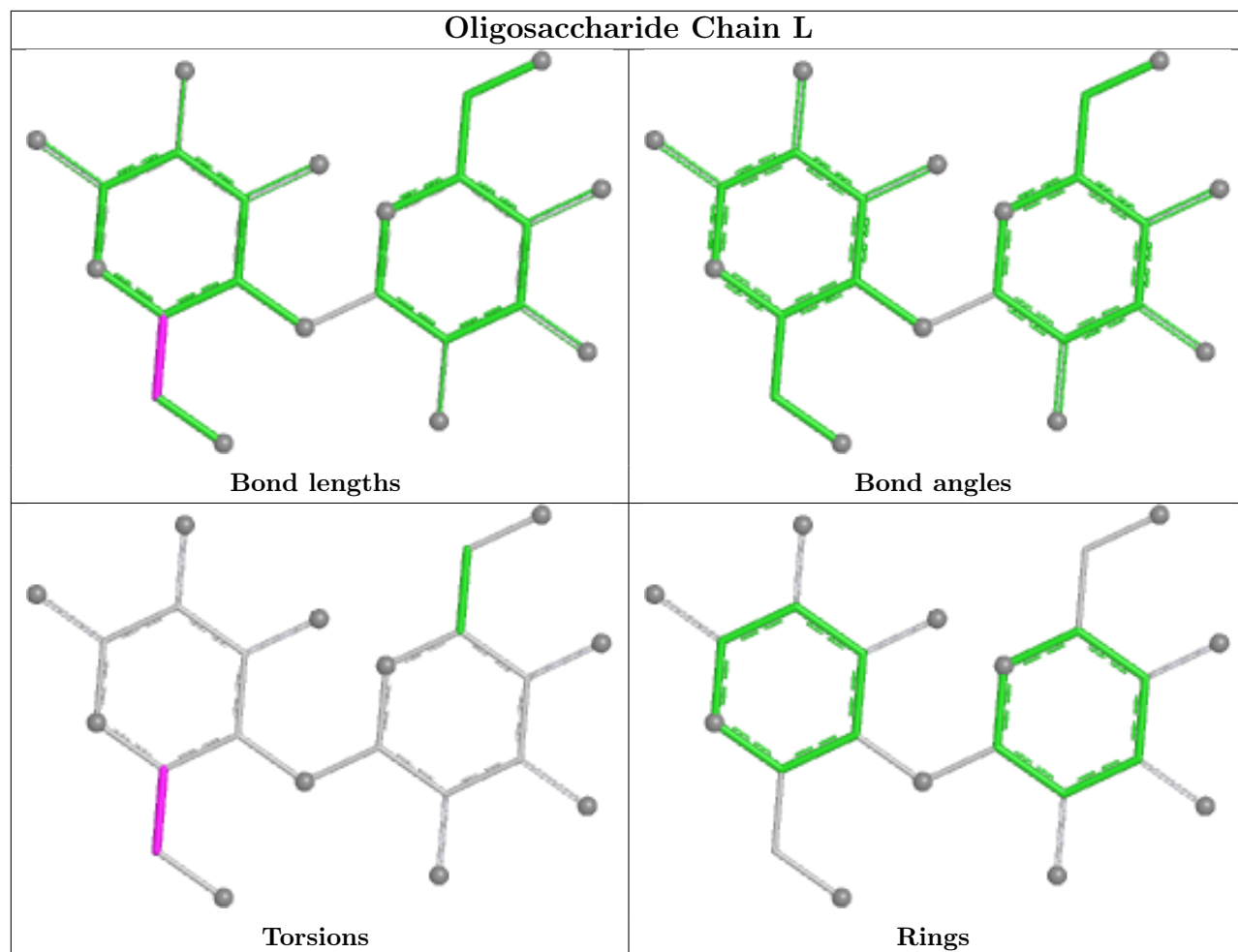
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.