



wwPDB EM Validation Summary Report ⓘ

Oct 16, 2024 – 12:23 AM JST

PDB ID : 8HWS
EMDB ID : EMD-35063
Title : The complex structure of Omicron BA.4 RBD with BD604, S309, and S304
Authors : He, Q.W.; Xu, Z.P.; Xie, Y.F.
Deposited on : 2023-01-02
Resolution : 2.36 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : **FAILED**
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

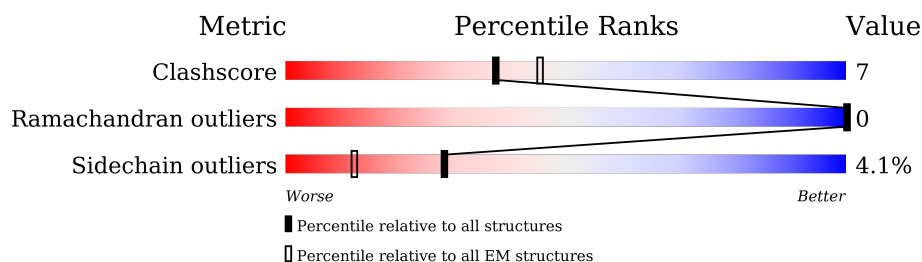
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.36 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	196	84% 15% .
2	H	229	74% 21% . .
3	L	215	79% 20% .
4	D	231	75% 19% . .
5	E	216	78% 20% ..
6	I	239	83% 12% . .
7	J	215	76% 22% .
8	B	3	67% 33%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 11511 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 Spike protein S2'.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	196	Total	C	N	O	S	0	0
			1563	1007	265	283	8		

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	452	ARG	LEU	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2
A	486	VAL	PHE	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2

- Molecule 2 is a protein called BD-604 Fab Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	219	Total	C	N	O	S	0	0
			1617	1016	272	322	7		

- Molecule 3 is a protein called BD-604 Fab Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	213	Total	C	N	O	S	0	0
			1632	1021	272	335	4		

- Molecule 4 is a protein called S304 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	222	Total	C	N	O	S	0	0
			1662	1048	276	331	7		

- Molecule 5 is a protein called S304 Fab Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	214	Total	C	N	O	S	0	0
			1634	1025	271	332	6		

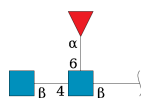
- Molecule 6 is a protein called S309 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	I	230	Total	C	N	O	S	1	0
			1732	1089	294	341	8		

- Molecule 7 is a protein called S309 Fab Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	J	214	Total	C	N	O	S	1	0
			1633	1016	278	333	6		

- Molecule 8 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.

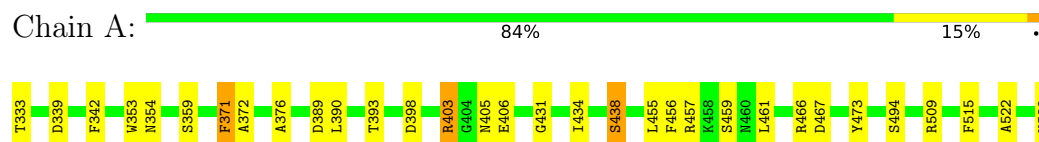


Mol	Chain	Residues	Atoms				AltConf	Trace
8	B	3	Total	C	N	O	0	0
			38	22	2	14		

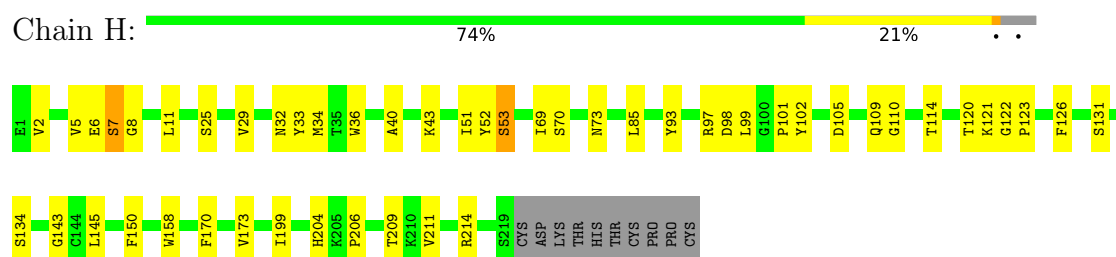
3 Residue-property plots

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.

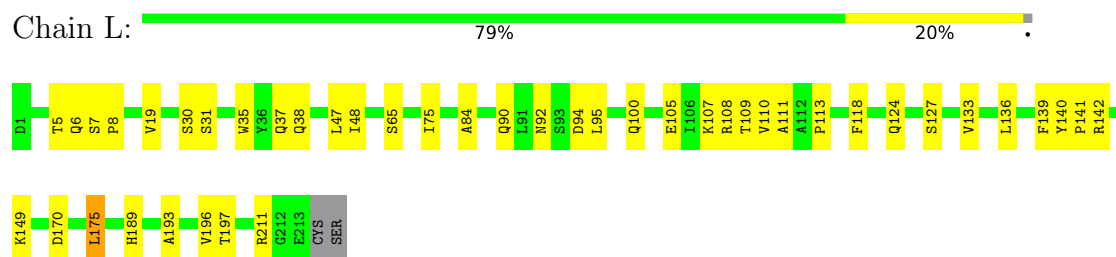
- Molecule 1: Spike protein S2'



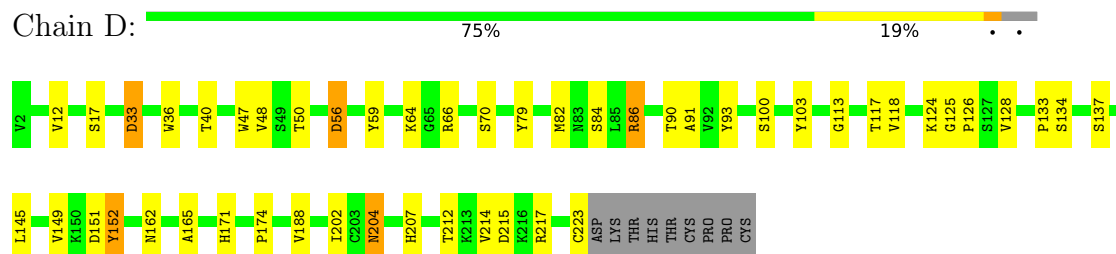
- Molecule 2: BD-604 Fab Heavy chain




- Molecule 3: BD-604 Fab Light chain

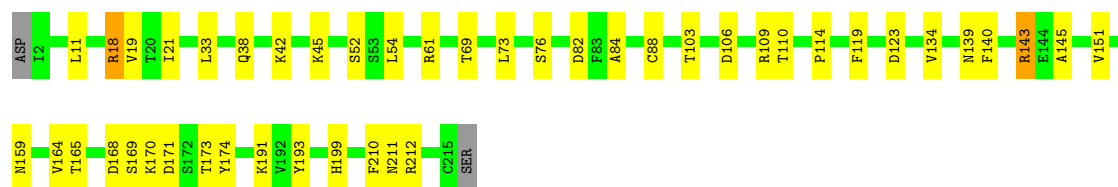


- Molecule 4: S304 Fab Heavy Chain




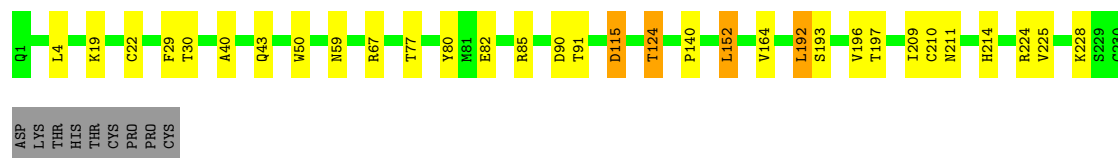
- Molecule 5: S304 Fab Light Chain

Chain E:  78% 20% ..



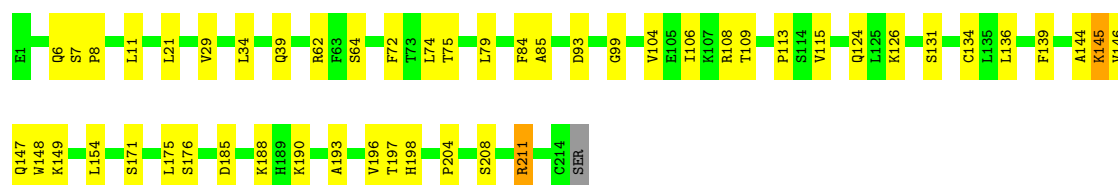
• Molecule 6: S309 Fab Heavy Chain

Chain I:  83% 12% . .



• Molecule 7: S309 Fab Light Chain

Chain J:  76% 22% .



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

Chain B:  67% 33%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	662631	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.31	0/1610	0.49	0/2191
2	H	0.28	0/1653	0.51	0/2252
3	L	0.27	0/1665	0.48	0/2261
4	D	0.28	0/1704	0.51	0/2321
5	E	0.28	0/1670	0.47	0/2269
6	I	0.30	0/1778	0.52	1/2422 (0.0%)
7	J	0.28	0/1669	0.52	0/2266
All	All	0.29	0/11749	0.50	1/15982 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	I	192	LEU	CA-CB-CG	5.04	126.89	115.30

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	1563	0	1489	17	0
2	H	1617	0	1596	27	0
3	L	1632	0	1586	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1662	0	1601	26	0
5	E	1634	0	1588	26	0
6	I	1732	0	1687	19	0
7	J	1633	0	1590	32	0
8	B	38	0	34	1	0
All	All	11511	0	11171	166	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

The worst 5 of 166 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)
4:D:90:THR:HG23	4:D:117:THR:HA	1.61	0.81
3:L:30:SER:OG	3:L:31:SER:N	2.17	0.75
3:L:136:LEU:HD11	3:L:196:VAL:HG21	1.73	0.70
2:H:5:VAL:HG13	2:H:109:GLN:HE22	1.57	0.68
6:I:209:ILE:HG12	6:I:224:ARG:HG2	1.75	0.68

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	194/196 (99%)	185 (95%)	9 (5%)	0	100	100
2	H	217/229 (95%)	207 (95%)	10 (5%)	0	100	100
3	L	211/215 (98%)	202 (96%)	9 (4%)	0	100	100
4	D	220/231 (95%)	214 (97%)	6 (3%)	0	100	100
5	E	212/216 (98%)	208 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	I	229/239 (96%)	225 (98%)	4 (2%)	0	100	100
7	J	213/215 (99%)	204 (96%)	9 (4%)	0	100	100
All	All	1496/1541 (97%)	1445 (97%)	51 (3%)	0	100	100

There are no Ramachandran outliers to report.

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 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	168/168 (100%)	159 (95%)	9 (5%)	18	21
2	H	183/193 (95%)	174 (95%)	9 (5%)	21	25
3	L	186/188 (99%)	181 (97%)	5 (3%)	40	50
4	D	185/194 (95%)	175 (95%)	10 (5%)	18	21
5	E	187/189 (99%)	180 (96%)	7 (4%)	29	38
6	I	193/201 (96%)	184 (95%)	9 (5%)	22	27
7	J	186/186 (100%)	182 (98%)	4 (2%)	47	59
All	All	1288/1319 (98%)	1235 (96%)	53 (4%)	28	33

5 of 53 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	D	100	SER
5	E	45	LYS
7	J	62	ARG
4	D	152	TYR
4	D	223	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
3	L	100	GLN
7	J	6	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 ⓘ

3 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	$\# Z > 2$
8	NAG	B	1	8,1	14,14,15	0.33	0	17,19,21	0.48	0
8	NAG	B	2	8	14,14,15	0.25	0	17,19,21	0.45	0
8	FUC	B	3	8	10,10,11	0.52	0	14,14,16	0.80	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
8	NAG	B	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	B	2	8	-	2/6/23/26	0/1/1/1
8	FUC	B	3	8	-	-	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

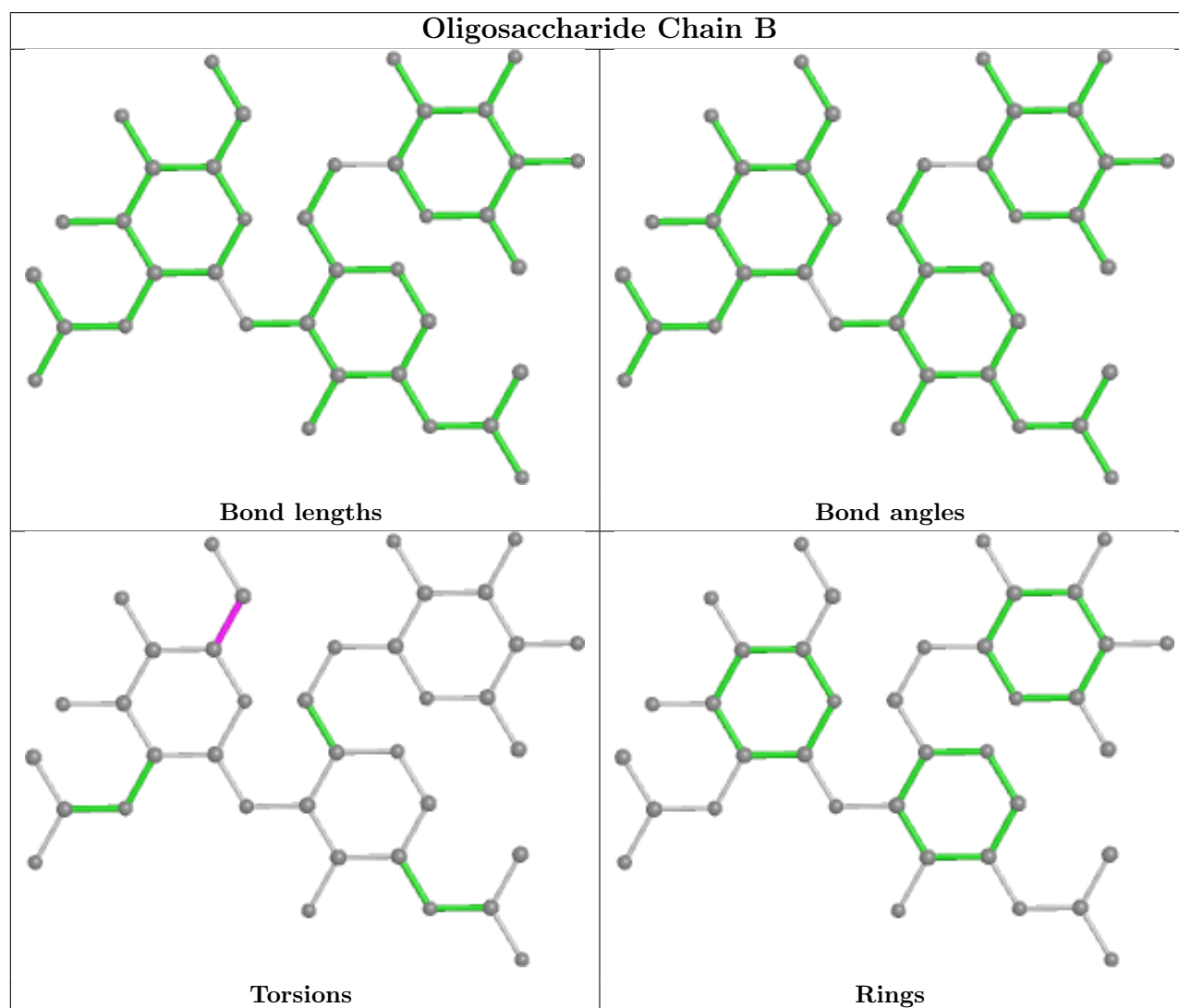
Mol	Chain	Res	Type	Atoms
8	B	2	NAG	O5-C5-C6-O6
8	B	2	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	1	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

There are no ligands in this entry.

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.