



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 17, 2026 – 09:18 PM UTC

PDB ID : 8ORP / pdb\_00008orp  
Title : Crystal structure of Drosophila melanogaster alpha-amylase in complex with the inhibitor acarbose  
Authors : Aghajari, N.; Haser, R.  
Deposited on : 2023-04-16  
Resolution : 2.00 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](mailto:validation@mail.wwpdb.org)

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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>.

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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  
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

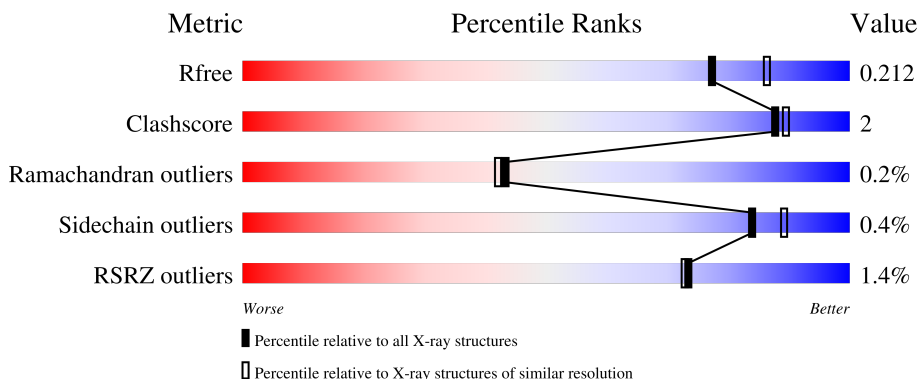
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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.



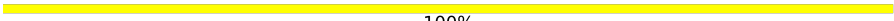
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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  $\geq 3$ , 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	494	 92% 5%
1	B	494	 91% 5%
2	C	3	 67% 33%
2	E	3	 67% 33%
3	D	2	 100%

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Mol	Chain	Length	Quality of chain
3	F	2	 100%

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	RY7	C	3	X	-	-	-
2	RY7	E	3	X	-	-	-
3	RY7	D	2	X	-	-	-
3	RY7	F	2	X	-	-	-

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 8503 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 Alpha-amylase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	476	Total	C	N	O	S	0	2	0
			3663	2281	642	723	17			
1	B	476	Total	C	N	O	S	0	0	0
			3657	2277	642	721	17			

- Molecule 2 is an oligosaccharide called 4,6-dideoxy-4- $\{[(1S,2S,3S,4R,5R)-2,3,4\text{-trihydroxy-5-(hydroxymethyl)cyclohexyl]amino}\}$ -alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	3	Total	C	N	O	0	0	0
			43	25	1	17			
2	E	3	Total	C	N	O	0	0	0
			43	25	1	17			

- Molecule 3 is an oligosaccharide called 4,6-dideoxy-4- $\{[(1S,2S,3S,4R,5R)-2,3,4\text{-trihydroxy-5-(hydroxymethyl)cyclohexyl]amino}\}$ -alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	D	2	Total	C	N	O	0	0	0
			32	19	1	12			
3	F	2	Total	C	N	O	0	0	0
			32	19	1	12			

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total	Cl	0
			1	1	
4	B	1	Total	Cl	0
			1	1	

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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

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

- Molecule 6 is STRONTIUM ION (CCD ID: SR) (formula: Sr) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total Sr 2 2	0	0
6	B	2	Total Sr 2 2	0	0

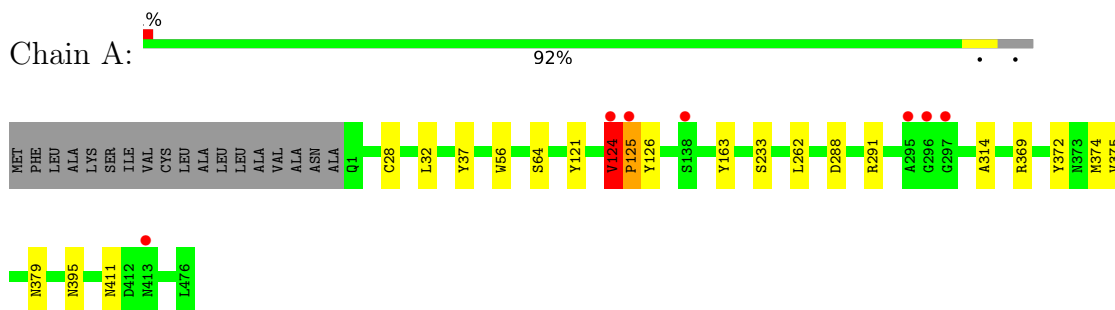
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	440	Total O 440 440	0	0
7	B	435	Total O 435 435	0	0

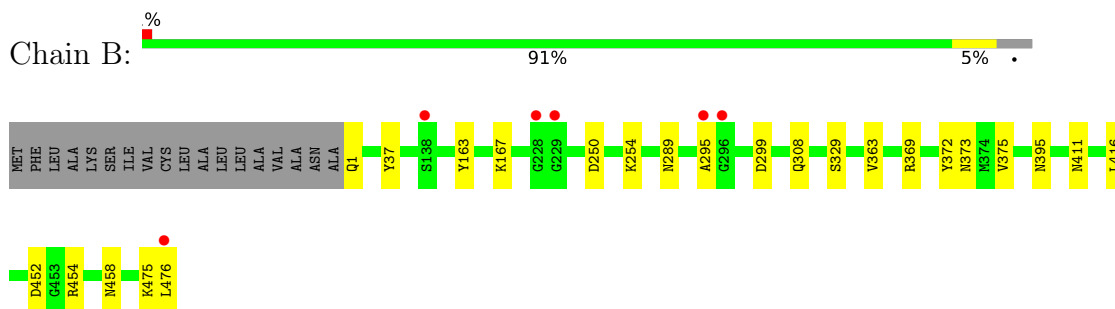
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

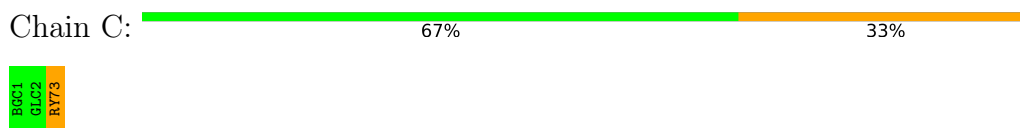
- Molecule 1: Alpha-amylase A



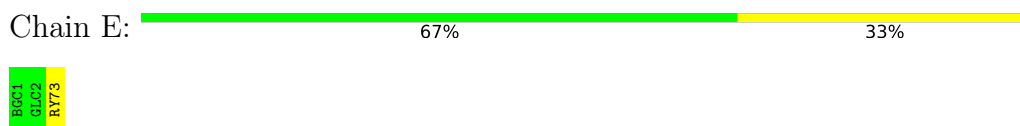
- Molecule 1: Alpha-amylase A



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,2S,3S,4R,5R)-2,3,4\text{-trihydroxy-5-(hydroxymethyl)cyclohexyl]amino}\}$ -alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,2S,3S,4R,5R)-2,3,4\text{-trihydroxy-5-(hydroxymethyl)cyclohexyl]amino}\}$ -alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose



- Molecule 3: 4,6-dideoxy-4-[(1S,2S,3S,4R,5R)-2,3,4-trihydroxy-5-(hydroxymethyl)cyclohexyl]amino}-alpha-D-glucofuranose-(1-4)-alpha-D-glucofuranose

Chain D:  100%

GLC1  
RY72

- Molecule 3: 4,6-dideoxy-4-[(1S,2S,3S,4R,5R)-2,3,4-trihydroxy-5-(hydroxymethyl)cyclohexyl]amino}-alpha-D-glucofuranose-(1-4)-alpha-D-glucofuranose

Chain F:  100%

GLC1  
RY72

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.08Å 72.54Å 98.21Å 90.00° 98.41° 90.00°	Depositor
Resolution (Å)	29.57 – 2.00 29.57 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.57-2.00) 99.7 (29.57-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 2.00Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.174 , 0.212 0.174 , 0.212	Depositor DCC
$R_{free}$ test set	3759 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.6	Xtrriage
Anisotropy	0.634	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 57.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8503	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.29% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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: CL, BGC, EDO, GLC, SR, RY7

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	0.16	0/3759	0.36	0/5100
1	B	0.16	0/3747	0.36	0/5084
All	All	0.16	0/7506	0.36	0/10184

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.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	124	VAL	Peptide

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3663	0	3422	14	0
1	B	3657	0	3412	16	0
2	C	43	0	17	1	0
2	E	43	0	17	0	0
3	D	32	0	9	0	0
3	F	32	0	9	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	72	0	108	4	0
5	B	80	0	118	3	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
7	A	440	0	0	0	0
7	B	435	0	0	0	0
All	All	8503	0	7112	30	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

The worst 5 of 30 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:295:ALA:HB3	5:B:515:EDO:H11	1.84	0.59
1:B:475:LYS:O	1:B:476:LEU:HB2	2.05	0.57
1:B:254:LYS:NZ	1:B:299:ASP:OD2	2.38	0.56
1:A:379:ASN:HA	5:A:515:EDO:H22	1.88	0.54
5:A:508:EDO:H21	1:B:454:ARG:HD2	1.89	0.53

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	476/494 (96%)	463 (97%)	11 (2%)	2 (0%)	30	27
1	B	474/494 (96%)	462 (98%)	12 (2%)	0	100	100
All	All	950/988 (96%)	925 (97%)	23 (2%)	2 (0%)	43	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	125	PRO
1	A	124	VAL

### 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	391/403 (97%)	390 (100%)	1 (0%)	86	91
1	B	389/403 (96%)	387 (100%)	2 (0%)	81	87
All	All	780/806 (97%)	777 (100%)	3 (0%)	84	89

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	233	SER
1	B	329	SER
1	B	363	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	458	ASN
1	B	413	ASN
1	A	458	ASN
1	A	417	ASN
1	B	205	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

10 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	BGC	C	1	2	11,11,12	0.16	0	16,16,17	0.19	0
2	GLC	C	2	2	11,11,12	0.72	0	15,15,17	0.77	0
2	RY7	C	3	3,2	21,22,23	1.38	4 (19%)	26,32,34	1.33	4 (15%)
3	GLC	D	1	3,2	11,11,12	0.56	0	15,15,17	1.11	1 (6%)
3	RY7	D	2	3	21,22,23	1.38	3 (14%)	26,32,34	1.54	3 (11%)
2	BGC	E	1	2	11,11,12	0.16	0	16,16,17	0.22	0
2	GLC	E	2	2	11,11,12	0.66	0	15,15,17	0.79	0
2	RY7	E	3	3,2	21,22,23	1.39	3 (14%)	26,32,34	1.49	4 (15%)
3	GLC	F	1	3,2	11,11,12	0.57	0	15,15,17	1.14	2 (13%)
3	RY7	F	2	3	21,22,23	1.40	3 (14%)	26,32,34	1.45	3 (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	BGC	C	1	2	-	-	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	RY7	C	3	3,2	1/1/9/11	3/6/43/46	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GLC	D	1	3,2	-	0/2/19/22	0/1/1/1
3	RY7	D	2	3	1/1/9/11	2/6/43/46	0/2/2/2
2	BGC	E	1	2	-	-	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	RY7	E	3	3,2	1/1/9/11	3/6/43/46	0/2/2/2
3	GLC	F	1	3,2	-	0/2/19/22	0/1/1/1
3	RY7	F	2	3	1/1/9/11	3/6/43/46	0/2/2/2

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	2	RY7	O5-C1	3.32	1.49	1.43
3	D	2	RY7	O5-C1	3.20	1.49	1.43
2	E	3	RY7	O5-C1	3.17	1.49	1.43
2	C	3	RY7	O5-C1	2.96	1.48	1.43
2	E	3	RY7	O5-C5	2.37	1.48	1.43

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2	RY7	C5A-C7A-C1A	5.10	115.12	108.53
3	F	2	RY7	C5A-C7A-C1A	4.79	114.72	108.53
2	E	3	RY7	O6A-C6A-C5A	-3.79	102.61	111.26
2	E	3	RY7	C5A-C7A-C1A	3.49	113.04	108.53
2	C	3	RY7	O6A-C6A-C5A	-3.25	103.85	111.26

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	3	RY7	C5A
2	E	3	RY7	C5A
3	D	2	RY7	C5A
3	F	2	RY7	C5A

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	3	RY7	C7A-C1A-N4-C4
2	C	3	RY7	C4A-C5A-C6A-O6A
2	C	3	RY7	C7A-C5A-C6A-O6A
2	E	3	RY7	C7A-C1A-N4-C4

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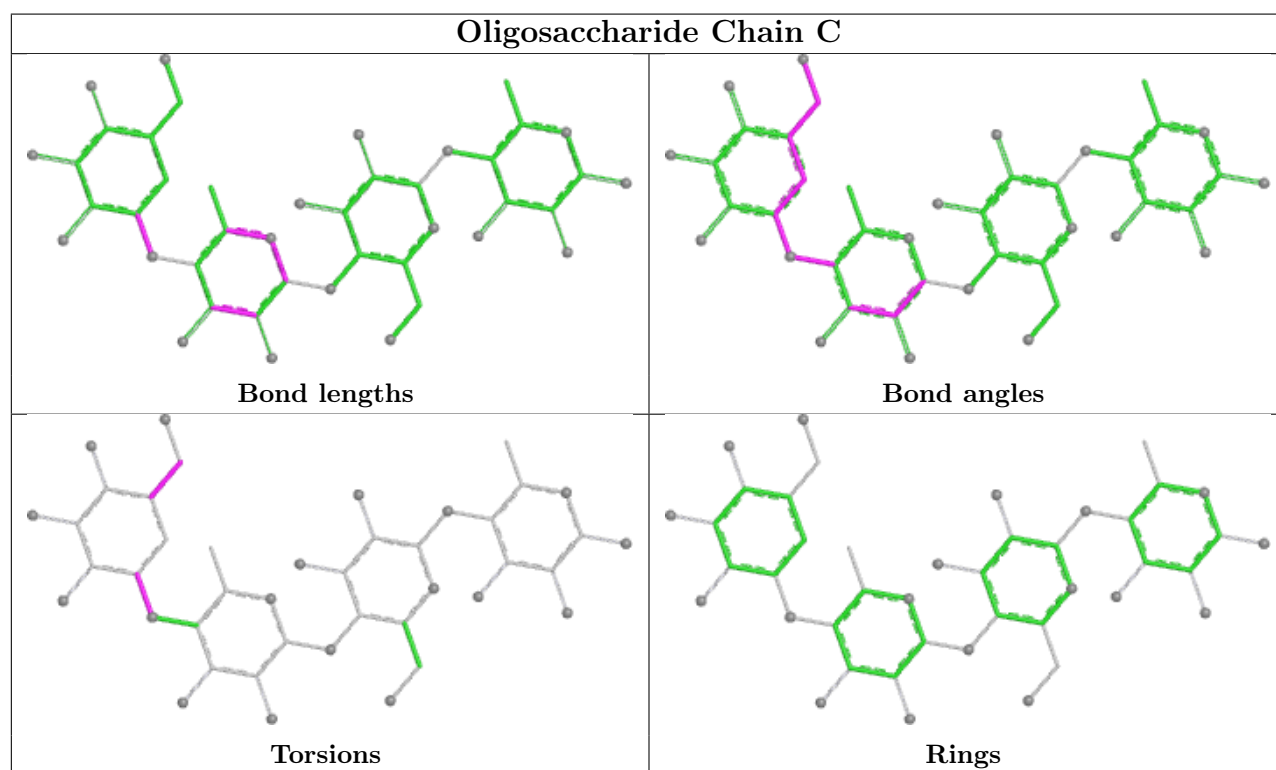
Mol	Chain	Res	Type	Atoms
2	E	3	RY7	C4A-C5A-C6A-O6A

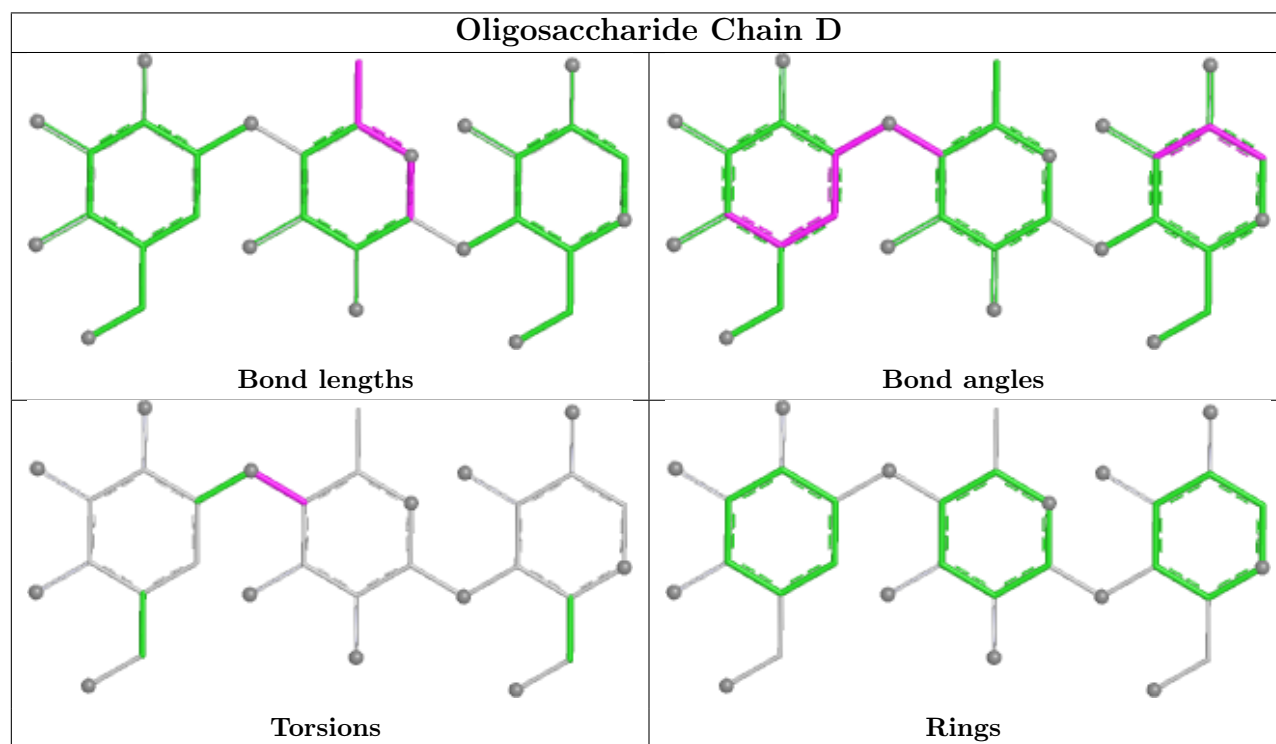
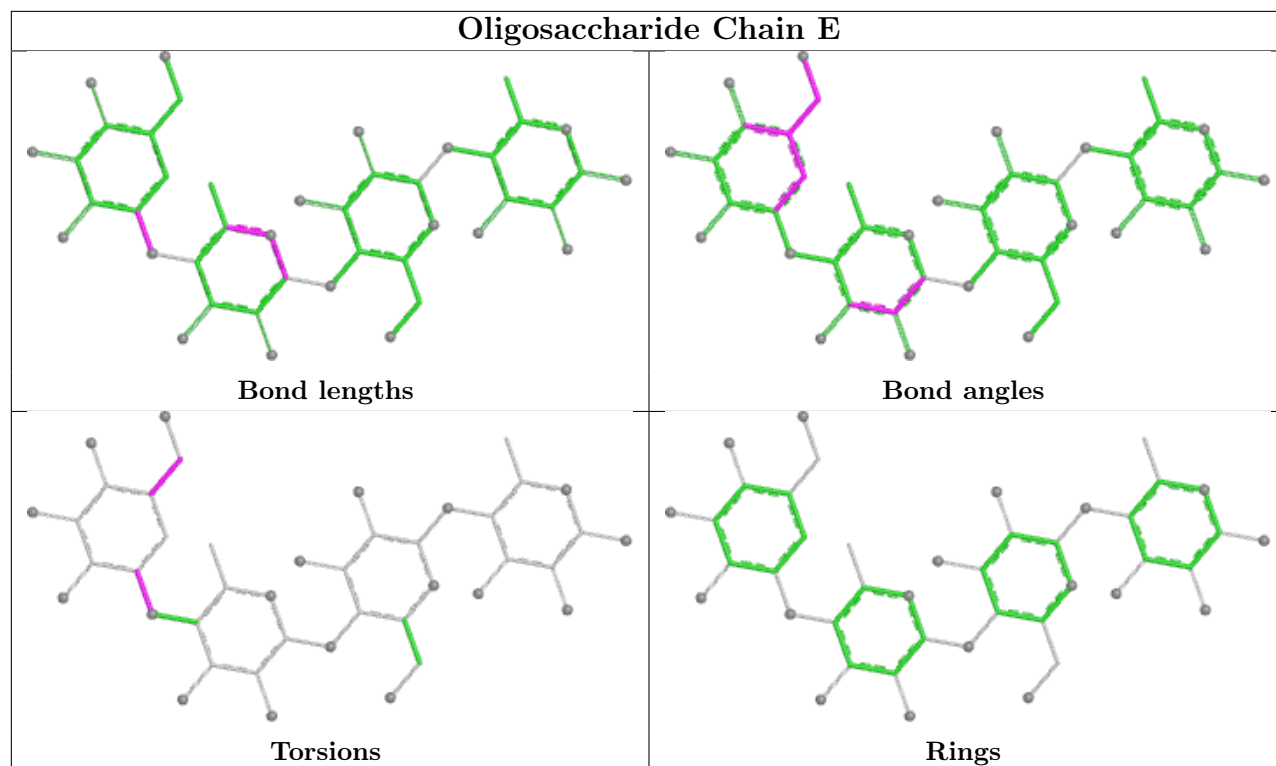
There are no ring outliers.

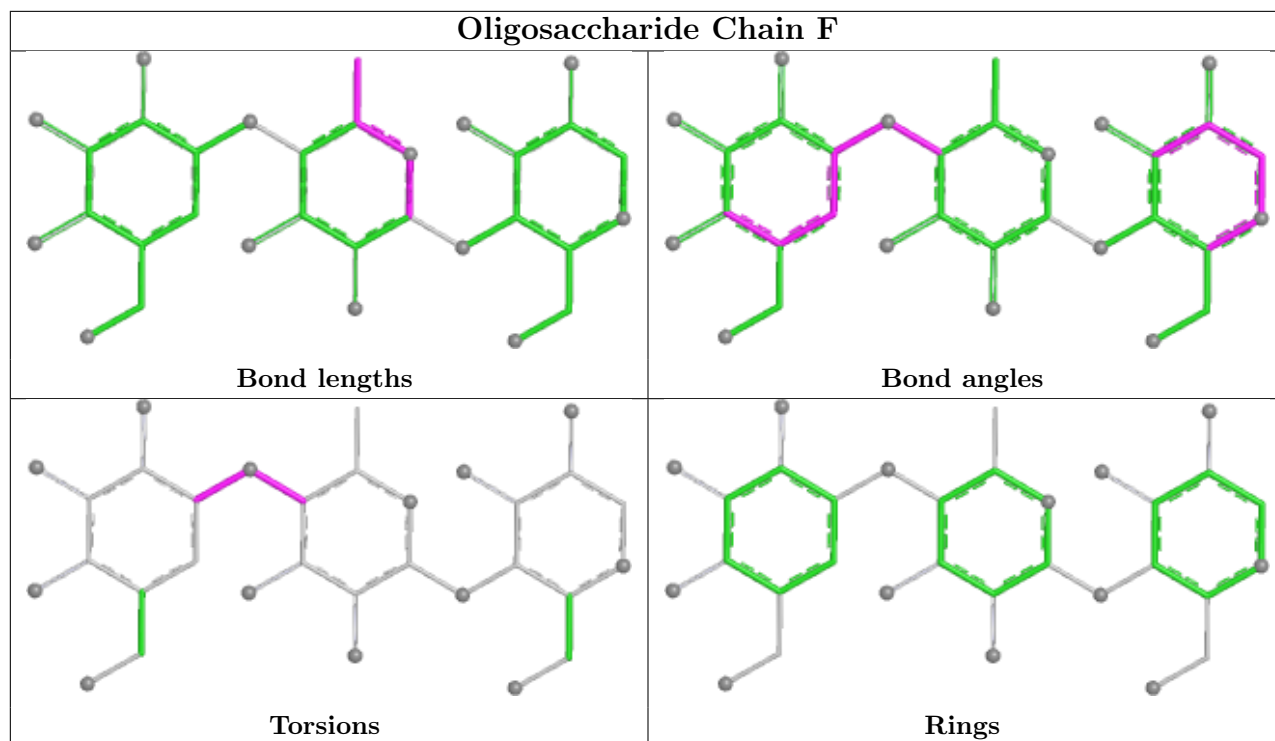
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	3	RY7	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 oligosaccharide.







## 5.6 Ligand geometry [i](#)

Of 44 ligands modelled in this entry, 6 are monoatomic - leaving 38 for Mogul analysis.

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	EDO	B	509	-	3,3,3	0.42	0	2,2,2	0.41	0
5	EDO	A	507	-	3,3,3	0.39	0	2,2,2	0.47	0
5	EDO	A	514	-	3,3,3	0.32	0	2,2,2	0.55	0
5	EDO	A	502	-	3,3,3	0.42	0	2,2,2	0.44	0
5	EDO	A	512	-	3,3,3	0.45	0	2,2,2	0.35	0
5	EDO	B	506	-	3,3,3	0.43	0	2,2,2	0.50	0
5	EDO	B	514	-	3,3,3	0.41	0	2,2,2	0.41	0
5	EDO	B	520	-	3,3,3	0.44	0	2,2,2	0.45	0
5	EDO	B	510	-	3,3,3	0.47	0	2,2,2	0.36	0
5	EDO	B	507	-	3,3,3	0.41	0	2,2,2	0.35	0
5	EDO	A	513	-	3,3,3	0.42	0	2,2,2	0.38	0
5	EDO	A	505	-	3,3,3	0.41	0	2,2,2	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	EDO	B	516	-	3,3,3	0.41	0	2,2,2	0.52	0
5	EDO	A	510	-	3,3,3	0.40	0	2,2,2	0.53	0
5	EDO	B	508	-	3,3,3	0.42	0	2,2,2	0.34	0
5	EDO	B	515	-	3,3,3	0.44	0	2,2,2	0.44	0
5	EDO	B	519	-	3,3,3	0.43	0	2,2,2	0.48	0
5	EDO	A	515	-	3,3,3	0.39	0	2,2,2	0.41	0
5	EDO	B	503	-	3,3,3	0.50	0	2,2,2	0.22	0
5	EDO	A	503	-	3,3,3	0.43	0	2,2,2	0.28	0
5	EDO	A	506	-	3,3,3	0.47	0	2,2,2	0.28	0
5	EDO	A	508	-	3,3,3	0.41	0	2,2,2	0.33	0
5	EDO	B	504	-	3,3,3	0.41	0	2,2,2	0.45	0
5	EDO	B	505	-	3,3,3	0.46	0	2,2,2	0.32	0
5	EDO	A	511	-	3,3,3	0.42	0	2,2,2	0.42	0
5	EDO	B	511	-	3,3,3	0.43	0	2,2,2	0.41	0
5	EDO	B	513	-	3,3,3	0.40	0	2,2,2	0.38	0
5	EDO	A	517	-	3,3,3	0.44	0	2,2,2	0.36	0
5	EDO	A	516	-	3,3,3	0.45	0	2,2,2	0.41	0
5	EDO	B	517	-	3,3,3	0.44	0	2,2,2	0.40	0
5	EDO	A	504	-	3,3,3	0.41	0	2,2,2	0.40	0
5	EDO	A	519	-	3,3,3	0.45	0	2,2,2	0.29	0
5	EDO	B	512	-	3,3,3	0.45	0	2,2,2	0.43	0
5	EDO	B	521	-	3,3,3	0.41	0	2,2,2	0.43	0
5	EDO	A	518	-	3,3,3	0.48	0	2,2,2	0.25	0
5	EDO	B	518	-	3,3,3	0.34	0	2,2,2	0.41	0
5	EDO	A	509	-	3,3,3	0.40	0	2,2,2	0.42	0
5	EDO	B	502	-	3,3,3	0.43	0	2,2,2	0.32	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
5	EDO	B	509	-	-	0/1/1/1	-
5	EDO	A	507	-	-	0/1/1/1	-
5	EDO	A	514	-	-	1/1/1/1	-
5	EDO	A	502	-	-	0/1/1/1	-
5	EDO	A	512	-	-	0/1/1/1	-
5	EDO	B	506	-	-	0/1/1/1	-
5	EDO	B	514	-	-	0/1/1/1	-
5	EDO	B	520	-	-	1/1/1/1	-
5	EDO	B	510	-	-	1/1/1/1	-
5	EDO	B	507	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	513	-	-	0/1/1/1	-
5	EDO	A	505	-	-	0/1/1/1	-
5	EDO	B	516	-	-	0/1/1/1	-
5	EDO	A	510	-	-	0/1/1/1	-
5	EDO	B	508	-	-	0/1/1/1	-
5	EDO	B	515	-	-	1/1/1/1	-
5	EDO	B	519	-	-	0/1/1/1	-
5	EDO	A	515	-	-	0/1/1/1	-
5	EDO	B	503	-	-	0/1/1/1	-
5	EDO	A	503	-	-	0/1/1/1	-
5	EDO	A	506	-	-	0/1/1/1	-
5	EDO	A	508	-	-	0/1/1/1	-
5	EDO	B	504	-	-	0/1/1/1	-
5	EDO	B	505	-	-	0/1/1/1	-
5	EDO	A	511	-	-	0/1/1/1	-
5	EDO	B	511	-	-	0/1/1/1	-
5	EDO	B	513	-	-	0/1/1/1	-
5	EDO	A	517	-	-	1/1/1/1	-
5	EDO	A	516	-	-	0/1/1/1	-
5	EDO	B	517	-	-	1/1/1/1	-
5	EDO	A	504	-	-	0/1/1/1	-
5	EDO	A	519	-	-	0/1/1/1	-
5	EDO	B	512	-	-	0/1/1/1	-
5	EDO	B	521	-	-	1/1/1/1	-
5	EDO	A	518	-	-	0/1/1/1	-
5	EDO	B	518	-	-	0/1/1/1	-
5	EDO	A	509	-	-	0/1/1/1	-
5	EDO	B	502	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	517	EDO	O1-C1-C2-O2
5	B	517	EDO	O1-C1-C2-O2
5	B	521	EDO	O1-C1-C2-O2
5	B	510	EDO	O1-C1-C2-O2
5	A	514	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	514	EDO	1	0
5	B	515	EDO	2	0
5	A	515	EDO	1	0
5	A	508	EDO	3	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	476/494 (96%)	-0.38	7 (1%) 72 71	10, 16, 28, 41	2 (0%)
1	B	476/494 (96%)	-0.44	6 (1%) 75 74	11, 16, 27, 38	0
All	All	952/988 (96%)	-0.41	13 (1%) 73 73	10, 16, 27, 41	2 (0%)

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124	VAL	5.3
1	A	296	GLY	3.9
1	B	296	GLY	3.4
1	B	476	LEU	3.4
1	A	125	PRO	2.8

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

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

### 6.3 Carbohydrates [i](#)

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, 95<sup>th</sup> 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.

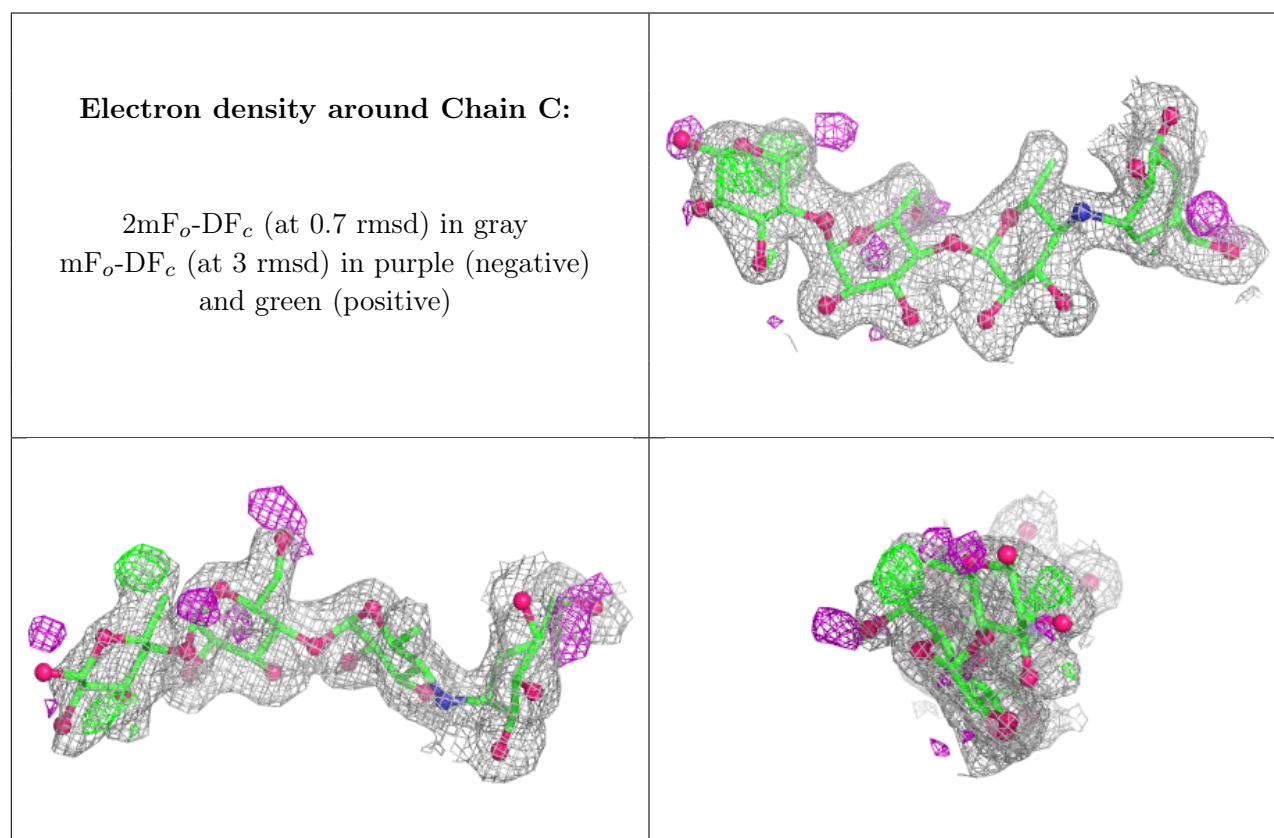
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	BGC	C	1	11/12	0.71	0.18	26,39,52,57	0
2	BGC	E	1	11/12	0.76	0.15	32,45,49,59	0
2	GLC	C	2	11/12	0.90	0.11	17,21,32,33	0
3	RY7	F	2	21/22	0.90	0.10	17,23,33,38	0
2	GLC	E	2	11/12	0.91	0.09	16,23,29,31	0

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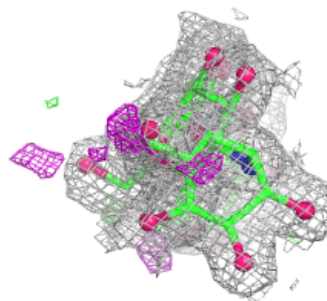
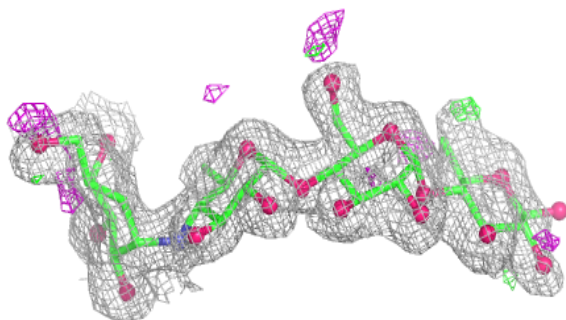
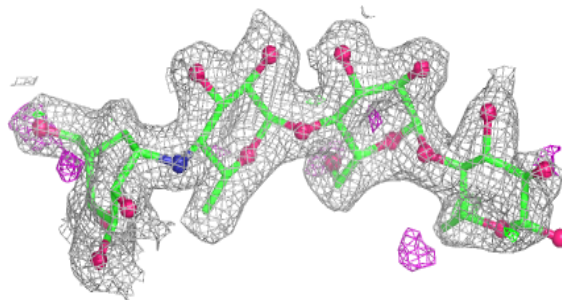
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	RY7	D	2	21/22	0.91	0.09	18,23,33,38	0
2	RY7	C	3	21/22	0.91	0.09	15,18,21,23	0
2	RY7	E	3	21/22	0.93	0.08	13,17,19,21	0
3	GLC	F	1	11/12	0.96	0.05	14,17,18,20	0
3	GLC	D	1	11/12	0.97	0.05	15,17,19,19	0

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.

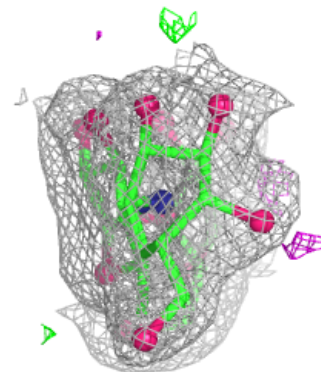
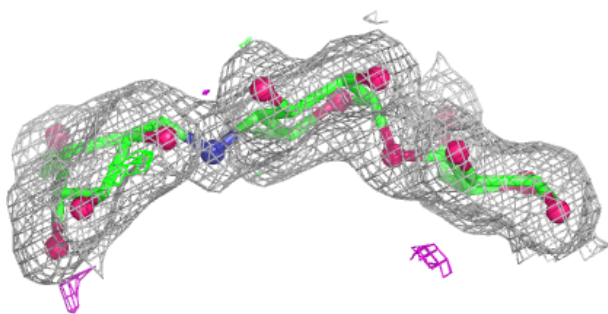
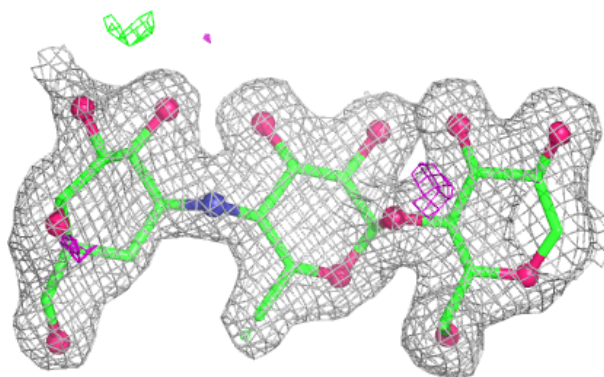


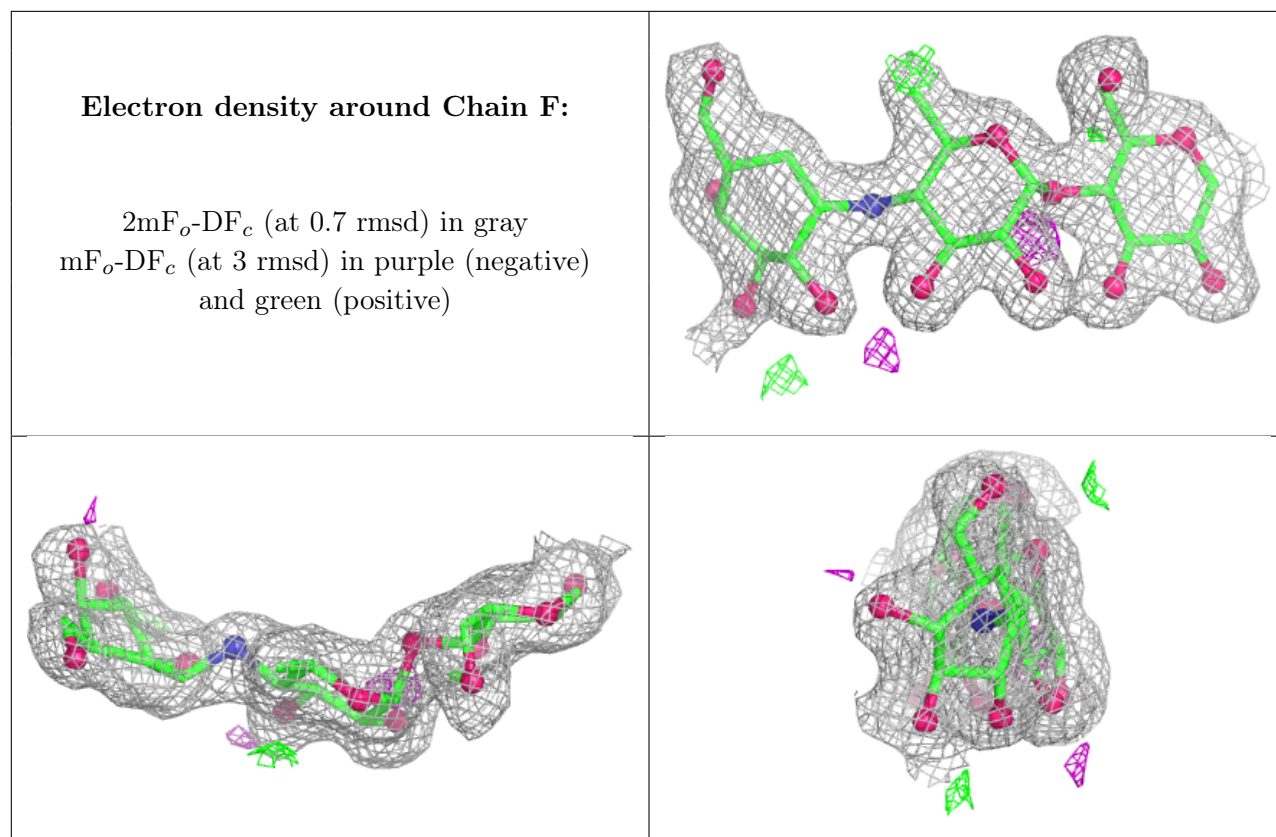
**Electron density around Chain E:**

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

**Electron density around Chain D:**

$2mF_o-DF_c$  (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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
5	EDO	B	506	4/4	0.67	0.17	22,34,35,39	0
5	EDO	B	514	4/4	0.73	0.18	22,27,27,31	4
5	EDO	B	519	4/4	0.74	0.19	26,28,37,40	0
5	EDO	A	506	4/4	0.75	0.15	24,26,26,30	4
5	EDO	B	516	4/4	0.76	0.16	28,31,37,45	0
5	EDO	B	502	4/4	0.77	0.17	21,22,22,23	4
5	EDO	B	510	4/4	0.78	0.17	26,28,29,44	0
5	EDO	B	518	4/4	0.78	0.17	23,30,31,42	0
5	EDO	B	503	4/4	0.78	0.19	26,26,30,42	0
5	EDO	A	502	4/4	0.79	0.13	29,35,38,42	0
5	EDO	B	520	4/4	0.79	0.16	21,27,31,43	0
5	EDO	B	505	4/4	0.80	0.18	32,33,34,40	0
5	EDO	A	508	4/4	0.80	0.24	24,26,35,43	0
5	EDO	A	515	4/4	0.81	0.15	17,18,26,28	4

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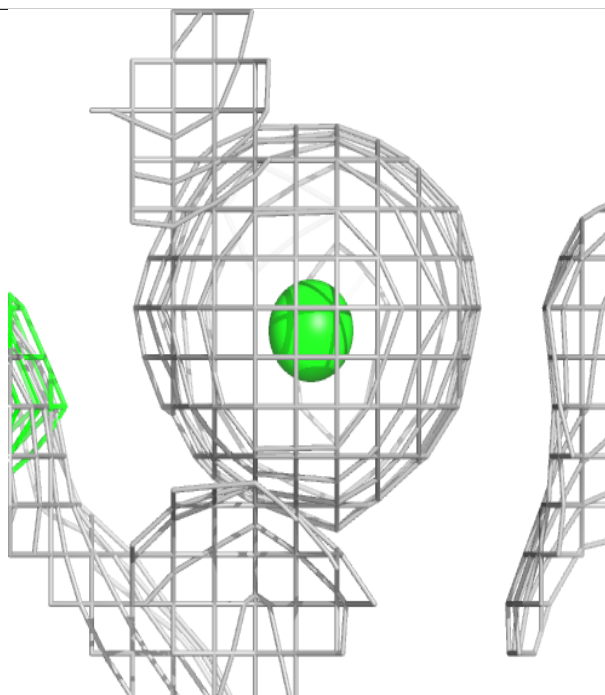
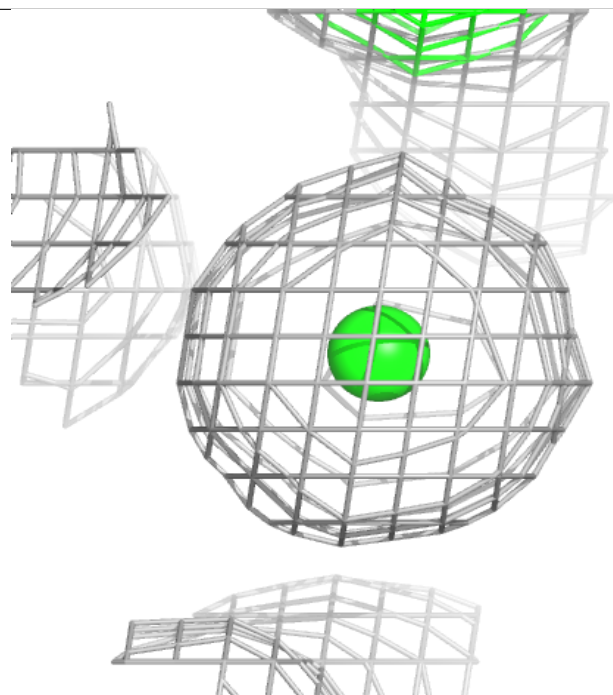
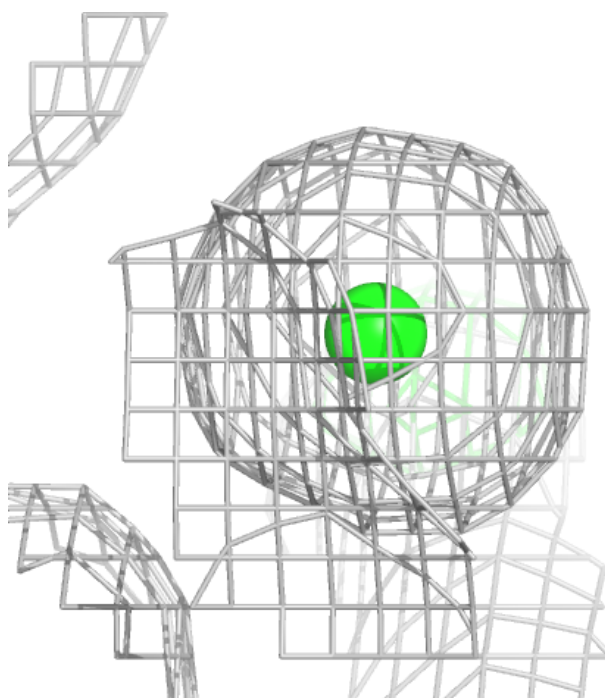
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	B	517	4/4	0.82	0.17	29,32,36,41	0
5	EDO	A	510	4/4	0.82	0.15	29,30,30,37	0
5	EDO	B	521	4/4	0.82	0.14	22,26,31,36	4
5	EDO	A	503	4/4	0.83	0.14	23,23,25,27	4
5	EDO	A	505	4/4	0.83	0.12	26,30,31,43	0
5	EDO	B	515	4/4	0.83	0.15	21,22,27,31	4
5	EDO	A	513	4/4	0.84	0.12	28,29,34,44	0
5	EDO	A	519	4/4	0.85	0.13	30,31,34,35	4
5	EDO	A	516	4/4	0.85	0.14	32,34,35,36	4
5	EDO	B	509	4/4	0.86	0.11	24,27,30,36	0
5	EDO	A	512	4/4	0.87	0.11	23,26,26,27	4
5	EDO	B	513	4/4	0.87	0.11	21,26,27,29	4
5	EDO	A	511	4/4	0.89	0.12	20,25,33,38	0
5	EDO	B	504	4/4	0.89	0.10	16,20,21,35	0
5	EDO	B	511	4/4	0.89	0.09	20,23,26,27	4
5	EDO	A	507	4/4	0.91	0.10	17,21,27,37	4
5	EDO	A	517	4/4	0.91	0.09	29,34,40,42	0
5	EDO	A	504	4/4	0.92	0.09	19,23,27,27	0
5	EDO	A	509	4/4	0.92	0.09	14,19,19,31	4
5	EDO	A	514	4/4	0.92	0.09	17,18,20,24	0
5	EDO	B	507	4/4	0.93	0.07	16,17,18,28	4
5	EDO	B	508	4/4	0.93	0.09	20,25,31,39	0
5	EDO	B	512	4/4	0.93	0.09	18,18,20,21	0
5	EDO	A	518	4/4	0.93	0.10	20,29,29,29	0
4	CL	B	501	1/1	0.99	0.02	13,13,13,13	0
6	SR	A	520	1/1	0.99	0.01	19,19,19,19	0
4	CL	A	501	1/1	1.00	0.02	14,14,14,14	0
6	SR	A	521	1/1	1.00	0.02	23,23,23,23	0
6	SR	B	522	1/1	1.00	0.01	18,18,18,18	0
6	SR	B	523	1/1	1.00	0.01	23,23,23,23	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.

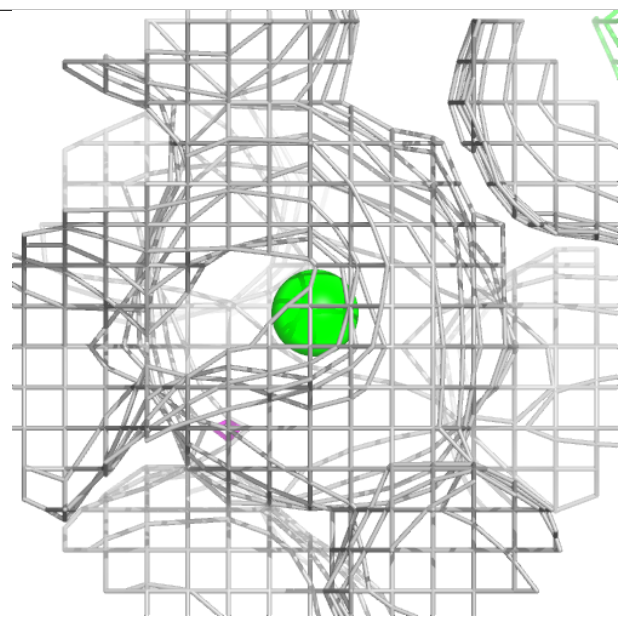
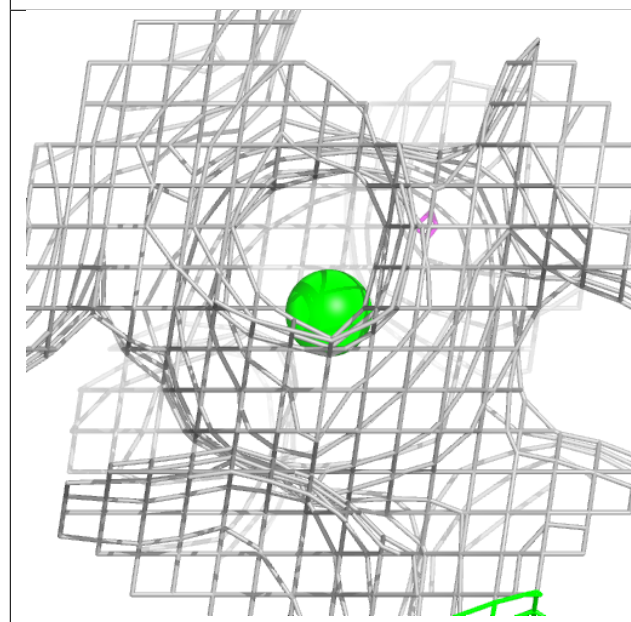
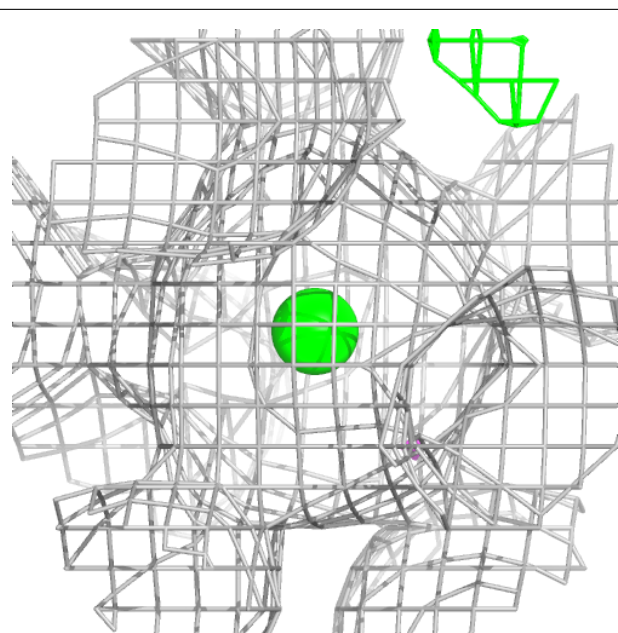
**Electron density around CL B 501:**

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



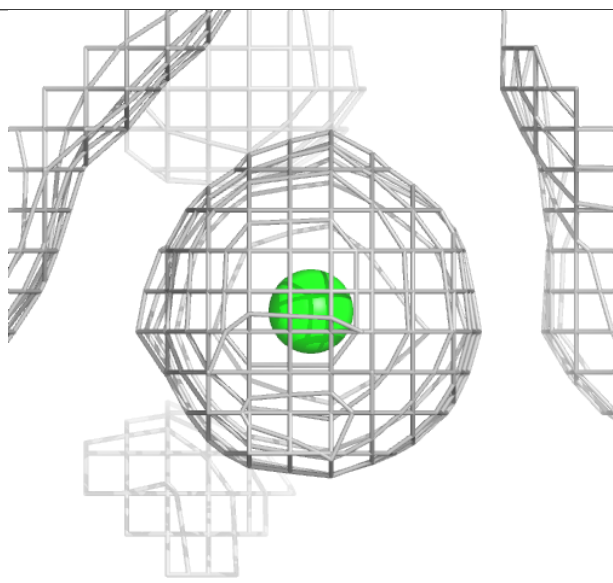
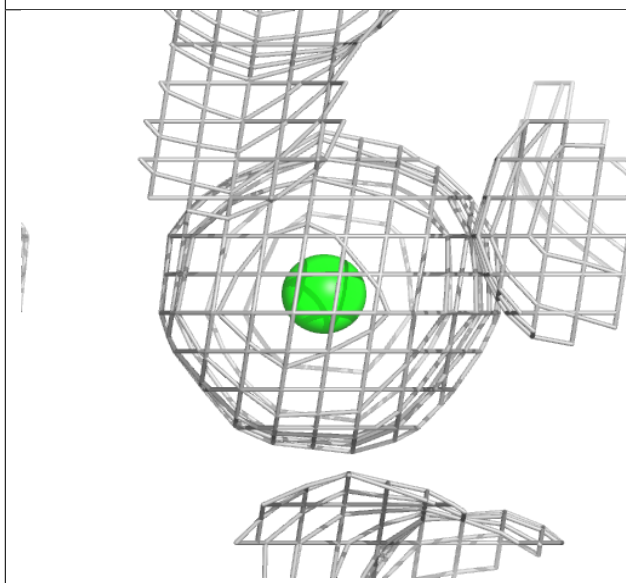
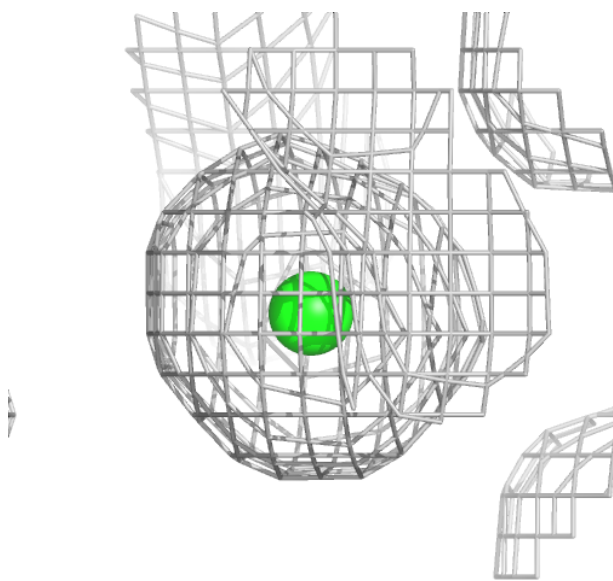
**Electron density around SR A 520:**

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



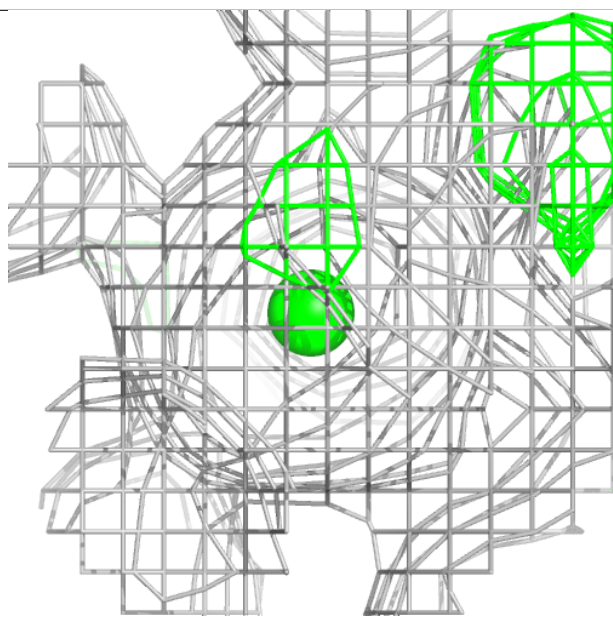
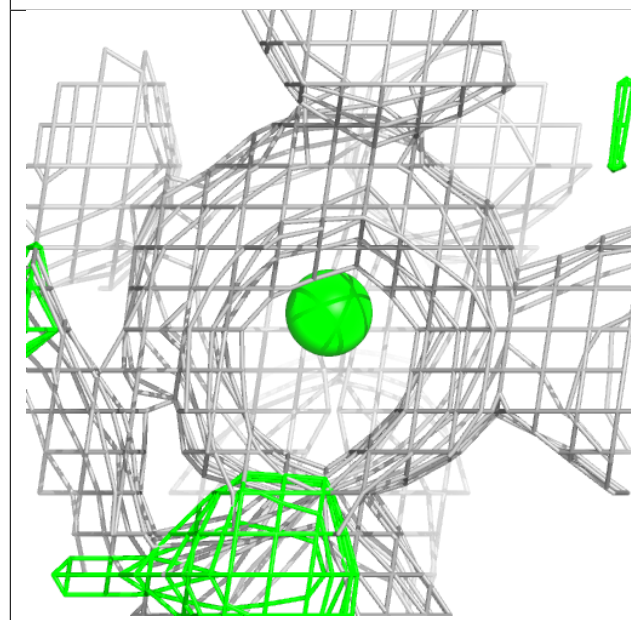
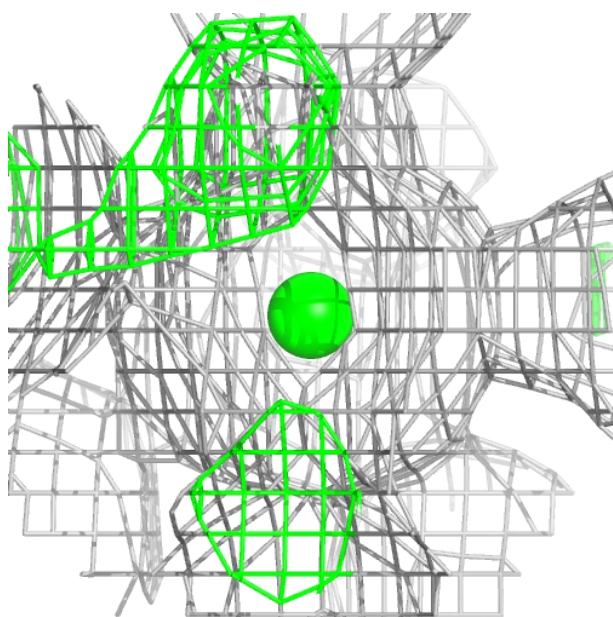
**Electron density around CL A 501:**

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



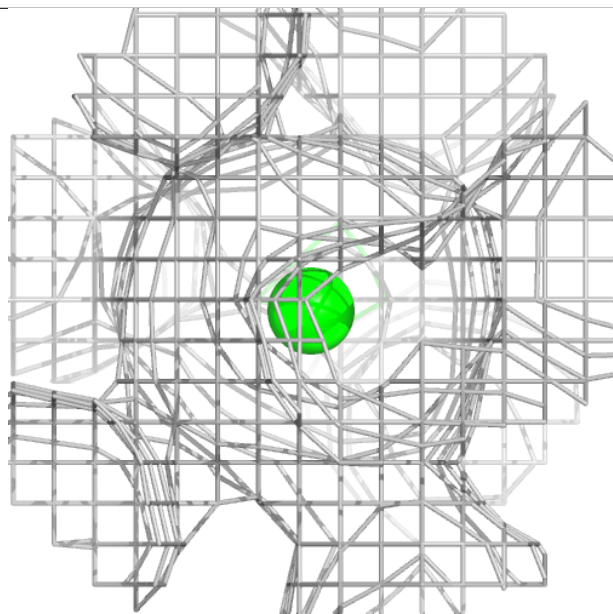
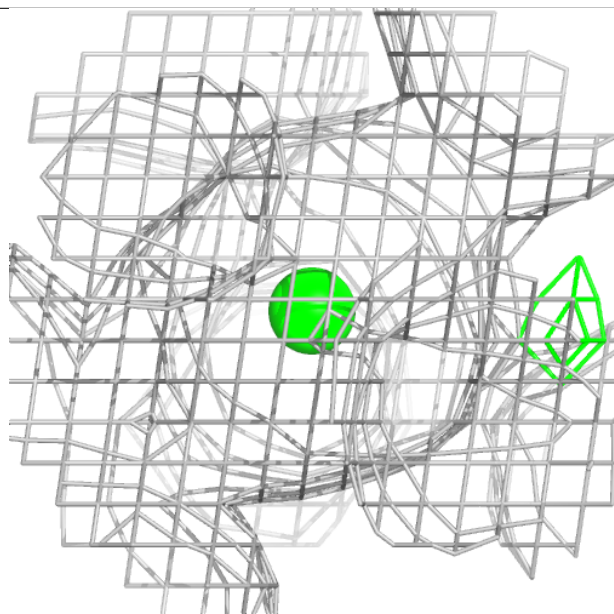
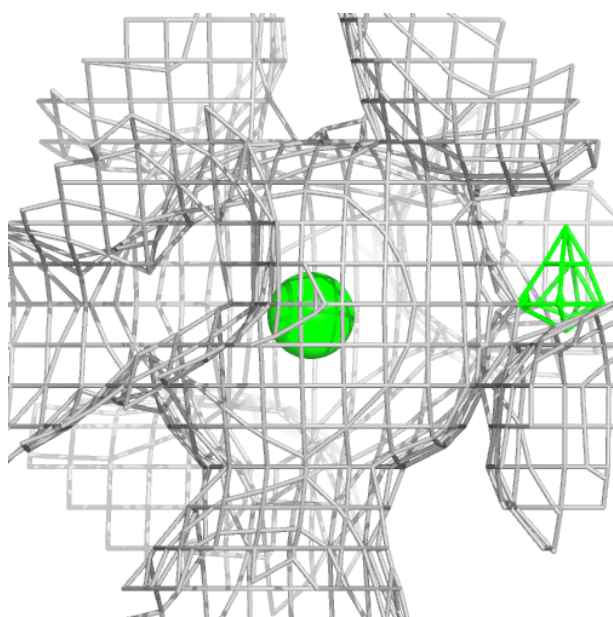
**Electron density around SR A 521:**

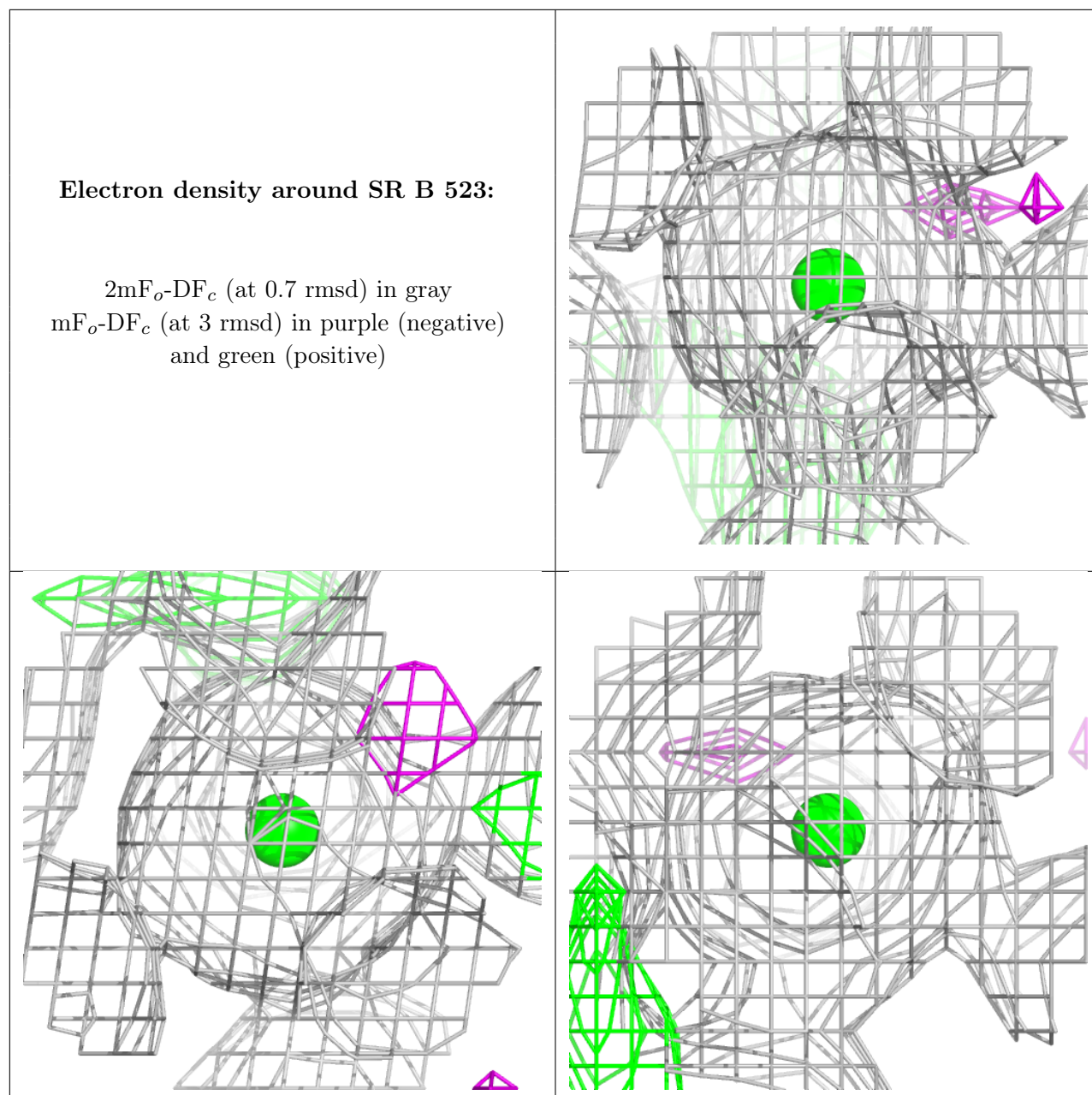
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around SR B 522:**

$2mF_o-DF_c$  (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.