



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 14, 2024 – 02:24 am GMT

PDB ID : 6TRV
Title : Structure of SapL1 lectin in complex with alpha methyl fucoside
Authors : Martinez Alarcon, D.; Varrot, A.
Deposited on : 2019-12-19
Resolution : 2.40 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

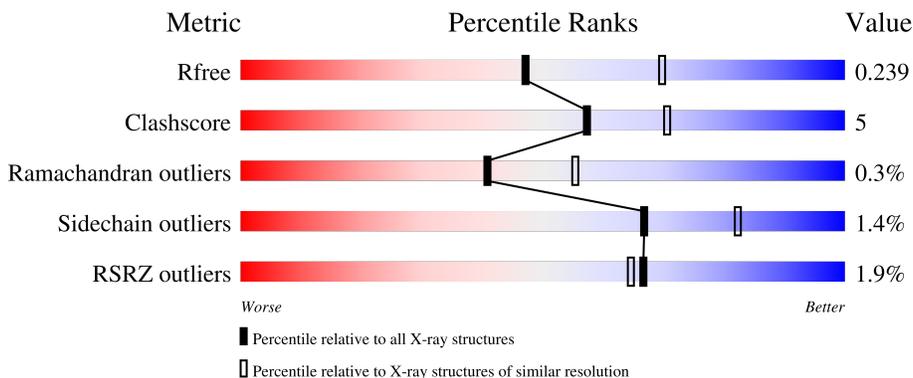
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AAA	297	 89% 10% .
1	BBB	297	 3% 83% 16% .

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4777 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.

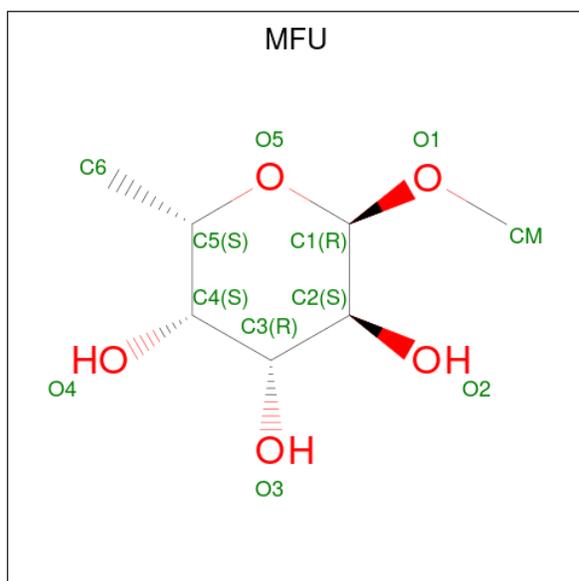
- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	AAA	294	2222	1416	386	419	1	0	0	0
1	BBB	294	2220	1415	385	419	1	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

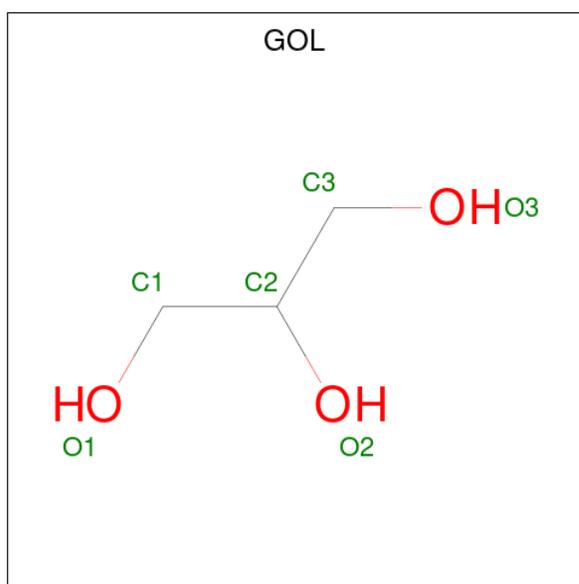
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-1	GLY	-	expression tag	UNP A0A084FYP2
AAA	0	HIS	-	expression tag	UNP A0A084FYP2
BBB	-1	GLY	-	expression tag	UNP A0A084FYP2
BBB	0	HIS	-	expression tag	UNP A0A084FYP2

- Molecule 2 is methyl alpha-L-fucopyranoside (three-letter code: MFU) (formula: C₇H₁₄O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	AAA	1	Total C O 12 7 5	0	0
2	AAA	1	Total C O 12 7 5	0	0
2	AAA	1	Total C O 12 7 5	0	0
2	AAA	1	Total C O 12 7 5	0	0
2	AAA	1	Total C O 12 7 5	0	0
2	BBB	1	Total C O 12 7 5	0	0
2	BBB	1	Total C O 12 7 5	0	0
2	BBB	1	Total C O 12 7 5	0	0
2	BBB	1	Total C O 12 7 5	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



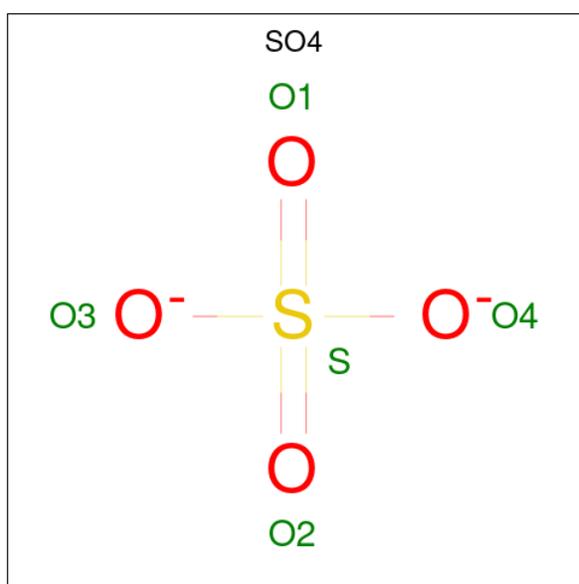
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	AAA	1	Total C O 6 3 3	0	0
3	AAA	1	Total C O 6 3 3	0	0
3	AAA	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	AAA	1	Total C O 6 3 3	0	0
3	BBB	1	Total C O 6 3 3	0	0
3	BBB	1	Total C O 6 3 3	0	0
3	BBB	1	Total C O 6 3 3	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	AAA	1	Total O S 5 4 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	AAA	111	Total O 111 111	0	2
5	BBB	68	Total O 69 69	0	1

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 and electron 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

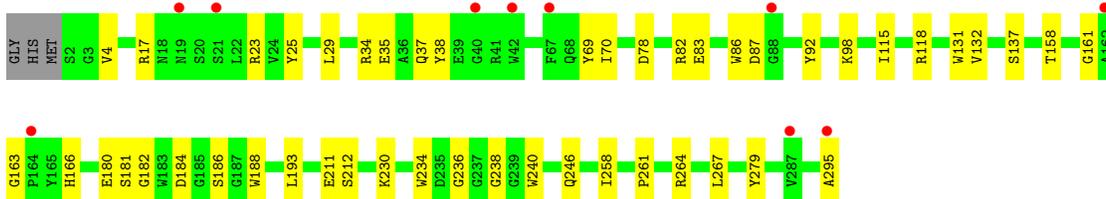
- Molecule 1: Uncharacterized protein

Chain AAA: 



- Molecule 1: Uncharacterized protein

Chain BBB: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.06Å 45.66Å 83.48Å 90.00° 105.05° 90.00°	Depositor
Resolution (Å)	39.99 – 2.40 38.78 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.9 (39.99-2.40) 97.9 (38.78-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.175 , 0.240 0.183 , 0.239	Depositor DCC
R_{free} test set	1089 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtrriage
Anisotropy	0.183	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 33.7	EDS
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
F_o, F_c correlation	0.96	EDS
Total number of atoms	4777	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: MFU, SO4, GOL

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	AAA	0.93	4/2287 (0.2%)	1.08	1/3125 (0.0%)
1	BBB	0.91	1/2285 (0.0%)	1.03	0/3123
All	All	0.92	5/4572 (0.1%)	1.05	1/6248 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	BBB	180	GLU	CD-OE1	-9.09	1.15	1.25
1	AAA	114	ASN	C-O	6.18	1.35	1.23
1	AAA	180	GLU	CD-OE2	-5.44	1.19	1.25
1	AAA	181	SER	CA-CB	-5.34	1.45	1.52
1	AAA	111	LEU	C-O	5.05	1.32	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	17	ARG	NE-CZ-NH1	5.38	122.99	120.30

There are no chirality outliers.

There are no planarity outliers.

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	AAA	2222	0	2107	15	0
1	BBB	2220	0	2100	27	0
2	AAA	60	0	67	2	0
2	BBB	48	0	56	3	0
3	AAA	24	0	25	0	0
3	BBB	18	0	24	2	0
4	AAA	5	0	0	0	0
5	AAA	111	0	0	2	0
5	BBB	69	0	0	0	0
All	All	4777	0	4379	44	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.

The worst 5 of 44 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)
2:AAA:301:MFU:HM3	2:AAA:301:MFU:O2	1.75	0.86
1:BBB:158:THR:OG1	3:BBB:307:GOL:H2	1.83	0.78
1:AAA:213:GLY:O	1:AAA:236:GLY:HA2	1.93	0.68
1:AAA:115:ILE:N	1:AAA:115:ILE:HD13	2.11	0.65
1:AAA:193:LEU:HD23	1:AAA:240:TRP:CD2	2.38	0.59

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	AAA	292/297 (98%)	280 (96%)	12 (4%)	0	100	100
1	BBB	292/297 (98%)	277 (95%)	13 (4%)	2 (1%)	22	32
All	All	584/594 (98%)	557 (95%)	25 (4%)	2 (0%)	41	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	236	GLY
1	BBB	163	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	AAA	216/221 (98%)	214 (99%)	2 (1%)	78	90
1	BBB	215/221 (97%)	211 (98%)	4 (2%)	57	75
All	All	431/442 (98%)	425 (99%)	6 (1%)	67	82

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	BBB	181	SER
1	BBB	186	SER
1	BBB	212	SER
1	AAA	176	ARG
1	AAA	64	SER

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](#)

17 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
2	MFU	BBB	301	-	12,12,12	1.35	1 (8%)	17,17,17	1.21	2 (11%)
2	MFU	BBB	302	-	12,12,12	1.31	1 (8%)	17,17,17	1.02	0
3	GOL	BBB	307	-	5,5,5	0.40	0	5,5,5	0.52	0
2	MFU	AAA	302	-	12,12,12	1.28	2 (16%)	17,17,17	1.01	2 (11%)
2	MFU	AAA	301	-	12,12,12	1.03	0	17,17,17	1.03	1 (5%)
3	GOL	AAA	306	-	5,5,5	0.22	0	5,5,5	0.42	0
4	SO4	AAA	310	-	4,4,4	2.44	2 (50%)	6,6,6	0.20	0
3	GOL	AAA	308	-	5,5,5	0.51	0	5,5,5	0.86	0
3	GOL	AAA	307	-	5,5,5	0.49	0	5,5,5	0.31	0
2	MFU	AAA	304	-	12,12,12	1.73	3 (25%)	17,17,17	1.30	2 (11%)
2	MFU	AAA	303	-	12,12,12	0.94	0	17,17,17	1.10	3 (17%)
3	GOL	AAA	309	-	5,5,5	0.42	0	5,5,5	0.78	0
3	GOL	BBB	306	-	5,5,5	0.39	0	5,5,5	0.21	0
2	MFU	BBB	304	-	12,12,12	2.84	3 (25%)	17,17,17	1.16	1 (5%)
2	MFU	AAA	305	-	12,12,12	2.56	5 (41%)	17,17,17	1.18	1 (5%)
2	MFU	BBB	303	-	12,12,12	1.31	1 (8%)	17,17,17	1.51	3 (17%)
3	GOL	BBB	305	-	5,5,5	0.24	0	5,5,5	0.21	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
2	MFU	BBB	301	-	-	0/2/22/22	0/1/1/1
2	MFU	BBB	302	-	-	0/2/22/22	0/1/1/1
3	GOL	BBB	307	-	-	2/4/4/4	-
2	MFU	AAA	302	-	-	0/2/22/22	0/1/1/1
2	MFU	AAA	301	-	-	2/2/22/22	0/1/1/1
3	GOL	AAA	306	-	-	2/4/4/4	-
3	GOL	AAA	308	-	-	4/4/4/4	-
3	GOL	AAA	307	-	-	2/4/4/4	-
2	MFU	AAA	304	-	-	0/2/22/22	0/1/1/1
2	MFU	AAA	303	-	-	0/2/22/22	0/1/1/1
3	GOL	AAA	309	-	-	4/4/4/4	-
3	GOL	BBB	306	-	-	3/4/4/4	-
2	MFU	BBB	304	-	-	0/2/22/22	0/1/1/1
2	MFU	AAA	305	-	-	2/2/22/22	0/1/1/1
2	MFU	BBB	303	-	-	2/2/22/22	0/1/1/1
3	GOL	BBB	305	-	-	1/4/4/4	-

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	BBB	304	MFU	O1-C1	7.92	1.53	1.40
2	AAA	305	MFU	O1-C1	6.00	1.50	1.40
2	AAA	305	MFU	C4-C5	4.12	1.62	1.52
2	BBB	304	MFU	O1-CM	3.61	1.55	1.42
2	AAA	304	MFU	O1-C1	3.49	1.46	1.40

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	303	MFU	CM-O1-C1	-4.58	106.20	113.27
2	AAA	304	MFU	O1-C1-C2	-4.15	103.29	108.15
2	AAA	301	MFU	CM-O1-C1	-3.33	108.14	113.27
2	AAA	305	MFU	C6-C5-C4	2.68	118.02	113.07
2	BBB	303	MFU	O1-C1-C2	-2.61	105.08	108.15

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	301	MFU	C2-C1-O1-CM

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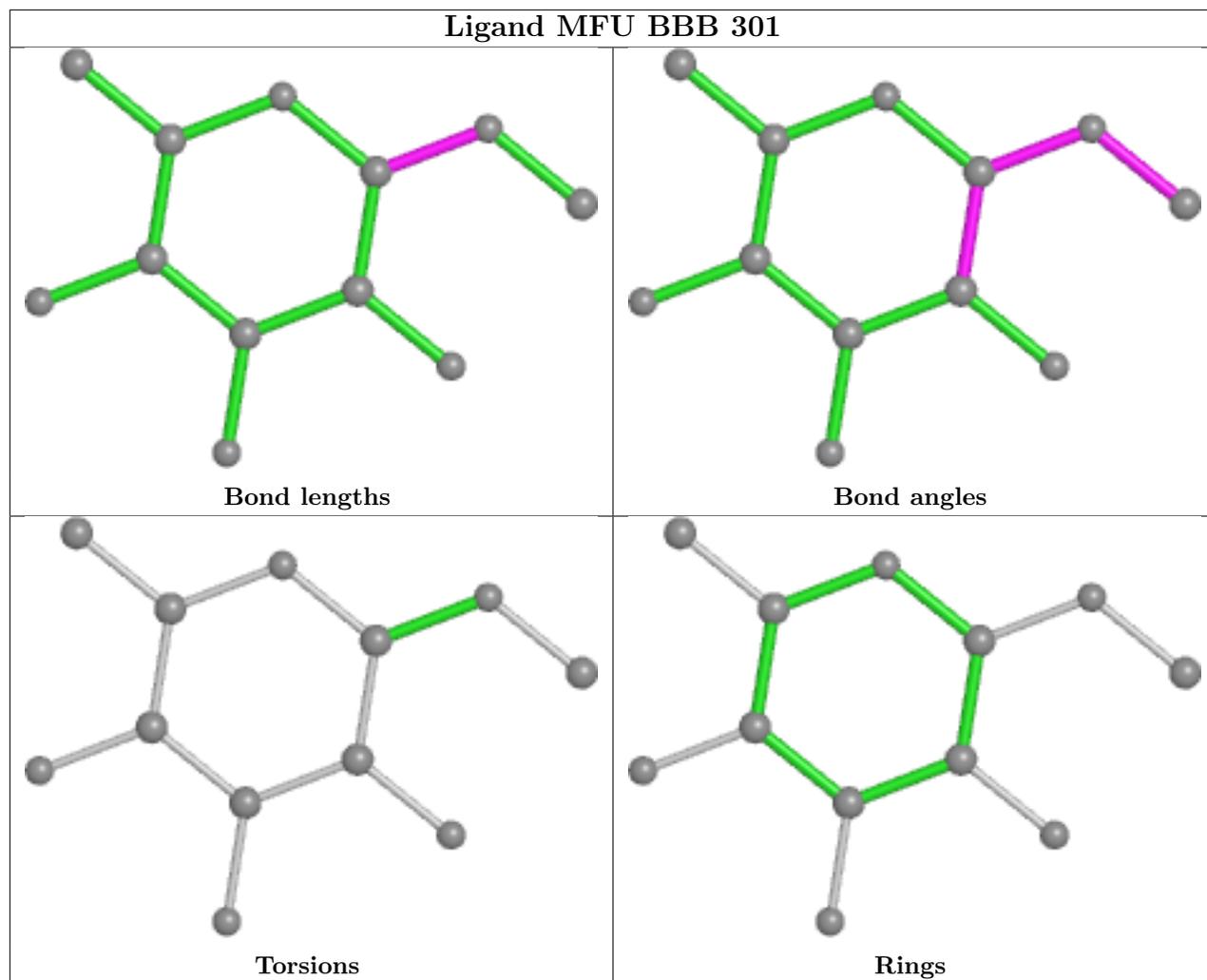
Mol	Chain	Res	Type	Atoms
2	AAA	301	MFU	O5-C1-O1-CM
3	AAA	307	GOL	O1-C1-C2-O2
3	AAA	307	GOL	O1-C1-C2-C3
3	AAA	308	GOL	O1-C1-C2-C3

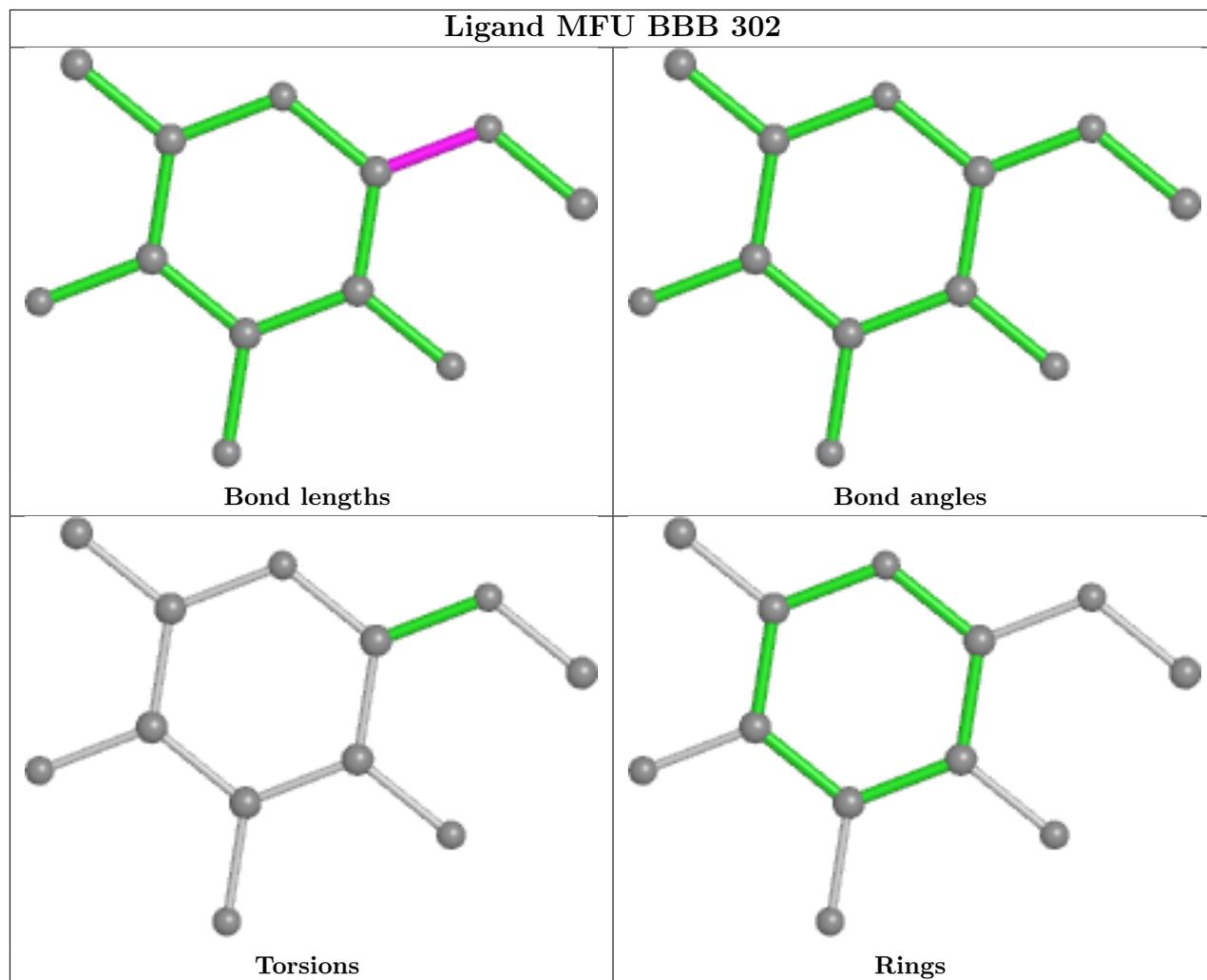
There are no ring outliers.

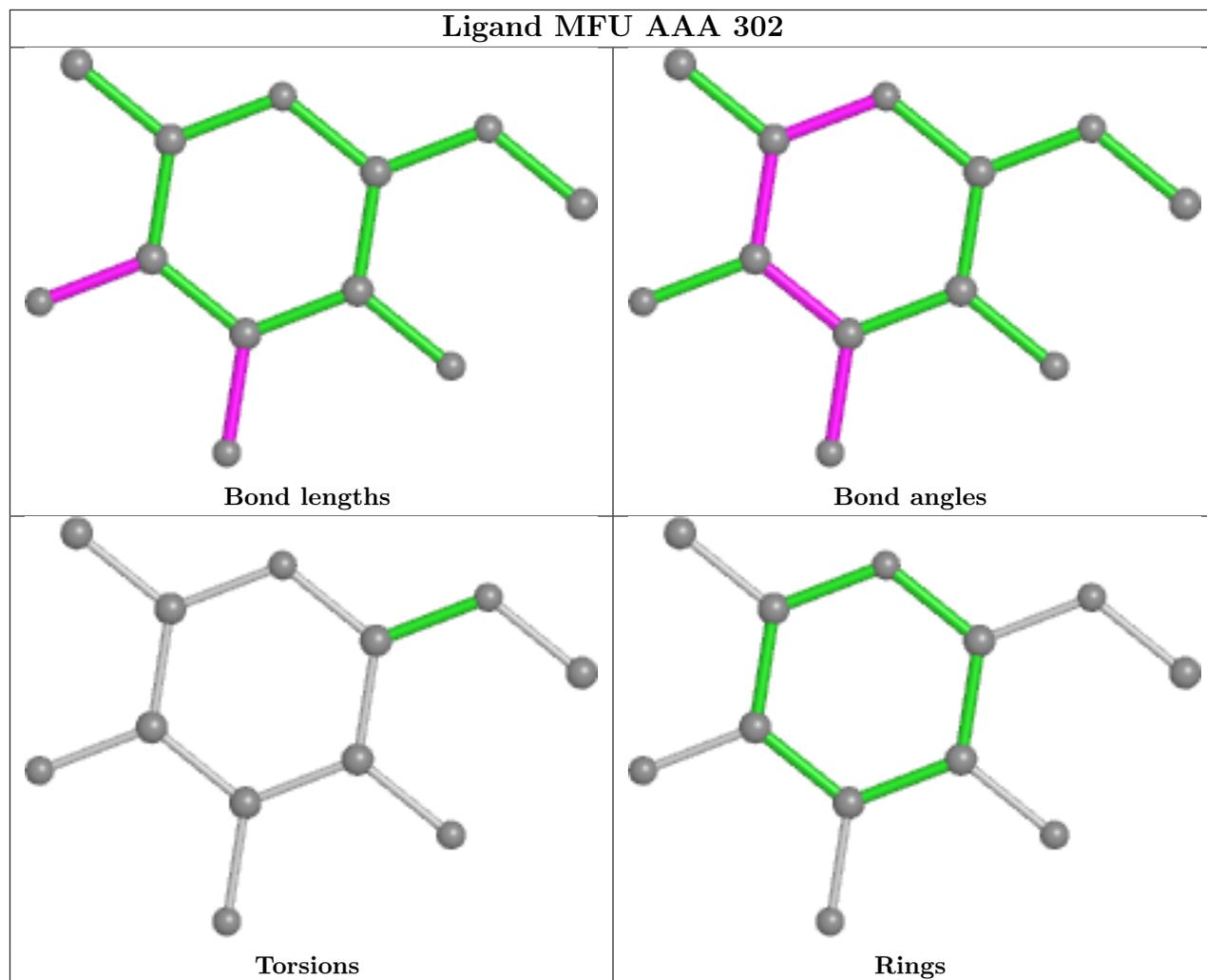
6 monomers are involved in 7 short contacts:

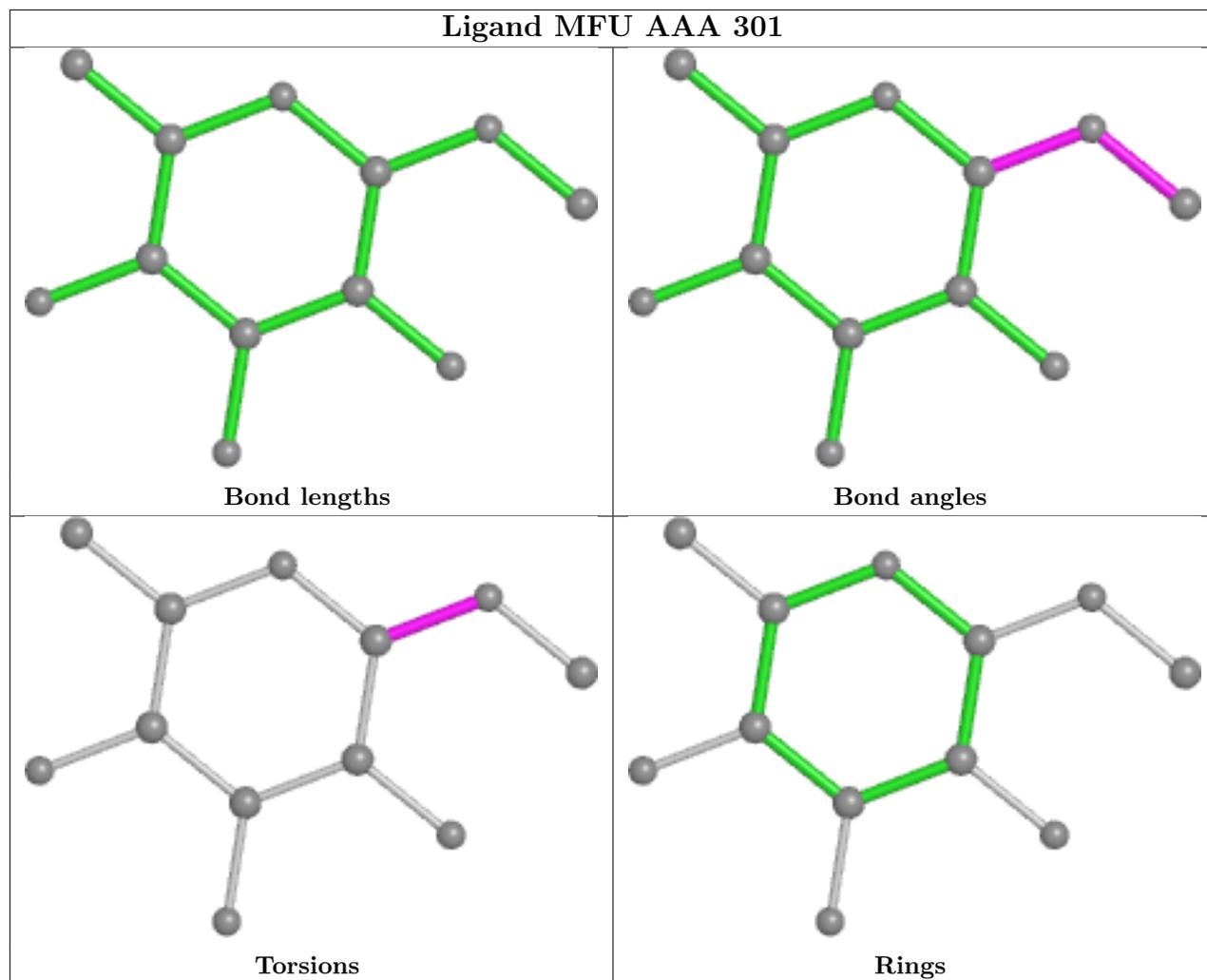
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	BBB	301	MFU	1	0
2	BBB	302	MFU	1	0
3	BBB	307	GOL	1	0
2	AAA	301	MFU	2	0
2	BBB	304	MFU	1	0
3	BBB	305	GOL	1	0

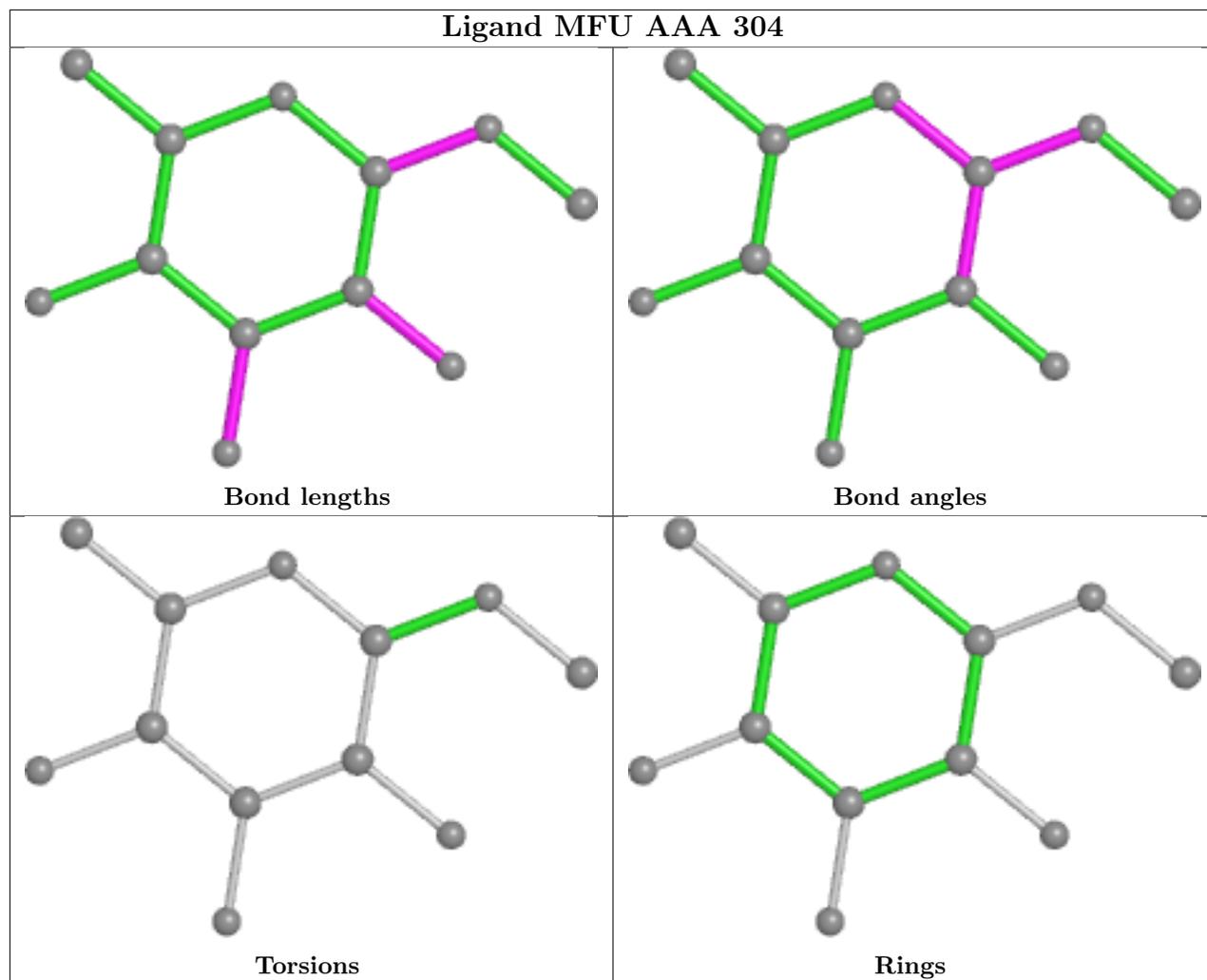
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

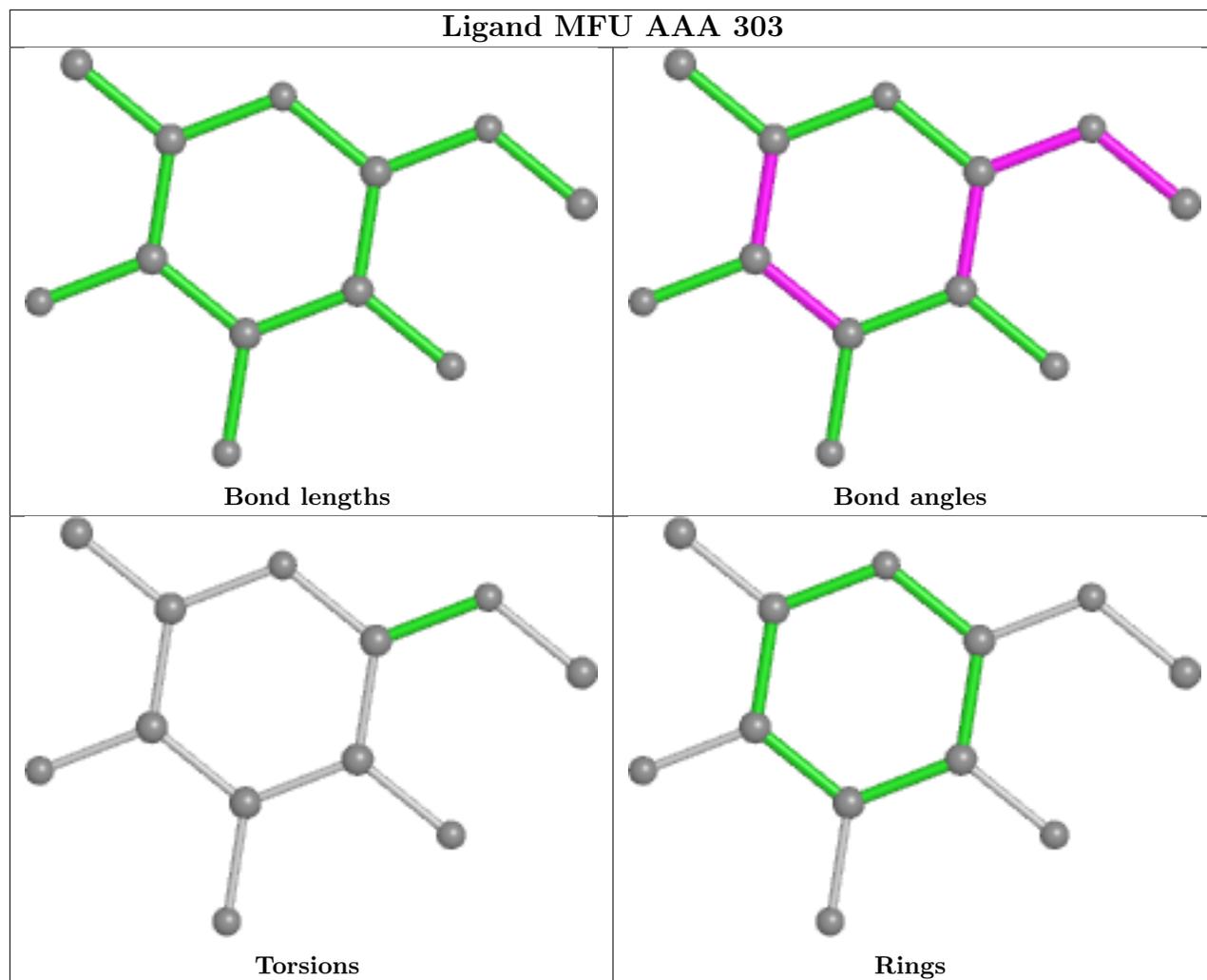


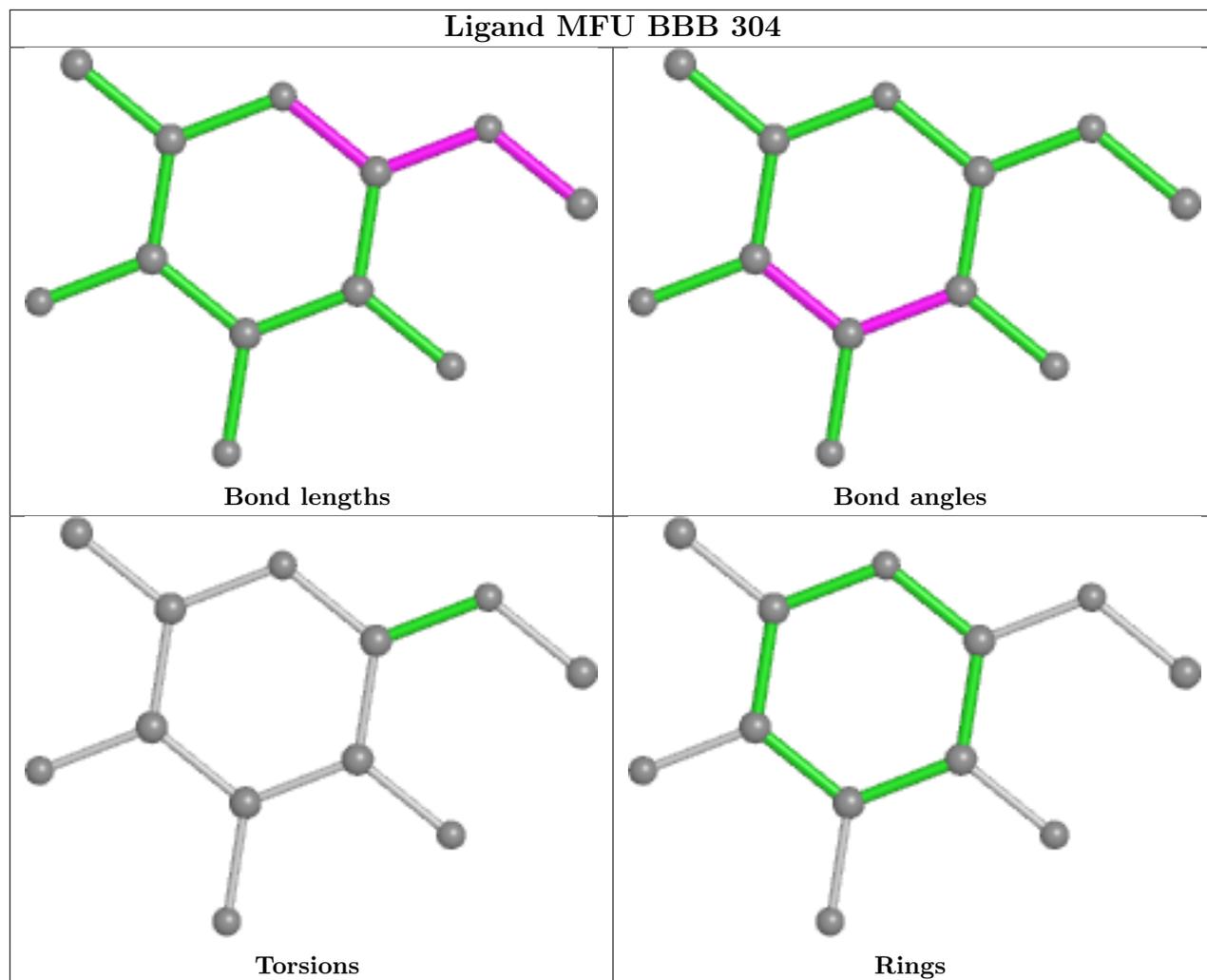


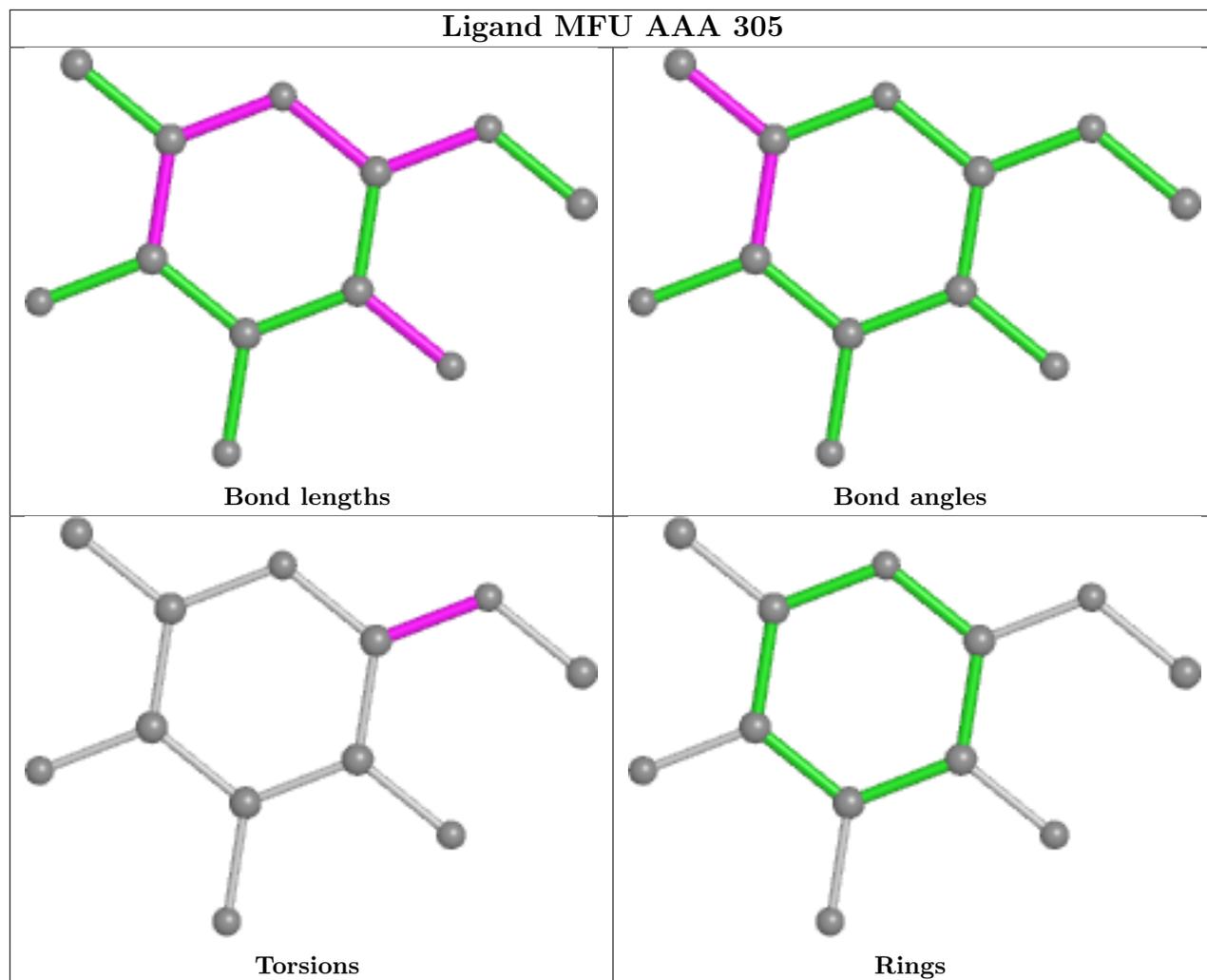


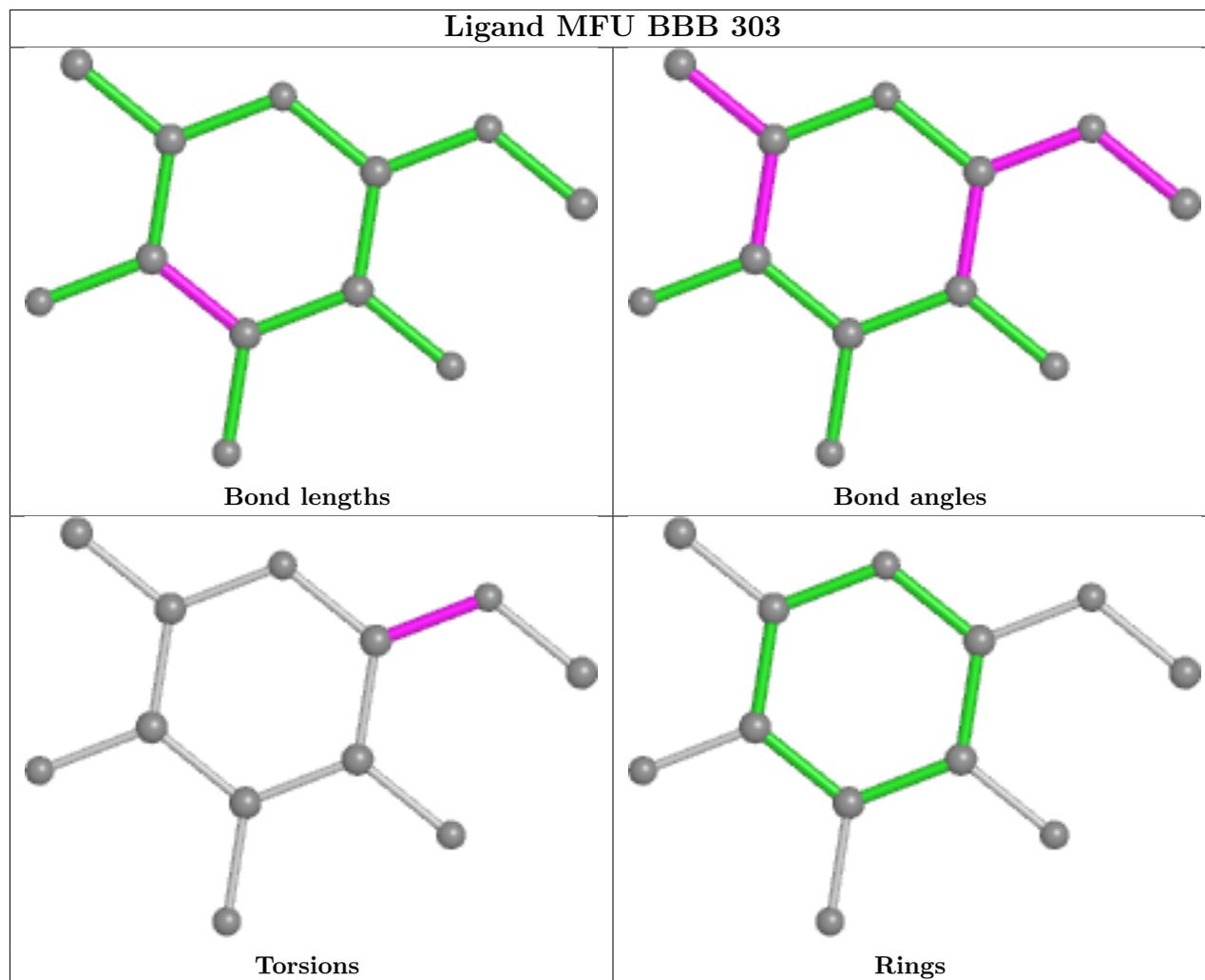












5.7 Other polymers [i](#)

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 [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	AAA	294/297 (98%)	-0.15	1 (0%) 94 93	22, 31, 44, 65	0
1	BBB	294/297 (98%)	0.13	10 (3%) 45 44	29, 39, 56, 74	0
All	All	588/594 (98%)	-0.01	11 (1%) 66 64	22, 35, 53, 74	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	295	ALA	6.0
1	BBB	19	ASN	3.4
1	BBB	88	GLY	3.2
1	BBB	40	GLY	2.9
1	AAA	295	ALA	2.8

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

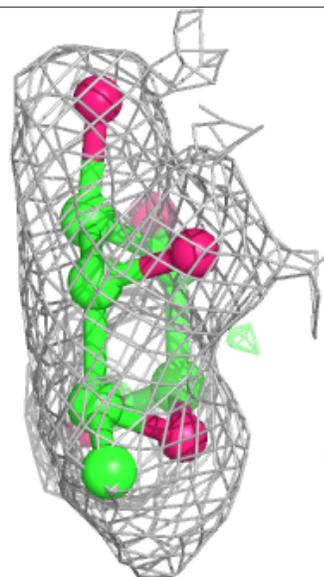
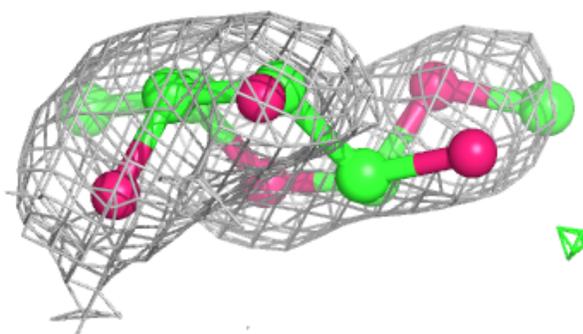
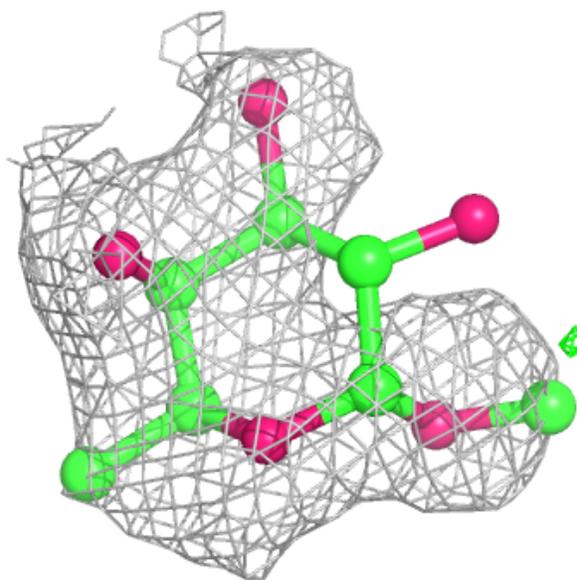
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	BBB	307	6/6	0.87	0.17	42,46,47,49	0
3	GOL	AAA	308	6/6	0.89	0.18	35,36,39,45	0
3	GOL	AAA	307	6/6	0.90	0.15	27,34,35,38	0
3	GOL	AAA	309	6/6	0.91	0.19	33,35,36,42	0
2	MFU	AAA	301	12/12	0.92	0.22	28,31,37,37	12
3	GOL	BBB	305	6/6	0.92	0.26	34,45,47,48	0
2	MFU	BBB	304	12/12	0.92	0.41	31,40,56,59	0
3	GOL	AAA	306	6/6	0.93	0.16	31,32,36,37	6
3	GOL	BBB	306	6/6	0.93	0.17	31,32,35,37	0
2	MFU	AAA	305	12/12	0.93	0.31	25,38,41,42	0
2	MFU	BBB	301	12/12	0.94	0.16	29,31,36,38	0
2	MFU	BBB	302	12/12	0.94	0.17	27,30,35,40	0
2	MFU	BBB	303	12/12	0.96	0.15	27,28,33,46	0
2	MFU	AAA	303	12/12	0.96	0.13	21,22,27,33	0
2	MFU	AAA	302	12/12	0.96	0.12	23,26,28,30	0
4	SO4	AAA	310	5/5	0.96	0.33	37,38,40,46	0
2	MFU	AAA	304	12/12	0.98	0.12	24,26,28,30	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

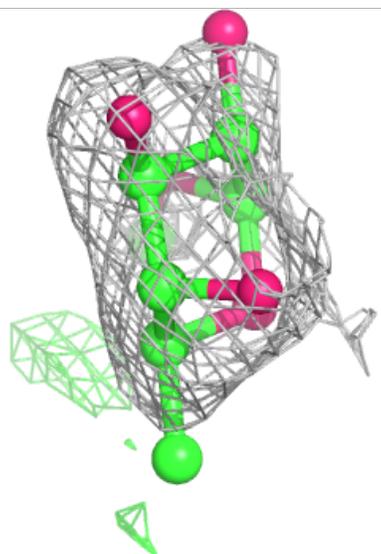
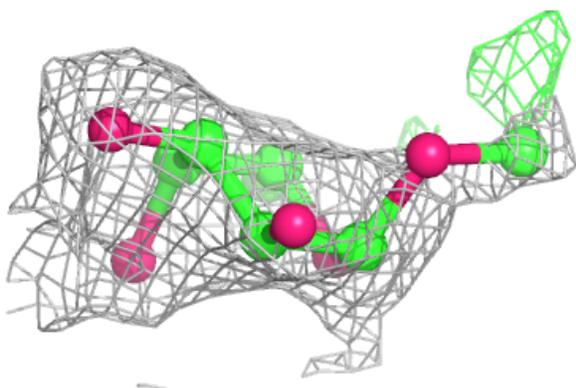
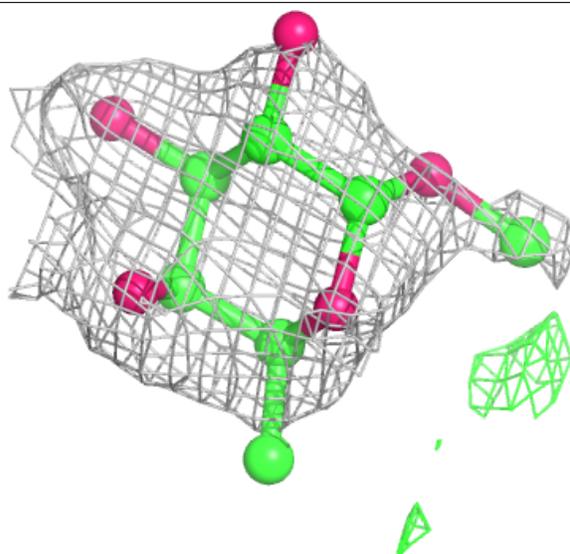
Electron density around MFU AAA 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



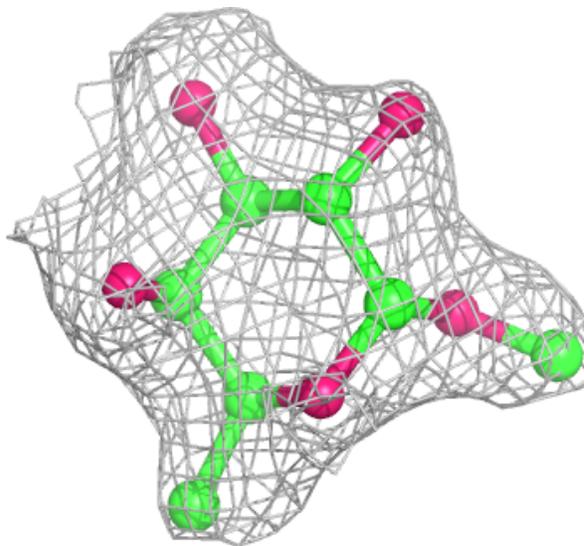
Electron density around MFU BBB 304:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

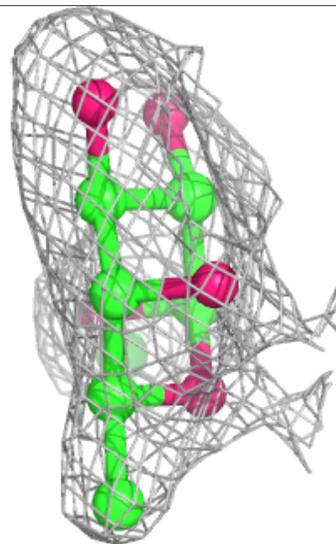
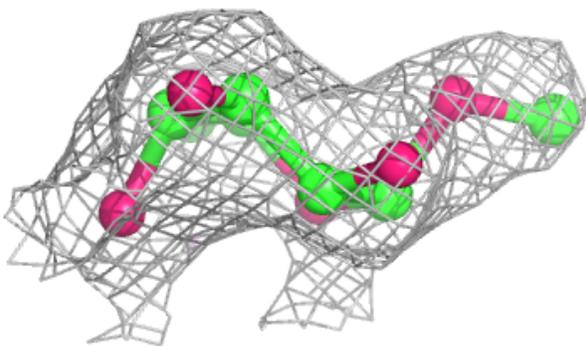


Electron density around MFU AAA 305:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



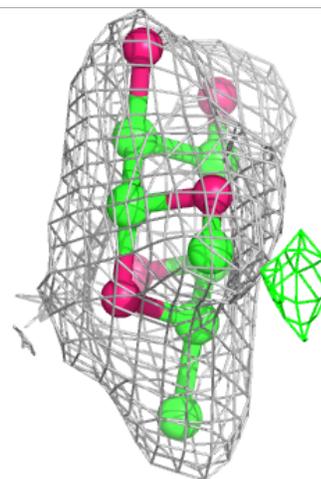
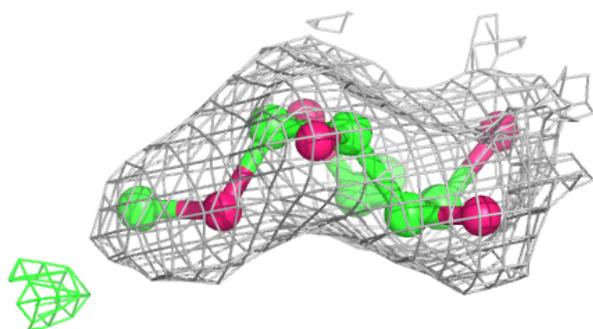
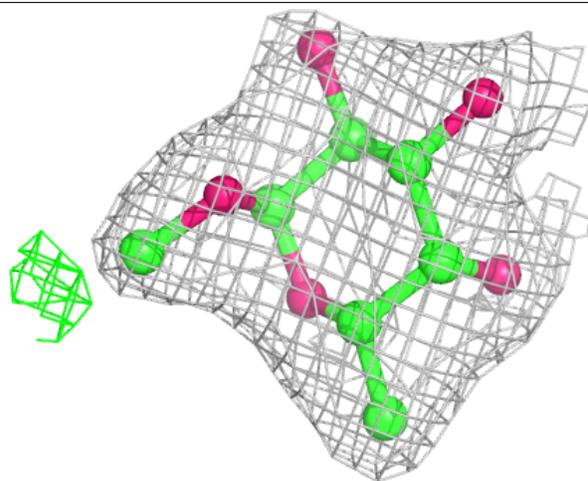
4



4

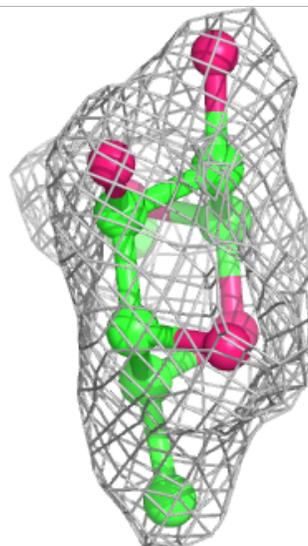
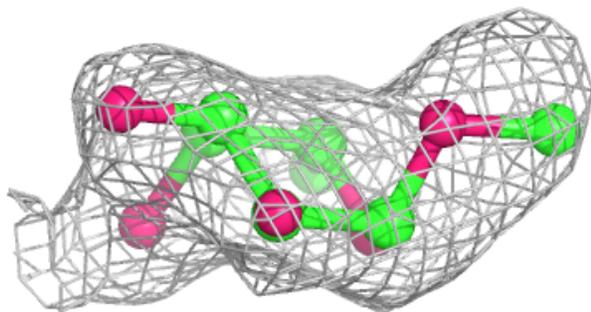
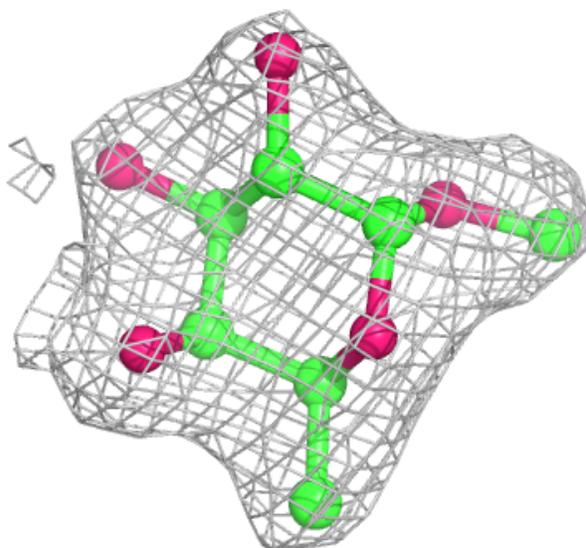
Electron density around MFU BBB 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



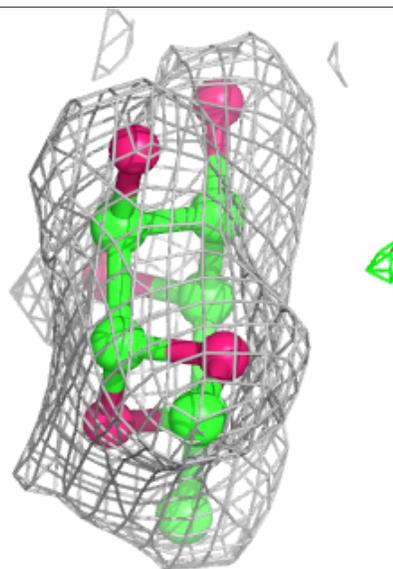
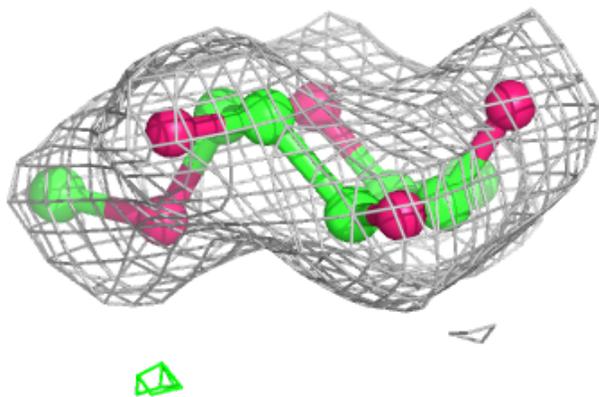
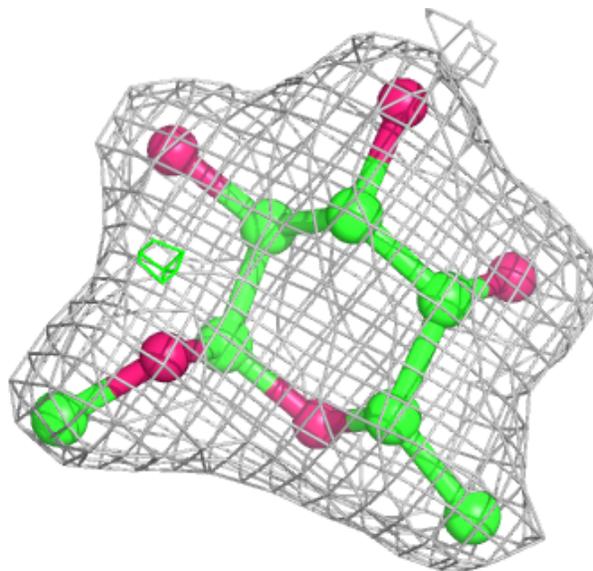
Electron density around MFU BBB 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



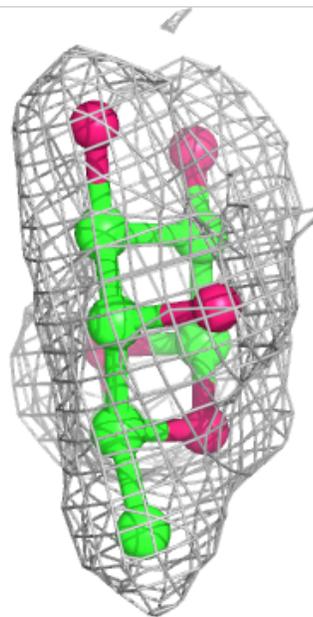
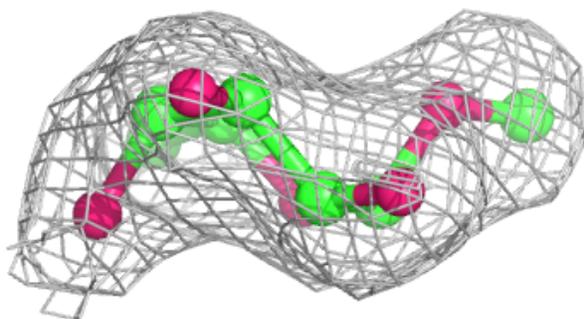
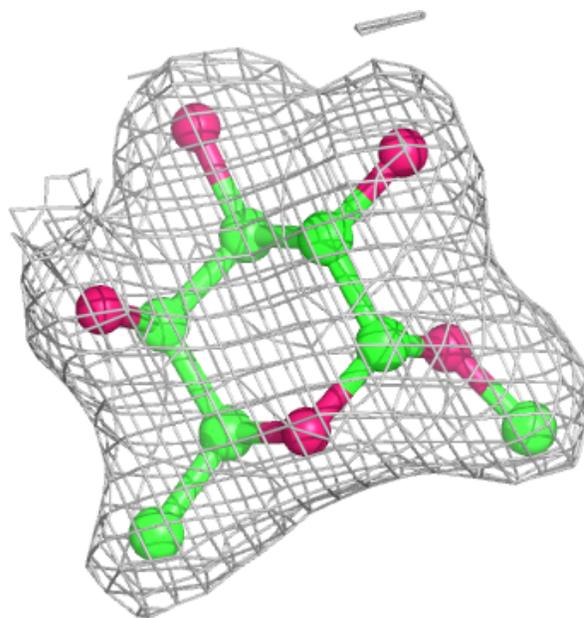
Electron density around MFU BBB 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



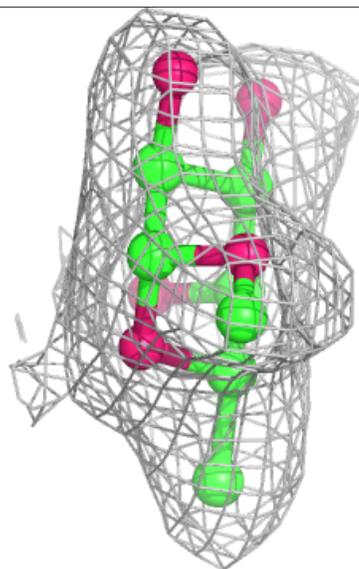
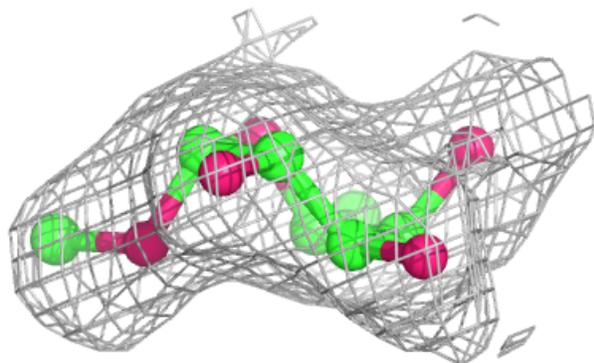
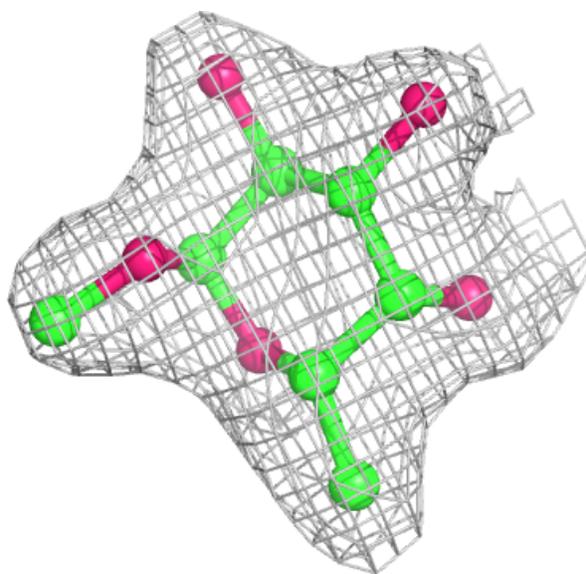
Electron density around MFU AAA 303:

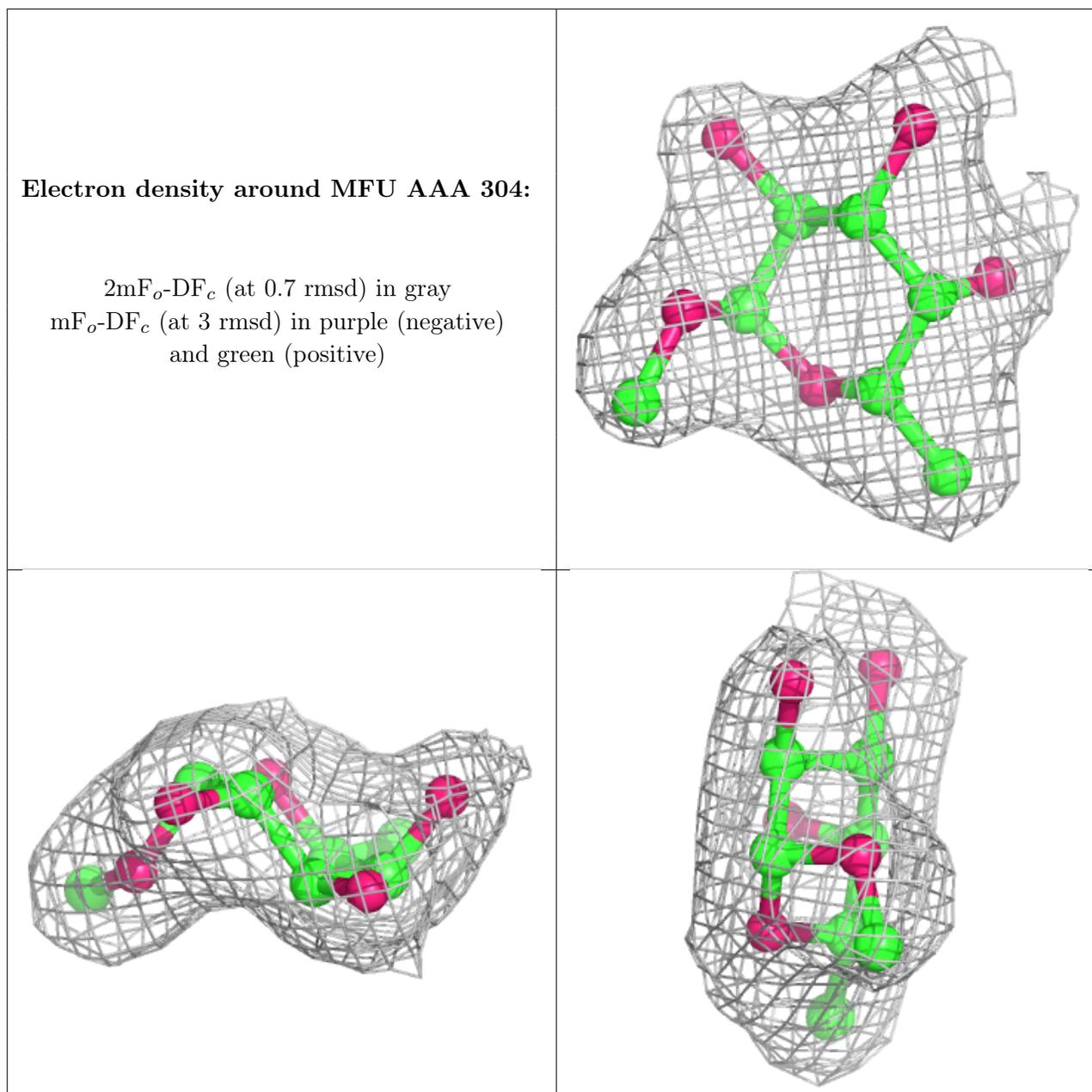
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MFU AAA 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.