



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 04:59 AM UTC

PDB ID : 5O47 / pdb_00005o47
Title : Structure of D80A-fructofuranosidase from Xanthophyllomyces dendrorhous complexed with fructosyl-hydroxytyrosol
Authors : Ramirez-Escudero, M.; Sanz-Aparicio, J.
Deposited on : 2017-05-26
Resolution : 1.91 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

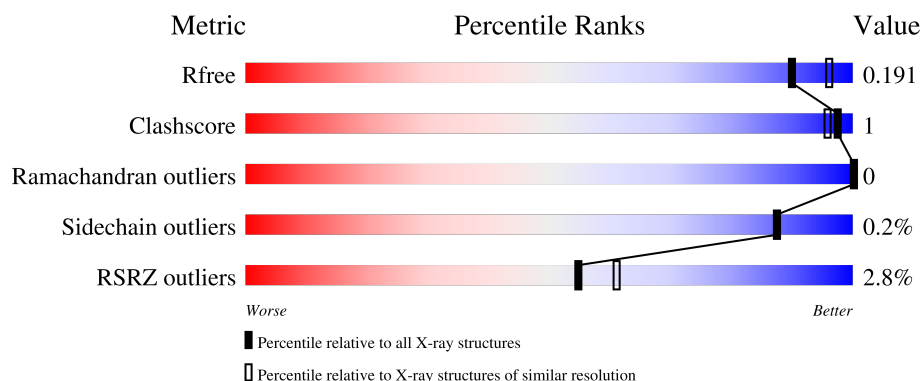
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

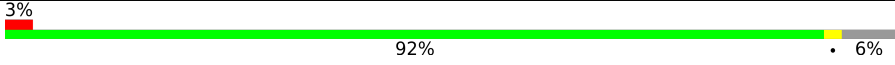
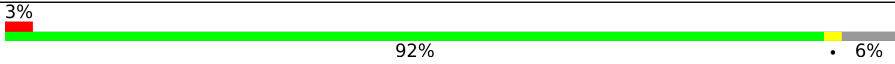
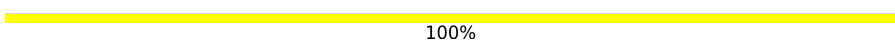
The reported resolution of this entry is 1.91 Å.

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.



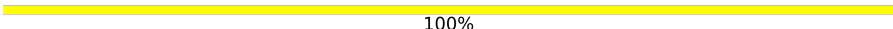
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	665	
1	B	665	
2	C	6	
2	F	6	
3	D	10	

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Mol	Chain	Length	Quality of chain
3	G	10	 80%20%
4	E	2	 50%50%
4	H	2	 100%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 12016 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.

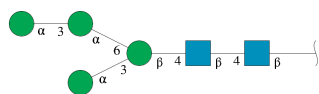
- Molecule 1 is a protein called Beta-fructofuranosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	624	Total	C	N	O	S	0	3	0
			4812	3063	786	956	7			
1	B	624	Total	C	N	O	S	0	3	0
			4815	3065	787	956	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	ILE	conflict	UNP J7HDY4
A	80	ALA	ASP	engineered mutation	UNP J7HDY4
A	663	ALA	SER	conflict	UNP J7HDY4
A	665	TYR	ARG	conflict	UNP J7HDY4
B	2	VAL	ILE	conflict	UNP J7HDY4
B	80	ALA	ASP	engineered mutation	UNP J7HDY4
B	663	ALA	SER	conflict	UNP J7HDY4
B	665	TYR	ARG	conflict	UNP J7HDY4

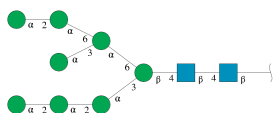
- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			72	40	2	30			
2	F	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyran

ose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



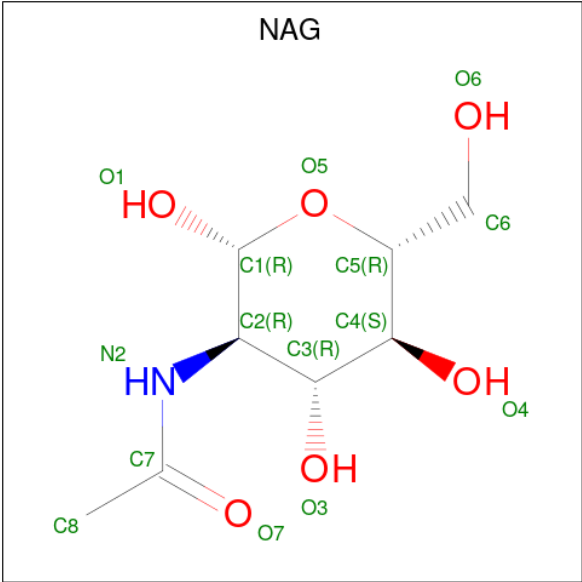
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	10	Total	C	N	O	0	0	0
			116	64	2	50			
3	G	10	Total	C	N	O	0	0	0
			116	64	2	50			

- Molecule 4 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
4	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



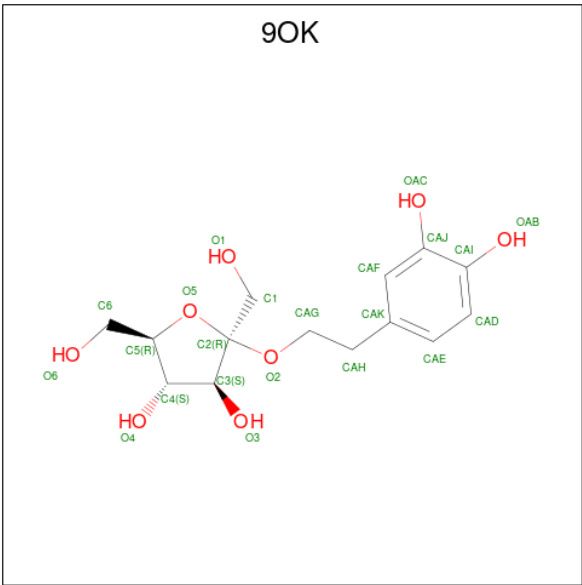
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1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

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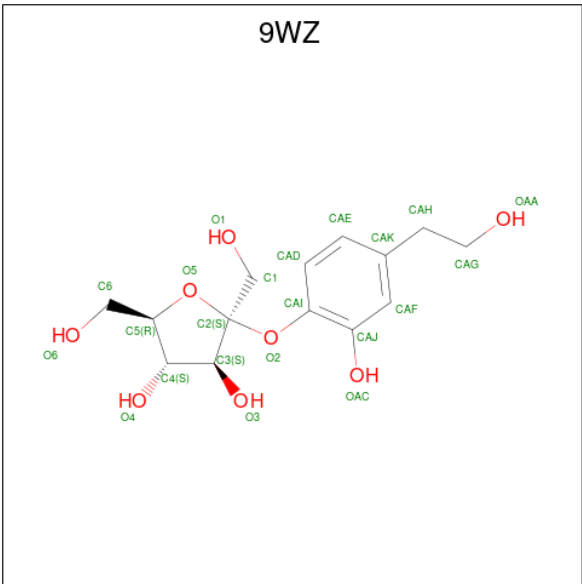
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1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

- Molecule 6 is 2-(3,4-dihydroxyphenyl)ethyl beta-D-fructofuranoside (CCD ID: 9OK) (formula: C₁₄H₂₀O₈).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			22	14	8		
6	B	1	Total	C	O	0	0
			22	14	8		

- Molecule 7 is 2-hydroxy-4-(2-hydroxyethyl)phenyl beta-D-fructofuranoside (CCD ID: 9WZ) (formula: C₁₄H₂₀O₈).



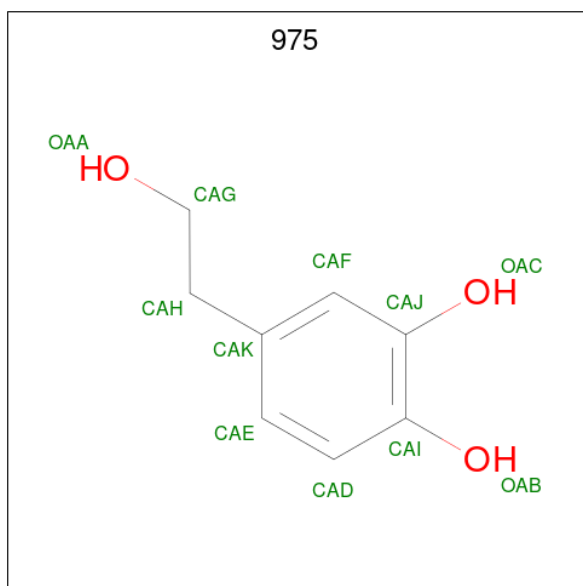
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			22	14	8		

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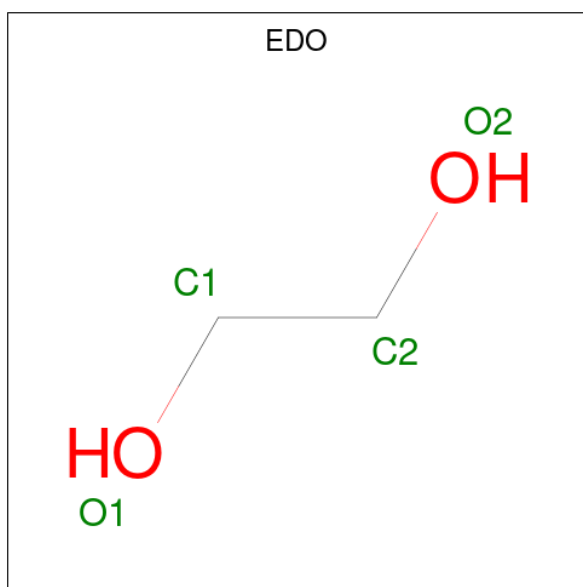
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			22	14	8		

- Molecule 8 is 4-(2-hydroxyethyl)benzene-1,2-diol (CCD ID: 975) (formula: $C_8H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			11	8	3		
8	B	1	Total	C	O	0	0
			11	8	3		

- Molecule 9 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	A	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		
9	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0
9	B	1	Total C O 4 2 2	0	0

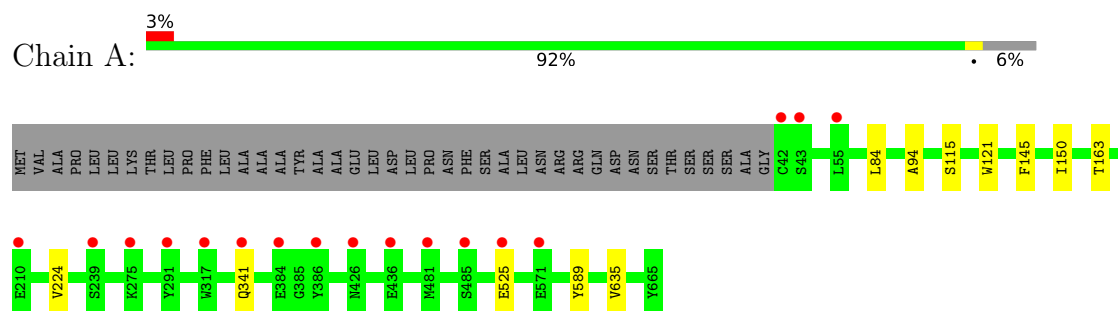
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	722	Total O 722 722	0	0
10	B	659	Total O 659 659	0	0

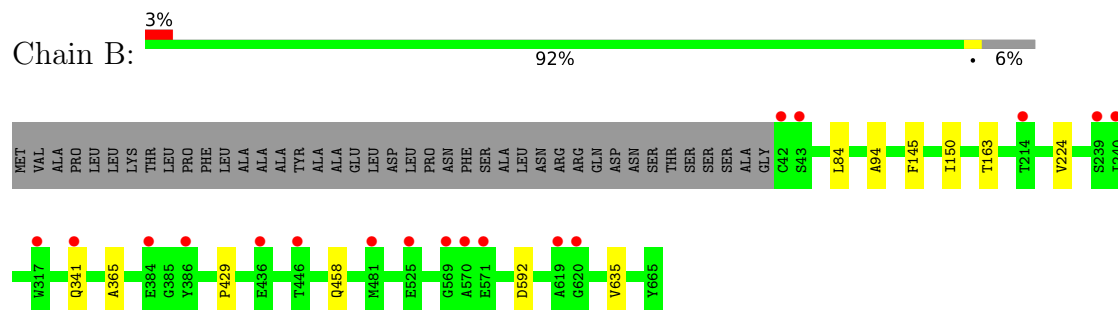
3 Residue-property plots

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.

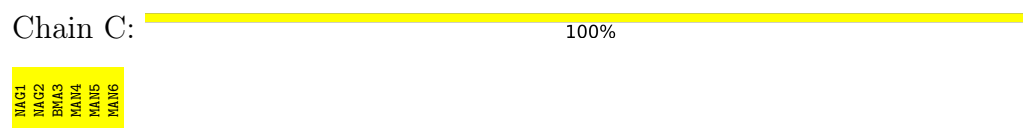
- Molecule 1: Beta-fructofuranosidase



- Molecule 1: Beta-fructofuranosidase



- Molecule 2: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain D:  70% 30%



- Molecule 3: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain G:  80% 20%

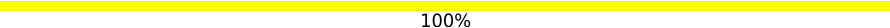


- Molecule 4: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain E:  50% 50%



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

Chain H:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	74.62Å 205.36Å 146.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.98 – 1.91 48.98 – 1.91	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.98-1.91) 100.0 (48.98-1.91)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.91 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.168 , 0.182 0.177 , 0.191	Depositor DCC
R_{free} test set	8691 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.710	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12016	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.25% 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: 975, NAG, MAN, 9OK, EDO, BMA, 9WZ

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	0/4946	0.70	0/6762
1	B	0.47	1/4949 (0.0%)	0.70	0/6766
All	All	0.47	1/9895 (0.0%)	0.70	0/13528

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	458	GLN	C-O	-5.10	1.18	1.23

There are no bond angle outliers.

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	A	4812	0	4504	7	0
1	B	4815	0	4510	8	0
2	C	72	0	61	0	0
2	F	72	0	61	0	0
3	D	116	0	97	0	1
3	G	116	0	97	0	0
4	E	28	0	25	0	0
4	H	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	210	0	195	2	0
5	B	168	0	156	1	0
6	A	22	0	0	0	0
6	B	22	0	0	0	0
7	A	22	0	0	1	0
7	B	22	0	0	2	0
8	A	11	0	0	0	0
8	B	11	0	0	0	0
9	A	44	0	66	1	0
9	B	44	0	66	2	0
10	A	722	0	0	0	1
10	B	659	0	0	0	0
All	All	12016	0	9863	16	2

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 1.

The worst 5 of 16 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:B:341:GLN:HG3	7:B:2003:9WZ:CAE	2.03	0.88
1:A:341:GLN:HG3	7:A:2003:9WZ:CAE	2.07	0.85
1:B:429:PRO:HB3	5:B:1444:NAG:H62	1.83	0.60
1:A:525:GLU:H	5:A:1606:NAG:H81	1.69	0.58
1:B:150:ILE:HG12	1:B:224:VAL:HG11	1.94	0.49

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:A:4100:HOH:O	10:A:4100:HOH:O[2_555]	1.90	0.30
3:D:8:MAN:O4	3:D:8:MAN:O4[2_555]	2.04	0.16

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	625/665 (94%)	597 (96%)	28 (4%)	0	100	100
1	B	625/665 (94%)	595 (95%)	30 (5%)	0	100	100
All	All	1250/1330 (94%)	1192 (95%)	58 (5%)	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	500/529 (94%)	499 (100%)	1 (0%)	87	87
1	B	500/529 (94%)	499 (100%)	1 (0%)	87	87
All	All	1000/1058 (94%)	998 (100%)	2 (0%)	87	87

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	635	VAL
1	B	635	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	HIS
1	A	342	ASN
1	B	46	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 ⓘ

36 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	0.29	0	17,19,21	1.11	2 (11%)
2	NAG	C	2	2	14,14,15	0.34	0	17,19,21	1.06	1 (5%)
2	BMA	C	3	2	11,11,12	0.26	0	15,15,17	1.10	2 (13%)
2	MAN	C	4	2	11,11,12	0.56	0	15,15,17	0.92	1 (6%)
2	MAN	C	5	2	11,11,12	0.32	0	15,15,17	0.95	1 (6%)
2	MAN	C	6	2	11,11,12	0.34	0	15,15,17	1.13	2 (13%)
3	NAG	D	1	3,1	14,14,15	0.42	0	17,19,21	0.81	0
3	MAN	D	10	3	11,11,12	0.26	0	15,15,17	0.58	0
3	NAG	D	2	3	14,14,15	0.29	0	17,19,21	0.90	0
3	BMA	D	3	3	11,11,12	0.28	0	15,15,17	0.77	0
3	MAN	D	4	3	11,11,12	0.26	0	15,15,17	1.47	2 (13%)
3	MAN	D	5	3	11,11,12	0.29	0	15,15,17	2.25	4 (26%)
3	MAN	D	6	3	11,11,12	0.33	0	15,15,17	0.73	0
3	MAN	D	7	3	11,11,12	0.29	0	15,15,17	0.87	0
3	MAN	D	8	3	11,11,12	0.31	0	15,15,17	0.67	0
3	MAN	D	9	3	11,11,12	0.40	0	15,15,17	0.94	0
4	NAG	E	1	1,4	14,14,15	0.30	0	17,19,21	0.89	0
4	NAG	E	2	4	14,14,15	0.25	0	17,19,21	1.29	2 (11%)
2	NAG	F	1	1,2	14,14,15	0.36	0	17,19,21	1.06	2 (11%)
2	NAG	F	2	2	14,14,15	0.33	0	17,19,21	0.76	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	F	3	2	11,11,12	0.26	0	15,15,17	1.36	2 (13%)
2	MAN	F	4	2	11,11,12	0.41	0	15,15,17	0.92	1 (6%)
2	MAN	F	5	2	11,11,12	0.30	0	15,15,17	0.76	1 (6%)
2	MAN	F	6	2	11,11,12	0.36	0	15,15,17	0.62	0
3	NAG	G	1	3,1	14,14,15	0.45	0	17,19,21	0.71	0
3	MAN	G	10	3	11,11,12	0.26	0	15,15,17	0.63	0
3	NAG	G	2	3	14,14,15	0.33	0	17,19,21	0.78	0
3	BMA	G	3	3	11,11,12	0.25	0	15,15,17	0.64	0
3	MAN	G	4	3	11,11,12	0.23	0	15,15,17	1.25	1 (6%)
3	MAN	G	5	3	11,11,12	0.32	0	15,15,17	1.04	1 (6%)
3	MAN	G	6	3	11,11,12	0.36	0	15,15,17	0.83	0
3	MAN	G	7	3	11,11,12	0.29	0	15,15,17	0.79	0
3	MAN	G	8	3	11,11,12	0.35	0	15,15,17	0.72	0
3	MAN	G	9	3	11,11,12	0.42	0	15,15,17	0.82	0
4	NAG	H	1	1,4	14,14,15	0.48	0	17,19,21	1.16	2 (11%)
4	NAG	H	2	4	14,14,15	0.46	0	17,19,21	1.15	2 (11%)

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	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	2/2/19/22	0/1/1/1
2	MAN	C	5	2	-	0/2/19/22	0/1/1/1
2	MAN	C	6	2	-	0/2/19/22	0/1/1/1
3	NAG	D	1	3,1	-	0/6/23/26	0/1/1/1
3	MAN	D	10	3	-	0/2/19/22	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
3	MAN	D	4	3	-	1/2/19/22	0/1/1/1
3	MAN	D	5	3	-	0/2/19/22	0/1/1/1
3	MAN	D	6	3	-	2/2/19/22	0/1/1/1
3	MAN	D	7	3	-	0/2/19/22	0/1/1/1
3	MAN	D	8	3	-	0/2/19/22	0/1/1/1
3	MAN	D	9	3	-	0/2/19/22	0/1/1/1
4	NAG	E	1	1,4	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	BMA	F	3	2	-	0/2/19/22	0/1/1/1
2	MAN	F	4	2	-	0/2/19/22	0/1/1/1
2	MAN	F	5	2	-	2/2/19/22	0/1/1/1
2	MAN	F	6	2	-	0/2/19/22	0/1/1/1
3	NAG	G	1	3,1	-	0/6/23/26	0/1/1/1
3	MAN	G	10	3	-	0/2/19/22	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	BMA	G	3	3	-	0/2/19/22	0/1/1/1
3	MAN	G	4	3	-	0/2/19/22	0/1/1/1
3	MAN	G	5	3	-	1/2/19/22	0/1/1/1
3	MAN	G	6	3	-	2/2/19/22	0/1/1/1
3	MAN	G	7	3	-	0/2/19/22	0/1/1/1
3	MAN	G	8	3	-	0/2/19/22	0/1/1/1
3	MAN	G	9	3	-	0/2/19/22	0/1/1/1
4	NAG	H	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	5	MAN	C1-O5-C5	7.48	122.21	112.19
3	D	4	MAN	C1-O5-C5	4.26	117.90	112.19
3	G	4	MAN	C1-O5-C5	3.98	117.52	112.19
2	F	3	BMA	C1-O5-C5	3.52	116.91	112.19
3	G	5	MAN	C1-O5-C5	3.33	116.66	112.19

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

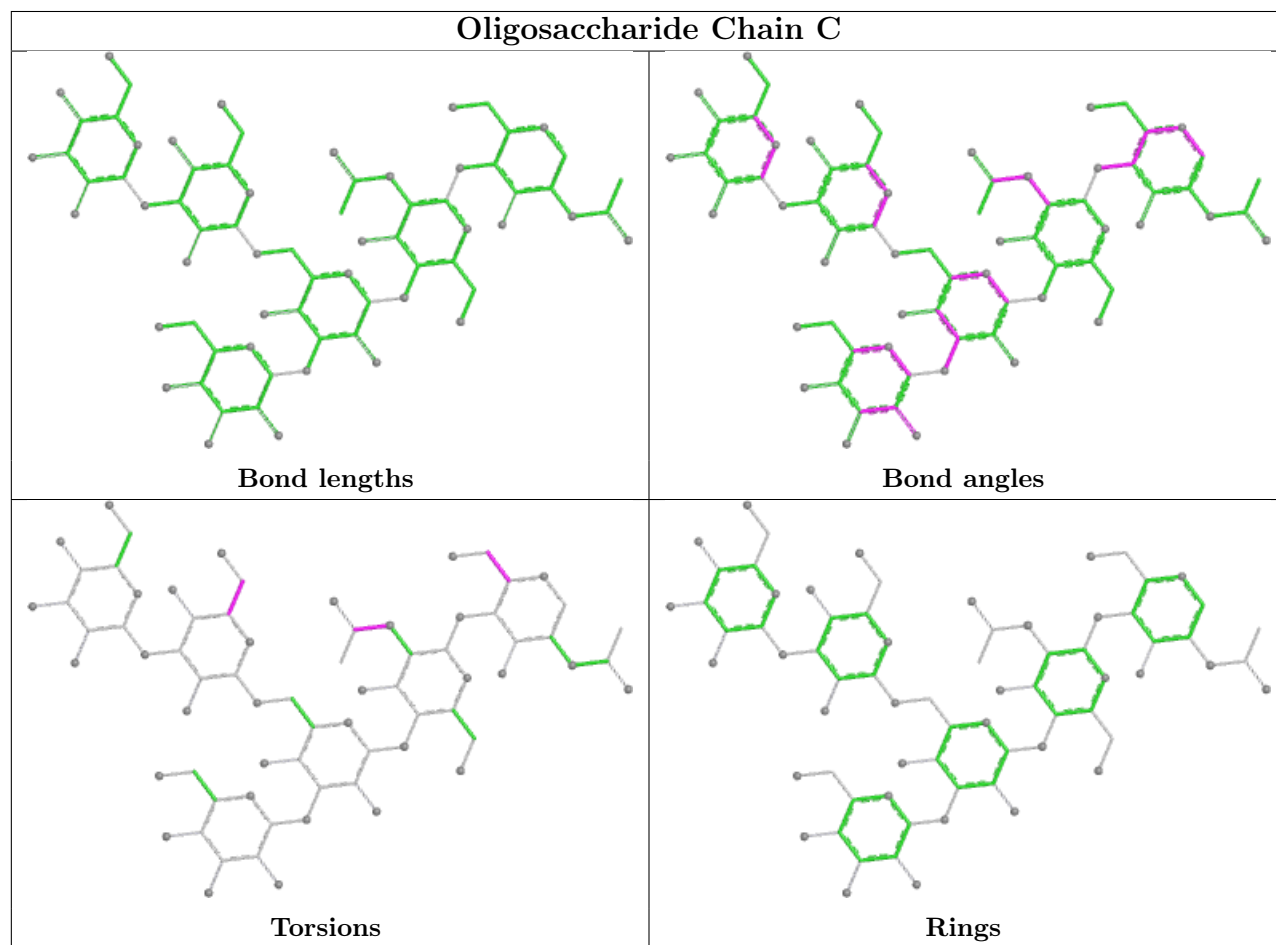
Mol	Chain	Res	Type	Atoms
4	H	2	NAG	C4-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
3	G	6	MAN	O5-C5-C6-O6
3	G	6	MAN	C4-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2

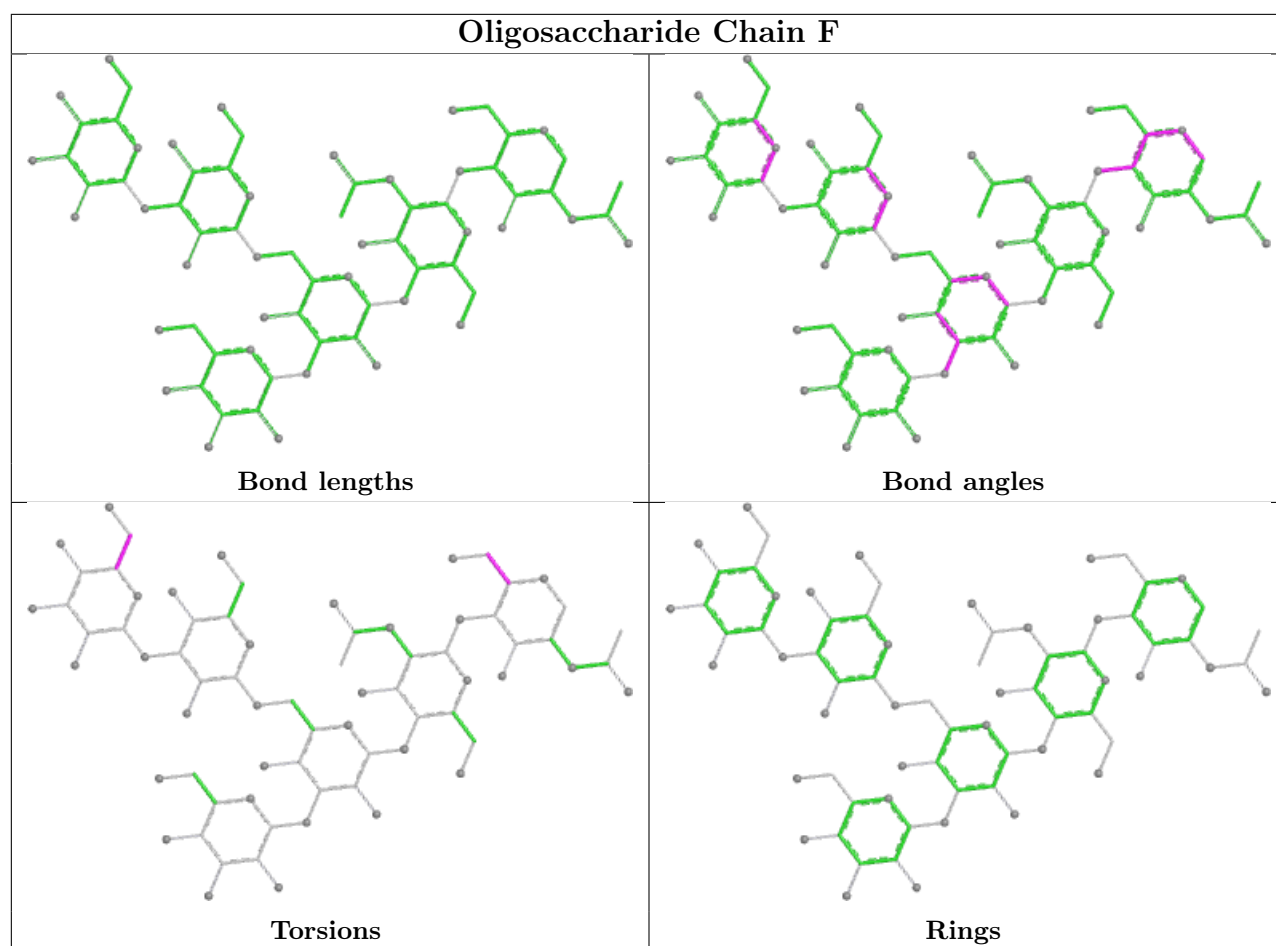
There are no ring outliers.

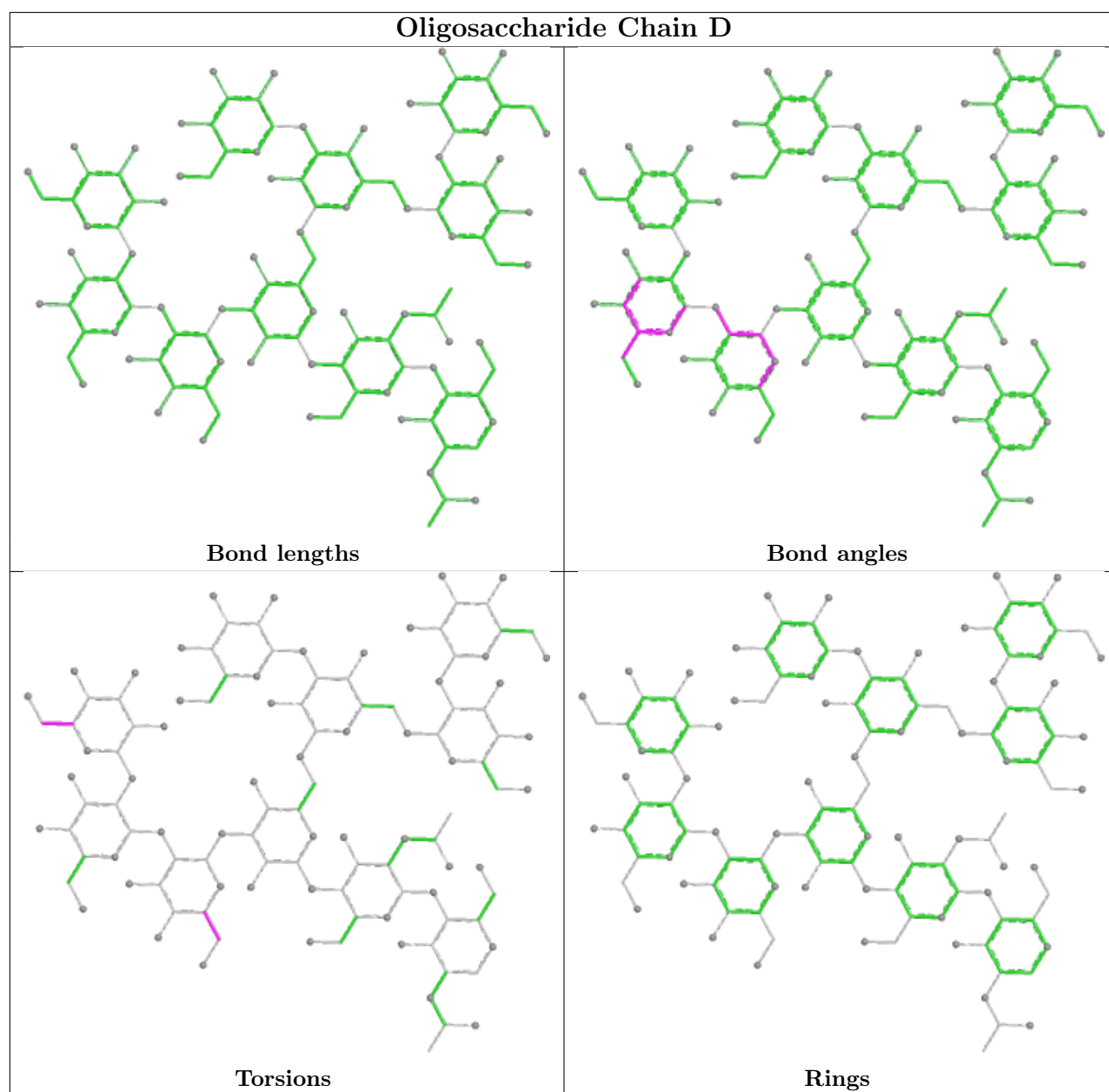
1 monomer is involved in 1 short contact:

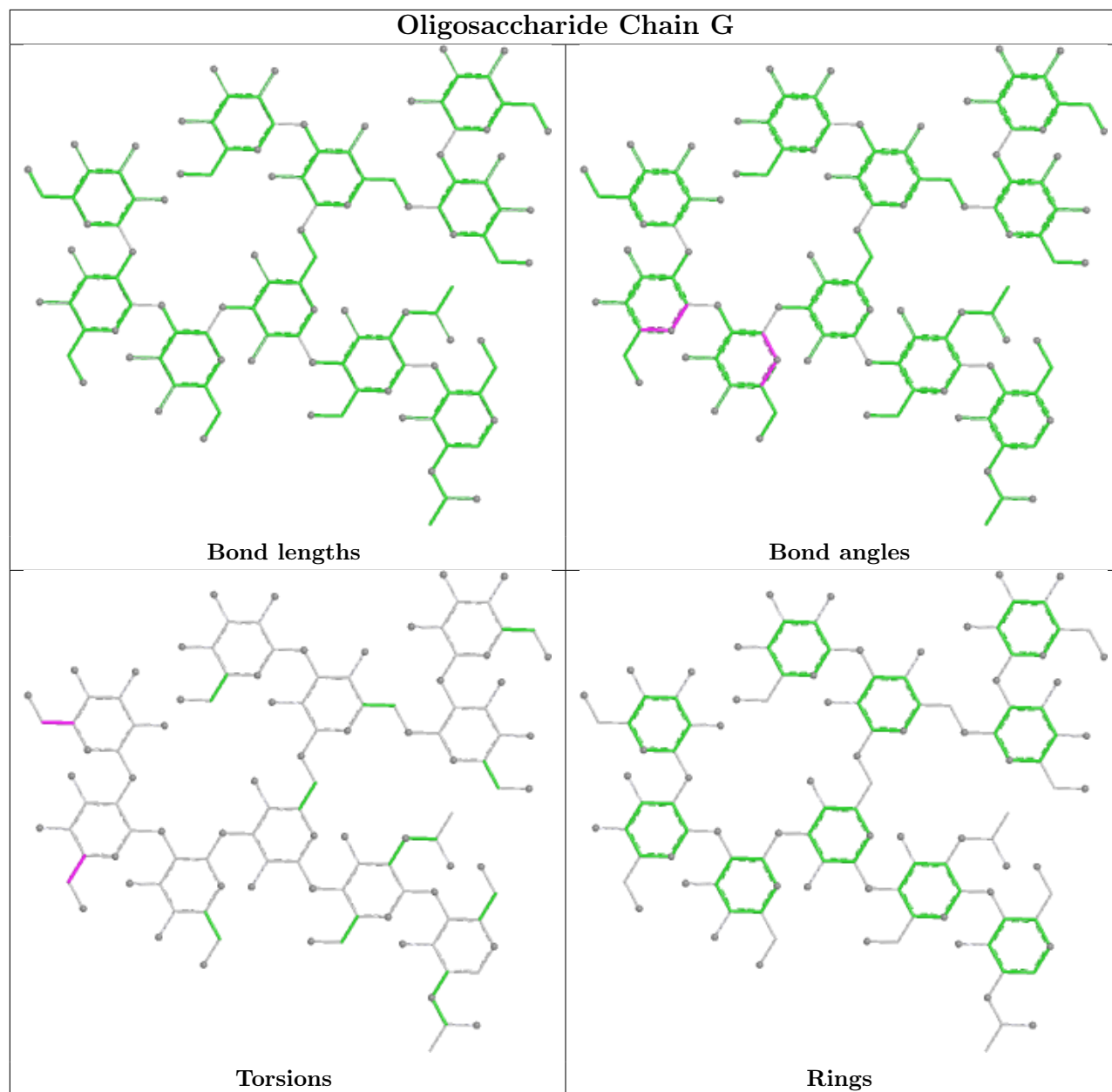
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	8	MAN	0	1

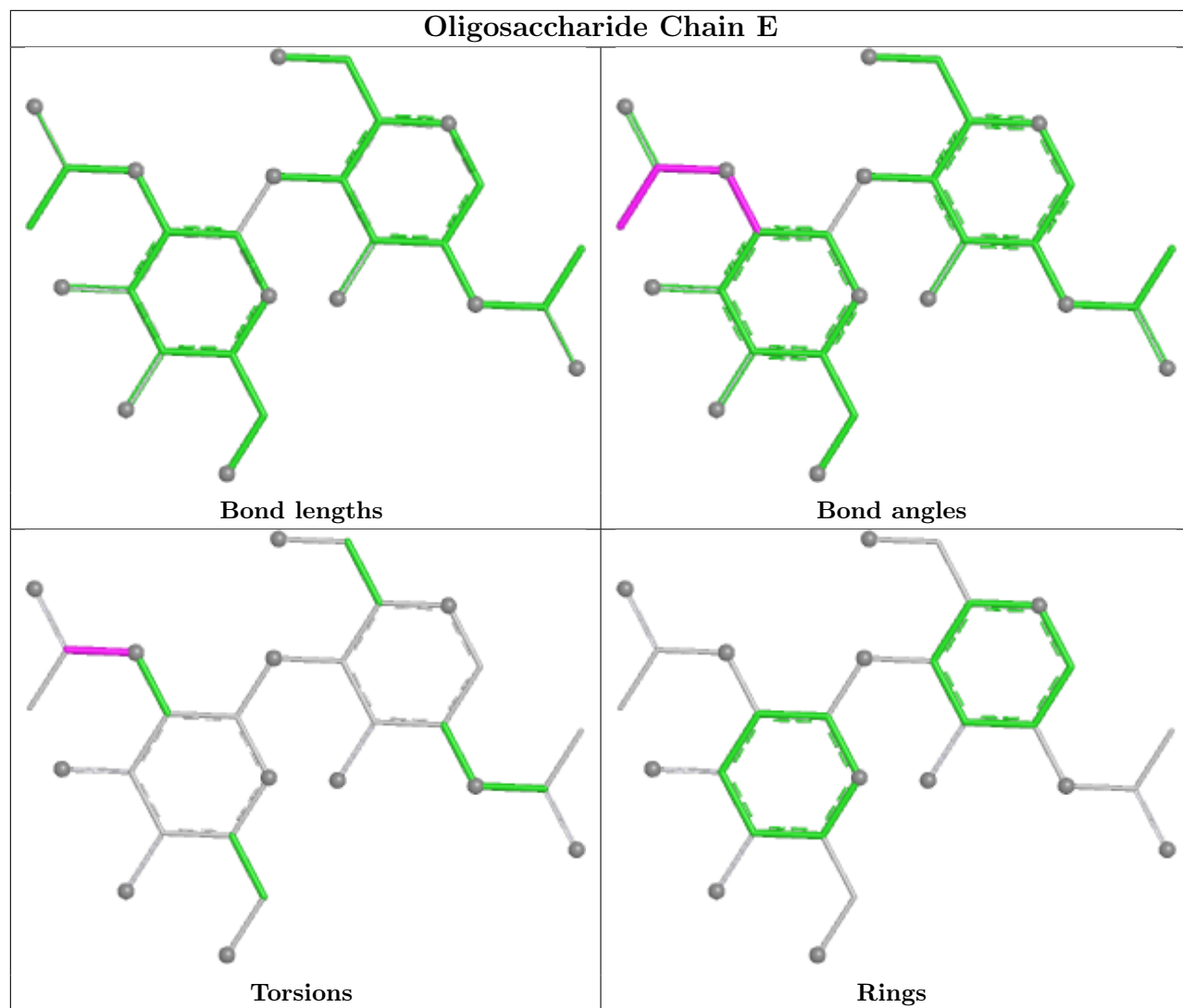
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

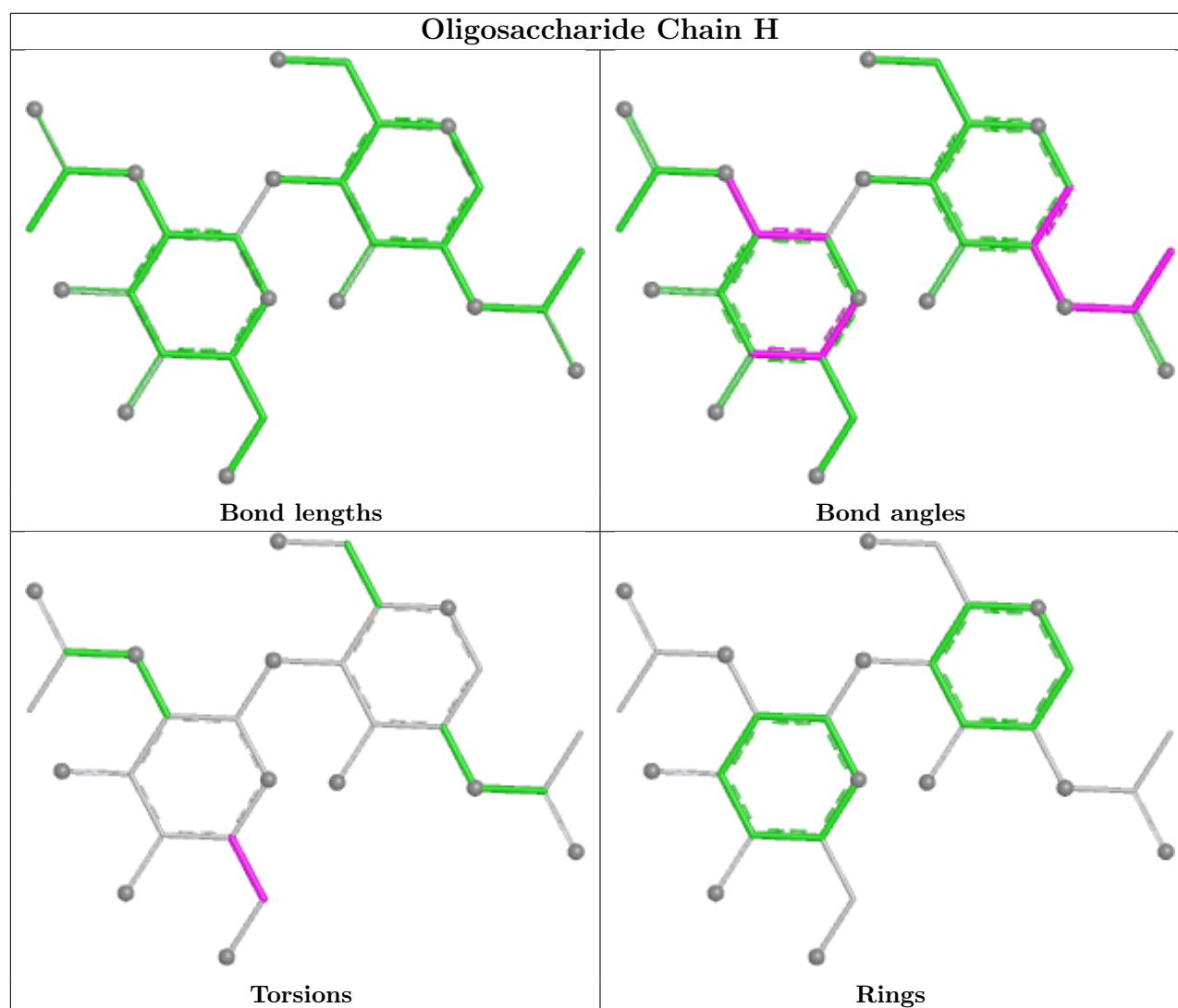












5.6 Ligand geometry [i](#)

55 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	$\# Z > 2$
5	NAG	B	1357	1	14,14,15	0.32	0	17,19,21	0.73	0
9	EDO	B	2014	-	3,3,3	0.45	0	2,2,2	0.27	0
5	NAG	A	1357	1	14,14,15	0.28	0	17,19,21	1.16	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	B	1125	1	14,14,15	0.50	0	17,19,21	2.16	5 (29%)
5	NAG	B	1644	1	14,14,15	0.30	0	17,19,21	0.64	0
6	9OK	A	2000	-	23,23,23	1.46	2 (8%)	32,33,33	1.32	3 (9%)
5	NAG	A	1242	1	14,14,15	0.30	0	17,19,21	1.13	2 (11%)
5	NAG	A	1483	1	14,14,15	0.29	0	17,19,21	0.93	1 (5%)
9	EDO	A	2019	-	3,3,3	0.58	0	2,2,2	0.38	0
9	EDO	B	2021	-	3,3,3	0.47	0	2,2,2	0.25	0
9	EDO	A	2016	-	3,3,3	0.50	0	2,2,2	0.12	0
9	EDO	B	2017	-	3,3,3	0.47	0	2,2,2	0.36	0
5	NAG	B	1555	1	14,14,15	0.31	0	17,19,21	0.82	0
5	NAG	A	1236	1	14,14,15	0.29	0	17,19,21	0.66	0
8	975	A	2004	-	11,11,11	1.61	1 (9%)	14,14,14	0.88	0
5	NAG	B	1242	1	14,14,15	0.35	0	17,19,21	1.57	4 (23%)
5	NAG	B	1319	1	14,14,15	0.35	0	17,19,21	1.49	1 (5%)
9	EDO	B	2013	-	3,3,3	0.53	0	2,2,2	0.28	0
5	NAG	B	1052	1	14,14,15	0.33	0	17,19,21	0.76	0
9	EDO	B	2020	-	3,3,3	0.50	0	2,2,2	0.20	0
8	975	B	2004	-	11,11,11	1.61	1 (9%)	14,14,14	1.12	1 (7%)
9	EDO	B	2006	-	3,3,3	0.47	0	2,2,2	0.26	0
9	EDO	A	2015	-	3,3,3	0.46	0	2,2,2	0.30	0
5	NAG	A	1052	1	14,14,15	0.24	0	17,19,21	1.16	3 (17%)
5	NAG	B	1215	1	14,14,15	0.30	0	17,19,21	1.43	1 (5%)
9	EDO	A	2009	-	3,3,3	0.49	0	2,2,2	0.12	0
7	9WZ	A	2003	-	22,23,23	2.14	2 (9%)	29,33,33	2.42	4 (13%)
9	EDO	B	2019	-	3,3,3	0.60	0	2,2,2	0.42	0
5	NAG	B	1539	1	14,14,15	0.54	0	17,19,21	1.60	2 (11%)
9	EDO	B	2008	-	3,3,3	0.49	0	2,2,2	0.29	0
9	EDO	A	2013	-	3,3,3	0.53	0	2,2,2	0.32	0
5	NAG	A	1444	1	14,14,15	0.28	0	17,19,21	0.76	1 (5%)
9	EDO	A	2007	-	3,3,3	0.44	0	2,2,2	0.20	0
5	NAG	A	1606	1	14,14,15	0.42	0	17,19,21	1.59	2 (11%)
5	NAG	A	1471	1	14,14,15	0.30	0	17,19,21	1.04	1 (5%)
5	NAG	B	1444	1	14,14,15	0.39	0	17,19,21	0.98	1 (5%)
9	EDO	B	2007	-	3,3,3	0.58	0	2,2,2	0.58	0
5	NAG	A	1512	1	14,14,15	0.42	0	17,19,21	2.64	6 (35%)
5	NAG	B	1471	1	14,14,15	0.30	0	17,19,21	0.64	0
5	NAG	A	1644	1	14,14,15	0.25	0	17,19,21	0.70	0
6	9OK	B	2000	-	23,23,23	1.36	2 (8%)	32,33,33	1.40	6 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	A	1215	1	14,14,15	0.31	0	17,19,21	1.41	1 (5%)
5	NAG	B	1483	1	14,14,15	0.31	0	17,19,21	1.25	2 (11%)
9	EDO	A	2010	-	3,3,3	0.46	0	2,2,2	0.32	0
5	NAG	A	1539	1	14,14,15	0.39	0	17,19,21	1.43	2 (11%)
5	NAG	A	1319	1	14,14,15	0.43	0	17,19,21	1.62	3 (17%)
9	EDO	A	2011	-	3,3,3	0.50	0	2,2,2	0.20	0
9	EDO	A	2006	-	3,3,3	0.48	0	2,2,2	0.30	0
9	EDO	A	2020	-	3,3,3	0.48	0	2,2,2	0.14	0
9	EDO	B	2016	-	3,3,3	0.48	0	2,2,2	0.24	0
5	NAG	A	1555	1	14,14,15	0.29	0	17,19,21	1.64	3 (17%)
7	9WZ	B	2003	-	22,23,23	2.10	2 (9%)	29,33,33	2.17	8 (27%)
9	EDO	B	2015	-	3,3,3	0.49	0	2,2,2	0.26	0
9	EDO	A	2012	-	3,3,3	0.51	0	2,2,2	0.26	0
5	NAG	A	1125	1	14,14,15	0.46	0	17,19,21	1.29	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
5	NAG	B	1357	1	-	0/6/23/26	0/1/1/1
9	EDO	B	2014	-	-	0/1/1/1	-
5	NAG	A	1357	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1125	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1644	1	-	1/6/23/26	0/1/1/1
6	9OK	A	2000	-	-	2/12/31/31	0/2/2/2
5	NAG	A	1242	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1483	1	-	0/6/23/26	0/1/1/1
9	EDO	A	2019	-	-	1/1/1/1	-
9	EDO	B	2021	-	-	0/1/1/1	-
9	EDO	A	2016	-	-	1/1/1/1	-
9	EDO	B	2017	-	-	0/1/1/1	-
5	NAG	B	1555	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1236	1	-	0/6/23/26	0/1/1/1
8	975	A	2004	-	-	0/3/3/3	0/1/1/1
5	NAG	B	1242	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1319	1	-	2/6/23/26	0/1/1/1
9	EDO	B	2013	-	-	1/1/1/1	-
5	NAG	B	1052	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	EDO	B	2020	-	-	1/1/1/1	-
8	975	B	2004	-	-	1/3/3/3	0/1/1/1
9	EDO	B	2006	-	-	0/1/1/1	-
9	EDO	A	2015	-	-	0/1/1/1	-
5	NAG	A	1052	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1215	1	-	0/6/23/26	0/1/1/1
9	EDO	A	2009	-	-	1/1/1/1	-
7	9WZ	A	2003	-	-	9/13/32/32	0/2/2/2
9	EDO	B	2019	-	-	1/1/1/1	-
5	NAG	B	1539	1	-	2/6/23/26	0/1/1/1
9	EDO	B	2008	-	-	1/1/1/1	-
9	EDO	A	2013	-	-	1/1/1/1	-
5	NAG	A	1444	1	-	0/6/23/26	0/1/1/1
9	EDO	A	2007	-	-	0/1/1/1	-
5	NAG	A	1606	1	-	5/6/23/26	0/1/1/1
5	NAG	A	1471	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1444	1	-	0/6/23/26	0/1/1/1
9	EDO	B	2007	-	-	1/1/1/1	-
5	NAG	A	1512	1	-	4/6/23/26	0/1/1/1
5	NAG	B	1471	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1644	1	-	0/6/23/26	0/1/1/1
6	9OK	B	2000	-	-	2/12/31/31	0/2/2/2
5	NAG	A	1215	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1483	1	-	0/6/23/26	0/1/1/1
9	EDO	A	2010	-	-	0/1/1/1	-
5	NAG	A	1539	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1319	1	-	2/6/23/26	0/1/1/1
9	EDO	A	2011	-	-	1/1/1/1	-
9	EDO	A	2006	-	-	0/1/1/1	-
9	EDO	A	2020	-	-	0/1/1/1	-
9	EDO	B	2016	-	-	1/1/1/1	-
5	NAG	A	1555	1	-	4/6/23/26	0/1/1/1
7	9WZ	B	2003	-	-	9/13/32/32	0/2/2/2
9	EDO	B	2015	-	-	1/1/1/1	-
9	EDO	A	2012	-	-	1/1/1/1	-
5	NAG	A	1125	1	-	2/6/23/26	0/1/1/1

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	2003	9WZ	O2-CAI	-8.07	1.19	1.39
7	B	2003	9WZ	O2-CAI	-8.03	1.19	1.39
6	B	2000	9OK	CAH-CAK	-5.38	1.36	1.51
7	A	2003	9WZ	CAH-CAK	-5.23	1.36	1.51
8	A	2004	975	CAH-CAK	-5.03	1.37	1.51

The worst 5 of 68 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	2003	9WZ	O2-CAI-CAD	-10.86	88.68	122.11
5	A	1512	NAG	C1-O5-C5	8.57	123.67	112.19
7	B	2003	9WZ	O2-CAI-CAD	-8.29	96.59	122.11
5	A	1215	NAG	C1-O5-C5	5.09	119.01	112.19
5	B	1319	NAG	C1-O5-C5	4.94	118.80	112.19

There are no chirality outliers.

5 of 69 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	2000	9OK	CAH-CAG-O2-C2
6	B	2000	9OK	CAH-CAG-O2-C2
7	A	2003	9WZ	C1-C2-O2-CAI
7	A	2003	9WZ	OAA-CAG-CAH-CAK
7	B	2003	9WZ	C1-C2-O2-CAI

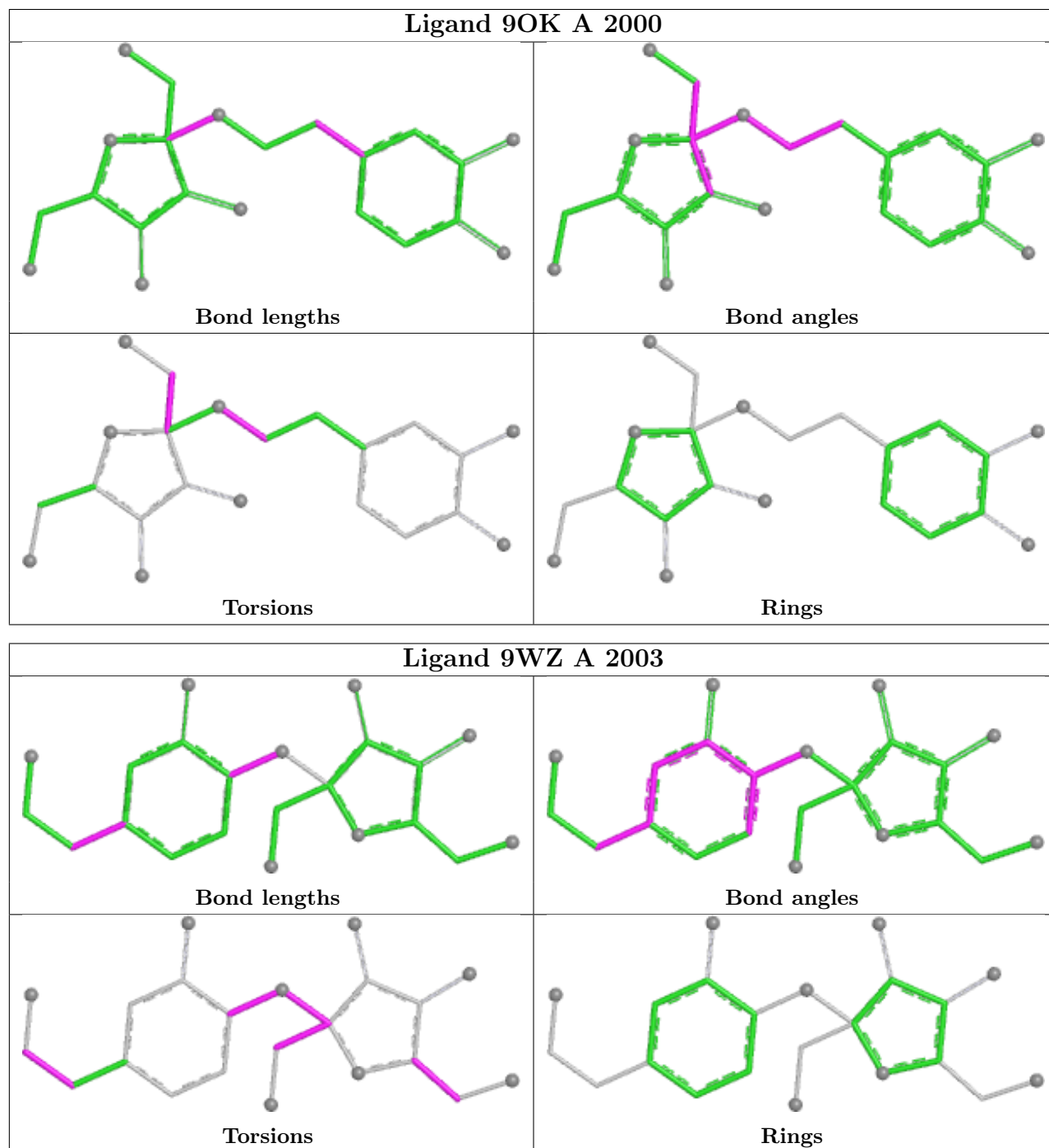
There are no ring outliers.

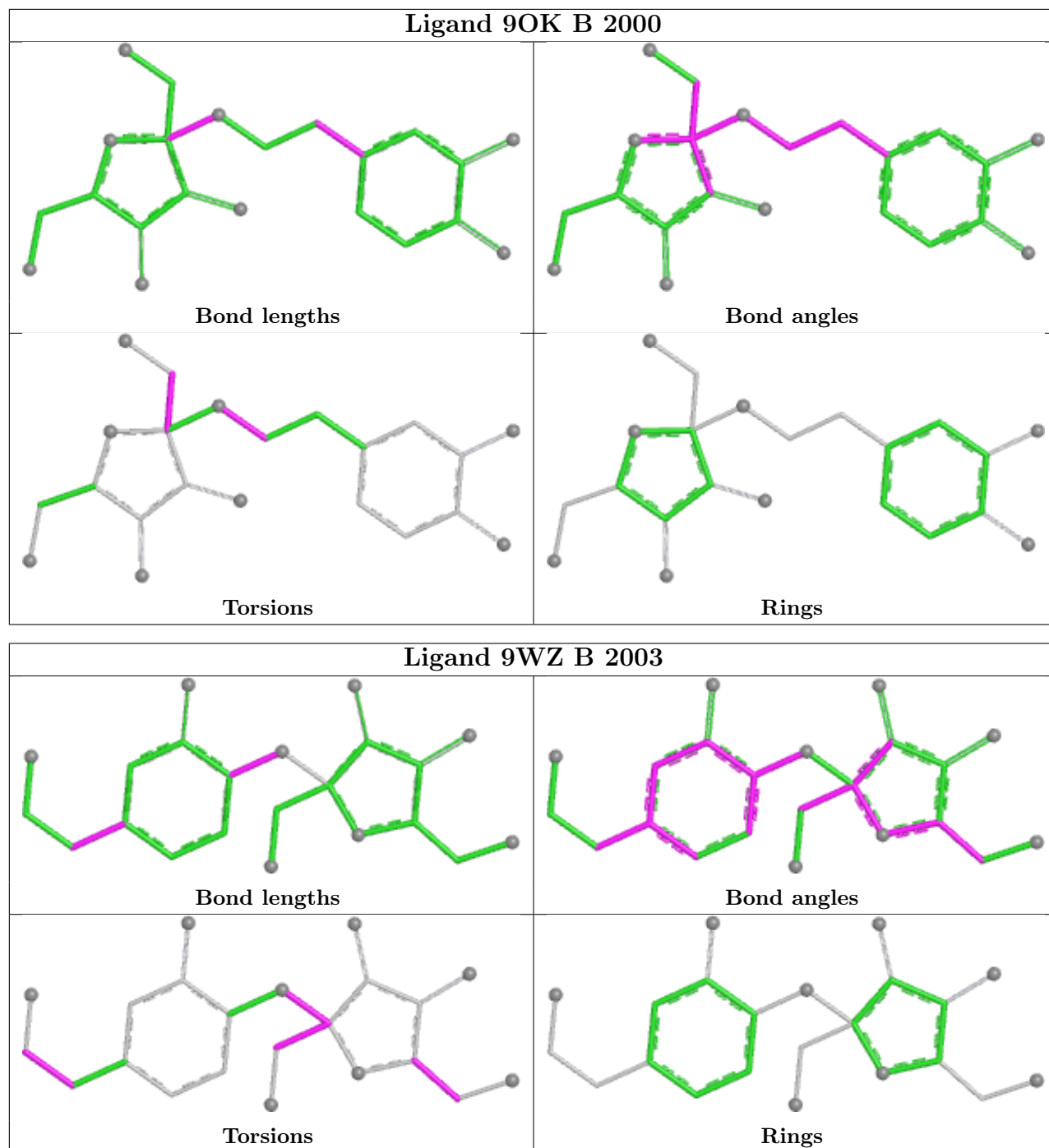
7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	2019	EDO	1	0
7	A	2003	9WZ	1	0
9	B	2019	EDO	1	0
5	A	1606	NAG	2	0
5	B	1444	NAG	1	0
9	B	2007	EDO	1	0
7	B	2003	9WZ	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	624/665 (93%)	-0.01	17 (2%) 56 62	9, 22, 33, 43	4 (0%)
1	B	624/665 (93%)	0.09	18 (2%) 53 60	10, 25, 37, 46	4 (0%)
All	All	1248/1330 (93%)	0.04	35 (2%) 55 61	9, 24, 36, 46	8 (0%)

The worst 5 of 35 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	291	TYR	4.1
1	A	341	GLN	3.7
1	B	481	MET	3.5
1	B	42	CYS	3.5
1	B	620	GLY	3.4

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
3	MAN	D	6	11/12	0.36	0.22	77,82,84,84	0
4	NAG	E	2	14/15	0.41	0.20	72,79,81,95	0
4	NAG	H	2	14/15	0.45	0.21	73,76,81,81	0
2	MAN	F	6	11/12	0.59	0.19	56,60,61,62	0
3	MAN	G	6	11/12	0.62	0.18	73,77,78,79	0

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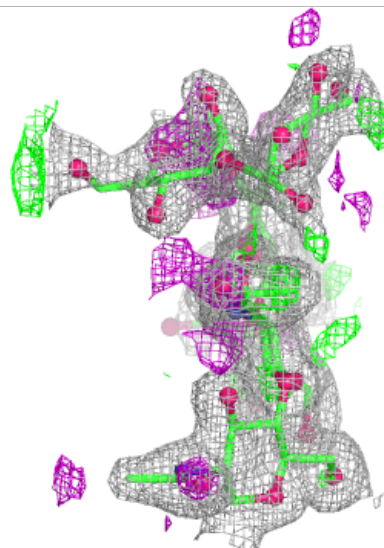
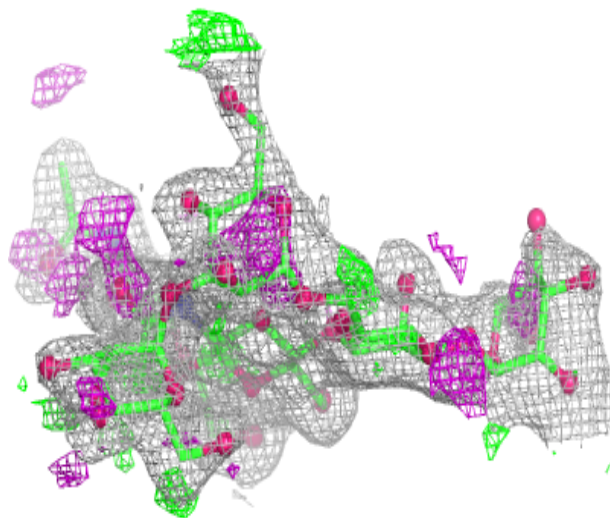
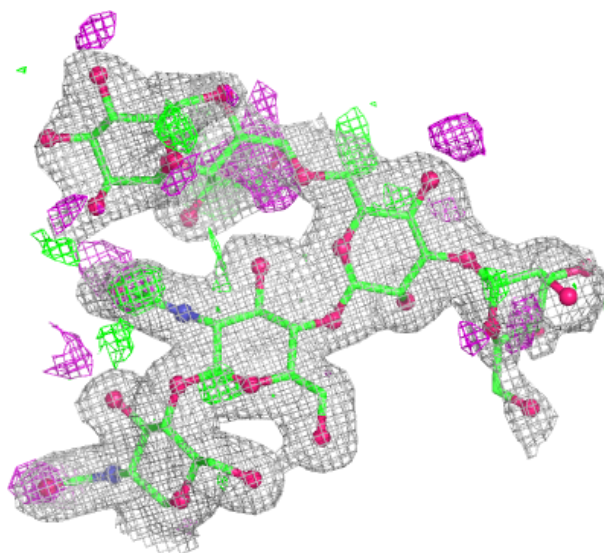
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MAN	G	9	11/12	0.64	0.21	55,59,59,60	0
2	MAN	C	6	11/12	0.66	0.18	57,63,64,69	0
3	MAN	D	9	11/12	0.70	0.19	55,59,60,61	0
2	MAN	C	4	11/12	0.73	0.20	48,50,52,52	0
3	MAN	G	5	11/12	0.80	0.15	56,59,60,66	0
3	MAN	G	4	11/12	0.81	0.14	39,42,44,49	0
3	MAN	D	5	11/12	0.81	0.14	56,59,61,68	0
4	NAG	H	1	14/15	0.82	0.14	47,51,57,65	0
2	MAN	C	5	11/12	0.82	0.18	50,50,51,51	0
3	MAN	D	8	11/12	0.83	0.15	39,41,43,49	0
3	MAN	D	4	11/12	0.84	0.13	38,42,43,48	0
2	MAN	F	5	11/12	0.84	0.15	56,56,56,57	0
2	MAN	F	4	11/12	0.84	0.14	52,54,55,56	0
3	MAN	G	8	11/12	0.88	0.12	41,44,46,50	0
2	BMA	C	3	11/12	0.88	0.14	34,40,45,48	0
4	NAG	E	1	14/15	0.88	0.11	44,48,54,63	0
2	BMA	F	3	11/12	0.89	0.13	38,43,48,49	0
2	NAG	C	2	14/15	0.92	0.11	24,25,26,30	0
3	BMA	D	3	11/12	0.94	0.07	23,24,26,31	0
3	MAN	G	7	11/12	0.94	0.08	22,24,30,35	0
2	NAG	C	1	14/15	0.95	0.08	21,22,25,25	0
3	BMA	G	3	11/12	0.95	0.08	25,25,28,33	0
3	NAG	D	2	14/15	0.96	0.07	18,19,20,21	0
3	MAN	D	7	11/12	0.96	0.07	20,23,29,33	0
2	NAG	F	2	14/15	0.96	0.07	26,27,29,33	0
3	MAN	G	10	11/12	0.97	0.06	18,19,19,20	0
3	NAG	G	2	14/15	0.97	0.06	20,21,22,23	0
2	NAG	F	1	14/15	0.97	0.06	23,25,27,27	0
3	NAG	D	1	14/15	0.97	0.06	16,17,17,17	0
3	NAG	G	1	14/15	0.97	0.06	18,18,19,19	0
3	MAN	D	10	11/12	0.98	0.05	17,17,18,18	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.

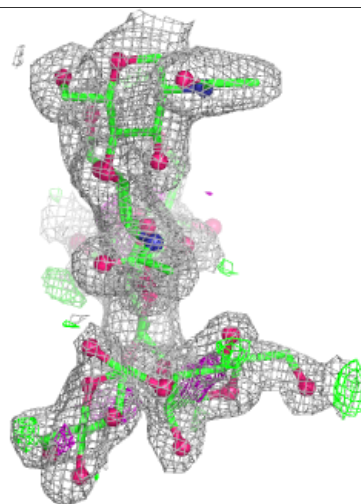
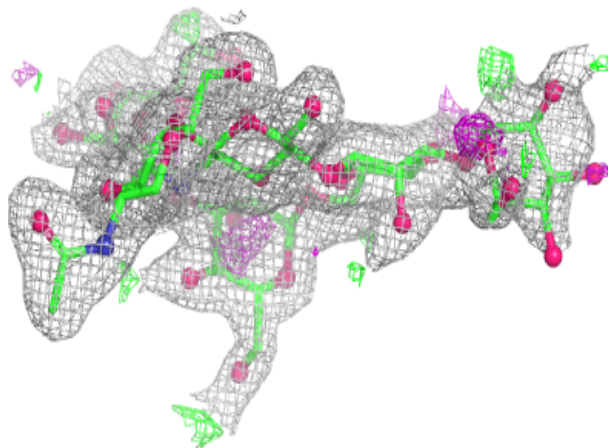
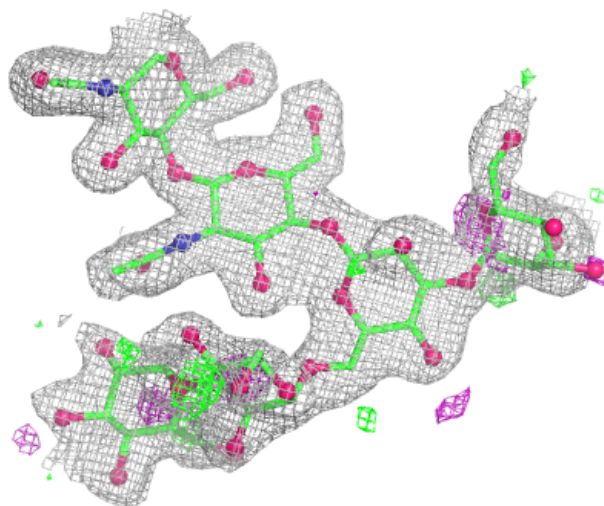
Electron density around Chain C:

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



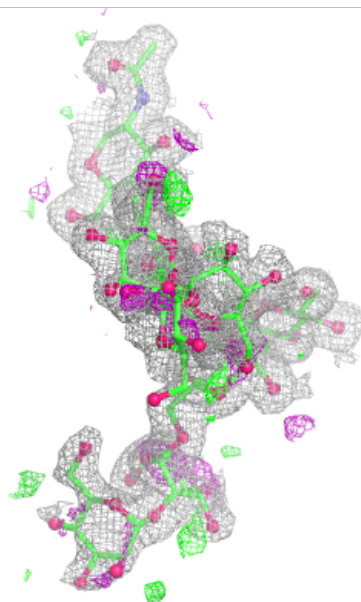
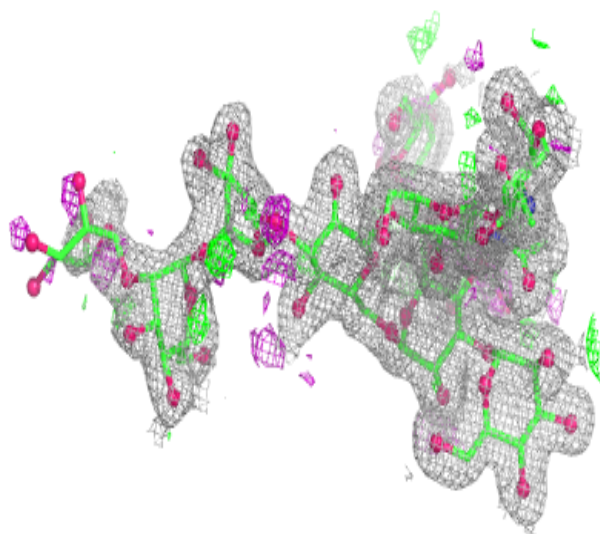
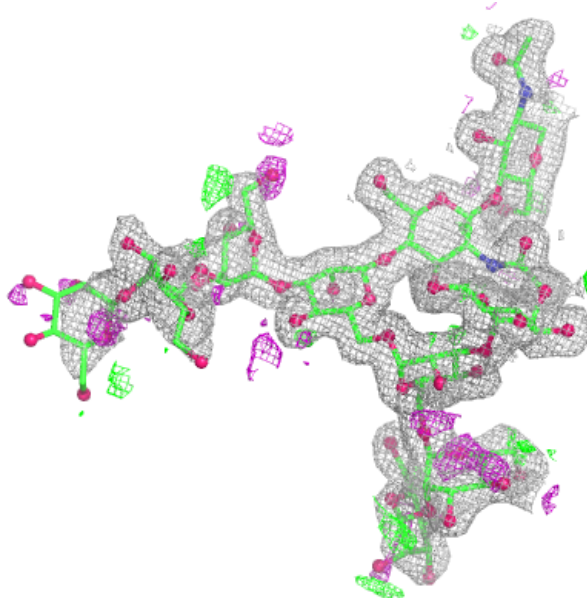
Electron density around Chain F:

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



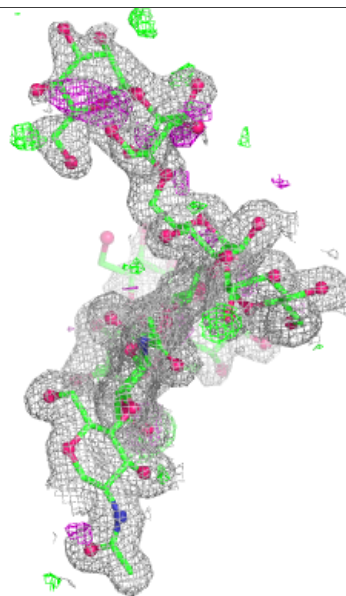
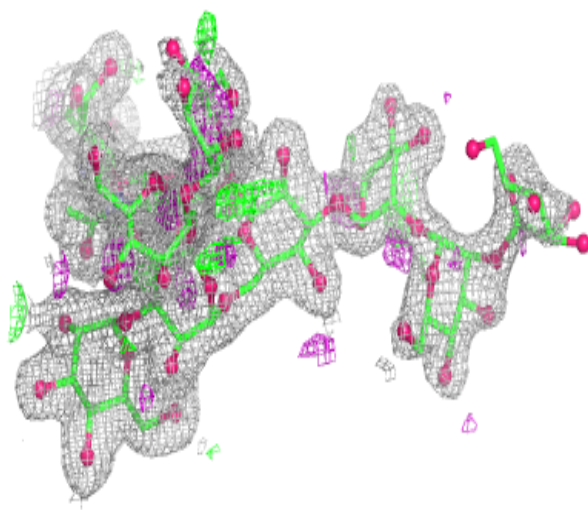
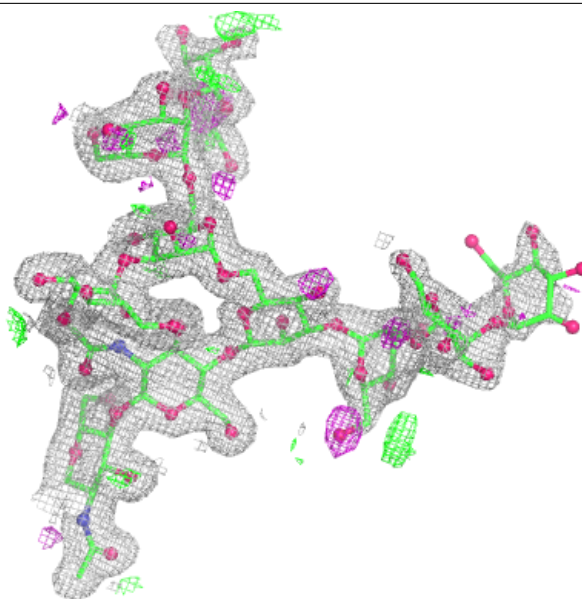
Electron density around Chain D:

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



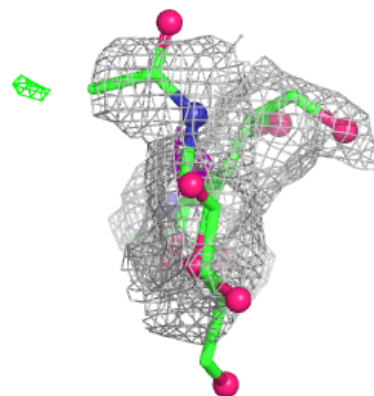
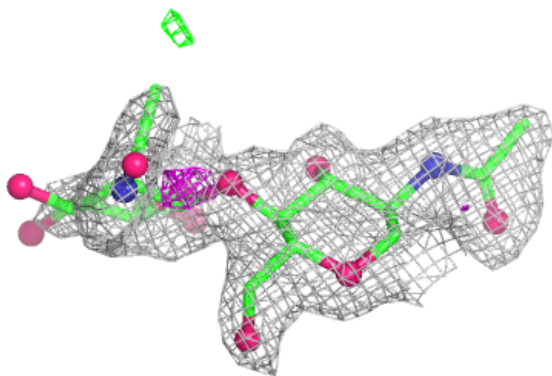
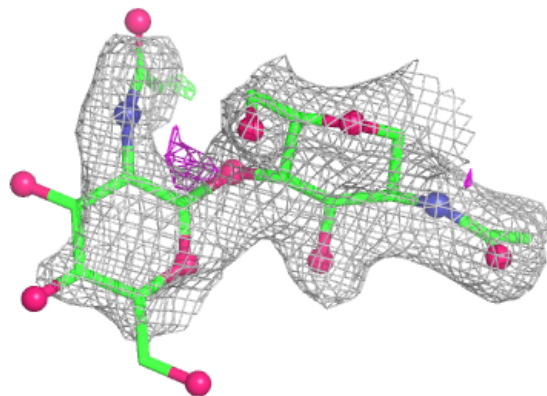
Electron density around Chain G:

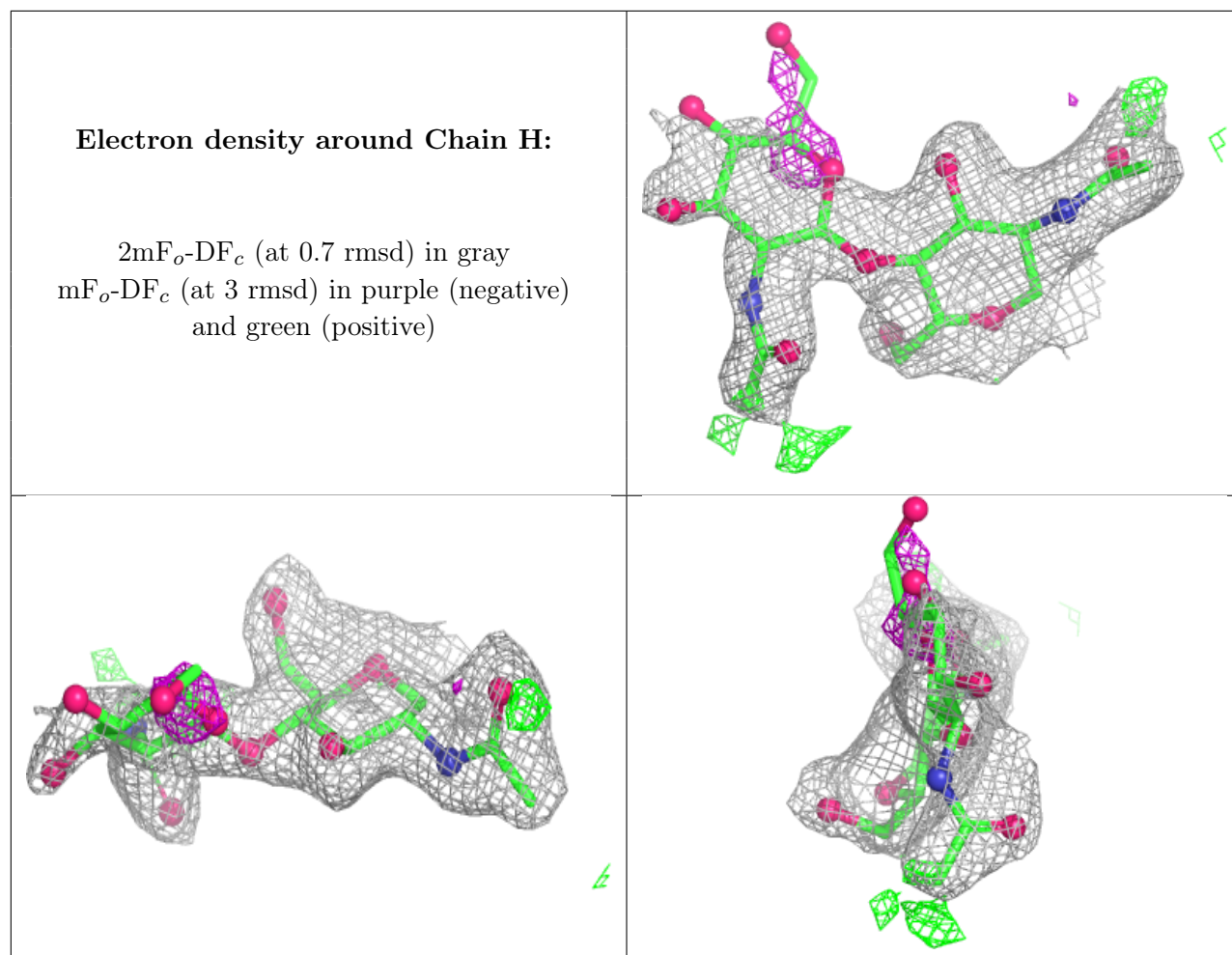
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain 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($\text{\AA}^2$)	Q<0.9
5	NAG	B	1319	14/15	0.52	0.20	64,68,71,71	0
5	NAG	B	1242	14/15	0.55	0.19	62,67,68,68	0
5	NAG	B	1215	14/15	0.60	0.20	59,65,67,68	0
5	NAG	A	1512	14/15	0.61	0.22	63,68,73,78	0
5	NAG	A	1555	14/15	0.62	0.19	39,42,43,44	0
5	NAG	B	1125	14/15	0.63	0.21	54,62,64,64	0
9	EDO	B	2008	4/4	0.66	0.29	49,49,50,50	0
5	NAG	A	1125	14/15	0.67	0.20	54,62,63,64	0
7	9WZ	B	2003	22/22	0.68	0.23	21,26,27,28	22
7	9WZ	A	2003	22/22	0.68	0.23	21,26,26,27	22

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	B	1052	14/15	0.70	0.18	55,59,62,62	0
5	NAG	A	1215	14/15	0.70	0.18	53,59,60,60	0
5	NAG	A	1242	14/15	0.71	0.17	51,55,57,57	0
9	EDO	A	2007	4/4	0.71	0.26	49,49,49,50	0
9	EDO	A	2010	4/4	0.71	0.21	48,49,49,50	0
9	EDO	A	2016	4/4	0.71	0.27	52,53,53,53	0
5	NAG	A	1606	14/15	0.71	0.17	44,47,48,48	0
9	EDO	A	2009	4/4	0.72	0.21	56,57,57,57	0
5	NAG	A	1052	14/15	0.73	0.17	48,52,53,54	0
9	EDO	B	2020	4/4	0.73	0.21	60,60,60,60	0
5	NAG	A	1319	14/15	0.74	0.17	55,59,61,62	0
9	EDO	A	2019	4/4	0.74	0.27	33,33,33,33	4
9	EDO	B	2021	4/4	0.74	0.23	56,56,56,56	0
9	EDO	B	2007	4/4	0.75	0.30	47,47,47,48	0
5	NAG	A	1539	14/15	0.75	0.17	44,49,51,52	0
8	975	B	2004	11/11	0.76	0.16	27,38,40,41	0
8	975	A	2004	11/11	0.76	0.16	32,37,38,39	0
9	EDO	B	2014	4/4	0.78	0.19	53,53,53,53	0
9	EDO	B	2019	4/4	0.78	0.25	35,35,35,35	4
9	EDO	B	2015	4/4	0.79	0.21	46,46,46,46	0
5	NAG	A	1483	14/15	0.80	0.15	41,44,45,46	0
9	EDO	B	2006	4/4	0.80	0.20	48,48,48,48	0
9	EDO	A	2006	4/4	0.80	0.18	48,48,48,48	0
5	NAG	B	1483	14/15	0.82	0.14	47,50,52,52	0
9	EDO	A	2011	4/4	0.83	0.22	39,40,40,41	0
9	EDO	A	2012	4/4	0.83	0.18	56,57,57,57	0
5	NAG	B	1539	14/15	0.83	0.15	46,51,53,53	0
9	EDO	B	2016	4/4	0.84	0.18	51,51,51,52	0
9	EDO	B	2017	4/4	0.86	0.23	49,49,49,50	0
9	EDO	A	2020	4/4	0.86	0.17	37,38,38,39	0
5	NAG	A	1236	14/15	0.86	0.12	32,33,35,35	0
9	EDO	B	2013	4/4	0.86	0.21	45,45,45,46	0
9	EDO	A	2015	4/4	0.87	0.16	45,45,45,46	0
5	NAG	B	1555	14/15	0.87	0.11	36,38,39,39	0
9	EDO	A	2013	4/4	0.87	0.27	44,44,44,44	0
5	NAG	B	1357	14/15	0.89	0.12	38,40,41,41	0
5	NAG	A	1357	14/15	0.89	0.11	33,34,35,36	0
5	NAG	B	1644	14/15	0.91	0.09	27,27,30,30	0
5	NAG	A	1471	14/15	0.92	0.09	31,32,33,35	0
6	9OK	B	2000	22/22	0.92	0.12	20,24,38,40	0
5	NAG	A	1444	14/15	0.92	0.09	27,27,28,28	0
5	NAG	B	1471	14/15	0.92	0.09	33,34,36,36	0

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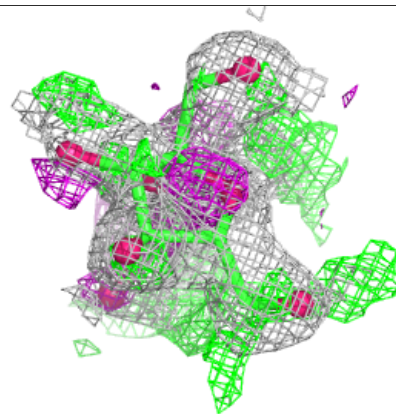
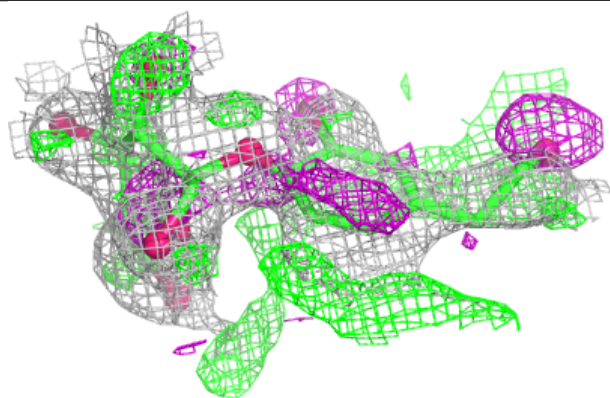
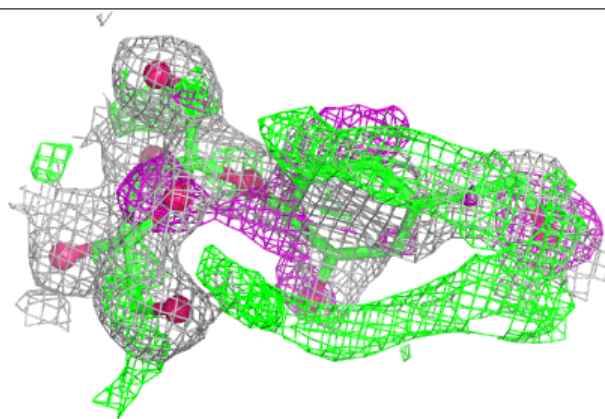
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	9OK	A	2000	22/22	0.93	0.12	18,22,37,39	0
5	NAG	A	1644	14/15	0.93	0.09	25,26,29,29	0
5	NAG	B	1444	14/15	0.95	0.08	28,29,31,31	0

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.

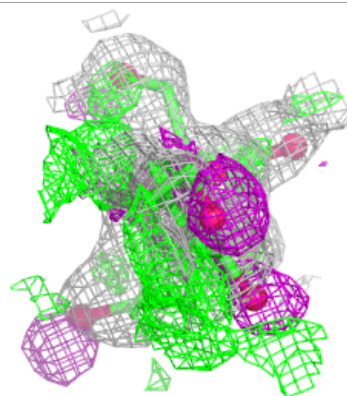
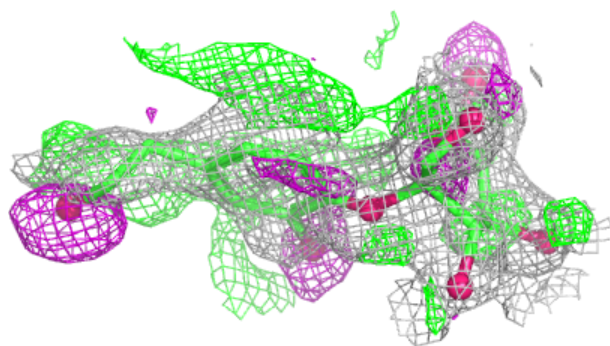
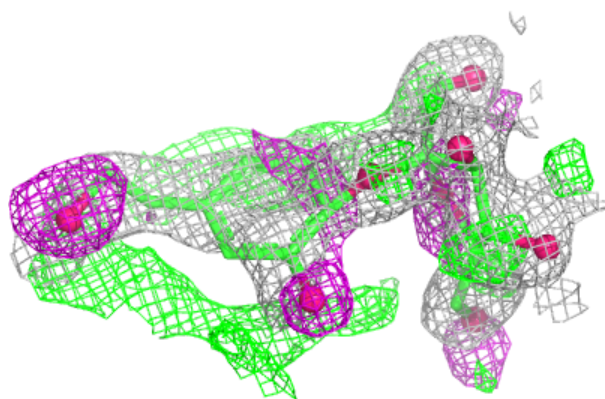
Electron density around 9WZ B 2003:

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

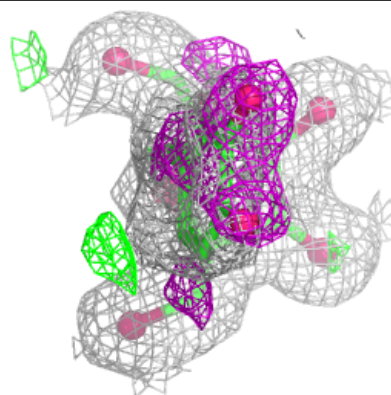
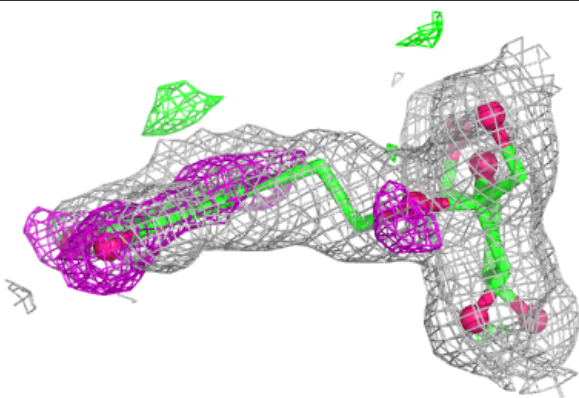
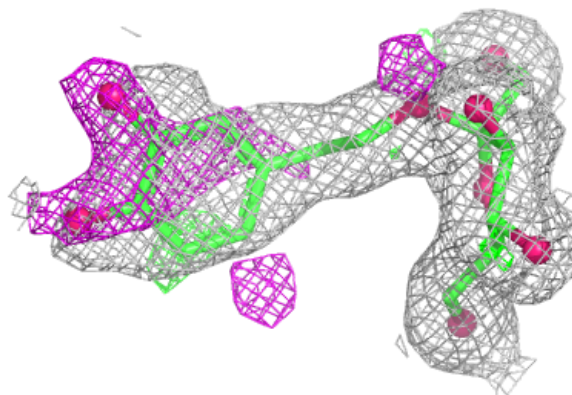


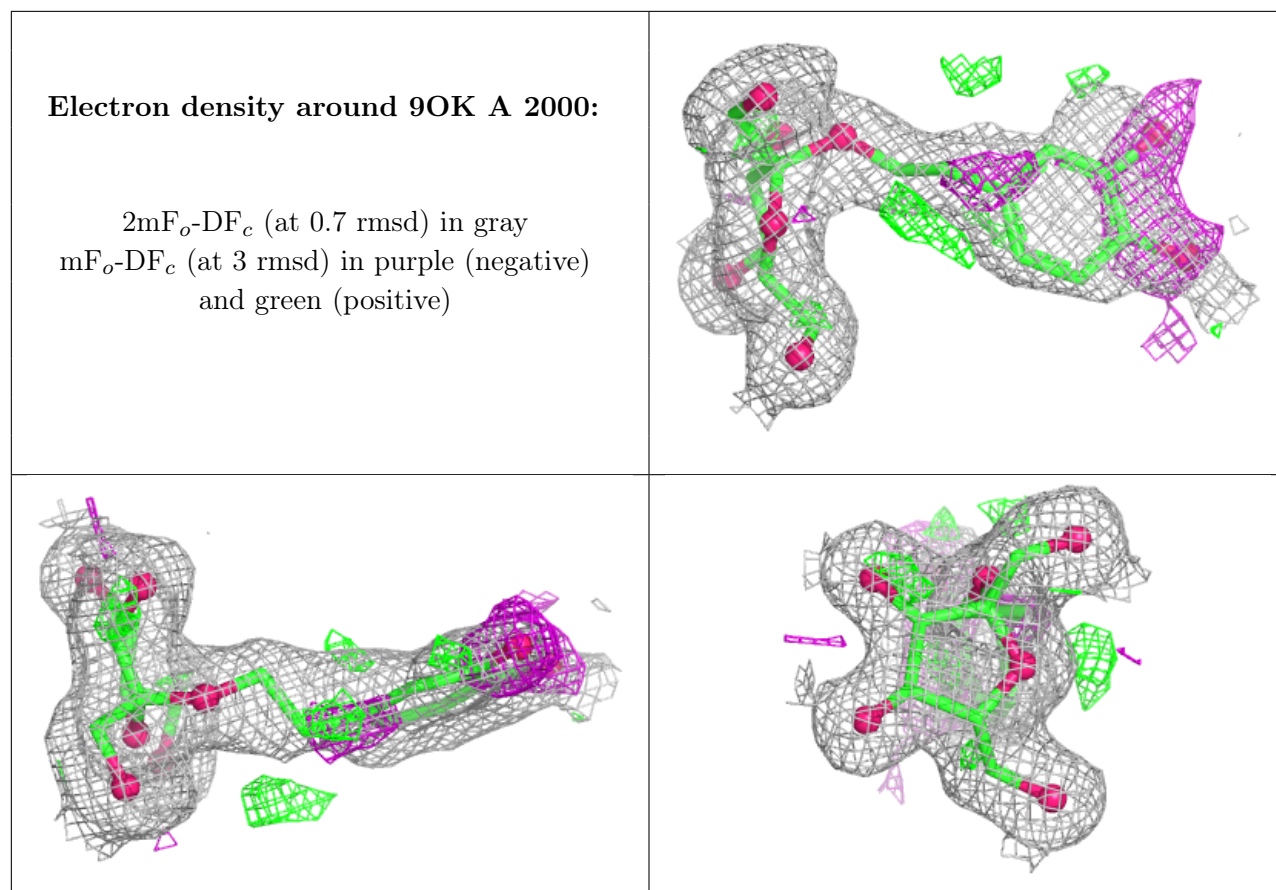
Electron density around 9WZ A 2003:

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

**Electron density around 9OK B 2000:**

$2mF_o-DF_c$ (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.