



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2021 – 10:41 am BST

PDB ID : 7BBC
Title : Joint X-ray/neutron room temperature structure of perdeuterated PLL lectin in complex with perdeuterated L-fucose
Authors : Gajdos, L.; Blakeley, M.P.; Kumar, A.; Wimmerova, M.; Haertlein, M.; Forsyth, V.T.; Imberty, A.; Devos, J.M.
Deposited on : 2020-12-17
Resolution : 1.84 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : **FAILED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.1

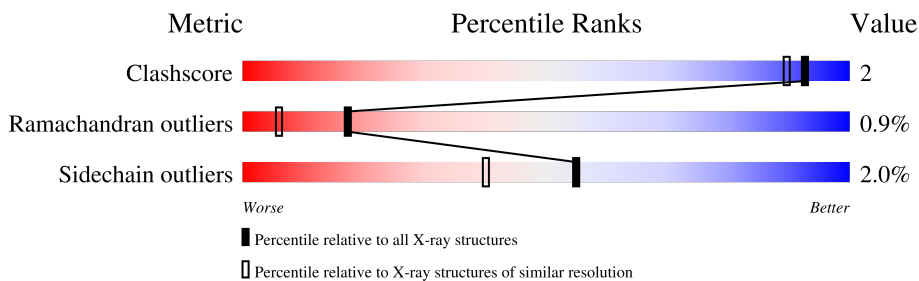
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION, NEUTRON DIFFRACTION

The reported resolution of this entry is 1.84 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	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)

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 failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	381	 87% 6% • 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6909 atoms, of which 0 are hydrogens and 3570 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.

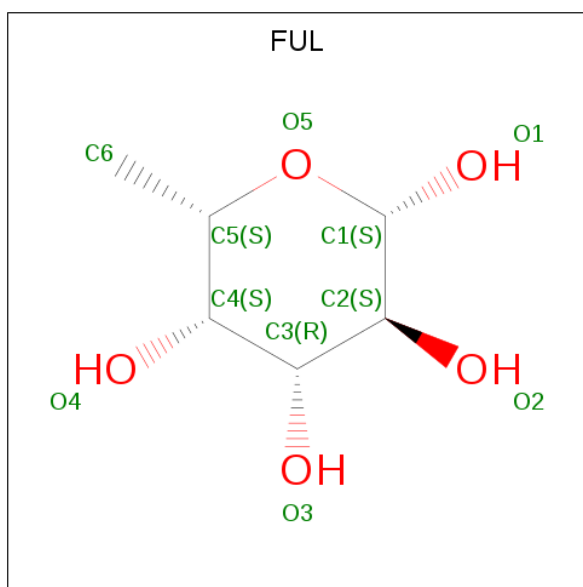
- Molecule 1 is a protein called PLL lectin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	D	N	O	S			
1	A	354	5769	1846	2830	521	566	6	0	42	0

There are 15 discrepancies between the modelled and reference sequences:

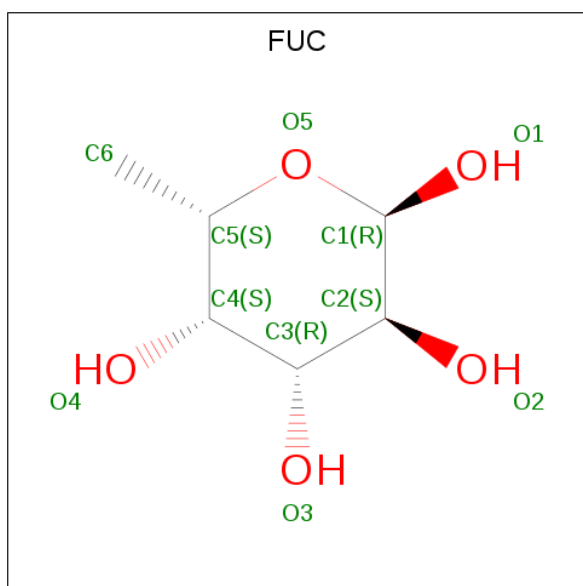
Chain	Residue	Modelled	Actual	Comment	Reference
A	10	HIS	TYR	conflict	UNP A0A329WTS5
A	139	VAL	ALA	conflict	UNP A0A329WTS5
A	369	LEU	-	expression tag	UNP A0A329WTS5
A	370	GLU	-	expression tag	UNP A0A329WTS5
A	371	HIS	-	expression tag	UNP A0A329WTS5
A	372	HIS	-	expression tag	UNP A0A329WTS5
A	373	HIS	-	expression tag	UNP A0A329WTS5
A	374	HIS	-	expression tag	UNP A0A329WTS5
A	375	HIS	-	expression tag	UNP A0A329WTS5
A	376	HIS	-	expression tag	UNP A0A329WTS5
A	377	TRP	-	expression tag	UNP A0A329WTS5
A	378	ARG	-	expression tag	UNP A0A329WTS5
A	379	SER	-	expression tag	UNP A0A329WTS5
A	380	GLY	-	expression tag	UNP A0A329WTS5
A	381	CYS	-	expression tag	UNP A0A329WTS5

- Molecule 2 is beta-L-fucopyranose (three-letter code: FUL) (formula: C₆H₁₂O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	D	O		
2	A	1	23	6	12	5	0	1
2	A	1	23	6	12	5	0	1
2	A	1	23	6	12	5	0	1

- Molecule 3 is alpha-L-fucopyranose (three-letter code: FUC) (formula: C₆H₁₂O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	D	O	0	1
			23	6	12	5		
3	A	1	Total	C	D	O	0	1
			23	6	12	5		
3	A	1	Total	C	D	O	0	1
			23	6	12	5		

- Molecule 4 is water.


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	334	Total	D	O	0	0
			1002	668	334		

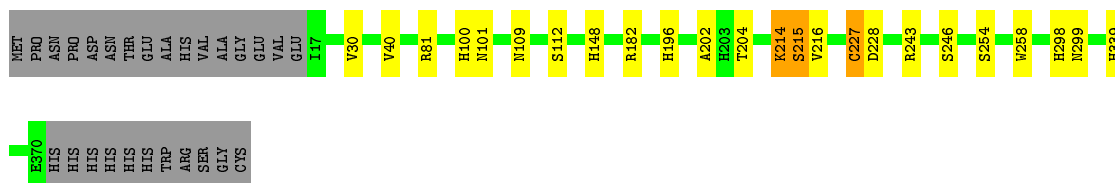
3 Residue-property plots [i](#)

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 failed to run properly.

- Molecule 1: PLL lectin

Chain A:  87% 6% • 7%



4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	72.67Å 89.21Å 159.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.00 – 1.84	Depositor
% Data completeness (in resolution range)	100.0 (33.00-1.84)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 1.84Å)	Xtrriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.121 , 0.140	Depositor
Wilson B-factor (Å ²)	19.1	Xtrriage
Anisotropy	0.186	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	6909	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.66	1/3203 (0.0%)	0.74	3/4391 (0.1%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	246	SER	CB-OG	-5.61	1.34	1.42

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	81	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	A	182	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	243	ARG	NE-CZ-NH2	-5.59	117.50	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	299	ASN	Peptide

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	A	5769	0	2683	9	0
2	A	69	0	35	0	0
3	A	69	0	34	0	0
4	A	1002	0	0	2	0
All	All	6909	0	2752	9	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 2.

All (9) 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:215[D]:SER:OG	1:A:216[D]:VAL:N	2.39	0.55
1:A:101[B]:ASN:OD1	1:A:112[B]:SER:O	2.25	0.55
1:A:202:ALA:O	1:A:204[B]:THR:HG23	2.05	0.52
1:A:214[D]:LYS:O	1:A:215[D]:SER:O	2.33	0.47
1:A:215[B]:SER:HA	4:A:600:HOH:O	2.10	0.47
1:A:227[B]:CYS:SG	1:A:298:HIS:NE2	2.83	0.44
1:A:30[B]:VAL:HG22	1:A:40:VAL:HG12	1.92	0.41
1:A:215[C]:SER:HB3	1:A:258:TRP:CE2	2.50	0.41
1:A:109:ASN:HB3	4:A:537:HOH:O	2.15	0.41

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	402/381 (106%)	384 (96%)	9 (2%)	9 (2%)	6 1

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	215[B]	SER
1	A	215[C]	SER
1	A	215[D]	SER
1	A	214[B]	LYS
1	A	214[C]	LYS
1	A	214[D]	LYS
1	A	228[A]	ASP
1	A	228[B]	ASP
1	A	228[C]	ASP

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	338/316 (107%)	329 (97%)	9 (3%)	44 28

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	HIS
1	A	148	HIS
1	A	196	HIS
1	A	227[A]	CYS
1	A	227[B]	CYS
1	A	227[C]	CYS
1	A	254[A]	SER
1	A	254[B]	SER
1	A	339	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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
3	FUC	A	406[A]	-	11,11,11	1.34	1 (9%)	15,16,16	1.15	1 (6%)
3	FUC	A	404[B]	-	11,11,11	1.26	1 (9%)	15,16,16	0.65	0
3	FUC	A	402[B]	-	11,11,11	1.20	1 (9%)	15,16,16	0.89	1 (6%)
2	FUL	A	405[B]	-	11,11,11	1.57	2 (18%)	15,16,16	1.65	3 (20%)
2	FUL	A	401[A]	-	11,11,11	1.17	1 (9%)	15,16,16	1.06	1 (6%)
2	FUL	A	403[A]	-	11,11,11	1.55	2 (18%)	15,16,16	1.21	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
3	FUC	A	406[A]	-	-	-	0/1/1/1
3	FUC	A	404[B]	-	-	-	0/1/1/1
3	FUC	A	402[B]	-	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FUL	A	405[B]	-	-	-	0/1/1/1
2	FUL	A	401[A]	-	-	-	0/1/1/1
2	FUL	A	403[A]	-	-	-	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	403[A]	FUL	O5-C1	3.53	1.51	1.42
3	A	406[A]	FUC	O5-C1	3.30	1.51	1.42
2	A	405[B]	FUL	O5-C1	3.24	1.51	1.42
3	A	404[B]	FUC	O5-C1	3.20	1.50	1.42
3	A	402[B]	FUC	O5-C1	3.13	1.50	1.42
2	A	401[A]	FUL	O5-C1	3.01	1.50	1.42
2	A	405[B]	FUL	O3-C3	2.10	1.47	1.43
2	A	403[A]	FUL	O2-C2	2.04	1.47	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	405[B]	FUL	O4-C4-C5	4.36	119.33	109.67
2	A	405[B]	FUL	O5-C5-C4	2.50	114.00	109.52
2	A	405[B]	FUL	O4-C4-C3	-2.31	105.00	110.35
2	A	401[A]	FUL	C1-C2-C3	-2.07	106.03	110.31
3	A	402[B]	FUC	C4-C3-C2	2.06	114.42	110.82
3	A	406[A]	FUC	O4-C4-C5	2.06	114.23	109.67

There are no chirality outliers.

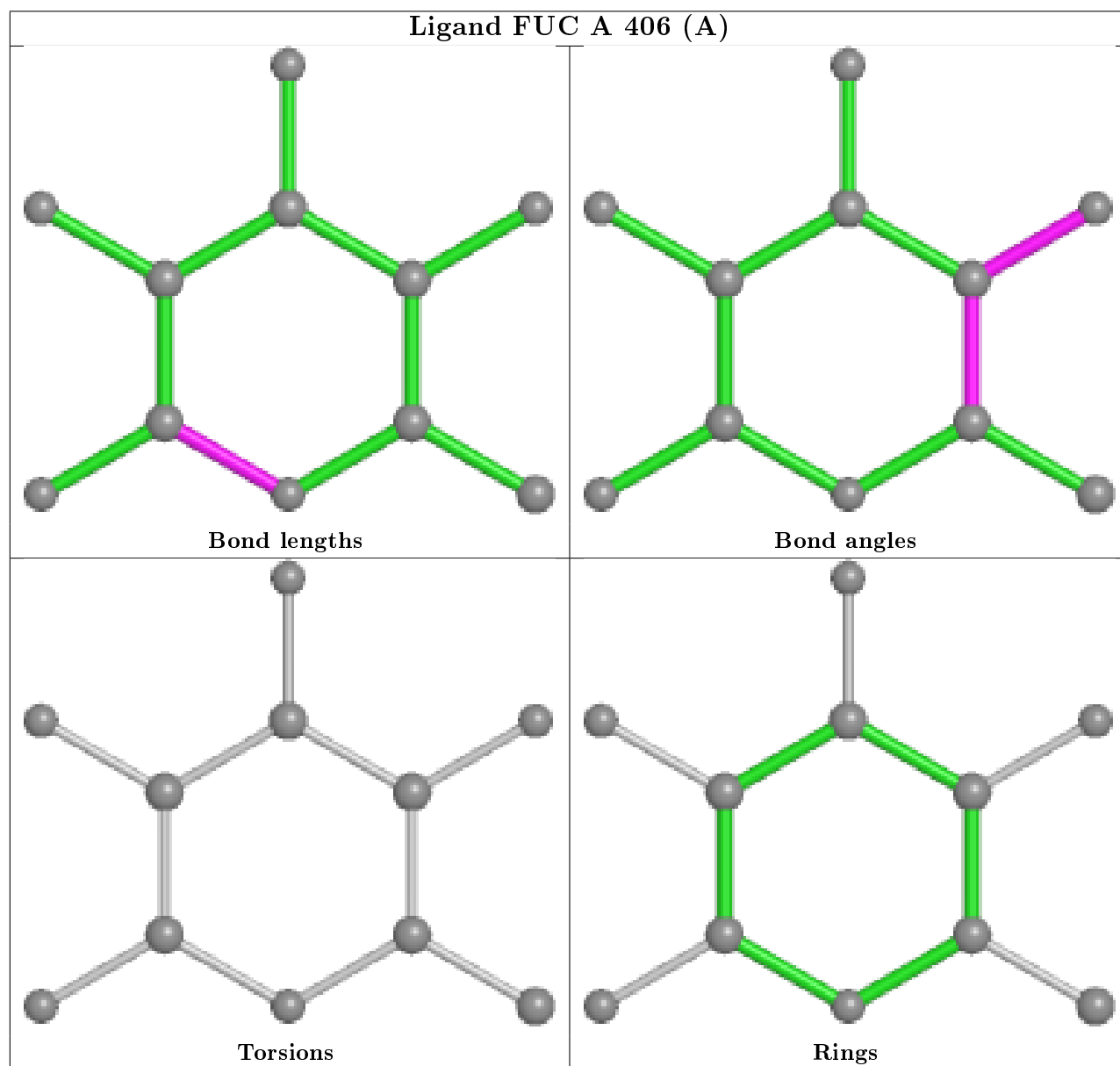
There are no torsion outliers.

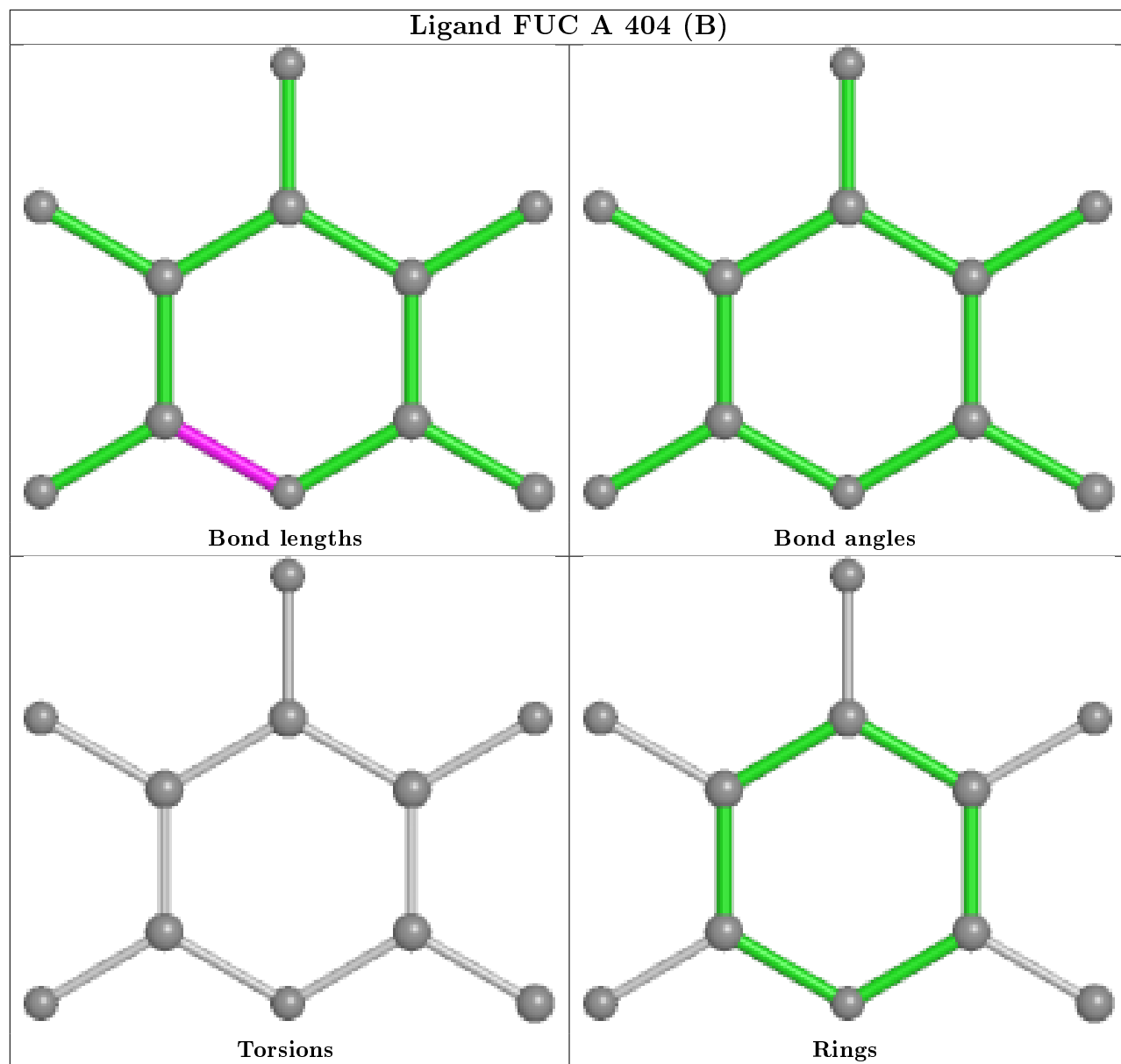
There are no ring outliers.

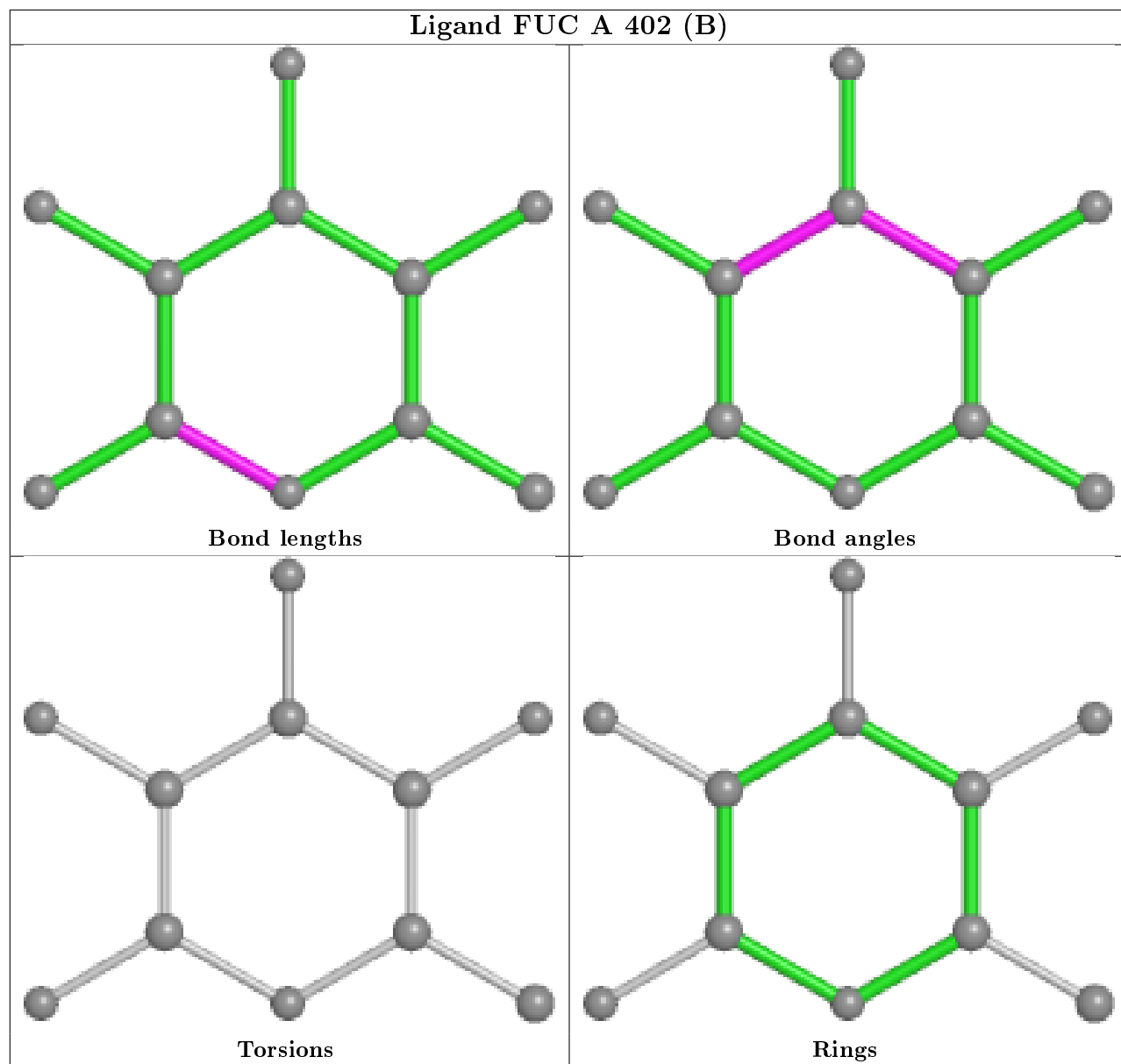
No monomer is involved in short contacts.

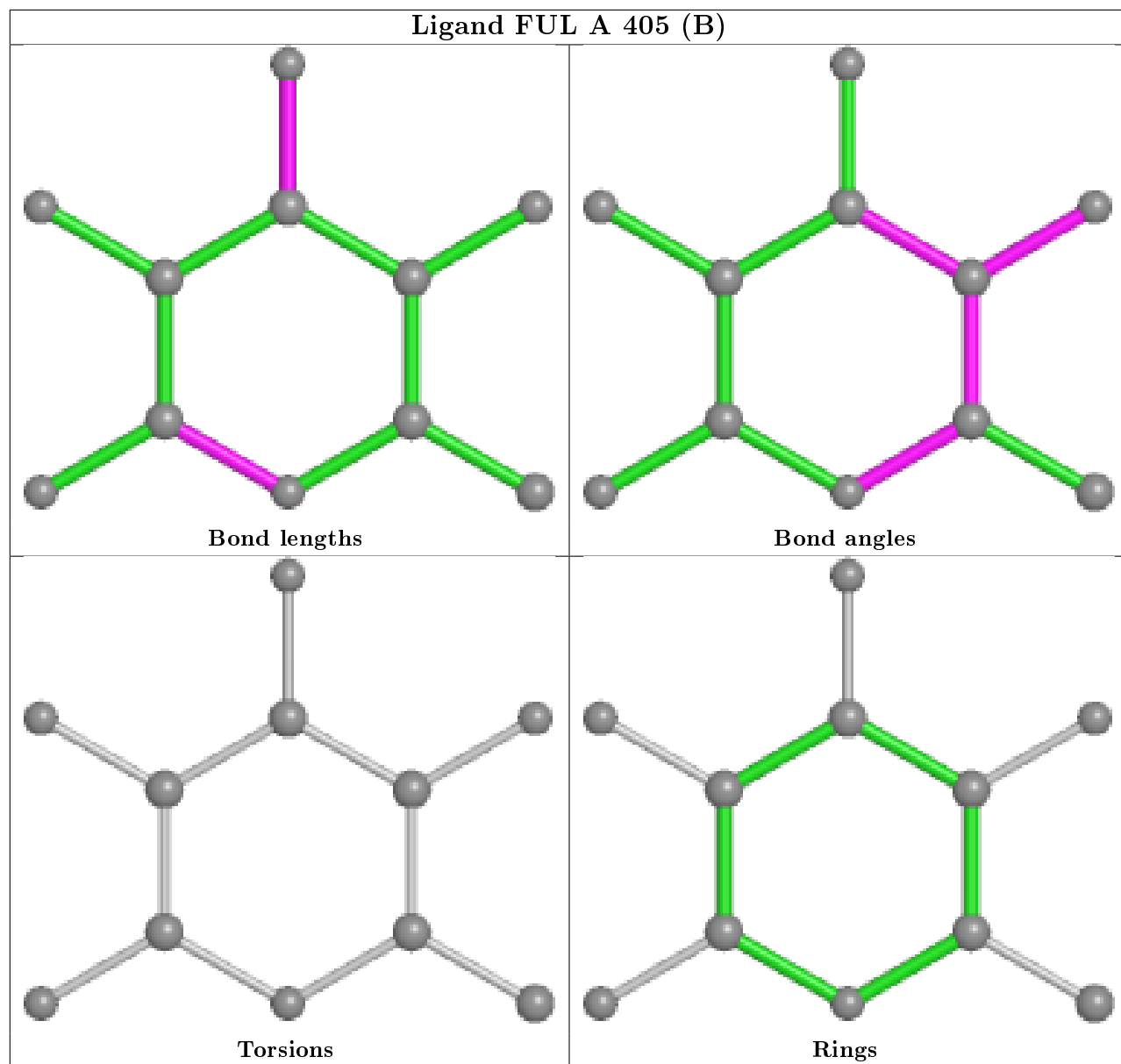
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

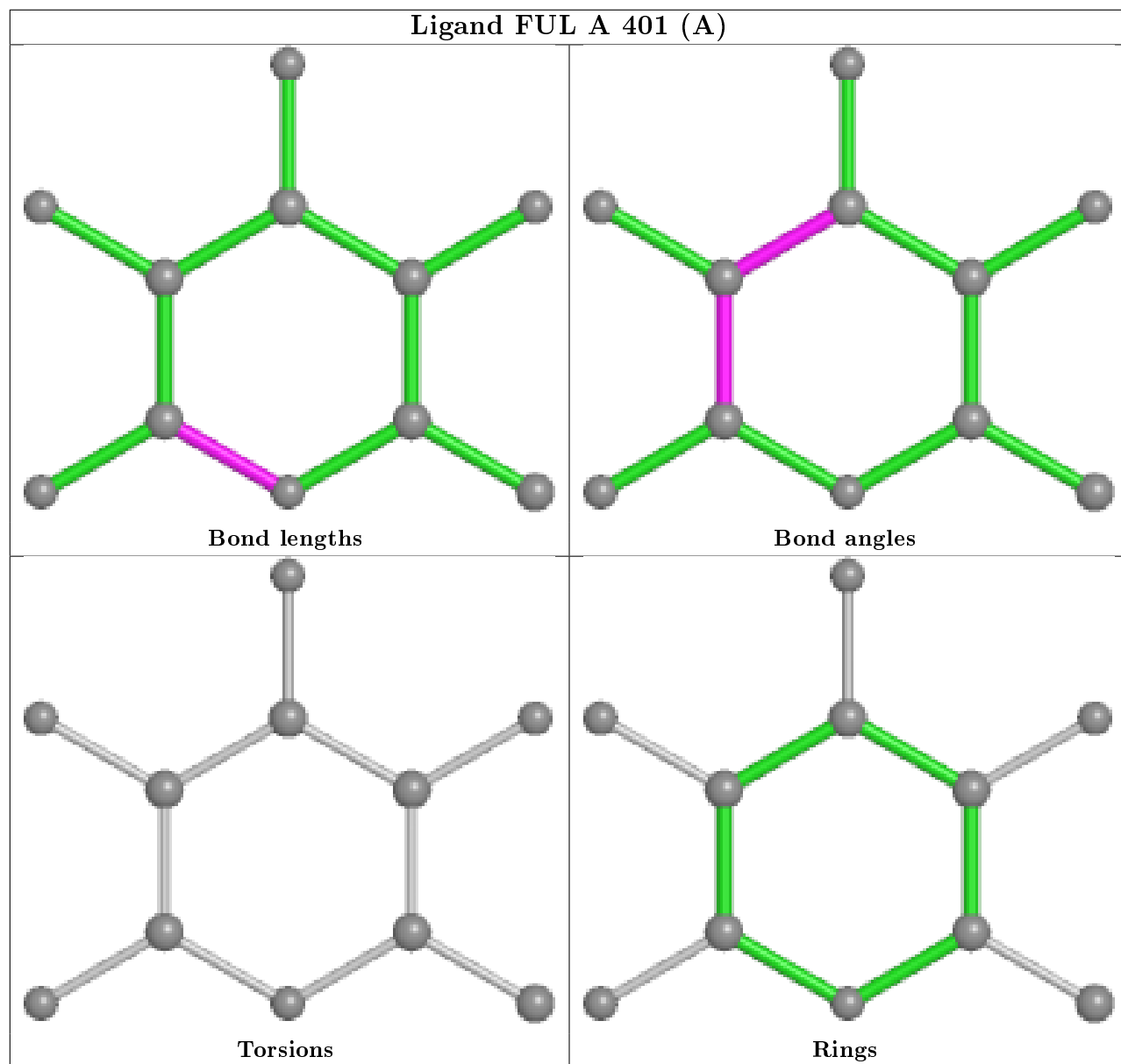
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

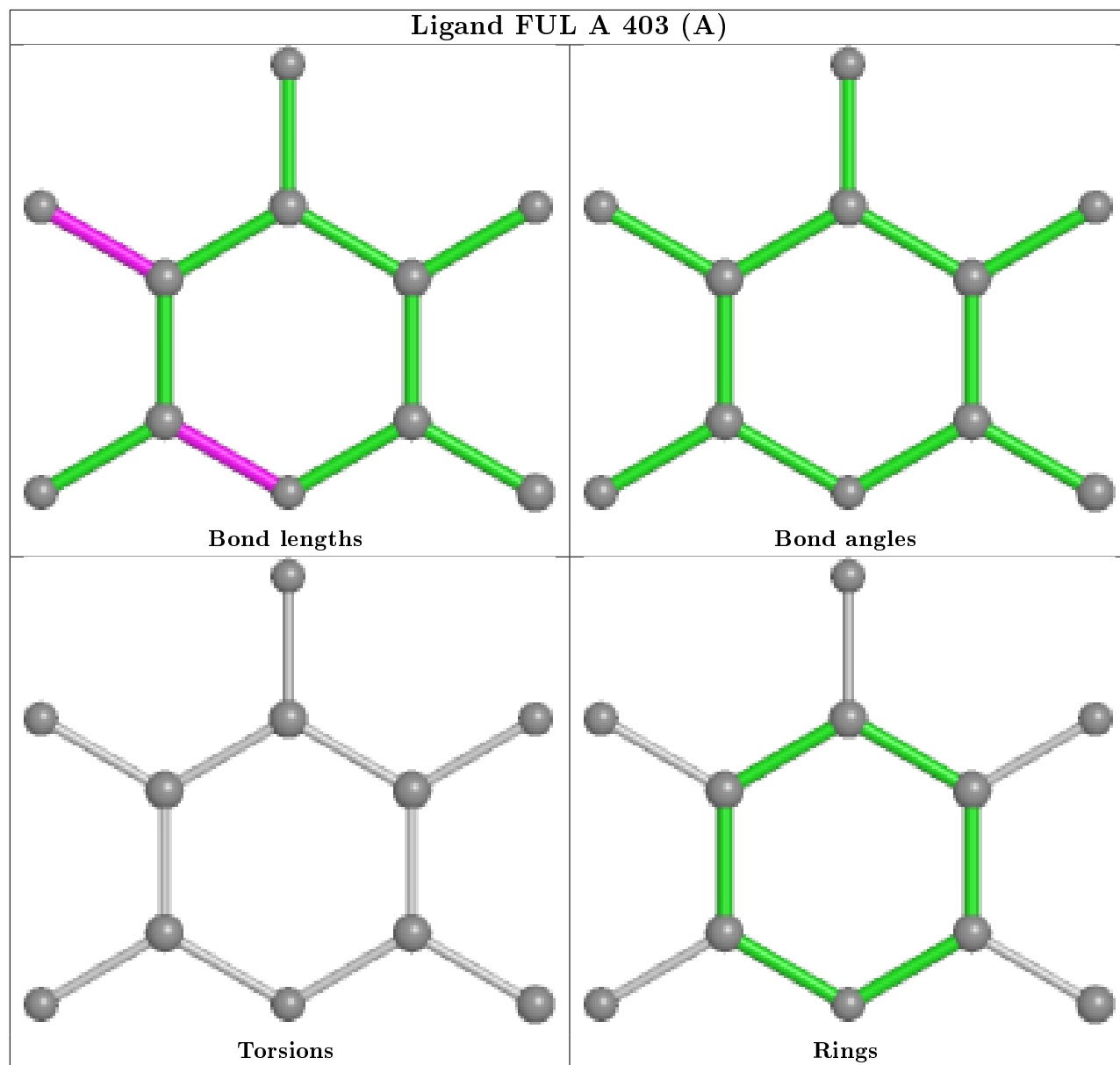












5.7 Other polymers [i](#)

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 failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

6.4 Ligands

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers

EDS failed to run properly - this section is therefore empty.