



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 13, 2024 – 12:16 AM JST

PDB ID : 7YO6
Title : Bifunctional xylosidase/glucosidase LXYL with intermediate substrate xylose for 1.5min
Authors : Yang, L.Y.
Deposited on : 2022-08-01
Resolution : 2.54 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

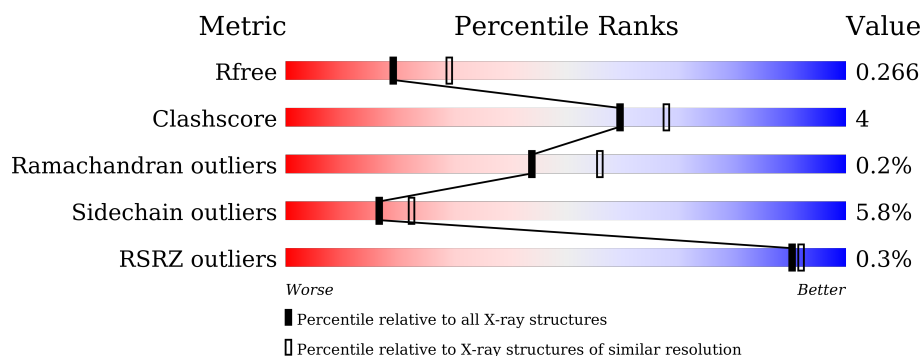
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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.



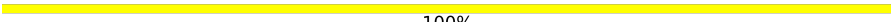
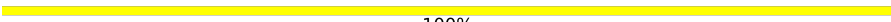

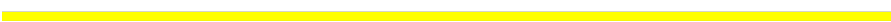
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	1004 (2.54-2.54)
Clashscore	180529	1055 (2.54-2.54)
Ramachandran outliers	177936	1048 (2.54-2.54)
Sidechain outliers	177891	1048 (2.54-2.54)
RSRZ outliers	164620	1004 (2.54-2.54)

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	A	803	
1	B	803	
2	C	4	
3	D	2	
3	G	2	
3	H	2	

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Mol	Chain	Length	Quality of chain
3	I	2	 100%
3	J	2	 100%
4	E	2	 100%
4	F	2	 100%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 12617 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 Beta-D-xylosidase/beta-D-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	756	Total	C	N	O	S	0	0	0
			5706	3615	954	1120	17			
1	B	757	Total	C	N	O	S	0	0	0
			5713	3619	955	1122	17			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-1)-alpha-D-mannopyranose-(3-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	0	0	0
			53	30	3	20			

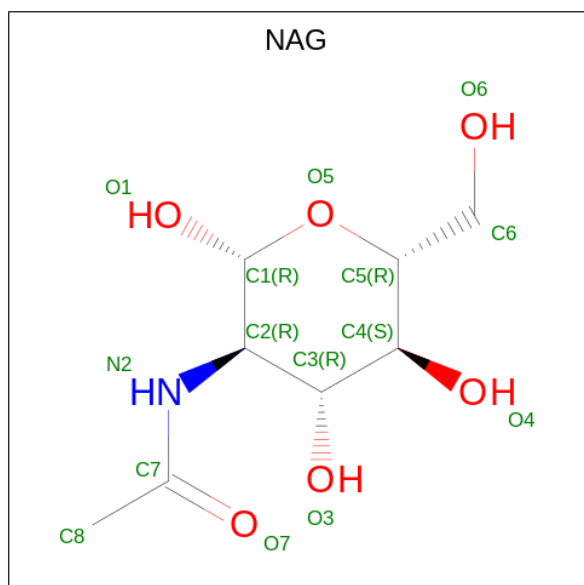
- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-alpha-D-mannopyranose.

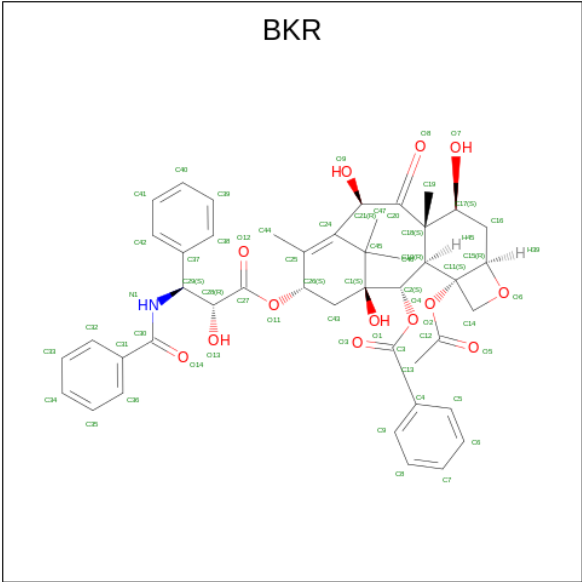
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	2	Total	C	N	O	0	0	0
			25	14	1	10			
4	F	2	Total	C	N	O	0	0	0
			26	14	1	11			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



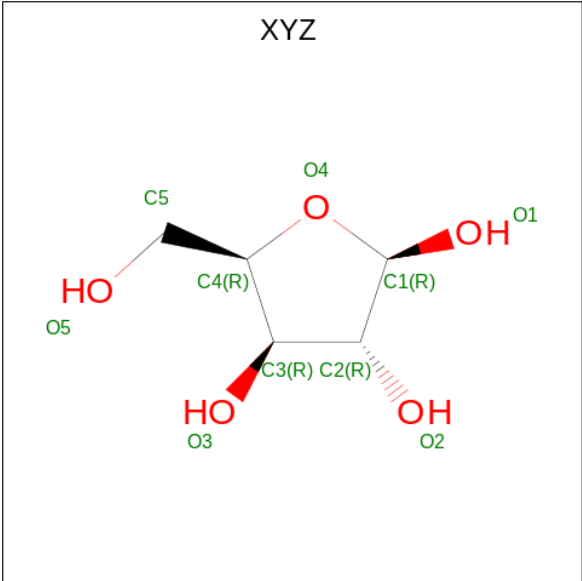
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			15	8	1	6		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is Deacetyltaxol (three-letter code: BKR) (formula: $C_{45}H_{49}NO_{13}$).



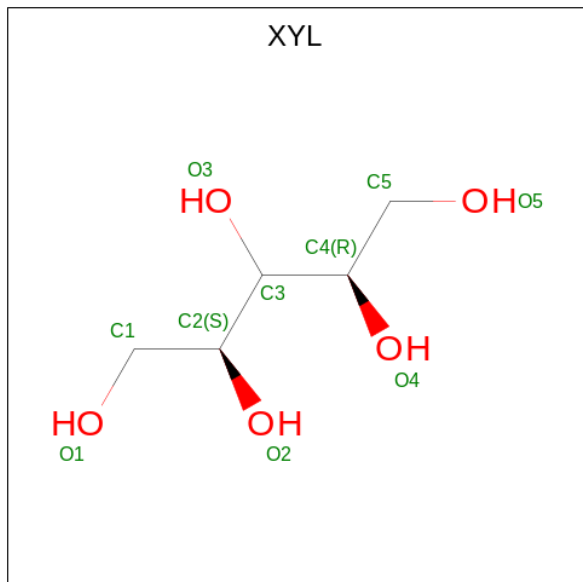
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			59	45	1	13		
6	B	1	Total	C	N	O	0	0
			59	45	1	13		

- Molecule 7 is beta-D-xylofuranose (three-letter code: XYZ) (formula: C₅H₁₀O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			10	5	5		

- Molecule 8 is Xylitol (three-letter code: XYL) (formula: $C_5H_{12}O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			10	5	5		

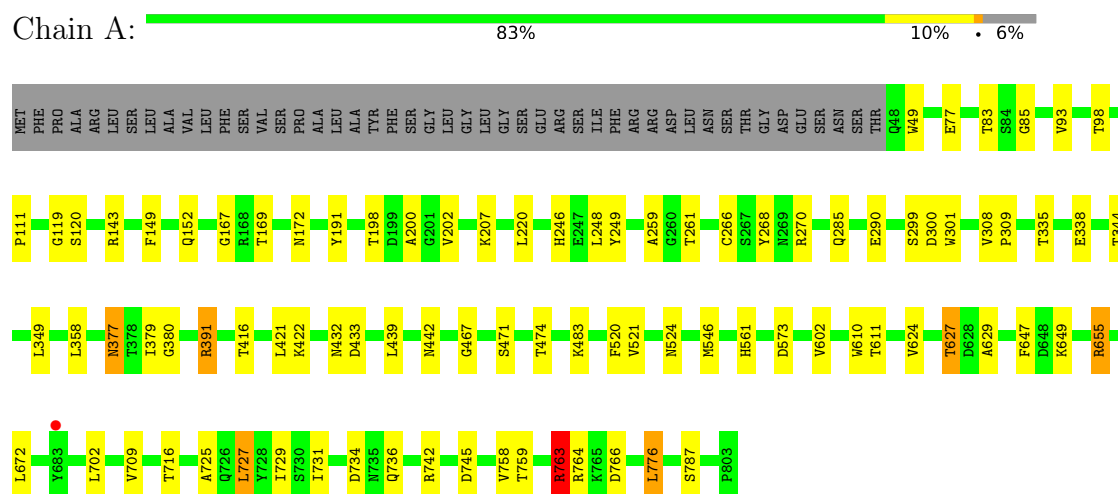
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	354	Total	O	0	0
			354	354		
9	B	377	Total	O	0	0
			377	377		

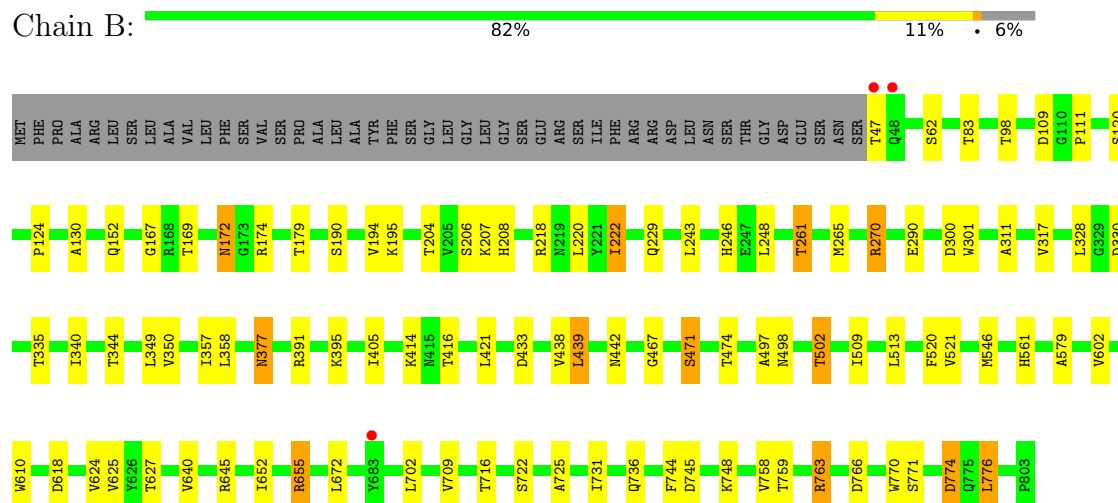
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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: Beta-D-xylosidase/beta-D-glucosidase



- Molecule 1: Beta-D-xylosidase/beta-D-glucosidase



- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-1)-alpha-D-mannopyranose-(3-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  100%



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-alpha-D-mannopyranose

Chain E:  100%



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-alpha-D-mannopyranose

Chain F:

100%

HAH1
HAG2

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	79.59Å 181.92Å 241.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	120.98 – 2.54 120.98 – 2.54	Depositor EDS
% Data completeness (in resolution range)	99.2 (120.98-2.54) 99.3 (120.98-2.54)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.53Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.191 , 0.265 0.198 , 0.266	Depositor DCC
R_{free} test set	2825 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12617	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 54.75 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.4922e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BKR, NAG, XYZ, MAN, XYL

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	0.65	0/5847	0.83	5/8000 (0.1%)
1	B	0.68	0/5854	0.84	10/8010 (0.1%)
All	All	0.66	0/11701	0.84	15/16010 (0.1%)

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	391	ARG	NE-CZ-NH1	7.57	124.09	120.30
1	A	763	ARG	NE-CZ-NH1	7.40	124.00	120.30
1	A	391	ARG	NE-CZ-NH1	6.86	123.73	120.30
1	B	270	ARG	NE-CZ-NH1	6.65	123.63	120.30
1	B	391	ARG	NE-CZ-NH2	-6.52	117.04	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5706	0	5494	35	0
1	B	5713	0	5502	55	0
2	C	53	0	46	2	0
3	D	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	28	0	25	0	0
3	H	28	0	25	0	0
3	I	28	0	25	0	0
3	J	28	0	25	0	0
4	E	25	0	22	3	0
4	F	26	0	24	0	0
5	A	57	0	54	2	0
5	B	28	0	26	2	0
6	A	59	0	0	1	0
6	B	59	0	0	5	0
7	A	10	0	7	2	0
8	B	10	0	12	3	0
9	A	354	0	0	1	0
9	B	377	0	0	7	0
All	All	12617	0	11312	97	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.

The worst 5 of 97 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:C:3:MAN:O4	4:E:1:MAN:C1	2.16	0.93
1:B:300:ASP:OD1	8:B:904:XYL:O4	1.94	0.83
1:A:300:ASP:OD1	7:A:906:XYZ:O2	1.95	0.82
1:A:246:HIS:HE1	1:A:290:GLU:OE2	1.76	0.68
1:B:438:VAL:HG23	1:B:439:LEU:HD13	1.75	0.67

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	754/803 (94%)	713 (95%)	39 (5%)	2 (0%)	37	46
1	B	755/803 (94%)	712 (94%)	42 (6%)	1 (0%)	48	61
All	All	1509/1606 (94%)	1425 (94%)	81 (5%)	3 (0%)	44	56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	167	GLY
1	A	422	LYS
1	A	167	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	A	607/648 (94%)	572 (94%)	35 (6%)	17	22
1	B	608/648 (94%)	573 (94%)	35 (6%)	17	22
All	All	1215/1296 (94%)	1145 (94%)	70 (6%)	17	22

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	602	VAL
1	B	624	VAL
1	B	709	VAL
1	A	655	ARG
1	A	649	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	377	ASN
1	B	676	HIS
1	B	740	GLN

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Mol	Chain	Res	Type
1	B	561	HIS
1	A	561	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

18 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.50	0	17,19,21	1.14	0
2	NAG	C	2	2	14,14,15	0.81	0	17,19,21	2.11	5 (29%)
2	MAN	C	3	2	11,11,12	1.08	1 (9%)	13,15,17	1.43	2 (15%)
2	NAG	C	4	2	14,14,15	1.25	2 (14%)	17,19,21	2.09	6 (35%)
3	NAG	D	1	3,1	14,14,15	0.55	0	17,19,21	1.41	3 (17%)
3	NAG	D	2	3	14,14,15	1.10	2 (14%)	17,19,21	1.91	5 (29%)
4	MAN	E	1	4	11,11,12	0.96	0	15,15,17	1.84	4 (26%)
4	NAG	E	2	4	14,14,15	0.69	0	17,19,21	2.27	7 (41%)
4	MAN	F	1	4	12,12,12	1.37	1 (8%)	17,17,17	1.84	5 (29%)
4	NAG	F	2	4	14,14,15	1.11	1 (7%)	17,19,21	2.39	6 (35%)
3	NAG	G	1	3,1	14,14,15	0.72	0	17,19,21	1.25	2 (11%)
3	NAG	G	2	3	14,14,15	0.93	0	17,19,21	1.05	1 (5%)
3	NAG	H	1	3,1	14,14,15	0.63	0	17,19,21	1.39	2 (11%)
3	NAG	H	2	3	14,14,15	0.87	0	17,19,21	1.26	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	I	1	3,1	14,14,15	0.46	0	17,19,21	1.47	3 (17%)
3	NAG	I	2	3	14,14,15	0.69	0	17,19,21	1.38	1 (5%)
3	NAG	J	1	3,1	14,14,15	0.48	0	17,19,21	1.33	3 (17%)
3	NAG	J	2	3	14,14,15	0.59	0	17,19,21	1.26	3 (17%)

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	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	MAN	C	3	2	-	2/2/18/22	0/1/1/1
2	NAG	C	4	2	-	2/6/23/26	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
4	MAN	E	1	4	-	0/2/19/22	0/1/1/1
4	NAG	E	2	4	-	4/6/23/26	0/1/1/1
4	MAN	F	1	4	-	0/2/22/22	0/1/1/1
4	NAG	F	2	4	-	3/6/23/26	0/1/1/1
3	NAG	G	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	NAG	I	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	NAG	J	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	2	NAG	C1-C2	3.28	1.57	1.52
4	F	1	MAN	O1-C1	2.63	1.48	1.39
3	D	2	NAG	O4-C4	2.55	1.49	1.43
3	D	2	NAG	O5-C5	2.32	1.48	1.43
2	C	3	MAN	C1-C2	2.08	1.54	1.52

The worst 5 of 61 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	2	NAG	C1-O5-C5	6.48	120.97	112.19
4	F	2	NAG	C2-N2-C7	4.90	129.89	122.90
4	F	2	NAG	C1-O5-C5	4.68	118.53	112.19
2	C	2	NAG	O4-C4-C5	4.64	120.81	109.30
4	F	2	NAG	C1-C2-N2	4.51	118.19	110.49

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	2	NAG	C1-C2-N2-C7
2	C	4	NAG	C4-C5-C6-O6
4	F	2	NAG	C4-C5-C6-O6
2	C	3	MAN	C4-C5-C6-O6
2	C	4	NAG	O5-C5-C6-O6

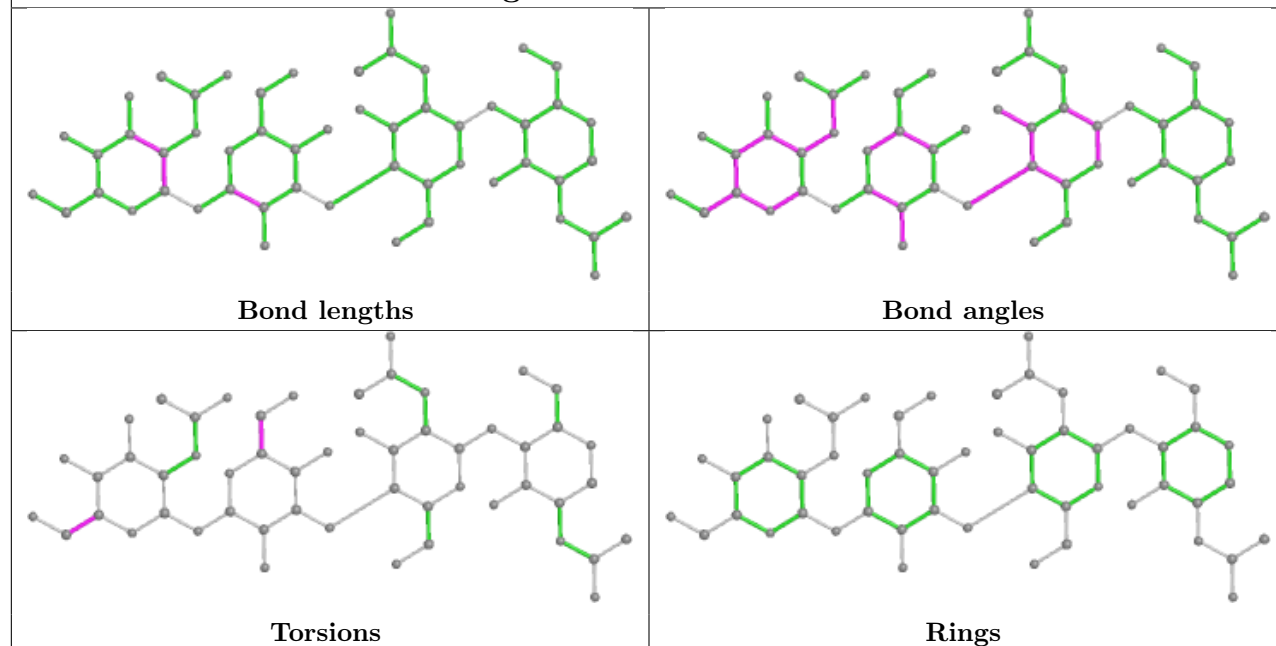
There are no ring outliers.

3 monomers are involved in 3 short contacts:

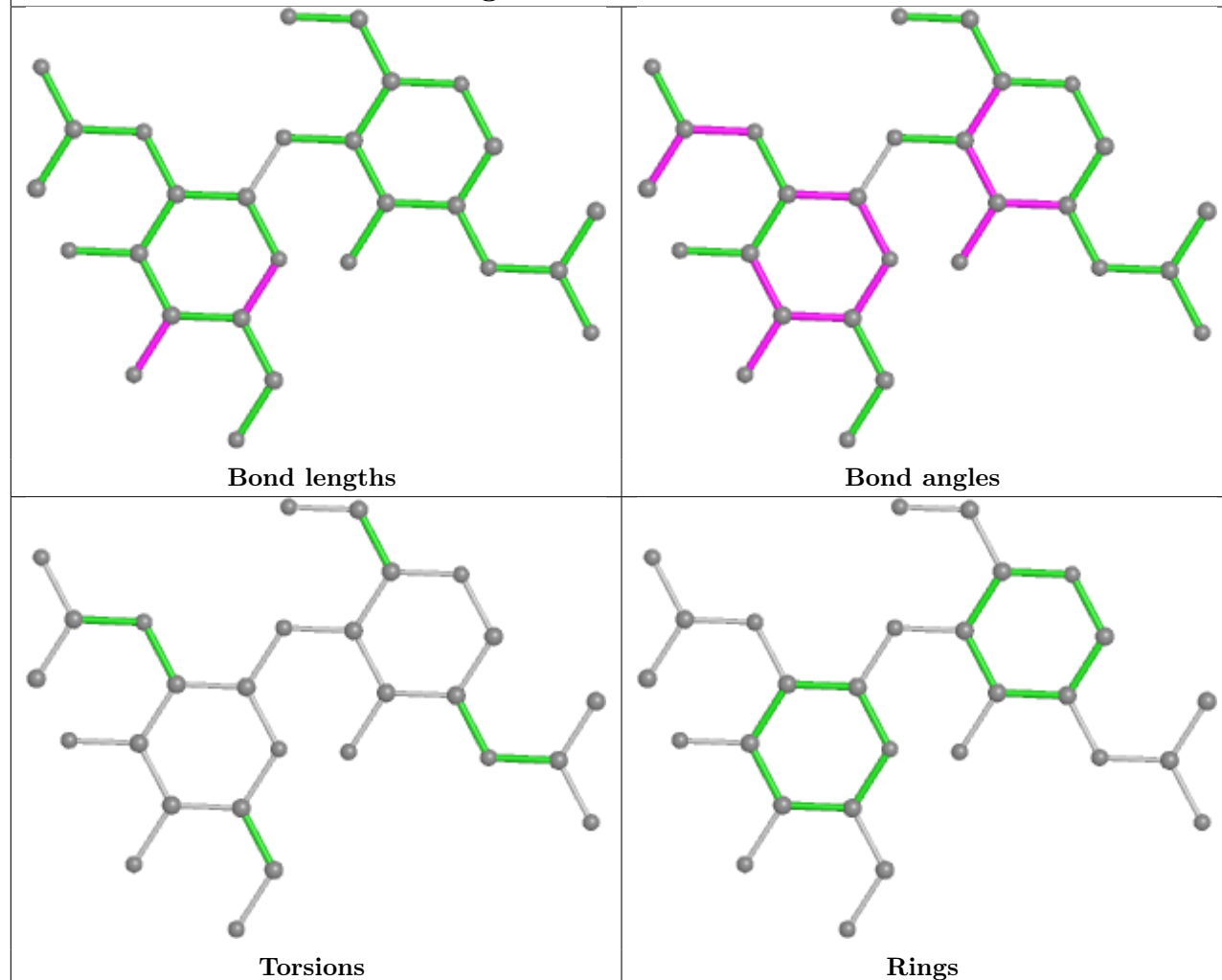
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	1	MAN	2	0
4	E	2	NAG	1	0
2	C	3	MAN	2	0

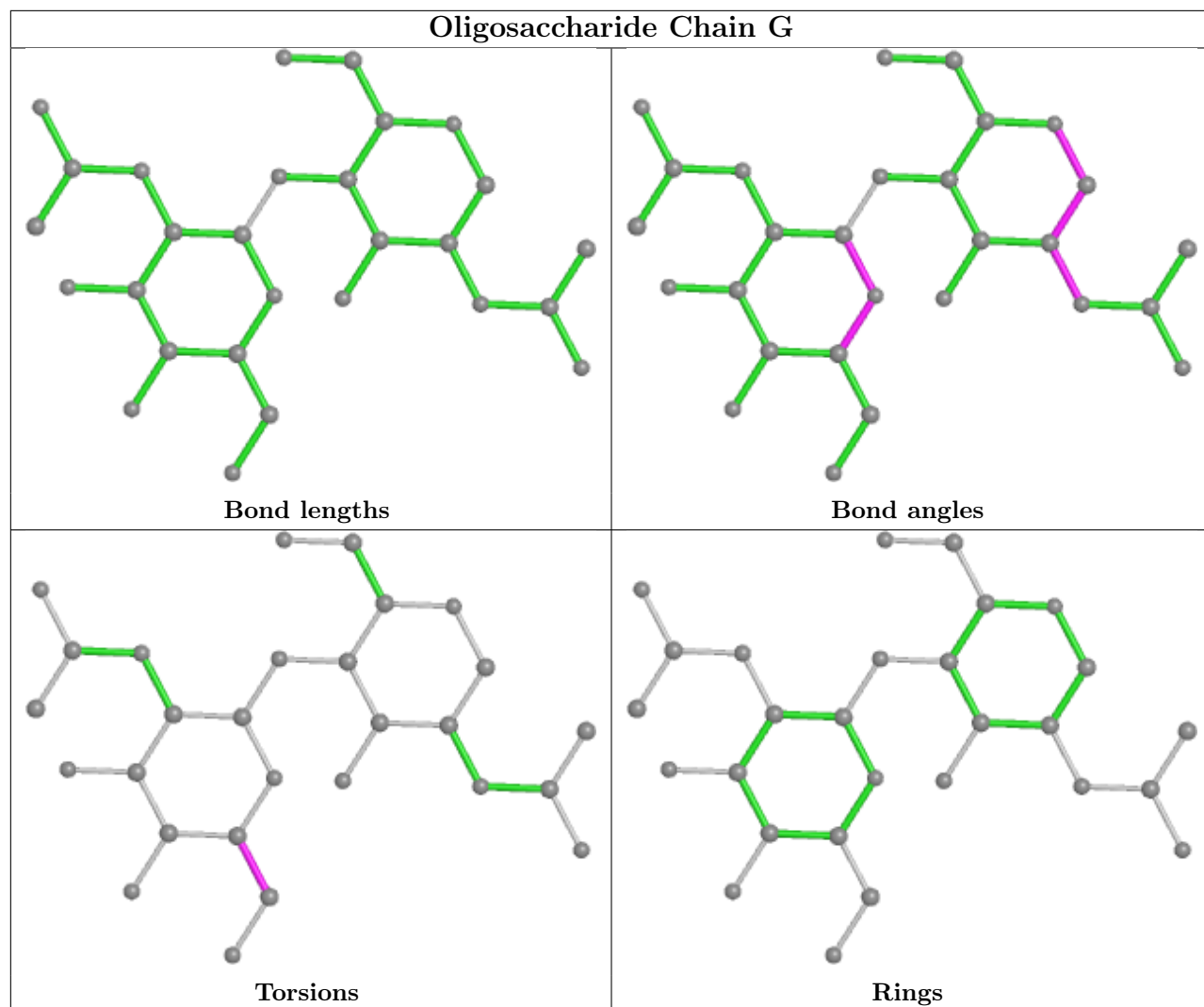
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

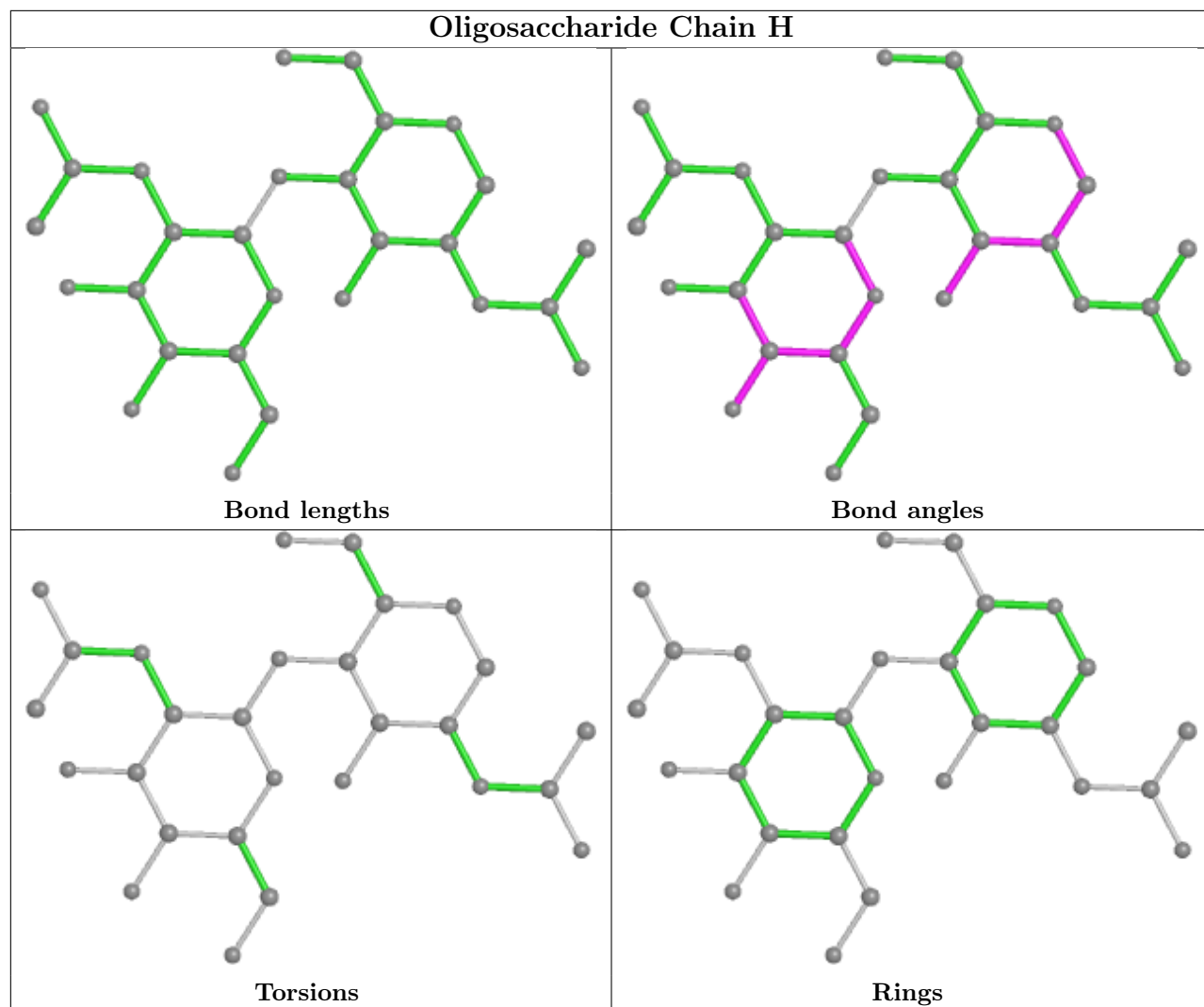
Oligosaccharide Chain C

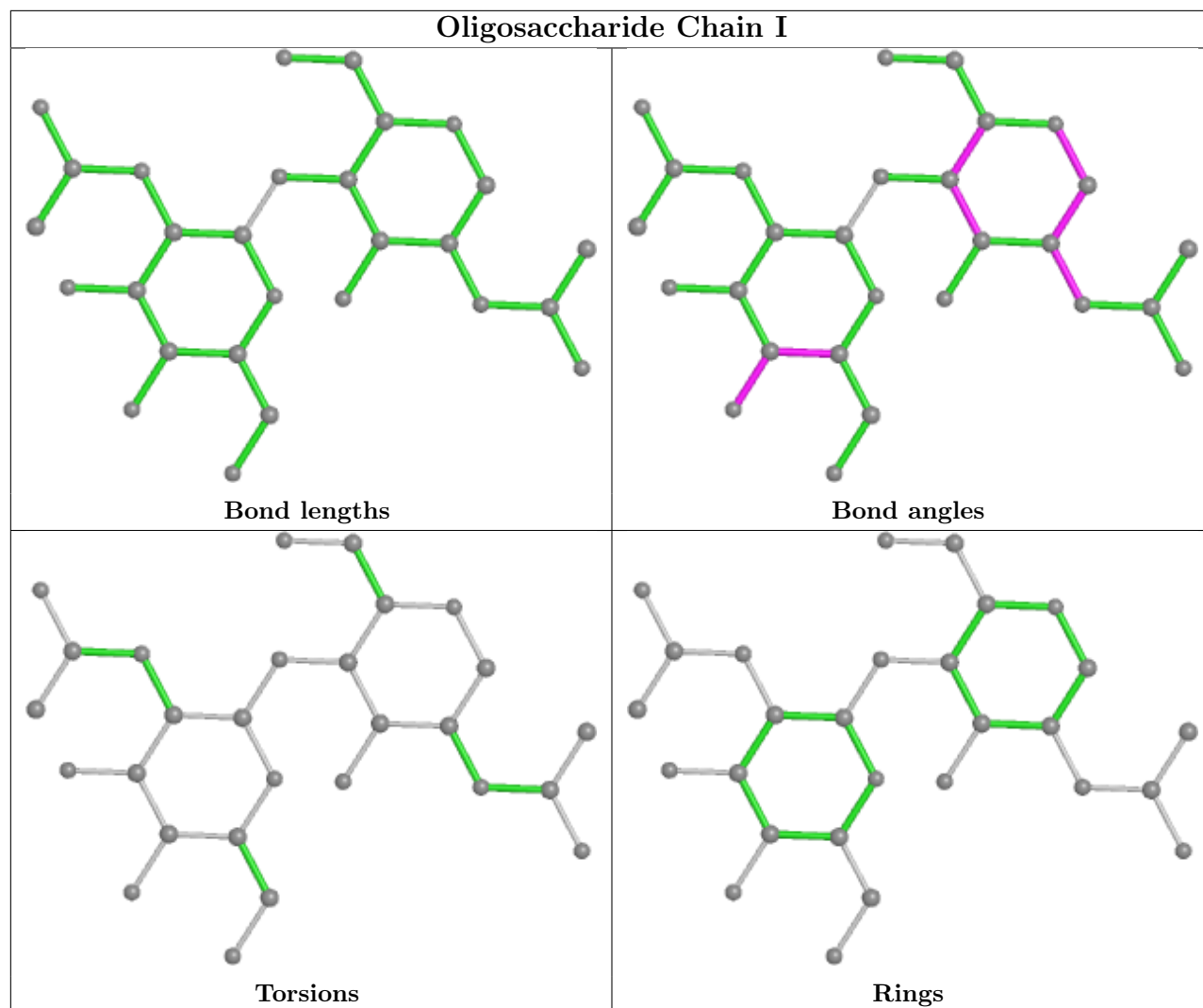


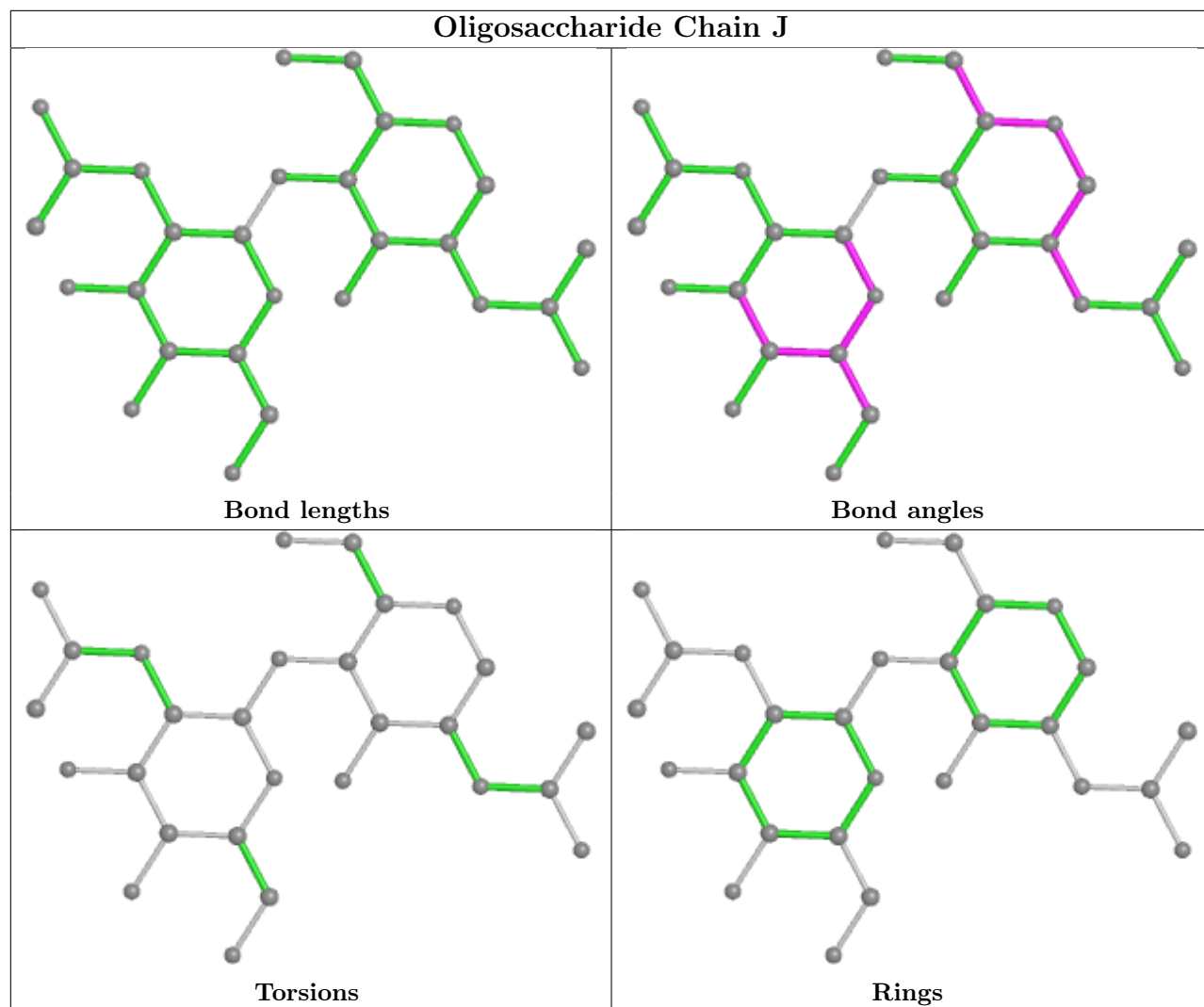
Oligosaccharide Chain D

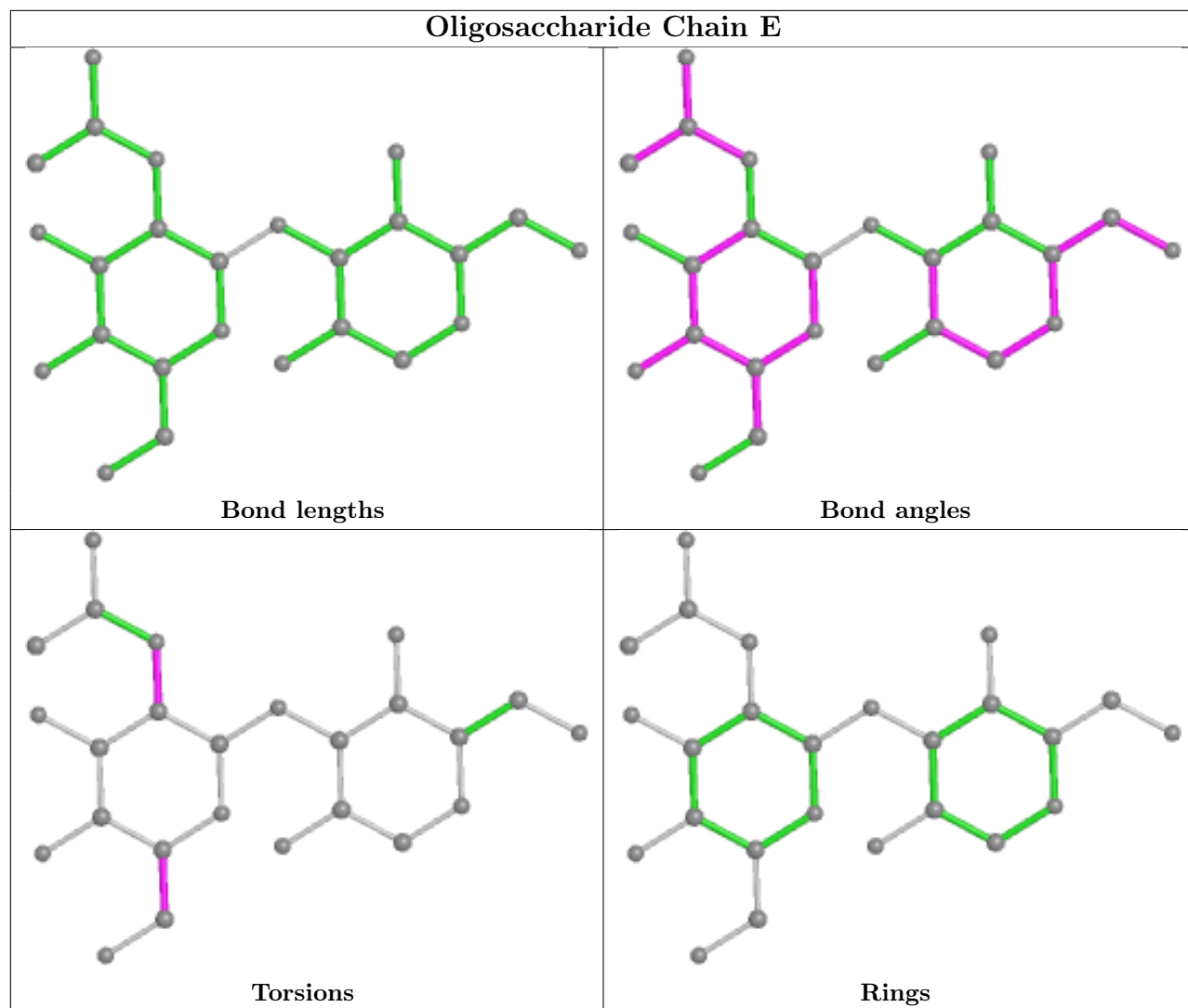


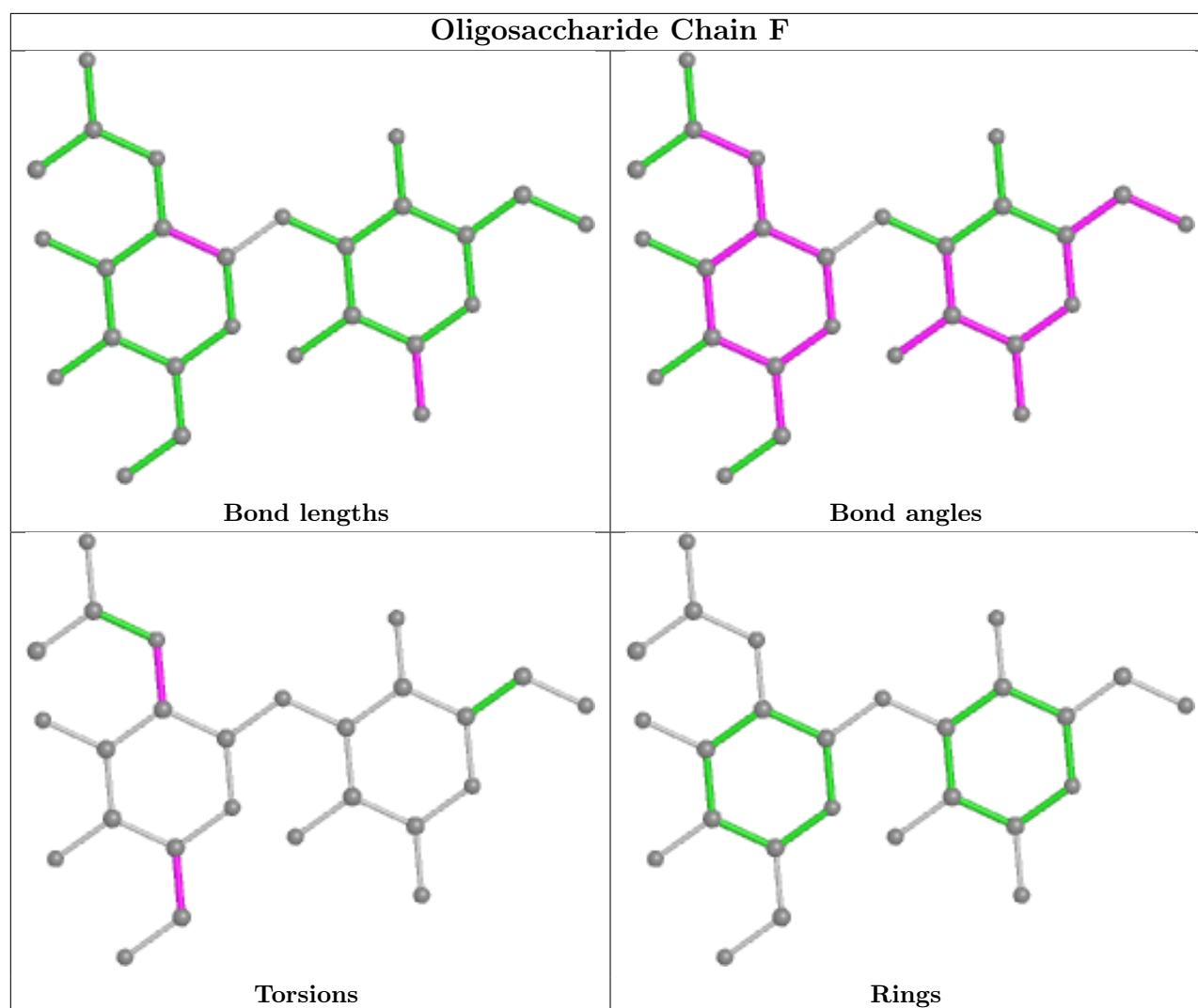












5.6 Ligand geometry [i](#)

10 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	$\# Z > 2$
6	BKR	A	903	-	65,65,65	1.35	5 (7%)	101,101,101	1.40	16 (15%)
5	NAG	A	905	1	14,14,15	0.64	0	17,19,21	1.54	4 (23%)
5	NAG	A	902	1	14,14,15	0.56	0	17,19,21	1.21	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	B	901	-	14,14,15	0.98	0	19,19,21	1.62	2 (10%)
8	XYL	B	904	-	9,9,9	0.27	0	11,11,11	0.56	0
6	BKR	B	903	-	65,65,65	1.37	6 (9%)	101,101,101	1.37	14 (13%)
5	NAG	A	904	-	15,15,15	1.55	2 (13%)	21,21,21	3.95	11 (52%)
5	NAG	B	902	1	14,14,15	0.41	0	17,19,21	1.23	1 (5%)
7	XYZ	A	906	-	10,10,10	0.95	0	13,14,14	3.60	8 (61%)
5	NAG	A	901	-	14,14,15	0.80	0	19,19,21	1.44	3 (15%)

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
6	BKR	A	903	-	-	4/37/123/123	0/7/7/7
5	NAG	A	905	1	-	0/6/23/26	0/1/1/1
5	NAG	A	902	1	-	0/6/23/26	0/1/1/1
5	NAG	B	901	-	-	2/6/22/26	0/1/1/1
8	XYL	B	904	-	-	2/12/12/12	-
6	BKR	B	903	-	-	9/37/123/123	0/7/7/7
5	NAG	A	904	-	-	5/6/26/26	0/1/1/1
5	NAG	B	902	1	-	0/6/23/26	0/1/1/1
7	XYZ	A	906	-	-	2/2/18/18	0/1/1/1
5	NAG	A	901	-	-	1/6/22/26	0/1/1/1

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	903	BKR	O2-C3	5.68	1.46	1.34
6	B	903	BKR	O2-C3	5.25	1.45	1.34
6	B	903	BKR	O11-C27	4.85	1.45	1.34
6	A	903	BKR	O11-C27	4.75	1.45	1.34
6	A	903	BKR	O4-C12	4.09	1.44	1.35

The worst 5 of 61 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	904	NAG	O5-C1-C2	9.55	119.11	109.52
5	A	904	NAG	C1-C2-N2	8.65	120.75	110.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	906	XYZ	C1-C2-C3	8.50	112.94	102.30
5	A	904	NAG	O5-C5-C4	-5.48	99.75	109.69
7	A	906	XYZ	O4-C1-C2	-5.31	97.93	104.46

There are no chirality outliers.

5 of 25 torsion outliers are listed below:

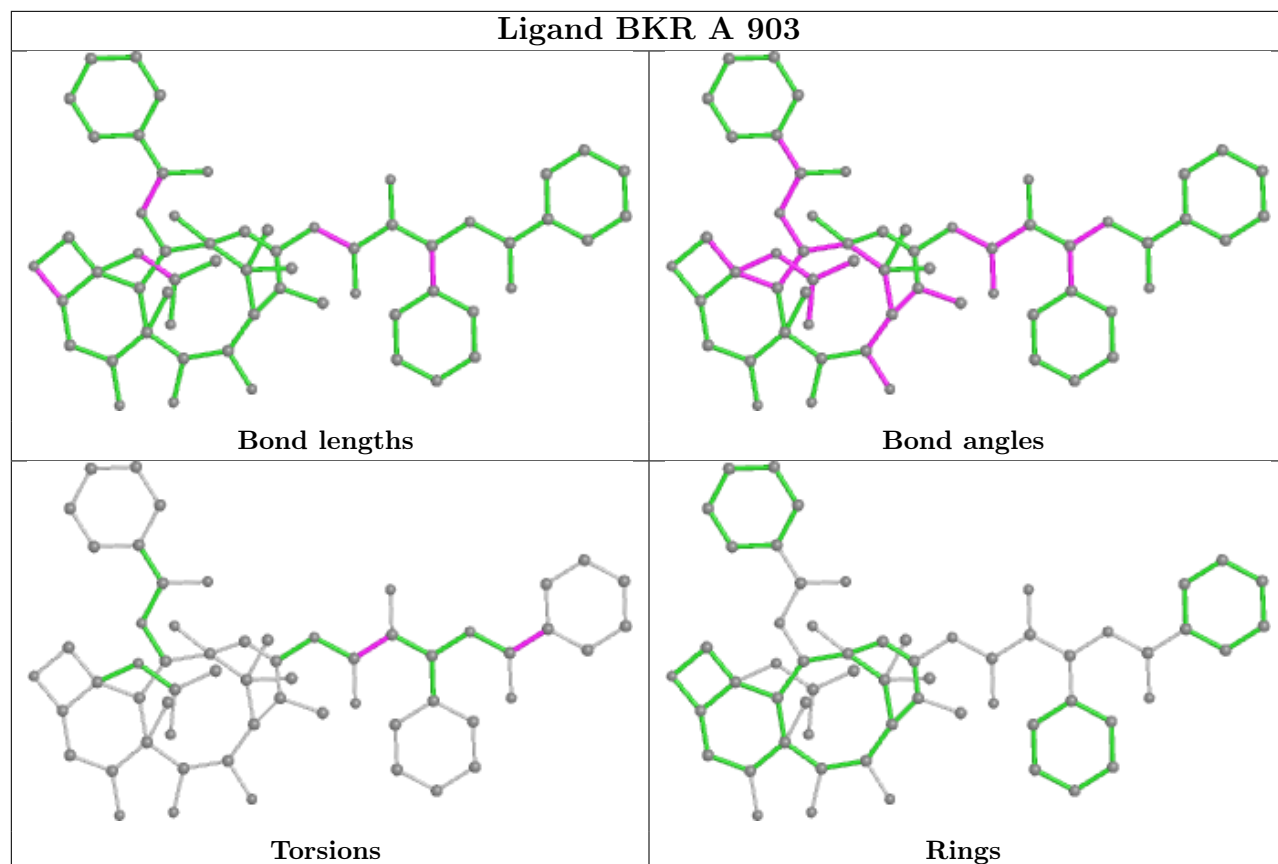
Mol	Chain	Res	Type	Atoms
5	B	901	NAG	C4-C5-C6-O6
5	B	901	NAG	O5-C5-C6-O6
5	A	904	NAG	O5-C5-C6-O6
5	A	904	NAG	C4-C5-C6-O6
5	A	904	NAG	C8-C7-N2-C2

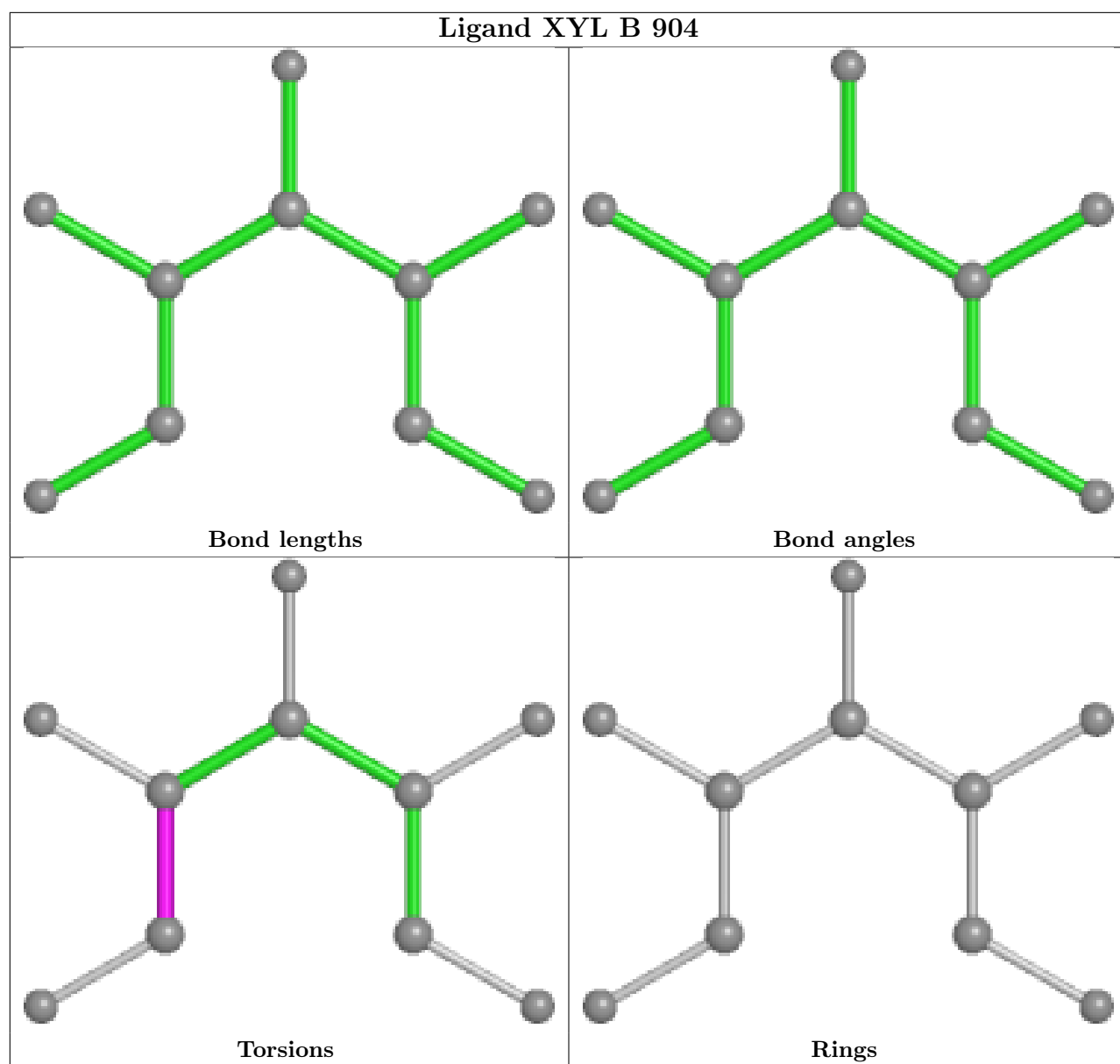
There are no ring outliers.

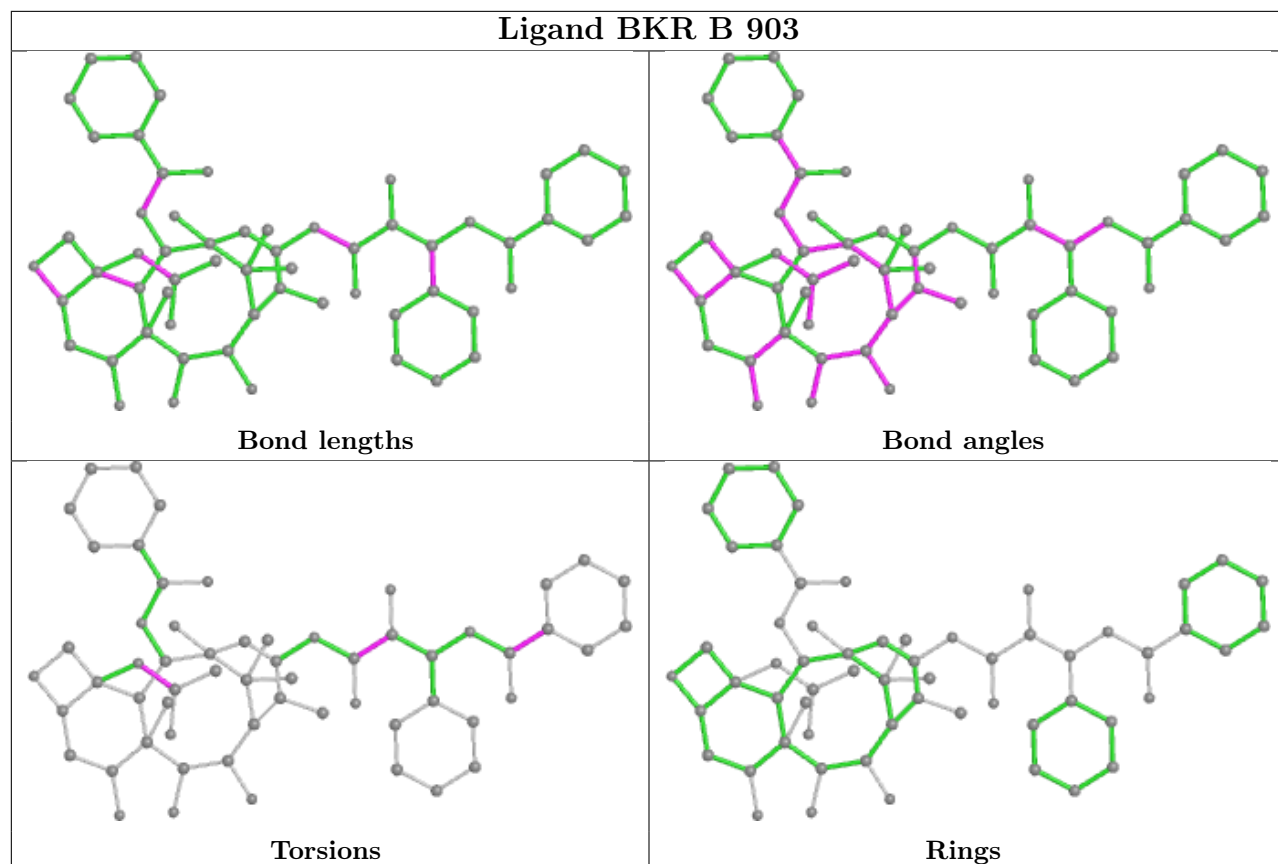
6 monomers are involved in 14 short contacts:

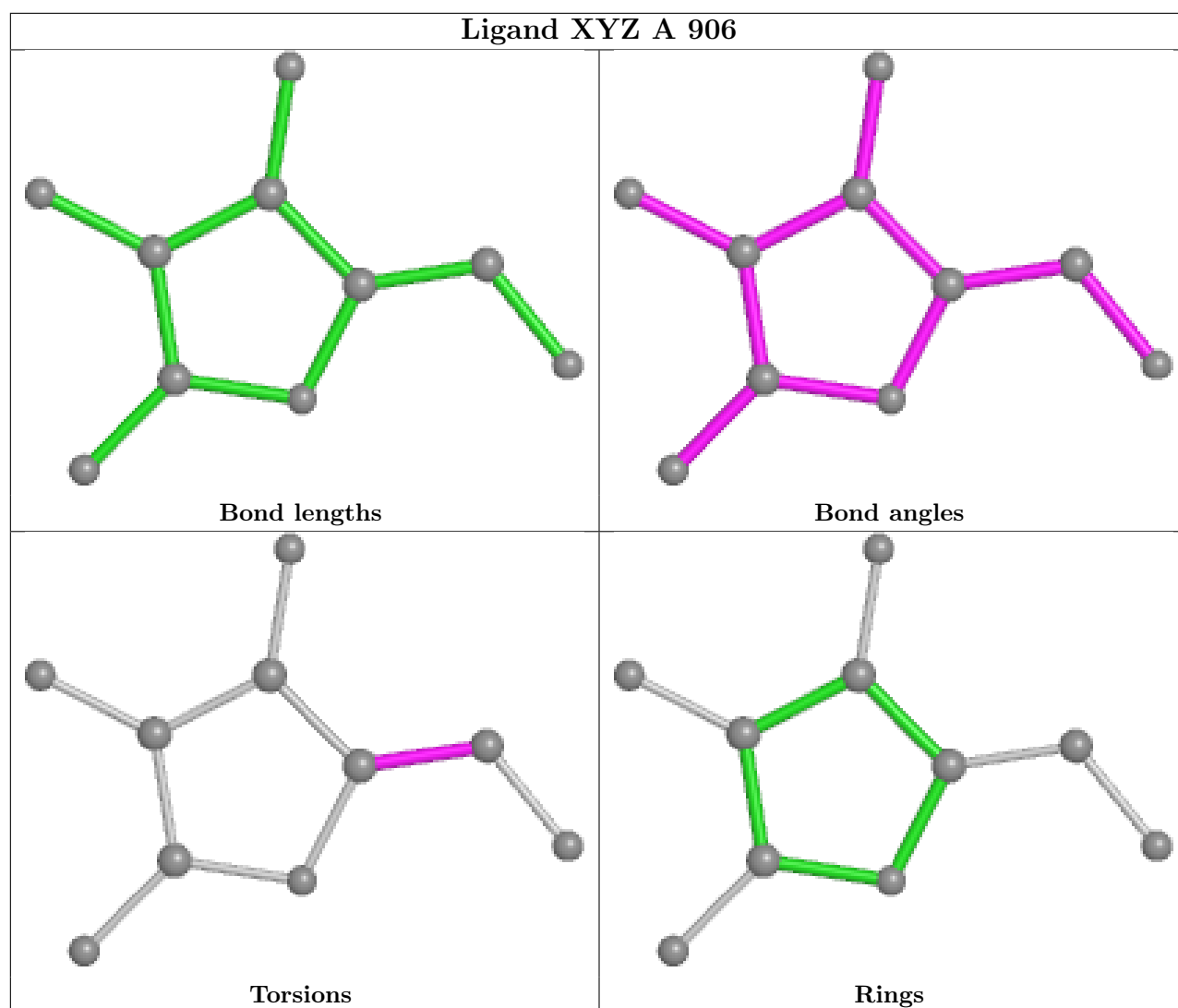
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	903	BKR	1	0
5	B	901	NAG	2	0
8	B	904	XYL	3	0
6	B	903	BKR	5	0
7	A	906	XYZ	2	0
5	A	901	NAG	2	0

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	A	756/803 (94%)	-0.43	1 (0%) 92 94	20, 31, 49, 82	0
1	B	757/803 (94%)	-0.38	3 (0%) 89 91	21, 32, 49, 93	0
All	All	1513/1606 (94%)	-0.40	4 (0%) 90 92	20, 32, 49, 93	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	683	TYR	3.8
1	B	47	THR	3.5
1	B	683	TYR	3.4
1	B	48	GLN	3.1

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](#)

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.

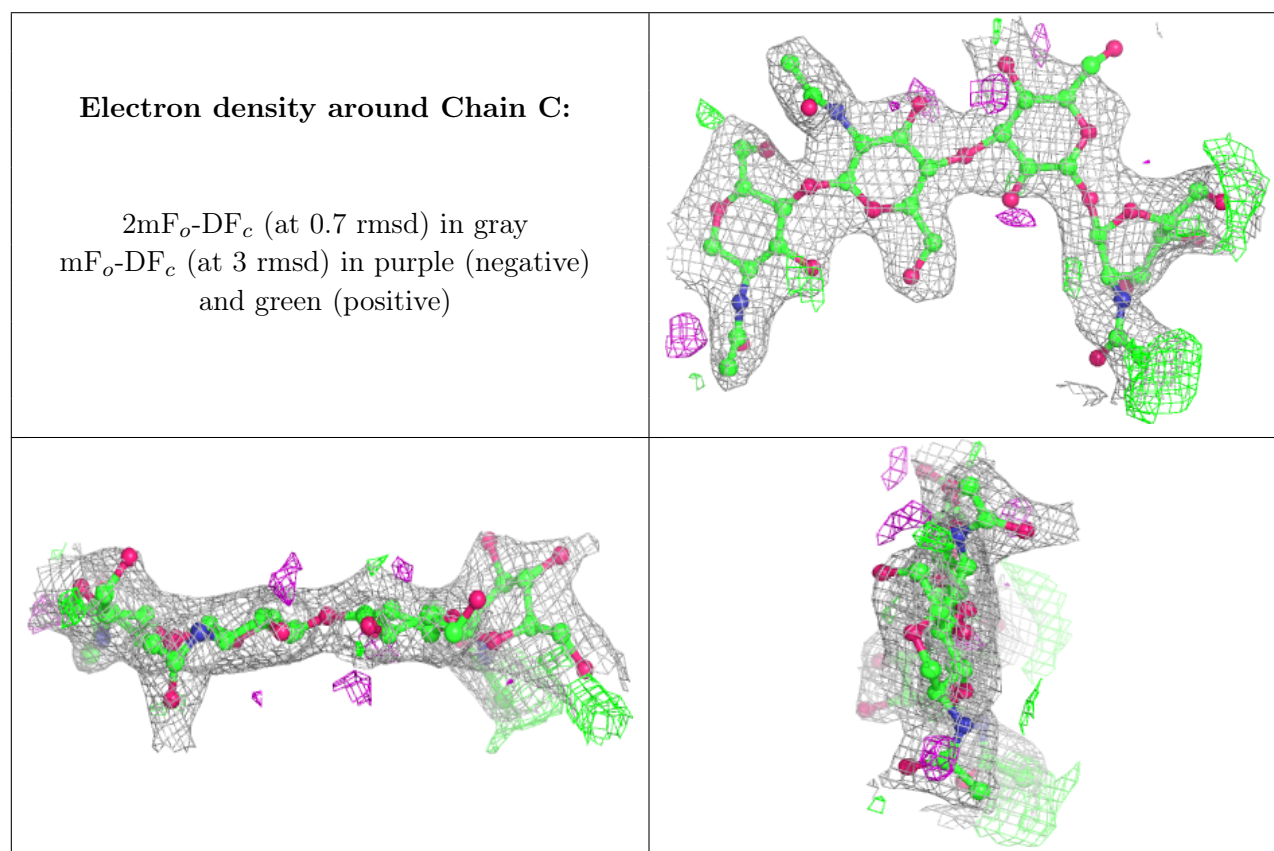
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	C	4	14/15	0.56	0.17	51,61,73,76	0
3	NAG	I	2	14/15	0.71	0.13	49,55,64,67	0
4	NAG	F	2	14/15	0.76	0.16	47,55,69,78	0
4	NAG	E	2	14/15	0.81	0.16	58,66,83,93	0
4	MAN	E	1	11/12	0.83	0.10	38,44,46,55	0

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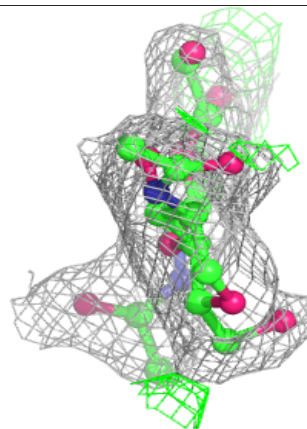
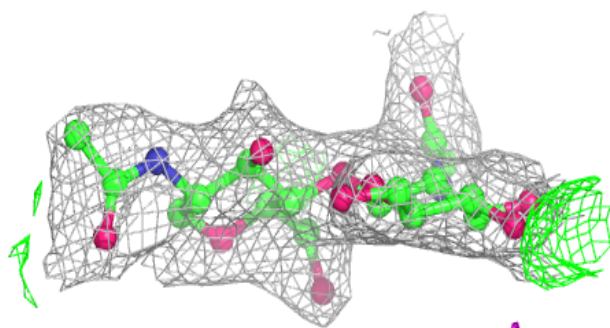
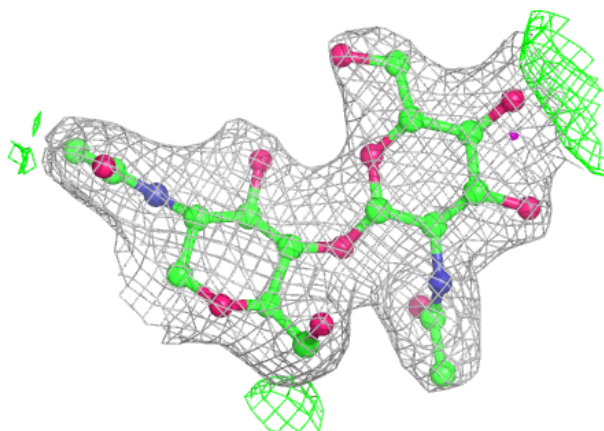
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	G	2	14/15	0.83	0.11	34,37,45,46	0
3	NAG	J	2	14/15	0.83	0.10	47,52,57,61	0
4	MAN	F	1	12/12	0.84	0.09	39,44,46,47	0
2	MAN	C	3	11/12	0.85	0.14	38,52,61,68	0
3	NAG	D	2	14/15	0.88	0.10	32,32,33,34	0
2	NAG	C	1	14/15	0.91	0.09	35,37,42,47	0
2	NAG	C	2	14/15	0.91	0.09	32,35,36,39	0
3	NAG	H	2	14/15	0.92	0.08	31,35,42,43	0
3	NAG	D	1	14/15	0.93	0.07	30,32,35,35	0
3	NAG	J	1	14/15	0.94	0.07	34,39,41,43	0
3	NAG	I	1	14/15	0.94	0.08	37,40,42,43	0
3	NAG	G	1	14/15	0.94	0.07	23,25,28,34	0
3	NAG	H	1	14/15	0.96	0.06	27,28,30,32	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



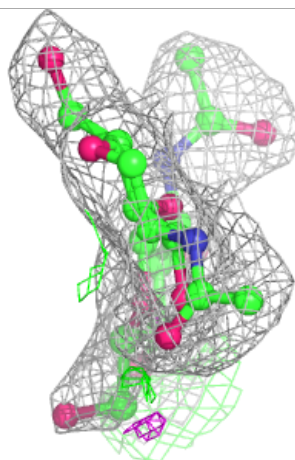
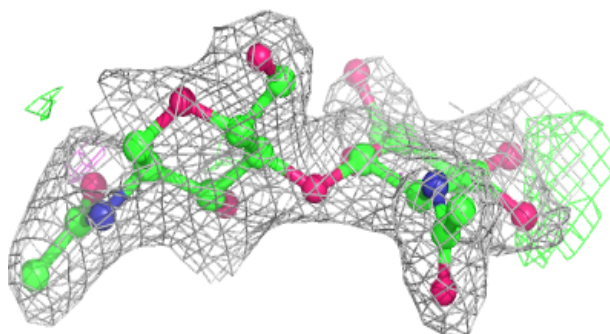
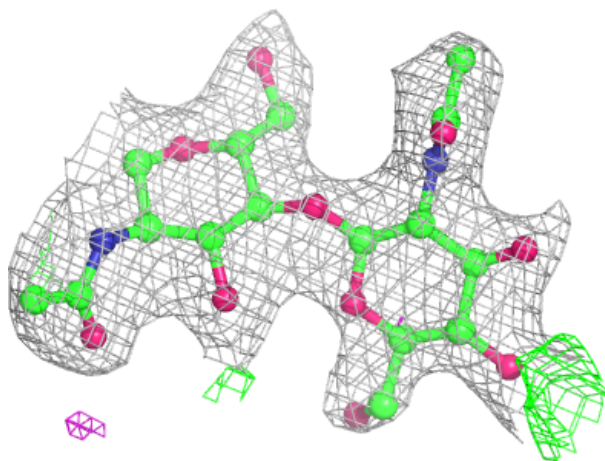
Electron density around Chain D:

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



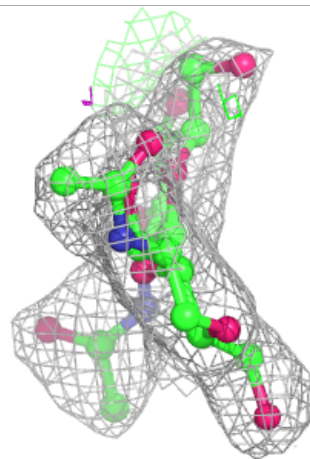
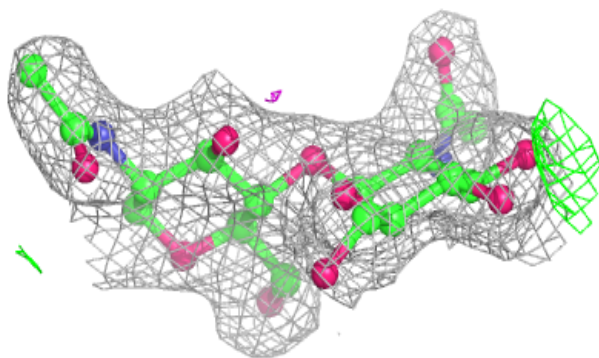
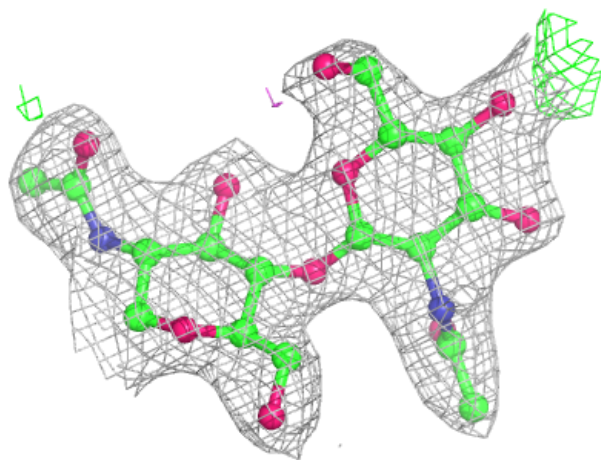
Electron density around Chain G:

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



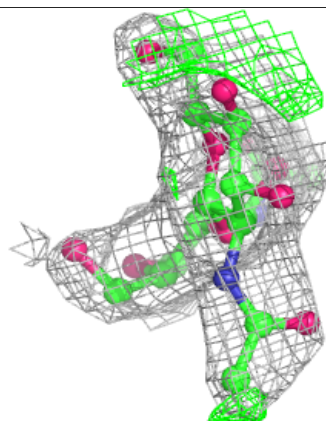
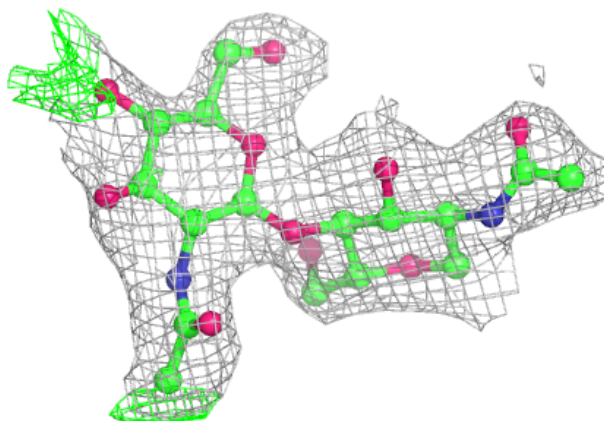
Electron density around Chain H:

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

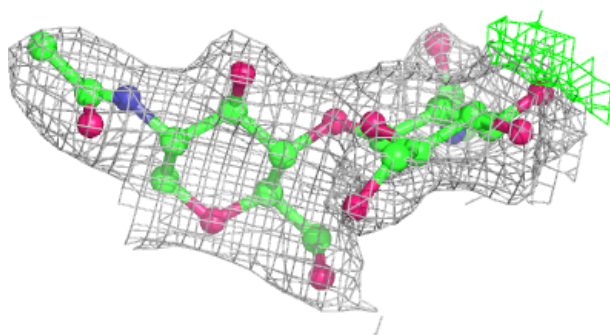
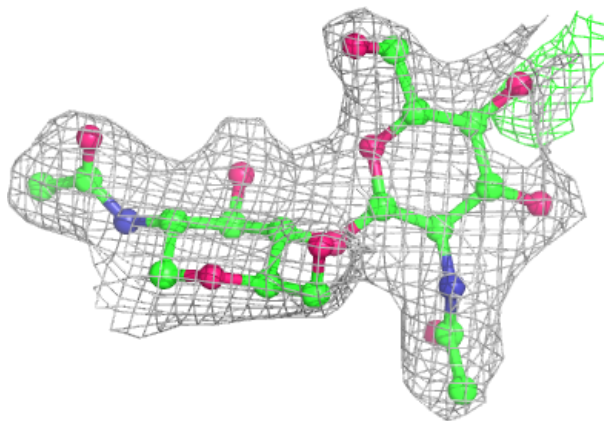


Electron density around Chain I:

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

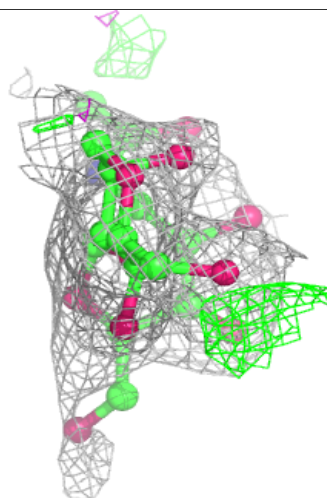
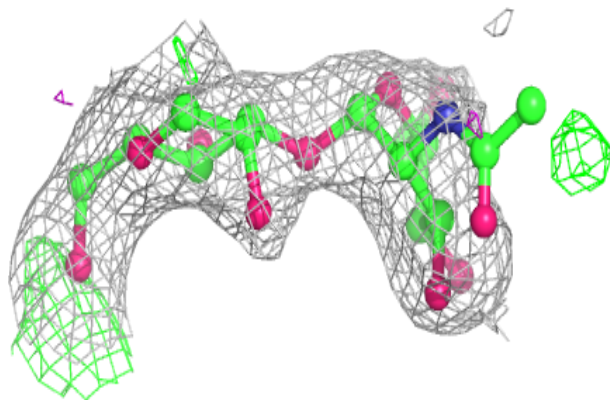
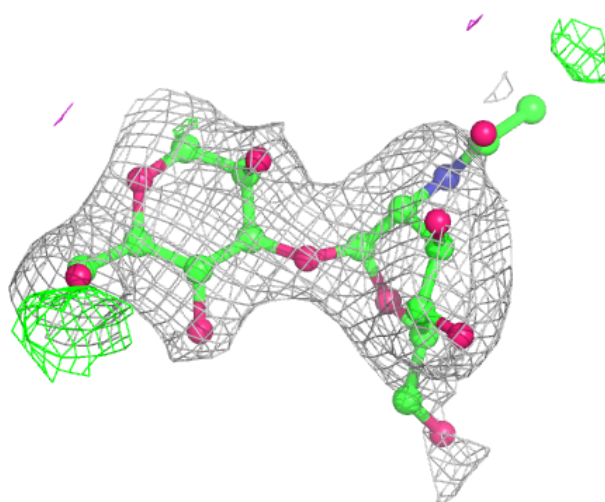
**Electron density around Chain J:**

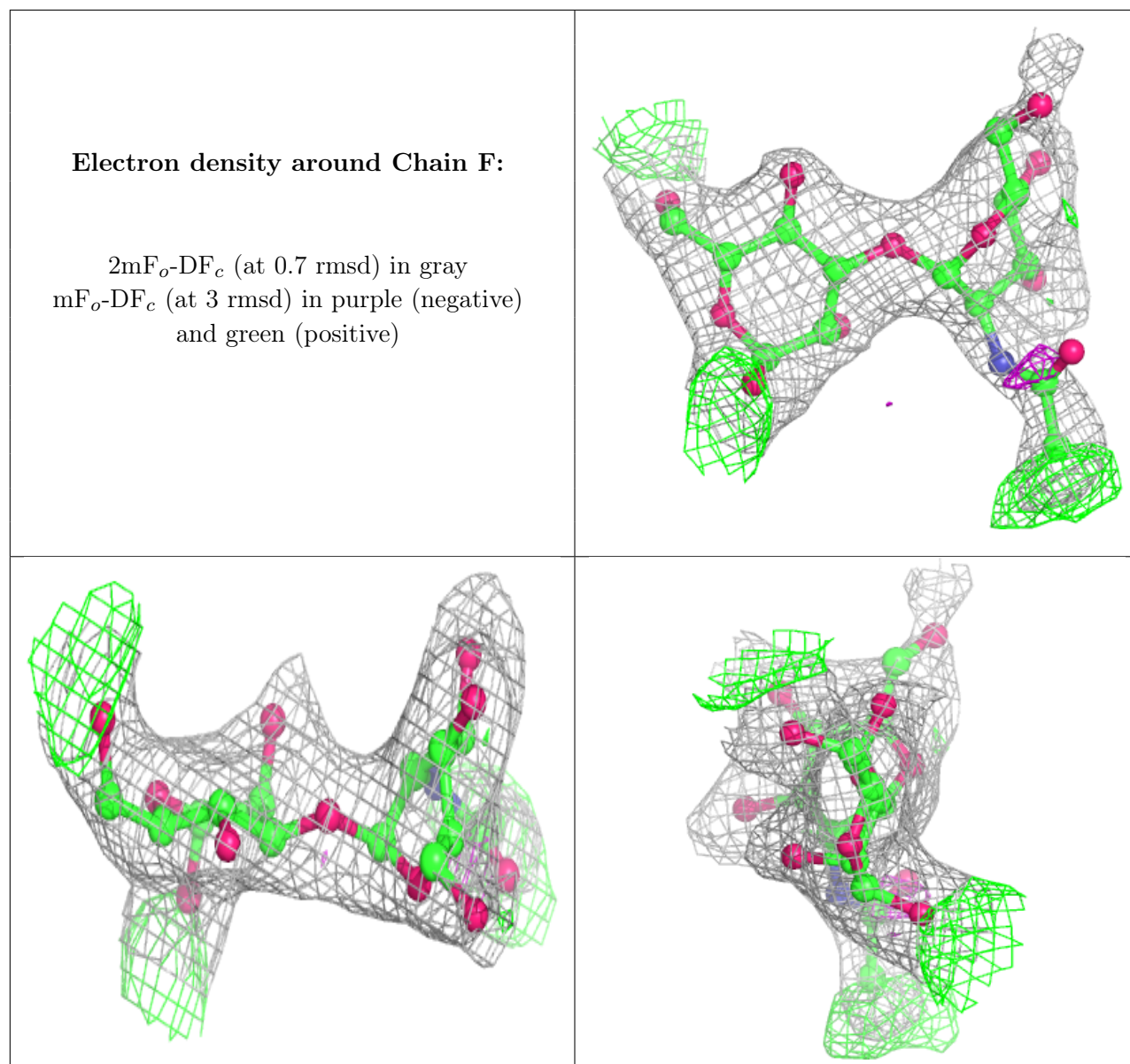
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain E:

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





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(Å ²)	Q<0.9
5	NAG	A	904	15/15	0.70	0.17	48,56,61,63	0
8	XYL	B	904	10/10	0.76	0.30	66,85,89,99	0
5	NAG	A	901	14/15	0.82	0.14	63,76,82,82	0
7	XYZ	A	906	10/10	0.84	0.15	36,47,49,54	0
5	NAG	B	901	14/15	0.84	0.12	60,65,67,71	0

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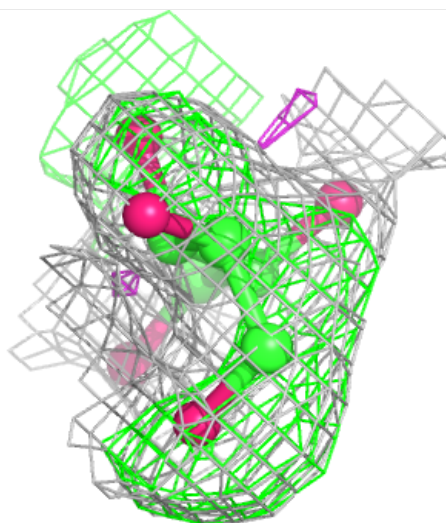
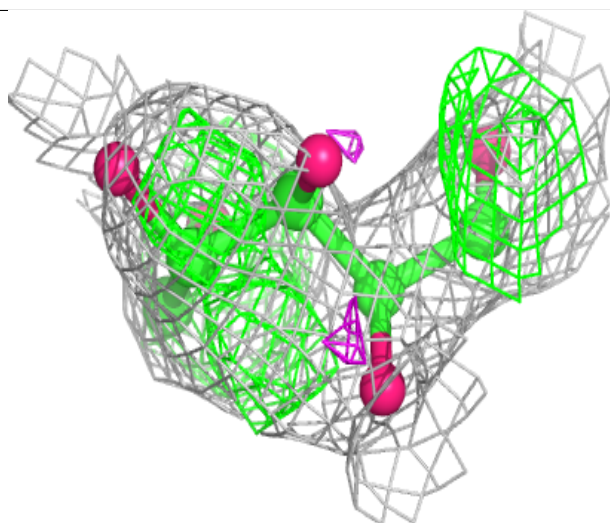
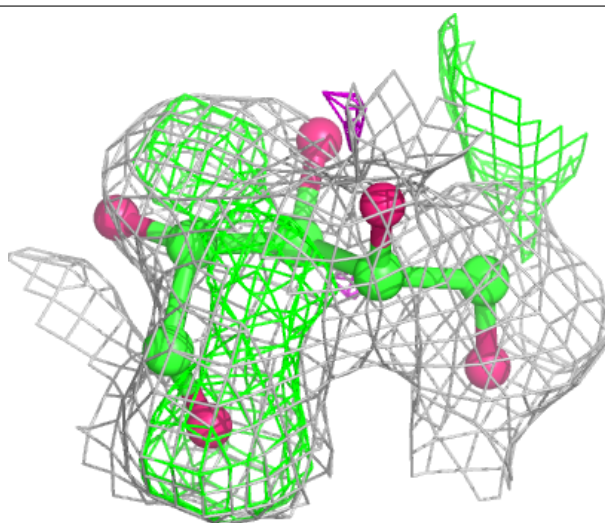
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	BKR	A	903	59/59	0.86	0.11	41,49,54,56	0
6	BKR	B	903	59/59	0.89	0.10	35,49,64,65	0
5	NAG	A	905	14/15	0.92	0.07	44,46,48,49	0
5	NAG	A	902	14/15	0.92	0.08	29,37,40,42	0
5	NAG	B	902	14/15	0.93	0.07	33,34,38,41	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.

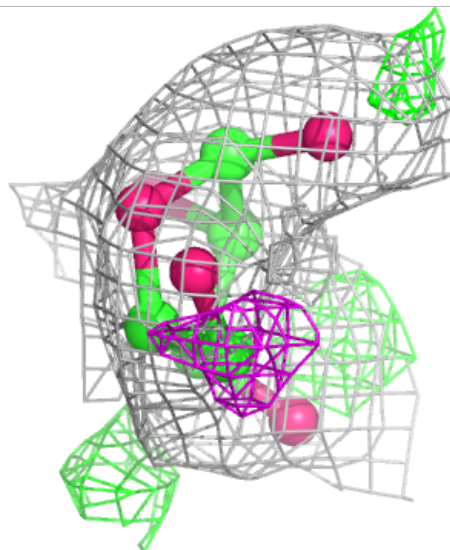
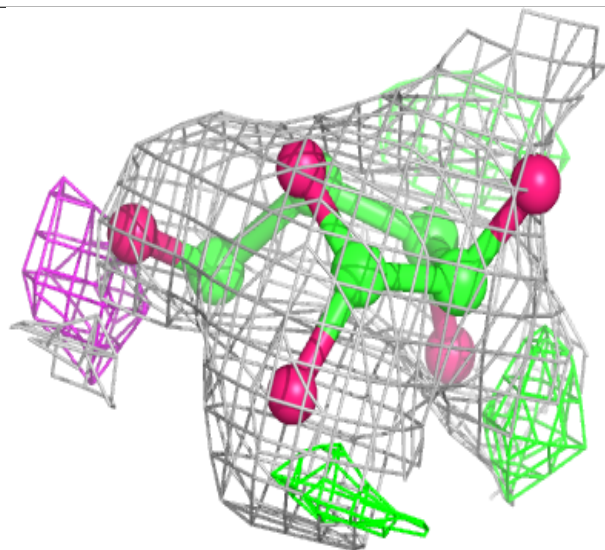
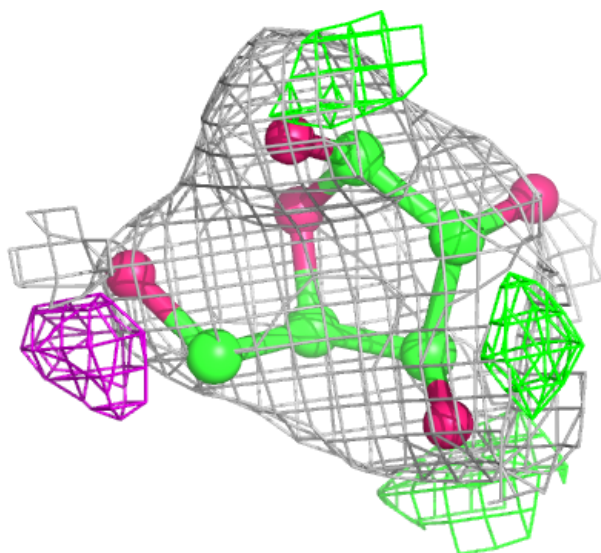
Electron density around XYL B 904:

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mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



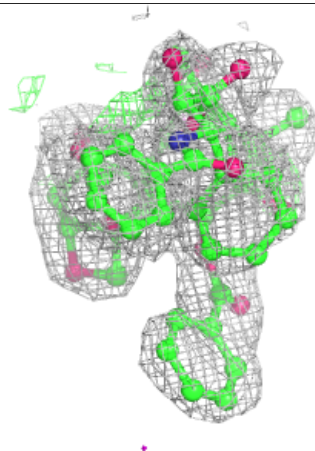
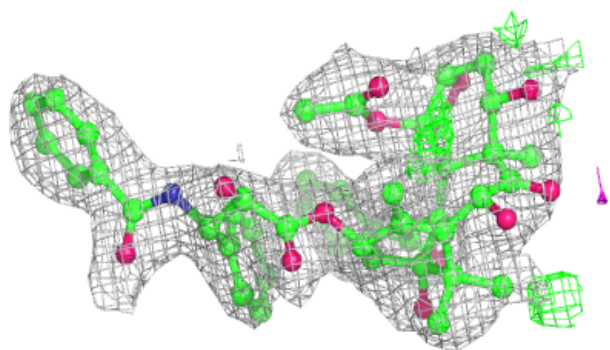
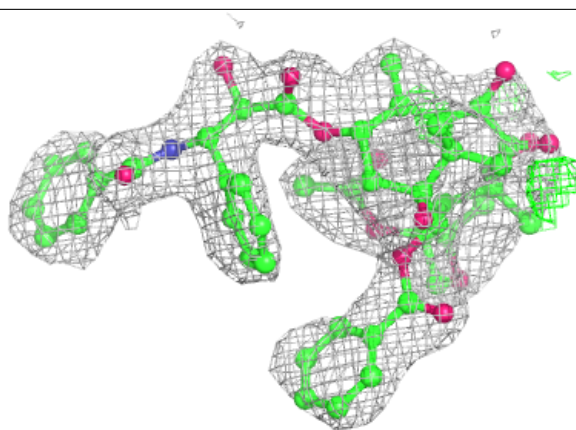
Electron density around XYZ A 906:

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



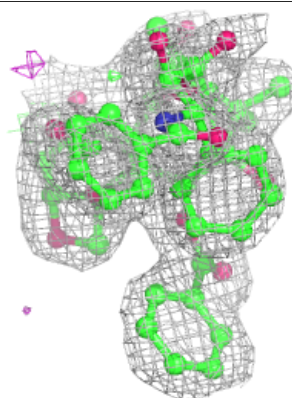
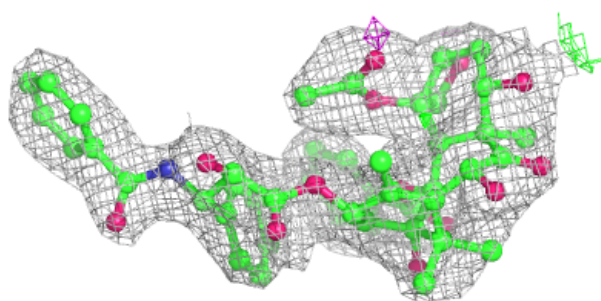
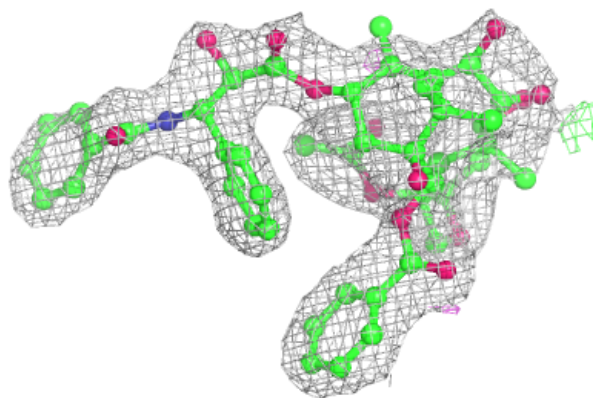
Electron density around BKR A 903:

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



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 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.