



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 03:09 am BST

PDB ID : 4BEI
Title : V. cholera biofilm scaffolding protein RbmA in complex with 18-crown- 6
Authors : Maestre-Reyna, M.; Wang, A.H.-J.
Deposited on : 2013-03-11
Resolution : 2.60 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 : 2.11
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) : 2.11

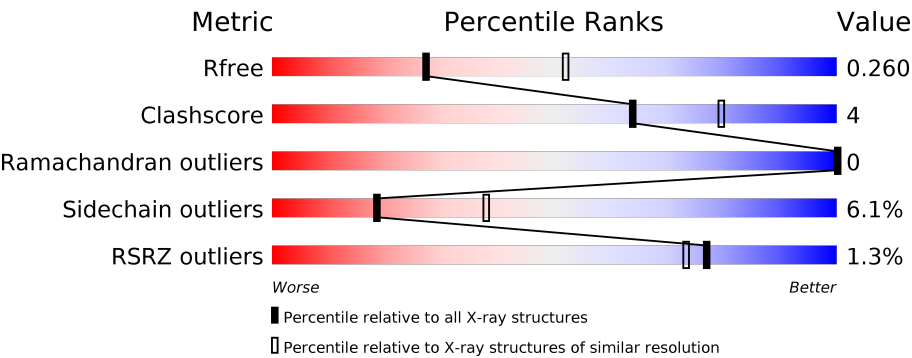
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 on 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	A	262	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>77%11%11%</div></div>
1	B	262	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>81%8%11%</div></div>
1	C	262	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>81%7%11%</div></div>
1	D	262	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>81%8%11%</div></div>
1	E	262	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>79%8%11%</div></div>
1	F	262	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>80%8%11%</div></div>

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Mol	Chain	Length	Quality of chain
1	G	262	<div> <div>%</div> <div> <div>81%8%11%</div> </div> </div>
1	H	262	<div> <div>%</div> <div> <div>81%8%11%</div> </div> </div>

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	O4B	A	1272	-	-	X	-
2	O4B	E	1272	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14574 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 RBMA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1772	1124	293	350	5			
1	B	233	Total	C	N	O	S	0	1	0
			1738	1107	290	338	3			
1	C	233	Total	C	N	O	S	0	2	0
			1749	1113	291	341	4			
1	D	234	Total	C	N	O	S	0	3	0
			1787	1137	294	351	5			
1	E	233	Total	C	N	O	S	0	3	0
			1782	1135	294	348	5			
1	F	233	Total	C	N	O	S	0	2	0
			1755	1115	295	342	3			
1	G	233	Total	C	N	O	S	0	1	0
			1743	1109	291	339	4			
1	H	234	Total	C	N	O	S	0	2	0
			1786	1134	294	353	5			

There are 168 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MET	-	expression tag	UNP C3NSJ9
A	11	GLY	-	expression tag	UNP C3NSJ9
A	12	SER	-	expression tag	UNP C3NSJ9
A	13	SER	-	expression tag	UNP C3NSJ9
A	14	HIS	-	expression tag	UNP C3NSJ9
A	15	HIS	-	expression tag	UNP C3NSJ9
A	16	HIS	-	expression tag	UNP C3NSJ9
A	17	HIS	-	expression tag	UNP C3NSJ9
A	18	HIS	-	expression tag	UNP C3NSJ9
A	19	HIS	-	expression tag	UNP C3NSJ9
A	20	SER	-	expression tag	UNP C3NSJ9
A	21	SER	-	expression tag	UNP C3NSJ9
A	22	GLY	-	expression tag	UNP C3NSJ9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	23	LEU	-	expression tag	UNP C3NSJ9
A	24	VAL	-	expression tag	UNP C3NSJ9
A	25	PRO	-	expression tag	UNP C3NSJ9
A	26	ARG	-	expression tag	UNP C3NSJ9
A	27	GLY	-	expression tag	UNP C3NSJ9
A	28	SER	-	expression tag	UNP C3NSJ9
A	29	HIS	-	expression tag	UNP C3NSJ9
A	30	MET	-	expression tag	UNP C3NSJ9
B	10	MET	-	expression tag	UNP C3NSJ9
B	11	GLY	-	expression tag	UNP C3NSJ9
B	12	SER	-	expression tag	UNP C3NSJ9
B	13	SER	-	expression tag	UNP C3NSJ9
B	14	HIS	-	expression tag	UNP C3NSJ9
B	15	HIS	-	expression tag	UNP C3NSJ9
B	16	HIS	-	expression tag	UNP C3NSJ9
B	17	HIS	-	expression tag	UNP C3NSJ9
B	18	HIS	-	expression tag	UNP C3NSJ9
B	19	HIS	-	expression tag	UNP C3NSJ9
B	20	SER	-	expression tag	UNP C3NSJ9
B	21	SER	-	expression tag	UNP C3NSJ9
B	22	GLY	-	expression tag	UNP C3NSJ9
B	23	LEU	-	expression tag	UNP C3NSJ9
B	24	VAL	-	expression tag	UNP C3NSJ9
B	25	PRO	-	expression tag	UNP C3NSJ9
B	26	ARG	-	expression tag	UNP C3NSJ9
B	27	GLY	-	expression tag	UNP C3NSJ9
B	28	SER	-	expression tag	UNP C3NSJ9
B	29	HIS	-	expression tag	UNP C3NSJ9
B	30	MET	-	expression tag	UNP C3NSJ9
C	10	MET	-	expression tag	UNP C3NSJ9
C	11	GLY	-	expression tag	UNP C3NSJ9
C	12	SER	-	expression tag	UNP C3NSJ9
C	13	SER	-	expression tag	UNP C3NSJ9
C	14	HIS	-	expression tag	UNP C3NSJ9
C	15	HIS	-	expression tag	UNP C3NSJ9
C	16	HIS	-	expression tag	UNP C3NSJ9
C	17	HIS	-	expression tag	UNP C3NSJ9
C	18	HIS	-	expression tag	UNP C3NSJ9
C	19	HIS	-	expression tag	UNP C3NSJ9
C	20	SER	-	expression tag	UNP C3NSJ9
C	21	SER	-	expression tag	UNP C3NSJ9
C	22	GLY	-	expression tag	UNP C3NSJ9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	23	LEU	-	expression tag	UNP C3NSJ9
C	24	VAL	-	expression tag	UNP C3NSJ9
C	25	PRO	-	expression tag	UNP C3NSJ9
C	26	ARG	-	expression tag	UNP C3NSJ9
C	27	GLY	-	expression tag	UNP C3NSJ9
C	28	SER	-	expression tag	UNP C3NSJ9
C	29	HIS	-	expression tag	UNP C3NSJ9
C	30	MET	-	expression tag	UNP C3NSJ9
D	10	MET	-	expression tag	UNP C3NSJ9
D	11	GLY	-	expression tag	UNP C3NSJ9
D	12	SER	-	expression tag	UNP C3NSJ9
D	13	SER	-	expression tag	UNP C3NSJ9
D	14	HIS	-	expression tag	UNP C3NSJ9
D	15	HIS	-	expression tag	UNP C3NSJ9
D	16	HIS	-	expression tag	UNP C3NSJ9
D	17	HIS	-	expression tag	UNP C3NSJ9
D	18	HIS	-	expression tag	UNP C3NSJ9
D	19	HIS	-	expression tag	UNP C3NSJ9
D	20	SER	-	expression tag	UNP C3NSJ9
D	21	SER	-	expression tag	UNP C3NSJ9
D	22	GLY	-	expression tag	UNP C3NSJ9
D	23	LEU	-	expression tag	UNP C3NSJ9
D	24	VAL	-	expression tag	UNP C3NSJ9
D	25	PRO	-	expression tag	UNP C3NSJ9
D	26	ARG	-	expression tag	UNP C3NSJ9
D	27	GLY	-	expression tag	UNP C3NSJ9
D	28	SER	-	expression tag	UNP C3NSJ9
D	29	HIS	-	expression tag	UNP C3NSJ9
D	30	MET	-	expression tag	UNP C3NSJ9
E	10	MET	-	expression tag	UNP C3NSJ9
E	11	GLY	-	expression tag	UNP C3NSJ9
E	12	SER	-	expression tag	UNP C3NSJ9
E	13	SER	-	expression tag	UNP C3NSJ9
E	14	HIS	-	expression tag	UNP C3NSJ9
E	15	HIS	-	expression tag	UNP C3NSJ9
E	16	HIS	-	expression tag	UNP C3NSJ9
E	17	HIS	-	expression tag	UNP C3NSJ9
E	18	HIS	-	expression tag	UNP C3NSJ9
E	19	HIS	-	expression tag	UNP C3NSJ9
E	20	SER	-	expression tag	UNP C3NSJ9
E	21	SER	-	expression tag	UNP C3NSJ9
E	22	GLY	-	expression tag	UNP C3NSJ9

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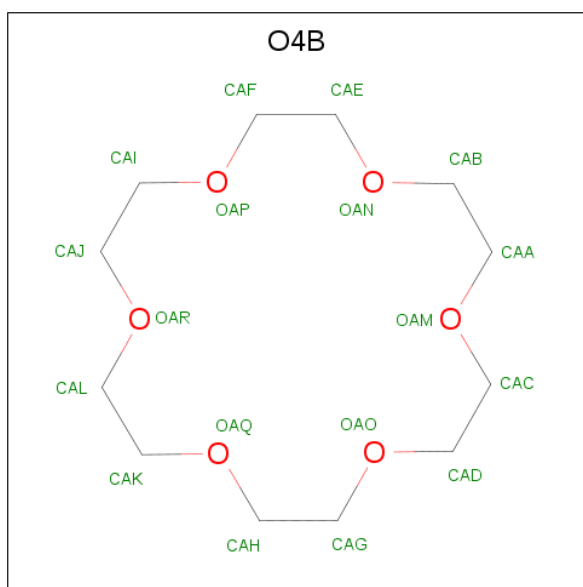
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E	24	VAL	-	expression tag	UNP C3NSJ9
E	25	PRO	-	expression tag	UNP C3NSJ9
E	26	ARG	-	expression tag	UNP C3NSJ9
E	27	GLY	-	expression tag	UNP C3NSJ9
E	28	SER	-	expression tag	UNP C3NSJ9
E	29	HIS	-	expression tag	UNP C3NSJ9
E	30	MET	-	expression tag	UNP C3NSJ9
F	10	MET	-	expression tag	UNP C3NSJ9
F	11	GLY	-	expression tag	UNP C3NSJ9
F	12	SER	-	expression tag	UNP C3NSJ9
F	13	SER	-	expression tag	UNP C3NSJ9
F	14	HIS	-	expression tag	UNP C3NSJ9
F	15	HIS	-	expression tag	UNP C3NSJ9
F	16	HIS	-	expression tag	UNP C3NSJ9
F	17	HIS	-	expression tag	UNP C3NSJ9
F	18	HIS	-	expression tag	UNP C3NSJ9
F	19	HIS	-	expression tag	UNP C3NSJ9
F	20	SER	-	expression tag	UNP C3NSJ9
F	21	SER	-	expression tag	UNP C3NSJ9
F	22	GLY	-	expression tag	UNP C3NSJ9
F	23	LEU	-	expression tag	UNP C3NSJ9
F	24	VAL	-	expression tag	UNP C3NSJ9
F	25	PRO	-	expression tag	UNP C3NSJ9
F	26	ARG	-	expression tag	UNP C3NSJ9
F	27	GLY	-	expression tag	UNP C3NSJ9
F	28	SER	-	expression tag	UNP C3NSJ9
F	29	HIS	-	expression tag	UNP C3NSJ9
F	30	MET	-	expression tag	UNP C3NSJ9
G	10	MET	-	expression tag	UNP C3NSJ9
G	11	GLY	-	expression tag	UNP C3NSJ9
G	12	SER	-	expression tag	UNP C3NSJ9
G	13	SER	-	expression tag	UNP C3NSJ9
G	14	HIS	-	expression tag	UNP C3NSJ9
G	15	HIS	-	expression tag	UNP C3NSJ9
G	16	HIS	-	expression tag	UNP C3NSJ9
G	17	HIS	-	expression tag	UNP C3NSJ9
G	18	HIS	-	expression tag	UNP C3NSJ9
G	19	HIS	-	expression tag	UNP C3NSJ9
G	20	SER	-	expression tag	UNP C3NSJ9
G	21	SER	-	expression tag	UNP C3NSJ9
G	22	GLY	-	expression tag	UNP C3NSJ9

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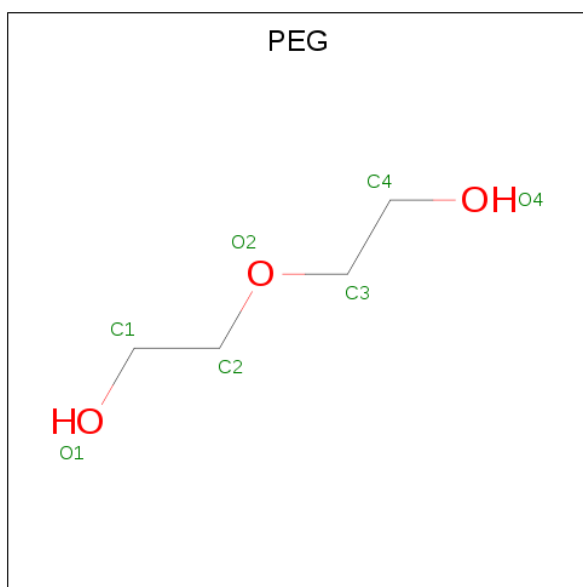
Chain	Residue	Modelled	Actual	Comment	Reference
G	23	LEU	-	expression tag	UNP C3NSJ9
G	24	VAL	-	expression tag	UNP C3NSJ9
G	25	PRO	-	expression tag	UNP C3NSJ9
G	26	ARG	-	expression tag	UNP C3NSJ9
G	27	GLY	-	expression tag	UNP C3NSJ9
G	28	SER	-	expression tag	UNP C3NSJ9
G	29	HIS	-	expression tag	UNP C3NSJ9
G	30	MET	-	expression tag	UNP C3NSJ9
H	10	MET	-	expression tag	UNP C3NSJ9
H	11	GLY	-	expression tag	UNP C3NSJ9
H	12	SER	-	expression tag	UNP C3NSJ9
H	13	SER	-	expression tag	UNP C3NSJ9
H	14	HIS	-	expression tag	UNP C3NSJ9
H	15	HIS	-	expression tag	UNP C3NSJ9
H	16	HIS	-	expression tag	UNP C3NSJ9
H	17	HIS	-	expression tag	UNP C3NSJ9
H	18	HIS	-	expression tag	UNP C3NSJ9
H	19	HIS	-	expression tag	UNP C3NSJ9
H	20	SER	-	expression tag	UNP C3NSJ9
H	21	SER	-	expression tag	UNP C3NSJ9
H	22	GLY	-	expression tag	UNP C3NSJ9
H	23	LEU	-	expression tag	UNP C3NSJ9
H	24	VAL	-	expression tag	UNP C3NSJ9
H	25	PRO	-	expression tag	UNP C3NSJ9
H	26	ARG	-	expression tag	UNP C3NSJ9
H	27	GLY	-	expression tag	UNP C3NSJ9
H	28	SER	-	expression tag	UNP C3NSJ9
H	29	HIS	-	expression tag	UNP C3NSJ9
H	30	MET	-	expression tag	UNP C3NSJ9

- Molecule 2 is 1,4,7,10,13,16-HEXA OXACYCLOOCTADECANE (three-letter code: O4B) (formula: C₁₂H₂₄O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			18	12	6		
2	B	1	Total	C	O	0	0
			18	12	6		
2	C	1	Total	C	O	0	0
			18	12	6		
2	D	1	Total	C	O	0	0
			18	12	6		
2	E	1	Total	C	O	0	0
			18	12	6		
2	F	1	Total	C	O	0	0
			18	12	6		
2	G	1	Total	C	O	0	0
			18	12	6		
2	H	1	Total	C	O	0	0
			18	12	6		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	D	1	Total	C	O	0	0
			7	4	3		
3	E	1	Total	C	O	0	0
			7	4	3		
3	F	1	Total	C	O	0	0
			7	4	3		


- Molecule 4 is water.

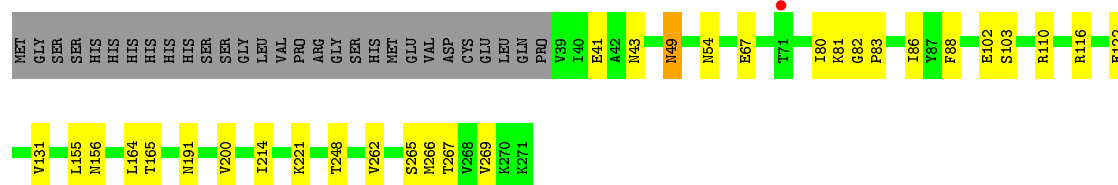
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	52	Total	O	0	0
			52	52		
4	B	28	Total	O	0	0
			28	28		
4	C	33	Total	O	0	0
			33	33		
4	D	28	Total	O	0	0
			28	28		
4	E	45	Total	O	0	0
			45	45		
4	F	21	Total	O	0	0
			21	21		
4	G	26	Total	O	0	1
			27	27		
4	H	56	Total	O	0	0
			56	56		

3 Residue-property plots


These plots are drawn for all protein, RNA and DNA 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.

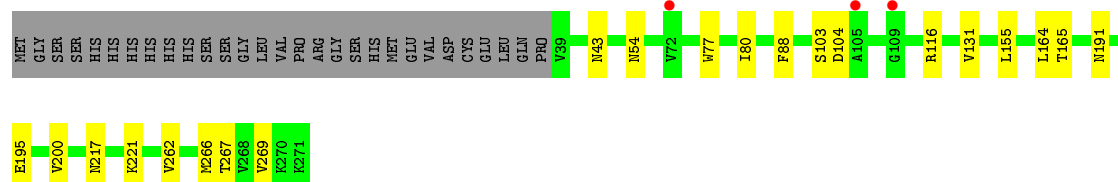
• Molecule 1: RBMA

Chain A: 




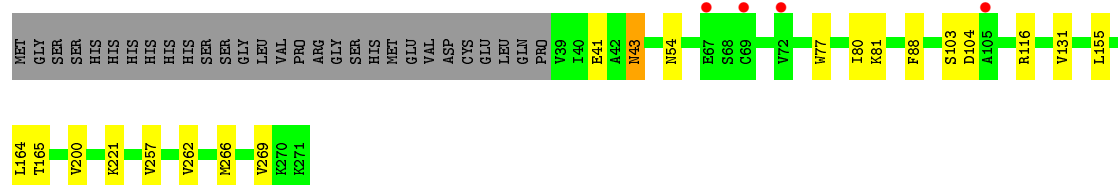
• Molecule 1: RBMA

Chain B: 




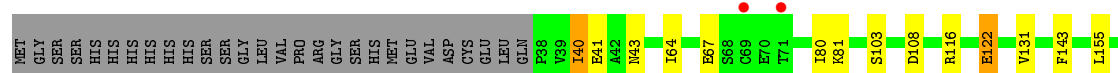
• Molecule 1: RBMA

Chain C: 



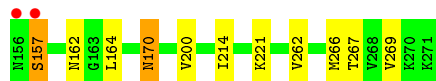
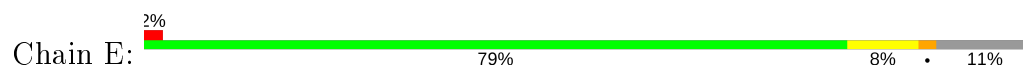
• Molecule 1: RBMA

Chain D: 

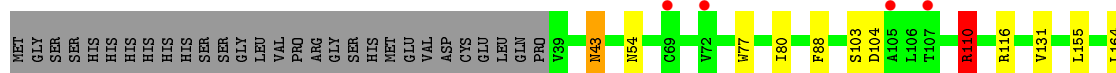
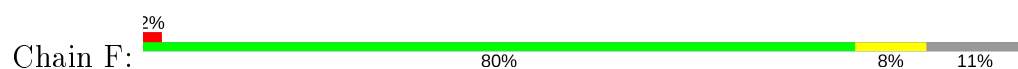




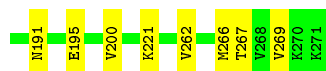
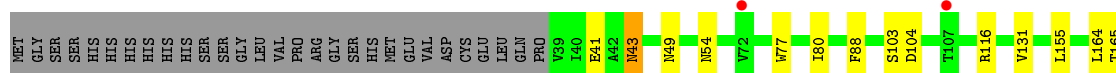
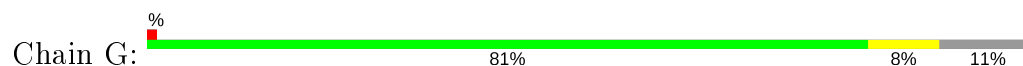
- Molecule 1: RBMA



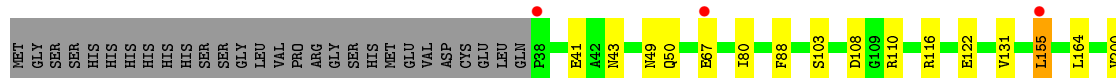
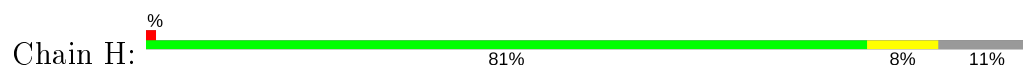
- Molecule 1: RBMA



- Molecule 1: RBMA



- Molecule 1: RBMA



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	136.88Å 136.88Å 116.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	21.96 – 2.60 21.95 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (21.96-2.60) 100.0 (21.95-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.222 , 0.261 0.224 , 0.260	Depositor DCC
R_{free} test set	3780 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	54.3	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 27.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.470 for -h,-k,l 0.487 for h,-h-k,-l 0.470 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14574	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: O4B, PEG

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	0.47	0/1805	0.73	0/2451
1	B	0.45	0/1774	0.70	0/2415
1	C	0.46	0/1785	0.70	0/2429
1	D	0.47	0/1825	0.71	0/2482
1	E	0.47	0/1819	0.71	0/2472
1	F	0.45	0/1791	0.72	1/2437 (0.0%)
1	G	0.46	0/1779	0.70	1/2420 (0.0%)
1	H	0.47	0/1826	0.73	0/2481
All	All	0.46	0/14404	0.71	2/19587 (0.0%)

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	110	ARG	NE-CZ-NH1	5.27	122.94	120.30
1	G	49	ASN	CB-CA-C	-5.01	100.38	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	ASN	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	1772	0	1709	32	0
1	B	1738	0	1656	9	0
1	C	1749	0	1666	11	0
1	D	1787	0	1717	18	0
1	E	1782	0	1720	24	0
1	F	1755	0	1674	13	0
1	G	1743	0	1671	11	0
1	H	1786	0	1726	16	0
2	A	18	0	24	9	0
2	B	18	0	24	0	0
2	C	18	0	24	2	0
2	D	18	0	24	1	0
2	E	18	0	24	10	0
2	F	18	0	24	1	0
2	G	18	0	24	0	0
2	H	18	0	24	0	0
3	A	7	0	10	0	0
3	D	7	0	10	0	0
3	E	7	0	10	0	0
3	F	7	0	10	0	0
4	A	52	0	0	2	0
4	B	28	0	0	3	0
4	C	33	0	0	3	0
4	D	28	0	0	1	0
4	E	45	0	0	1	0
4	F	21	0	0	3	0
4	G	27	0	0	4	0
4	H	56	0	0	1	0
All	All	14574	0	13771	119	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 4.

All (119) 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:E:81:LYS:HZ1	2:E:1272:O4B:HAI2	1.31	0.93
1:E:81:LYS:NZ	2:E:1272:O4B:HAE1	1.95	0.82
1:B:116:ARG:NH2	4:B:2007:HOH:O	2.12	0.81
1:F:116:ARG:NH2	4:F:2008:HOH:O	2.12	0.81
1:E:81:LYS:HZ1	2:E:1272:O4B:CAI	1.97	0.77
1:E:154:ASP:HB2	1:E:157:SER:OG	1.84	0.77
1:E:61[B]:GLN:NE2	1:E:115:THR:OG1	2.19	0.75
1:A:83:PRO:O	1:A:86:ILE:HG22	1.87	0.74
1:G:116:ARG:NH2	4:G:2004:HOH:O	2.21	0.73
1:D:41:GLU:OE1	1:D:67:GLU:OE2	2.07	0.73
1:D:116:ARG:NH2	4:D:2009:HOH:O	2.22	0.72
1:H:41:GLU:OE1	1:H:67:GLU:OE2	2.07	0.72
1:C:116:ARG:NH2	4:C:2006:HOH:O	2.24	0.71
1:C:257:VAL:O	1:H:110:ARG:NH1	2.27	0.68
1:A:81:LYS:HZ2	2:A:1272:O4B:HAJ2	1.61	0.66
1:A:110:ARG:NH1	1:F:257:VAL:O	2.27	0.66
1:A:41:GLU:OE1	1:A:67:GLU:OE2	2.14	0.66
1:E:81:LYS:HZ3	2:E:1272:O4B:HAE1	1.58	0.65
1:E:41:GLU:OE1	1:E:67:GLU:OE2	2.14	0.65
1:A:248:THR:HG21	2:A:1272:O4B:HAB2	1.79	0.65
1:H:49[B]:ASN:ND2	1:H:155:LEU:HD12	2.14	0.63
1:H:49[B]:ASN:ND2	1:H:155:LEU:CD1	2.62	0.62
1:H:116:ARG:NH2	4:H:2010:HOH:O	2.34	0.61
1:A:49:ASN:HB2	1:A:155:LEU:HD13	1.83	0.60
1:D:40:ILE:CG2	1:D:143:PHE:CD1	2.85	0.60
1:D:40:ILE:HG23	1:D:143:PHE:CD1	2.37	0.60
1:E:81:LYS:HZ1	2:E:1272:O4B:HAE1	1.68	0.59
1:F:43:ASN:ND2	4:F:2001:HOH:O	2.36	0.58
1:E:81:LYS:HZ2	2:E:1272:O4B:HAK1	1.68	0.57
1:H:49[B]:ASN:HD22	1:H:155:LEU:HD11	1.68	0.57
1:A:81:LYS:HZ2	2:A:1272:O4B:CAJ	2.18	0.57
1:A:165:THR:HG22	1:A:191:ASN:HB3	1.86	0.56
1:A:110:ARG:HH21	1:E:122:GLU:CD	2.08	0.56
1:B:165:THR:HG22	1:B:191:ASN:HB3	1.88	0.56
1:G:165:THR:HG23	4:G:2017:HOH:O	2.04	0.56
1:A:265:SER:OG	2:A:1272:O4B:HAC1	2.05	0.56
1:A:116:ARG:NH2	4:A:2011:HOH:O	2.39	0.56
1:G:165:THR:HG22	1:G:191:ASN:HB3	1.88	0.55
1:D:122:GLU:CD	1:H:110:ARG:HH21	2.10	0.55
1:D:40:ILE:HG13	1:D:64:ILE:HG23	1.89	0.55
1:C:41:GLU:CB	4:C:2001:HOH:O	2.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ASN:ND2	4:A:2003:HOH:O	2.35	0.55
1:C:81:LYS:HZ1	2:C:1272:O4B:HAG2	1.72	0.54
1:F:165:THR:HG22	1:F:191:ASN:HB3	1.88	0.54
1:F:265:SER:CB	2:F:1272:O4B:HAH1	2.37	0.54
1:D:122:GLU:OE1	1:H:110:ARG:NH2	2.42	0.53
1:A:82:GLY:C	1:A:86:ILE:HG22	2.27	0.53
1:A:49:ASN:HB2	1:A:155:LEU:CD1	2.39	0.53
1:A:110:ARG:NH2	1:E:122:GLU:OE1	2.42	0.52
1:E:116:ARG:NH2	4:E:2009:HOH:O	2.41	0.52
1:E:81:LYS:NZ	2:E:1272:O4B:HAI2	2.16	0.52
1:A:102:GLU:CD	1:F:195:GLU:H	2.13	0.52
1:H:49[B]:ASN:HD22	1:H:155:LEU:CD1	2.24	0.51
1:F:110:ARG:HH11	1:F:110:ARG:CG	2.25	0.50
1:F:165:THR:HG23	4:F:2016:HOH:O	2.12	0.50
1:A:80:ILE:HG13	1:A:88:PHE:CE1	2.47	0.50
1:B:80:ILE:HG13	1:B:88:PHE:CE1	2.47	0.50
1:G:80:ILE:HG13	1:G:88:PHE:CE1	2.47	0.49
1:F:80:ILE:HG13	1:F:88:PHE:CE1	2.48	0.49
1:E:164:LEU:CD2	1:E:200:VAL:HG11	2.43	0.48
1:B:195:GLU:CB	4:B:2023:HOH:O	2.61	0.48
1:C:80:ILE:HG13	1:C:88:PHE:CE1	2.48	0.48
1:E:80:ILE:HG13	1:E:88:PHE:CE1	2.48	0.48
1:H:80:ILE:HG13	1:H:88:PHE:CE1	2.47	0.48
1:A:164:LEU:CD2	1:A:200:VAL:HG11	2.44	0.48
1:A:83:PRO:O	1:A:86:ILE:CG2	2.58	0.48
1:E:170:ASN:HD22	1:E:170:ASN:C	2.17	0.48
1:D:170:ASN:HD22	1:D:170:ASN:C	2.18	0.47
1:D:164:LEU:CD2	1:D:200:VAL:HG11	2.43	0.47
1:E:81:LYS:HZ2	2:E:1272:O4B:CAK	2.28	0.47
1:G:164:LEU:CD2	1:G:200:VAL:HG11	2.44	0.47
1:H:164:LEU:CD2	1:H:200:VAL:HG11	2.44	0.47
1:B:164:LEU:CD2	1:B:200:VAL:HG11	2.44	0.47
1:F:164:LEU:CD2	1:F:200:VAL:HG11	2.44	0.47
1:C:164:LEU:CD2	1:C:200:VAL:HG11	2.45	0.47
1:D:40:ILE:HG22	1:D:143:PHE:CE1	2.49	0.46
1:A:110:ARG:NH2	1:E:122:GLU:OE2	2.48	0.46
1:A:265:SER:CB	2:A:1272:O4B:HAC1	2.45	0.46
1:A:82:GLY:C	1:A:86:ILE:CG2	2.84	0.46
1:D:40:ILE:HG23	1:D:143:PHE:CG	2.51	0.45
1:G:41:GLU:CB	4:G:2001:HOH:O	2.64	0.45
1:A:110:ARG:NH2	1:E:122:GLU:CD	2.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:80:ILE:HD12	1:G:131:VAL:HG22	1.99	0.45
1:A:82:GLY:CA	1:A:86:ILE:CG2	2.95	0.45
1:A:81:LYS:HD3	2:A:1272:O4B:HAK1	1.99	0.44
1:B:165:THR:HG23	4:B:2021:HOH:O	2.16	0.44
1:E:81:LYS:HZ2	2:E:1272:O4B:HAG2	1.81	0.44
1:D:122:GLU:OE2	1:H:110:ARG:NH2	2.51	0.44
1:B:80:ILE:HD12	1:B:131:VAL:HG22	1.99	0.44
1:C:80:ILE:HD12	1:C:131:VAL:HG22	1.99	0.44
1:C:81:LYS:HZ1	2:C:1272:O4B:HAK1	1.83	0.44
1:A:82:GLY:CA	1:A:86:ILE:HG23	2.48	0.43
1:E:81:LYS:NZ	2:E:1272:O4B:HAK1	2.31	0.43
1:A:81:LYS:NZ	2:A:1272:O4B:HAK2	2.34	0.43
1:D:80:ILE:HD12	1:D:131:VAL:HG22	2.00	0.43
1:E:80:ILE:HD12	1:E:131:VAL:HG22	2.01	0.43
1:A:80:ILE:HD12	1:A:131:VAL:HG22	2.01	0.43
1:H:80:ILE:HD12	1:H:131:VAL:HG22	2.00	0.43
1:D:122:GLU:CD	1:H:110:ARG:NH2	2.70	0.43
1:C:77:TRP:CZ2	1:D:214:ILE:HG12	2.54	0.43
1:F:80:ILE:HD12	1:F:131:VAL:HG22	1.99	0.43
1:A:214:ILE:HG12	1:B:77:TRP:CZ2	2.53	0.42
1:E:214:ILE:HG12	1:F:77:TRP:CZ2	2.53	0.42
1:G:77:TRP:CZ2	1:H:214:ILE:HG12	2.55	0.42
1:D:81:LYS:HD3	2:D:1272:O4B:HAK1	2.00	0.42
1:D:164:LEU:HD21	1:D:200:VAL:HG11	2.02	0.42
1:D:40:ILE:HG12	1:D:41:GLU:N	2.34	0.42
1:A:164:LEU:HD21	1:A:200:VAL:HG11	2.02	0.41
1:F:164:LEU:HD21	1:F:200:VAL:HG11	2.03	0.41
1:B:164:LEU:HD21	1:B:200:VAL:HG11	2.03	0.41
1:E:164:LEU:HD21	1:E:200:VAL:HG11	2.02	0.41
1:A:81:LYS:HZ3	2:A:1272:O4B:HAK2	1.85	0.41
1:A:81:LYS:HD3	2:A:1272:O4B:HAG2	2.03	0.41
1:G:195:GLU:OE1	1:G:195:GLU:N	2.54	0.41
1:G:43:ASN:ND2	4:G:2002:HOH:O	2.54	0.41
1:H:164:LEU:HD21	1:H:200:VAL:HG11	2.03	0.41
1:C:164:LEU:HD21	1:C:200:VAL:HG11	2.03	0.40
1:C:43:ASN:ND2	4:C:2002:HOH:O	2.54	0.40
1:G:164:LEU:HD21	1:G:200:VAL:HG11	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	231/262 (88%)	226 (98%)	5 (2%)	0	100	100
1	B	232/262 (88%)	227 (98%)	5 (2%)	0	100	100
1	C	233/262 (89%)	228 (98%)	5 (2%)	0	100	100
1	D	235/262 (90%)	231 (98%)	4 (2%)	0	100	100
1	E	234/262 (89%)	229 (98%)	5 (2%)	0	100	100
1	F	233/262 (89%)	228 (98%)	5 (2%)	0	100	100
1	G	232/262 (88%)	227 (98%)	5 (2%)	0	100	100
1	H	234/262 (89%)	229 (98%)	5 (2%)	0	100	100
All	All	1864/2096 (89%)	1825 (98%)	39 (2%)	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	A	187/221 (85%)	177 (95%)	10 (5%)	22	45
1	B	177/221 (80%)	166 (94%)	11 (6%)	18	37
1	C	179/221 (81%)	169 (94%)	10 (6%)	21	42
1	D	188/221 (85%)	176 (94%)	12 (6%)	17	35
1	E	187/221 (85%)	173 (92%)	14 (8%)	13	27
1	F	180/221 (81%)	169 (94%)	11 (6%)	18	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	180/221 (81%)	170 (94%)	10 (6%)	21	42
1	H	190/221 (86%)	179 (94%)	11 (6%)	20	40
All	All	1468/1768 (83%)	1379 (94%)	89 (6%)	18	38

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	49	ASN
1	A	54	ASN
1	A	103	SER
1	A	122	GLU
1	A	221	LYS
1	A	262	VAL
1	A	266	MET
1	A	267	THR
1	A	269	VAL
1	B	43	ASN
1	B	54	ASN
1	B	103	SER
1	B	104	ASP
1	B	155	LEU
1	B	217	ASN
1	B	221	LYS
1	B	262	VAL
1	B	266	MET
1	B	267	THR
1	B	269	VAL
1	C	43	ASN
1	C	54	ASN
1	C	103	SER
1	C	104	ASP
1	C	155	LEU
1	C	165	THR
1	C	221	LYS
1	C	262	VAL
1	C	266	MET
1	C	269	VAL
1	D	40	ILE
1	D	43	ASN
1	D	103	SER

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Mol	Chain	Res	Type
1	D	108	ASP
1	D	122	GLU
1	D	155	LEU
1	D	170	ASN
1	D	221	LYS
1	D	262	VAL
1	D	266	MET
1	D	267	THR
1	D	269	VAL
1	E	43	ASN
1	E	103	SER
1	E	108	ASP
1	E	116	ARG
1	E	122	GLU
1	E	155	LEU
1	E	157	SER
1	E	162	ASN
1	E	170	ASN
1	E	221	LYS
1	E	262	VAL
1	E	266	MET
1	E	267	THR
1	E	269	VAL
1	F	43	ASN
1	F	54	ASN
1	F	103	SER
1	F	104	ASP
1	F	110	ARG
1	F	155	LEU
1	F	221	LYS
1	F	262	VAL
1	F	266	MET
1	F	267	THR
1	F	269	VAL
1	G	43	ASN
1	G	54	ASN
1	G	103	SER
1	G	104	ASP
1	G	155	LEU
1	G	221	LYS
1	G	262	VAL
1	G	266	MET

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Mol	Chain	Res	Type
1	G	267	THR
1	G	269	VAL
1	H	43	ASN
1	H	50	GLN
1	H	103	SER
1	H	108	ASP
1	H	122	GLU
1	H	155	LEU
1	H	221	LYS
1	H	262	VAL
1	H	266	MET
1	H	267	THR
1	H	269	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	61	GLN
1	A	146	ASN
1	A	217	ASN
1	A	250	ASN
1	A	252	GLN
1	B	50	GLN
1	B	54	ASN
1	B	100	GLN
1	B	113	ASN
1	B	146	ASN
1	B	156	ASN
1	B	250	ASN
1	B	252	GLN
1	C	49	ASN
1	C	54	ASN
1	C	146	ASN
1	C	156	ASN
1	C	217	ASN
1	C	250	ASN
1	C	252	GLN
1	D	146	ASN
1	D	170	ASN
1	D	217	ASN
1	D	250	ASN

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Mol	Chain	Res	Type
1	D	252	GLN
1	E	146	ASN
1	E	162	ASN
1	E	170	ASN
1	E	217	ASN
1	E	250	ASN
1	E	252	GLN
1	F	49	ASN
1	F	54	ASN
1	F	113	ASN
1	F	146	ASN
1	F	156	ASN
1	F	250	ASN
1	F	252	GLN
1	G	54	ASN
1	G	146	ASN
1	G	156	ASN
1	G	217	ASN
1	G	250	ASN
1	G	252	GLN
1	G	264	GLN
1	H	54	ASN
1	H	146	ASN
1	H	217	ASN
1	H	250	ASN
1	H	252	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

12 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	O4B	E	1272	-	18,18,18	0.90	0	18,18,18	1.02	0
3	PEG	D	1273	-	6,6,6	0.51	0	5,5,5	0.35	0
2	O4B	C	1272	-	18,18,18	0.90	0	18,18,18	0.92	0
2	O4B	G	1272	-	18,18,18	1.01	1 (5%)	18,18,18	0.82	0
3	PEG	F	1273	-	6,6,6	0.59	0	5,5,5	0.58	0
2	O4B	A	1272	-	18,18,18	0.92	0	18,18,18	0.76	0
2	O4B	H	1272	-	18,18,18	0.94	1 (5%)	18,18,18	0.96	1 (5%)
3	PEG	A	1273	-	6,6,6	0.56	0	5,5,5	0.12	0
2	O4B	B	1272	-	18,18,18	0.68	0	18,18,18	0.82	0
3	PEG	E	1273	-	6,6,6	0.63	0	5,5,5	0.66	0
2	O4B	D	1272	-	18,18,18	0.87	0	18,18,18	0.86	0
2	O4B	F	1272	-	18,18,18	0.94	0	18,18,18	0.72	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	O4B	E	1272	-	-	10/18/18/18	1/1/1/1
3	PEG	D	1273	-	-	1/4/4/4	-
2	O4B	C	1272	-	-	9/18/18/18	0/1/1/1
2	O4B	G	1272	-	-	7/18/18/18	0/1/1/1
3	PEG	F	1273	-	-	3/4/4/4	-
2	O4B	A	1272	-	-	8/18/18/18	0/1/1/1
2	O4B	H	1272	-	-	7/18/18/18	0/1/1/1
3	PEG	A	1273	-	-	2/4/4/4	-
2	O4B	B	1272	-	-	2/18/18/18	0/1/1/1
3	PEG	E	1273	-	-	3/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	O4B	D	1272	-	-	6/18/18/18	0/1/1/1
2	O4B	F	1272	-	-	3/18/18/18	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	1272	O4B	OAM-CAA	2.25	1.51	1.42
2	H	1272	O4B	OAM-CAA	2.08	1.51	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1272	O4B	CAB-OAN-CAE	-2.24	103.56	113.29

There are no chirality outliers.

All (61) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1272	O4B	CAK-CAL-OAR-CAJ
2	E	1272	O4B	CAG-CAH-OAQ-CAK
2	A	1272	O4B	OAQ-CAK-CAL-OAR
2	A	1272	O4B	OAQ-CAK-CAL-OAR
2	E	1272	O4B	OAP-CAI-CAJ-OAR
2	E	1272	O4B	OAP-CAI-CAJ-OAR
2	C	1272	O4B	OAP-CAI-CAJ-OAR
2	G	1272	O4B	OAM-CAA-CAB-OAN
2	A	1272	O4B	OAQ-CAK-CAL-OAR
2	B	1272	O4B	OAQ-CAK-CAL-OAR
3	D	1273	PEG	O2-C3-C4-O4
2	E	1272	O4B	OAQ-CAK-CAL-OAR
2	A	1272	O4B	OAN-CAE-CAF-OAP
3	F	1273	PEG	O1-C1-C2-O2
3	A	1273	PEG	O2-C3-C4-O4
2	E	1272	O4B	OAQ-CAK-CAL-OAR
2	C	1272	O4B	OAQ-CAK-CAL-OAR
2	G	1272	O4B	OAQ-CAK-CAL-OAR
2	H	1272	O4B	OAM-CAA-CAB-OAN
2	E	1272	O4B	CAA-CAB-OAN-CAE
3	F	1273	PEG	O2-C3-C4-O4
3	E	1273	PEG	O1-C1-C2-O2
2	A	1272	O4B	CAG-CAH-OAQ-CAK
2	D	1272	O4B	OAM-CAA-CAB-OAN

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Mol	Chain	Res	Type	Atoms
2	G	1272	O4B	OAN-CAE-CAF-OAP
2	H	1272	O4B	OAQ-CAH-CAH-OAQ
2	D	1272	O4B	OAP-CAI-CAJ-OAR
2	F	1272	O4B	OAQ-CAK-CAL-OAR
2	F	1272	O4B	OAM-CAC-CAD-OAO
2	C	1272	O4B	OAM-CAA-CAB-OAN
2	F	1272	O4B	CAL-CAK-OAQ-CAH
2	G	1272	O4B	CAG-CAH-OAQ-CAK
2	G	1272	O4B	CAF-CAE-OAN-CAB
2	H	1272	O4B	CAG-CAH-OAQ-CAK
2	G	1272	O4B	CAB-CAA-OAM-CAC
2	A	1272	O4B	CAB-CAA-OAM-CAC
3	E	1273	PEG	C1-C2-O2-C3
2	E	1272	O4B	CAF-CAE-OAN-CAB
2	H	1272	O4B	CAC-CAD-OAO-CAG
2	H	1272	O4B	CAJ-CAI-OAP-CAF
3	F	1273	PEG	C1-C2-O2-C3
2	D	1272	O4B	CAB-CAA-OAM-CAC
2	C	1272	O4B	CAC-CAD-OAO-CAG
2	A	1272	O4B	CAF-CAE-OAN-CAB
2	E	1272	O4B	OAM-CAA-CAB-OAN
2	B	1272	O4B	OAM-CAC-CAD-OAO
2	G	1272	O4B	CAC-CAD-OAO-CAG
2	D	1272	O4B	CAC-CAD-OAO-CAG
2	D	1272	O4B	CAK-CAL-OAR-CAJ
2	C	1272	O4B	OAM-CAC-CAD-OAO
3	A	1273	PEG	C1-C2-O2-C3
2	D	1272	O4B	OAQ-CAH-CAH-OAQ
2	E	1272	O4B	CAK-CAL-OAR-CAJ
2	C	1272	O4B	CAG-CAH-OAQ-CAK
2	H	1272	O4B	CAD-CAC-OAM-CAA
2	E	1272	O4B	CAC-CAD-OAO-CAG
2	H	1272	O4B	OAM-CAC-CAD-OAO
2	C	1272	O4B	CAK-CAL-OAR-CAJ
2	E	1272	O4B	OAN-CAE-CAF-OAP
2	C	1272	O4B	CAI-CAJ-OAR-CAL
3	E	1273	PEG	C4-C3-O2-C2
2	C	1272	O4B	OAQ-CAK-CAL-OAR

All (1) ring outliers are listed below:

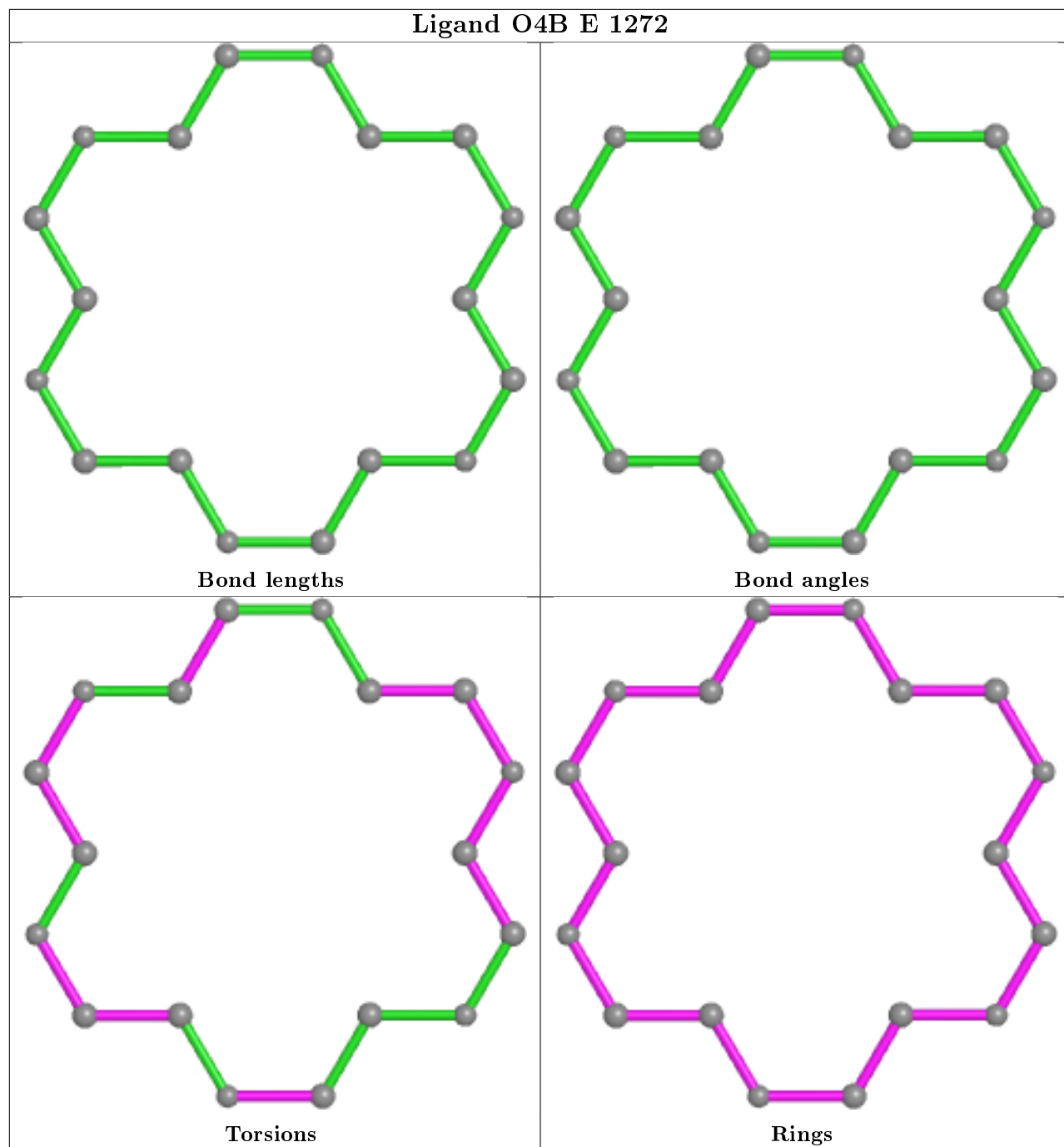
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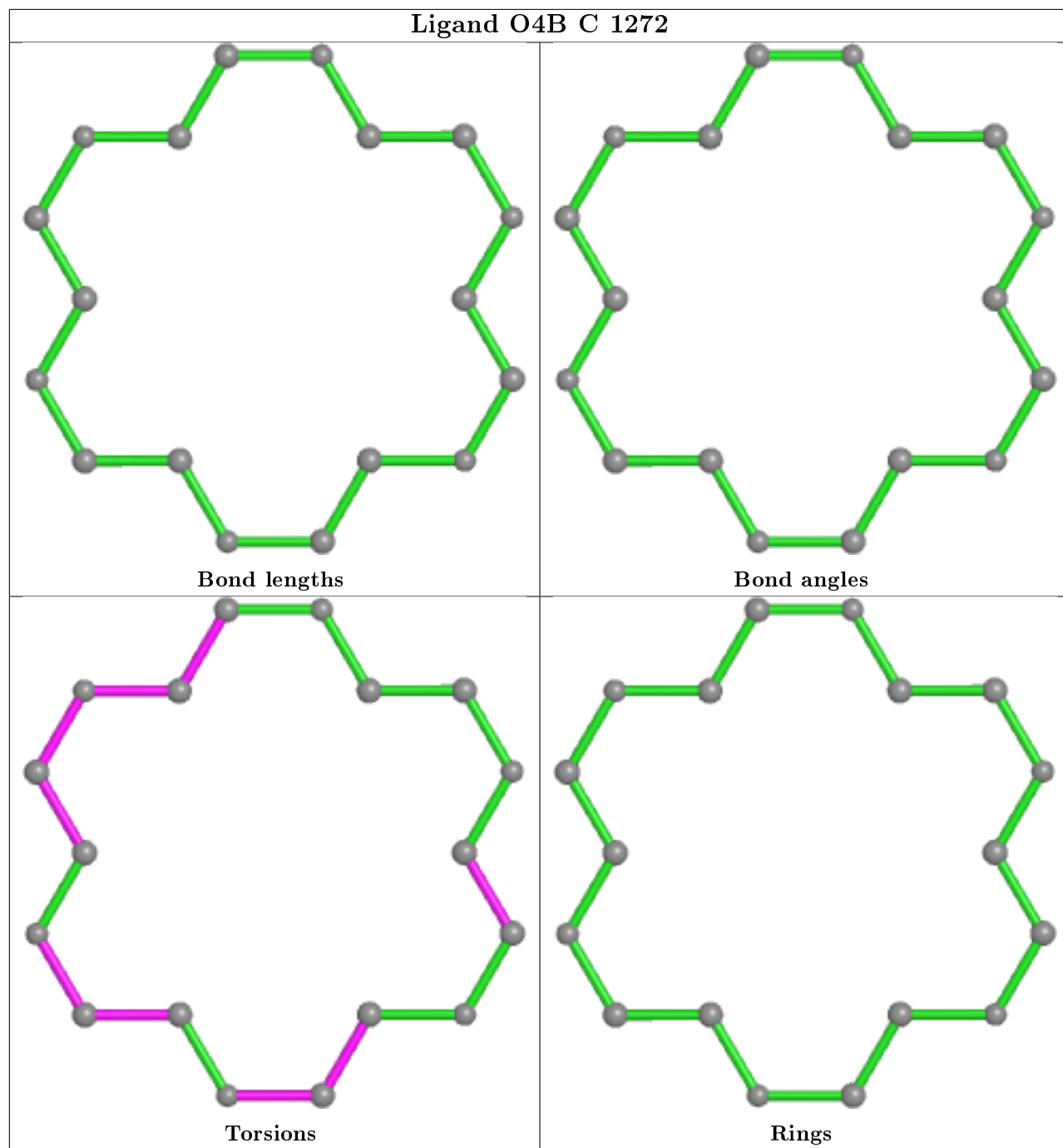
Mol	Chain	Res	Type	Atoms
2	E	1272	O4B	CAA-CAB-CAC-CAD-CAE-CAF-CAG-CAH-CAI-CAJ-CAK-CAL-OAM-O

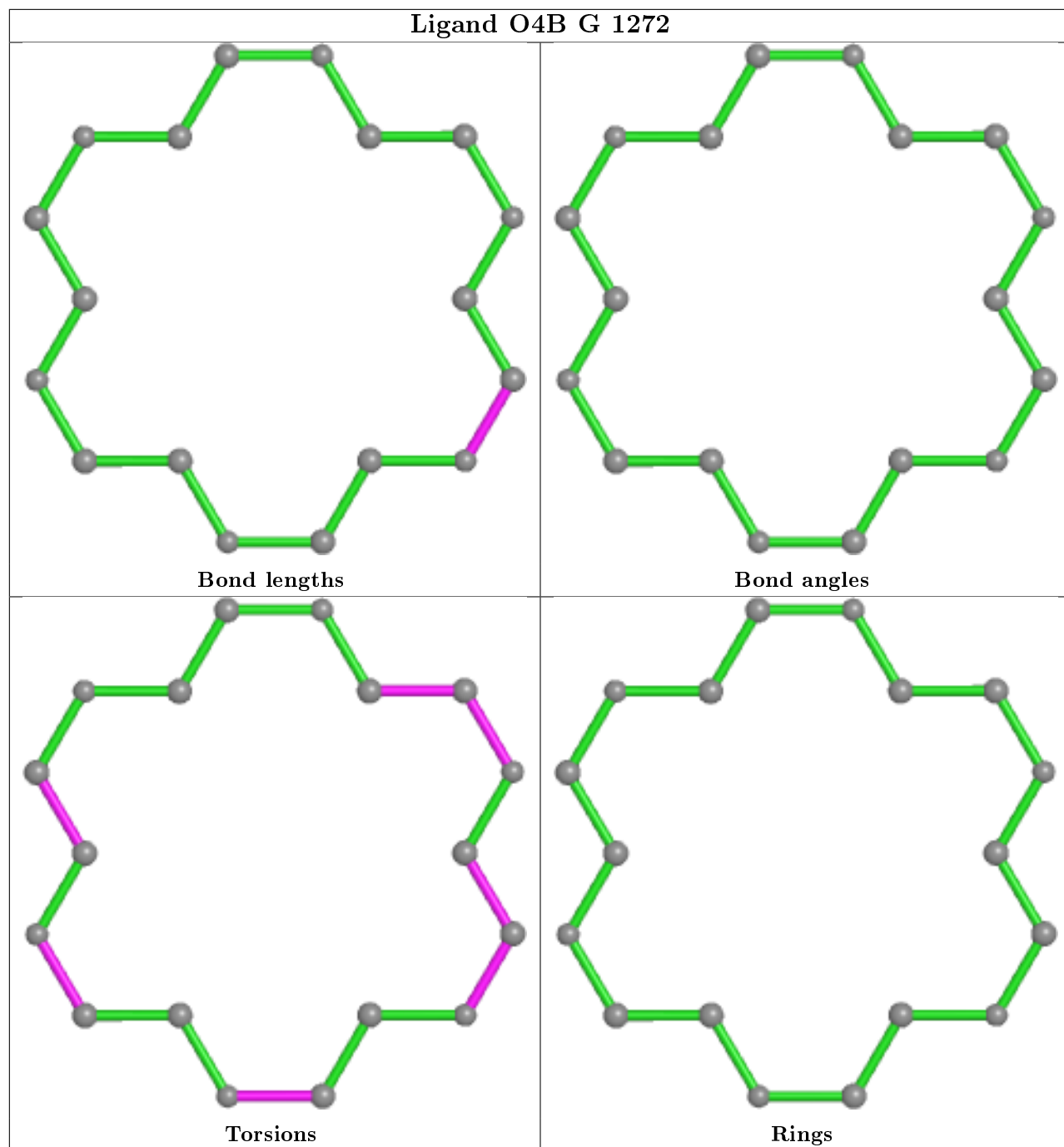
5 monomers are involved in 23 short contacts:

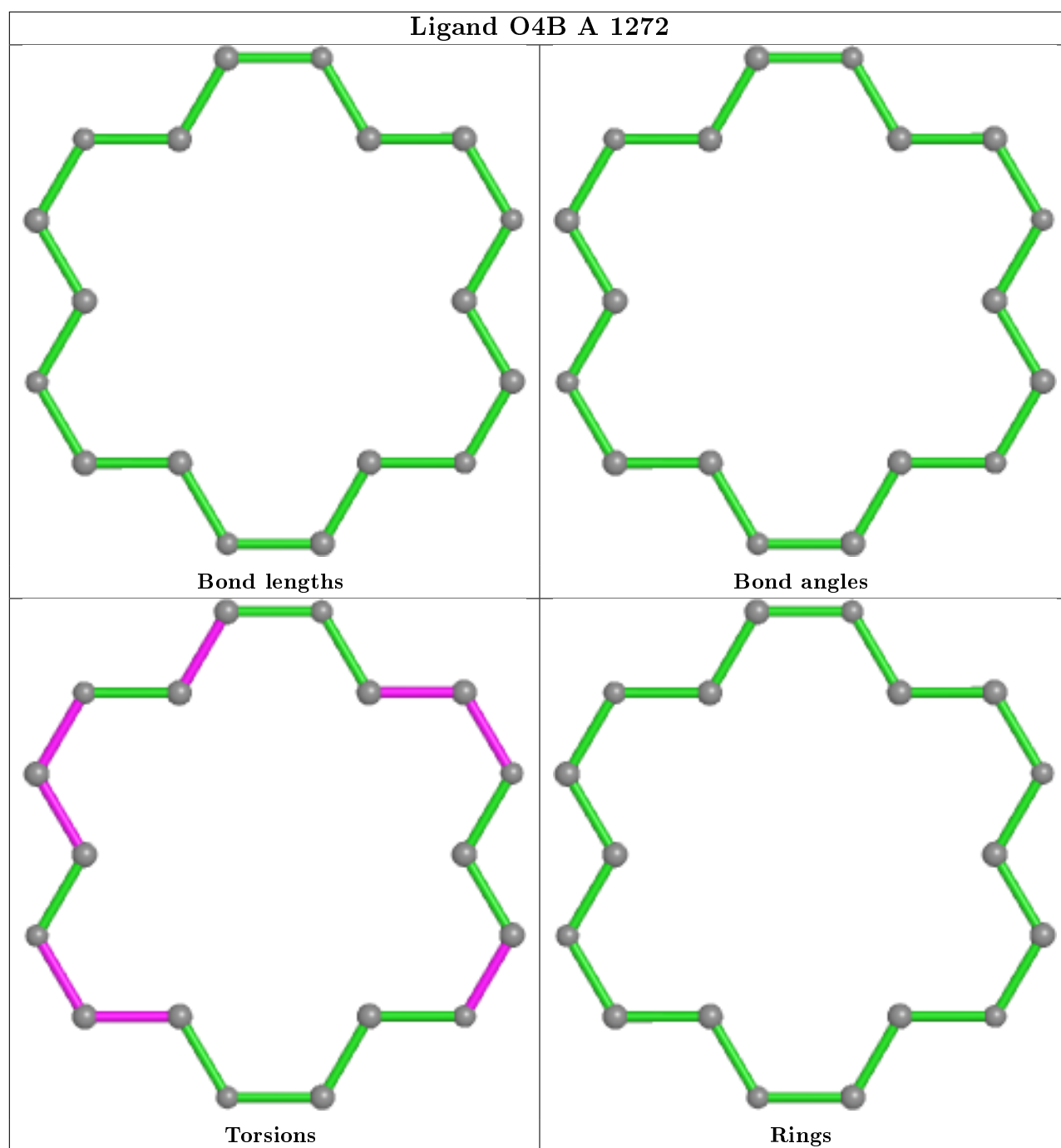
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1272	O4B	10	0
2	C	1272	O4B	2	0
2	A	1272	O4B	9	0
2	D	1272	O4B	1	0
2	F	1272	O4B	1	0

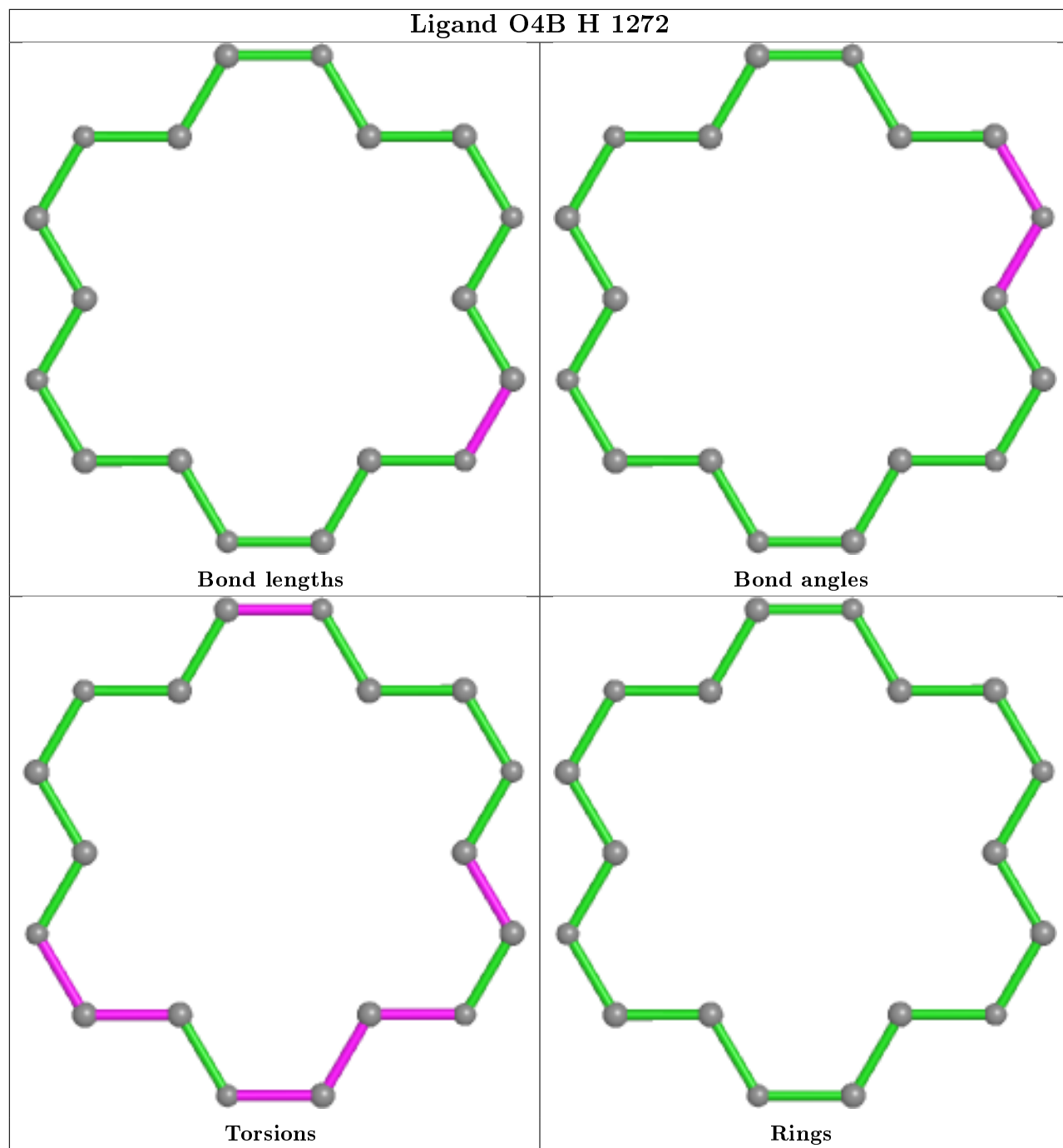
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

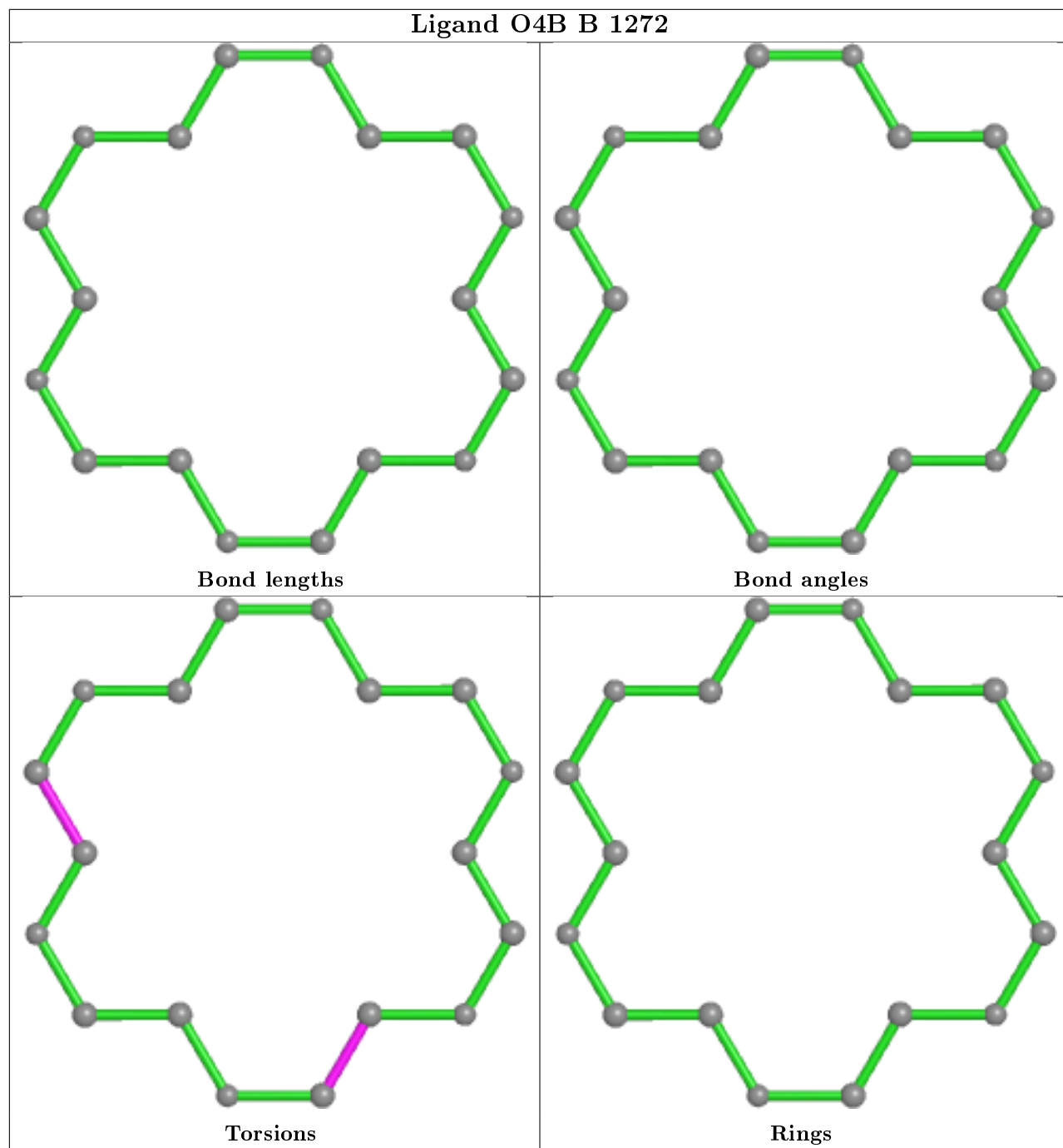


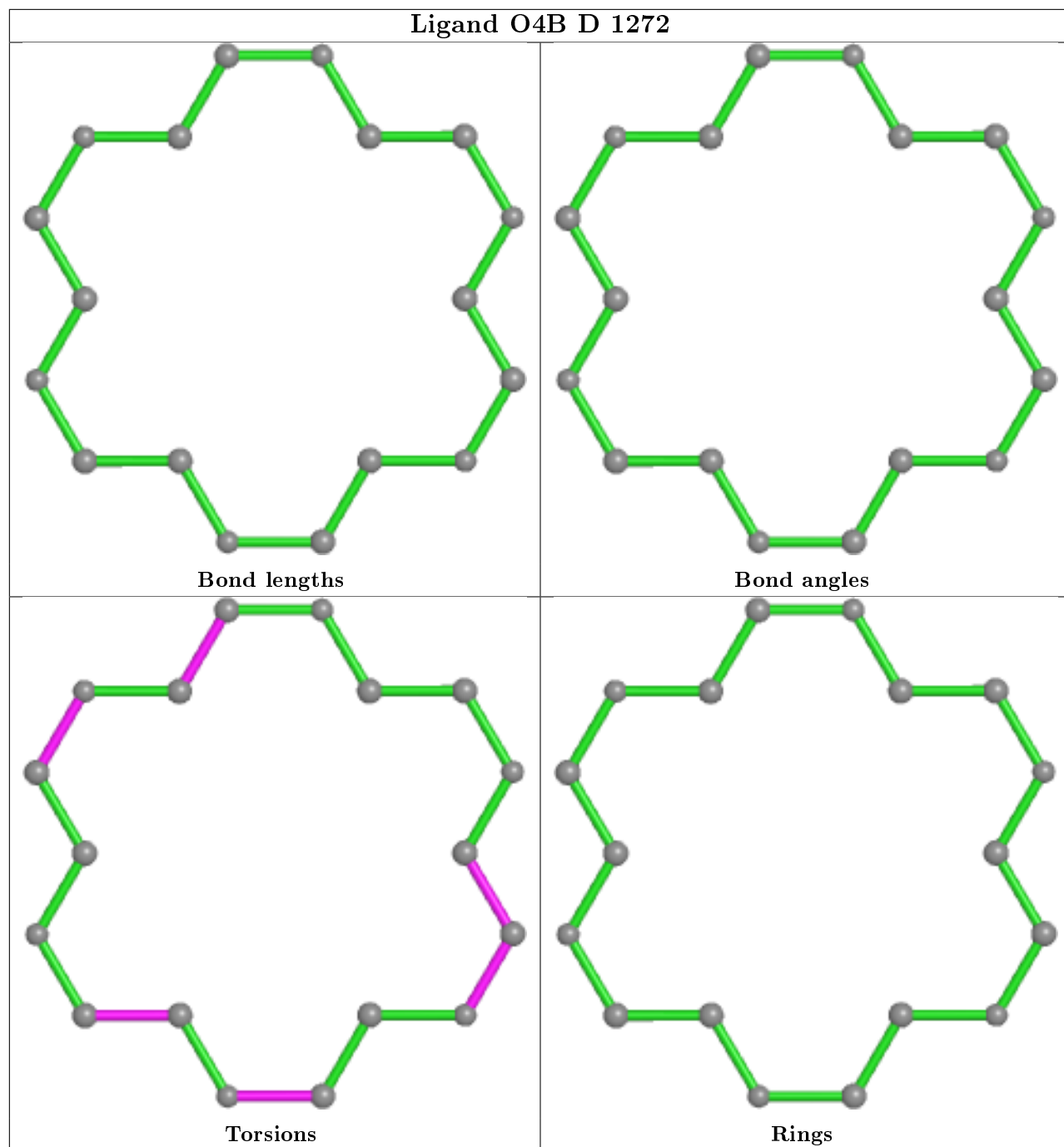


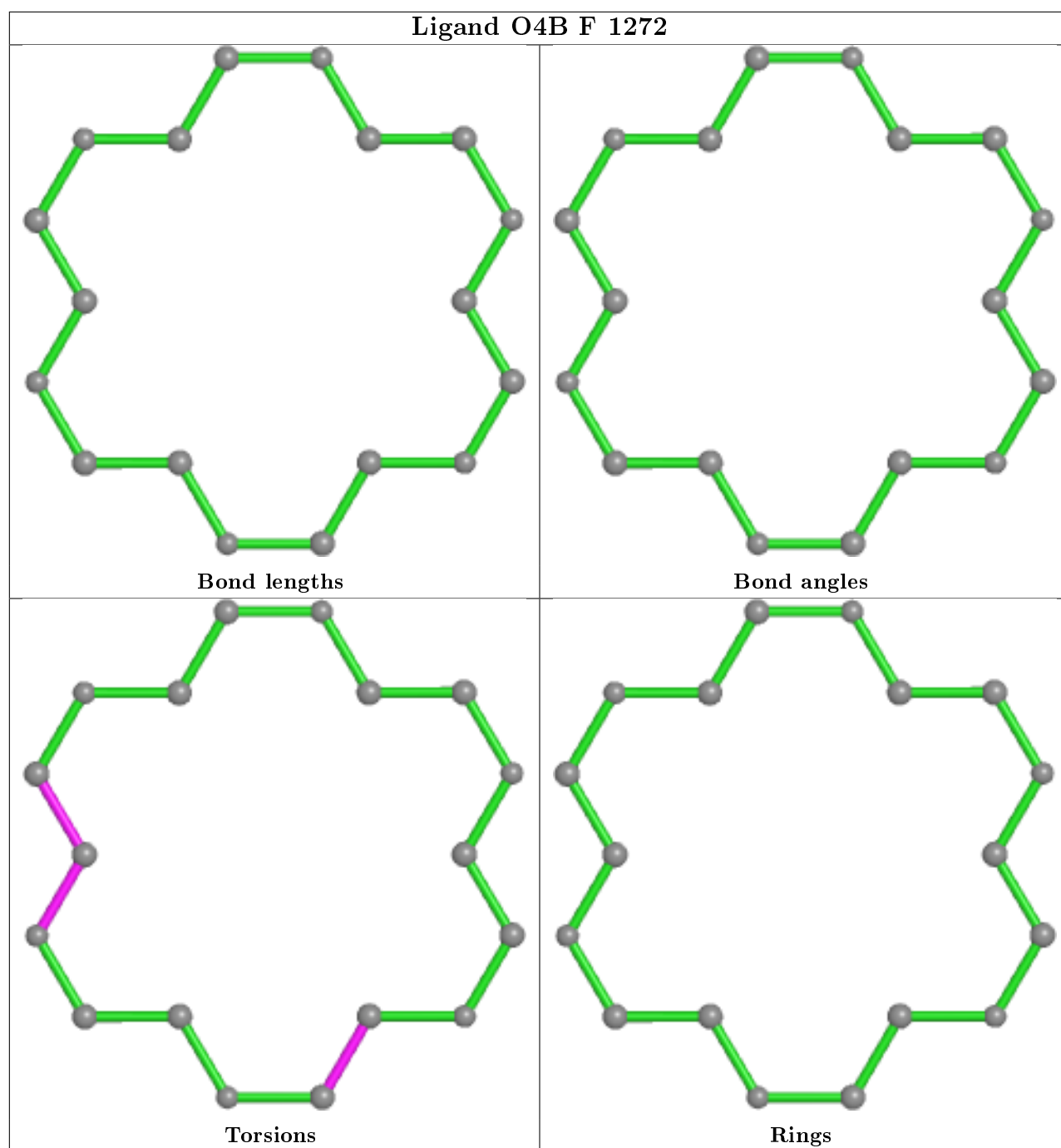












5.7 Other polymers [i](#)

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	A	233/262 (88%)	-0.06	1 (0%) 92 91	32, 54, 76, 89	7 (3%)
1	B	233/262 (88%)	0.06	3 (1%) 77 73	36, 57, 77, 87	10 (4%)
1	C	233/262 (88%)	0.04	4 (1%) 70 66	37, 55, 77, 96	11 (4%)
1	D	234/262 (89%)	-0.07	2 (0%) 84 82	31, 53, 80, 96	7 (2%)
1	E	233/262 (88%)	-0.08	4 (1%) 70 66	29, 53, 79, 96	7 (3%)
1	F	233/262 (88%)	0.03	5 (2%) 63 58	36, 55, 75, 92	12 (5%)
1	G	233/262 (88%)	0.03	2 (0%) 84 82	35, 56, 79, 91	8 (3%)
1	H	234/262 (89%)	-0.10	3 (1%) 77 73	31, 53, 77, 86	9 (3%)
All	All	1866/2096 (89%)	-0.02	24 (1%) 77 73	29, 55, 78, 96	71 (3%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	72	VAL	5.5
1	C	72	VAL	4.6
1	D	71	THR	4.3
1	B	72	VAL	4.3
1	F	105	ALA	3.5
1	E	71	THR	3.3
1	G	72	VAL	3.2
1	C	69	CYS	3.0
1	A	71	THR	2.6
1	F	107	THR	2.6
1	G	107	THR	2.5
1	H	38	PRO	2.5
1	H	67	GLU	2.4
1	C	105	ALA	2.3
1	B	105	ALA	2.3
1	B	109	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	67	GLU	2.2
1	E	157	SER	2.2
1	F	172	PHE	2.2
1	H	155	LEU	2.2
1	F	69	CYS	2.1
1	E	123	TYR	2.1
1	E	156	ASN	2.1
1	D	69	CYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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.

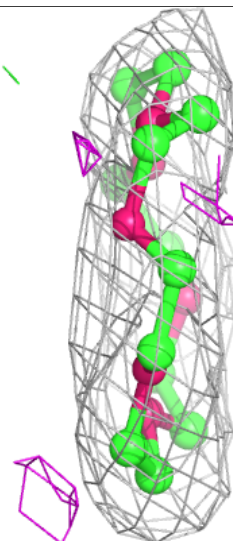
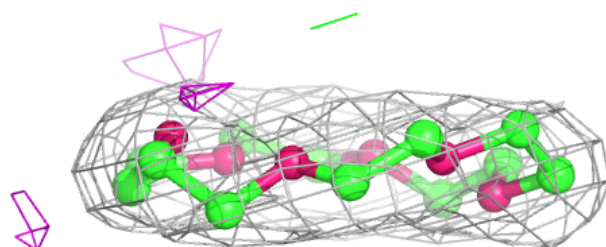
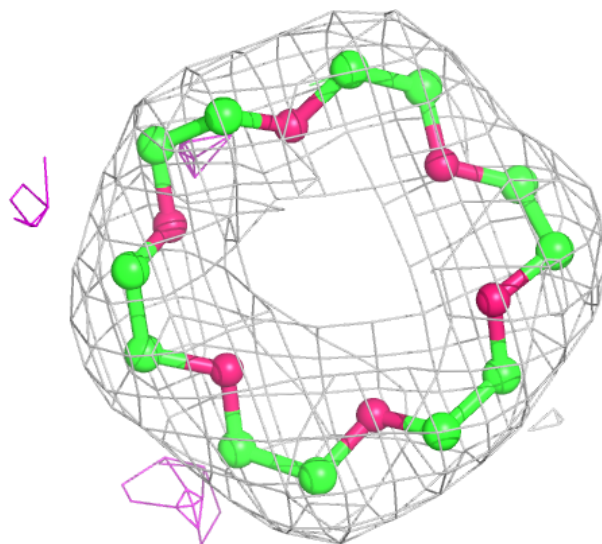
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PEG	D	1273	7/7	0.78	0.22	73,82,97,98	0
3	PEG	F	1273	7/7	0.82	0.27	78,81,87,89	0
3	PEG	E	1273	7/7	0.83	0.23	69,70,73,73	0
3	PEG	A	1273	7/7	0.85	0.28	37,40,41,42	7
2	O4B	F	1272	18/18	0.93	0.15	31,48,52,52	0
2	O4B	G	1272	18/18	0.94	0.19	38,47,50,51	0
2	O4B	C	1272	18/18	0.95	0.17	39,45,50,50	0
2	O4B	B	1272	18/18	0.95	0.19	38,46,50,51	0
2	O4B	A	1272	18/18	0.95	0.13	32,43,50,51	0
2	O4B	H	1272	18/18	0.95	0.15	37,40,44,44	0
2	O4B	D	1272	18/18	0.96	0.17	34,42,48,50	0
2	O4B	E	1272	18/18	0.96	0.14	37,41,48,49	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.

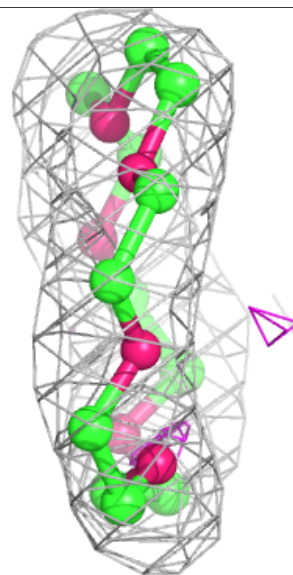
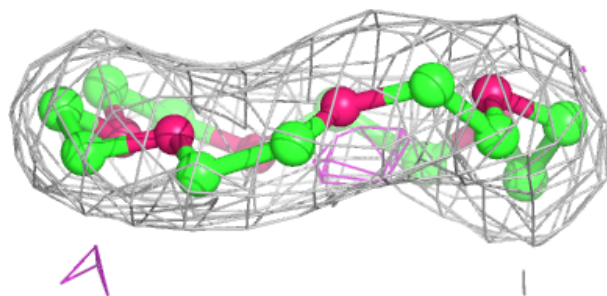
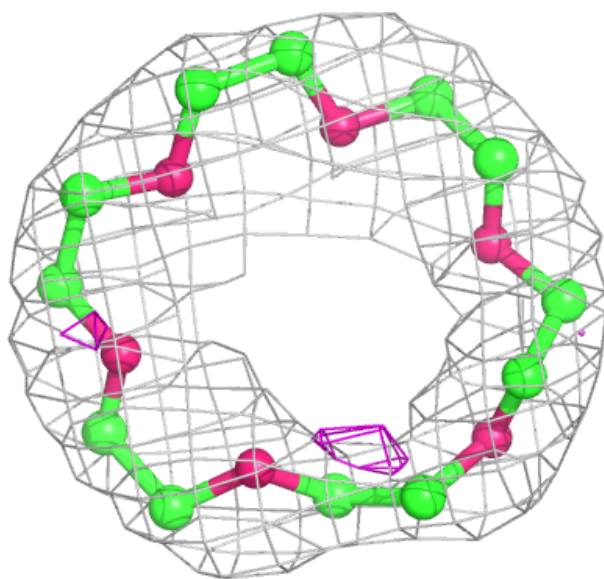
Electron density around O4B F 1272:

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



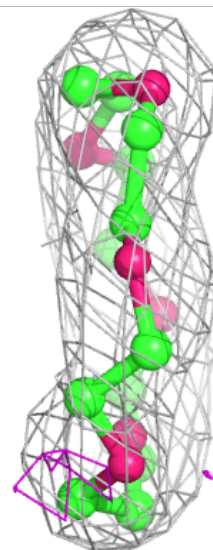
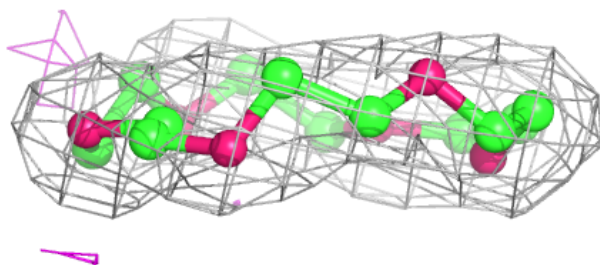
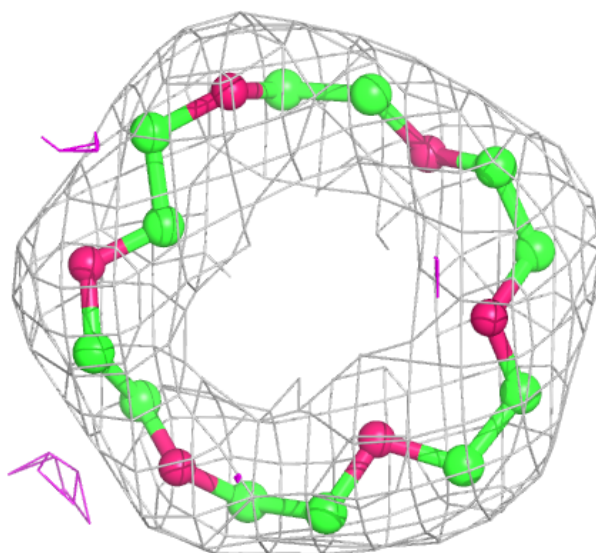
Electron density around O4B G 1272:

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



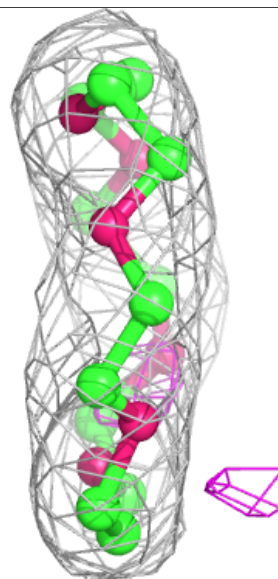
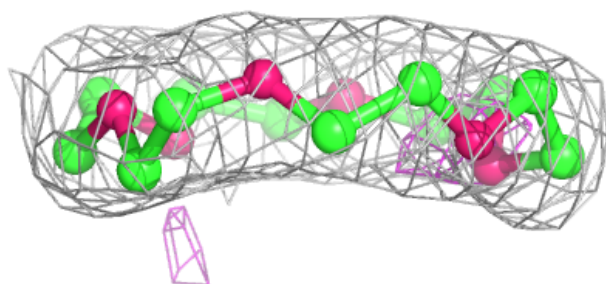
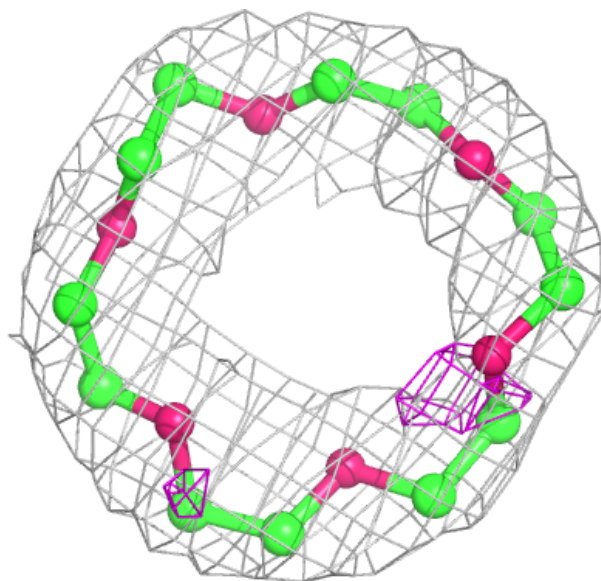
Electron density around O4B C 1272:

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



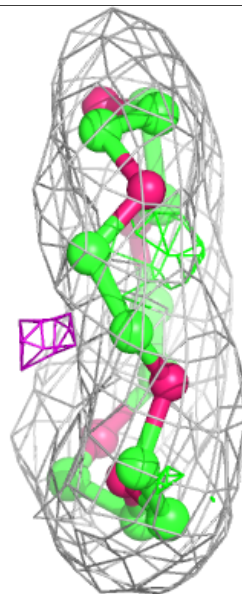
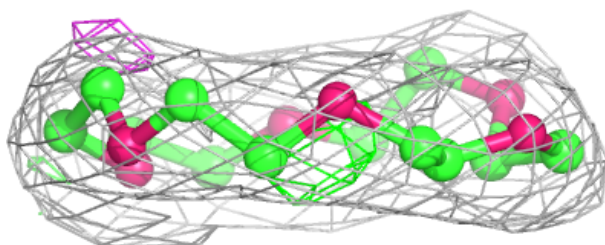
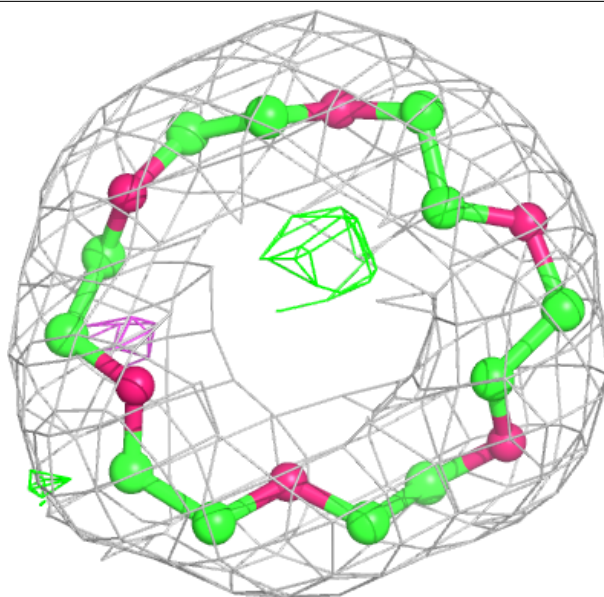
Electron density around O4B B 1272:

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



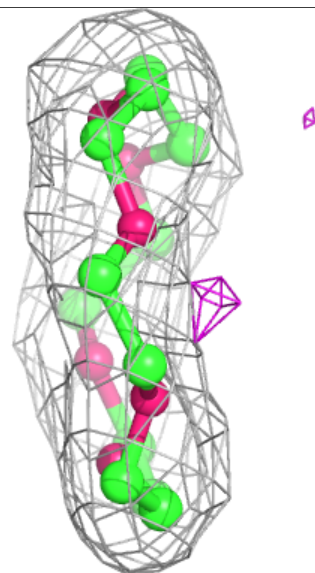
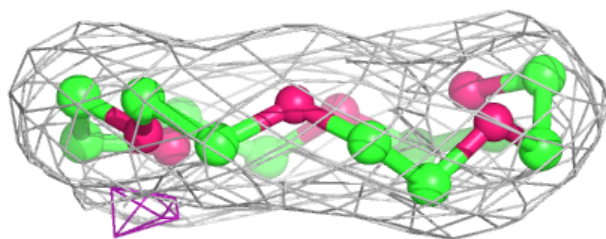
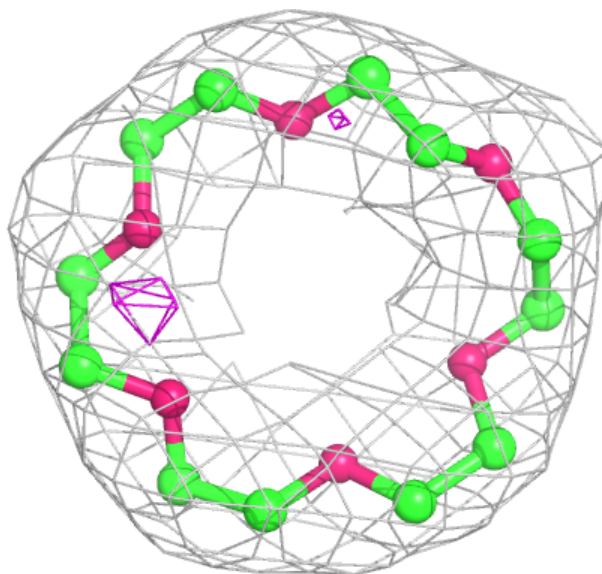
Electron density around O4B A 1272:

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



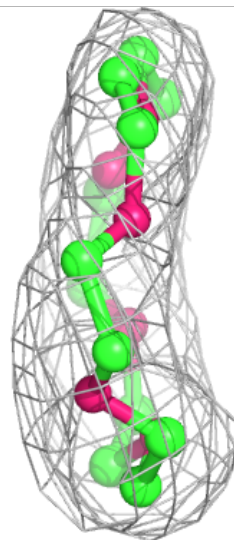
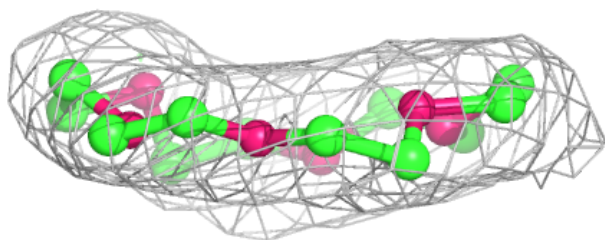
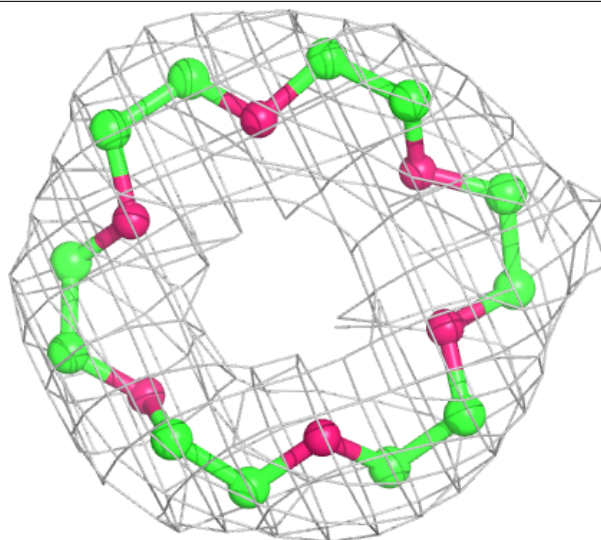
Electron density around O4B H 1272:

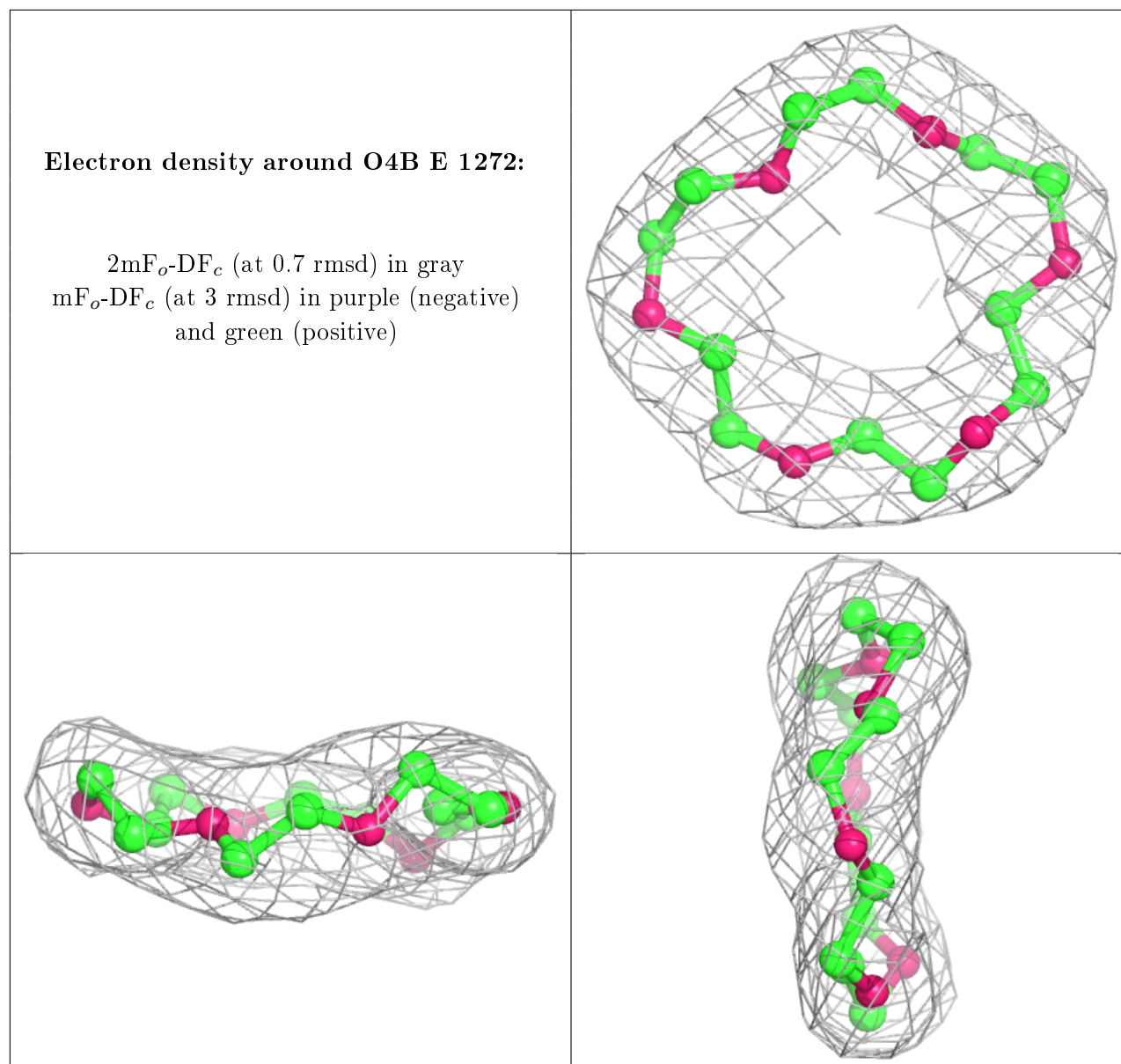
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around O4B D 1272:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.