wwPDB X-ray Structure Validation Summary Report (i)

Jun 13, 2024 – 11:00 AM EDT

PDB ID	:	1MXG
Title	:	Crystal Structure of a (Ca,Zn)-dependent alpha-amylase from the hyperthermophilic archaeon Pyrococcus woesei in complex with acarbose
Authors	:	Linden, A.; Mayans, O.; Meyer-Klaucke, W.; Antranikian, G.; Wilmanns, M.
Deposited on	:	2002-10-02
Resolution	:	1.60 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

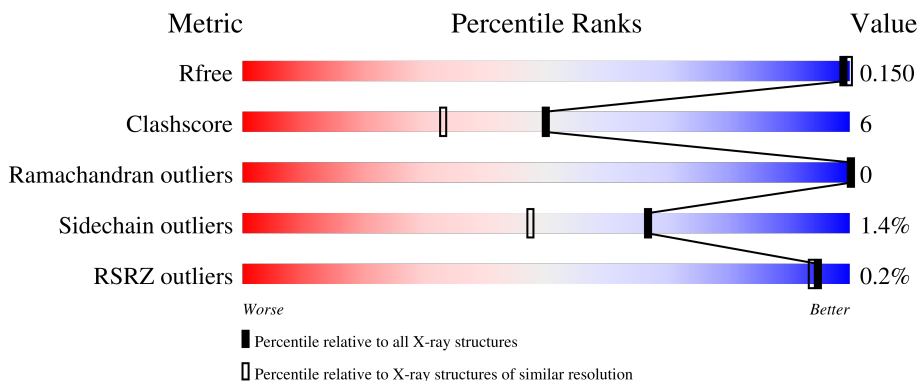|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.

Metric	Percentile Ranks	Value
Rfree		0.150
Clashscore		6
Ramachandran outliers		0
Sidechain outliers		1.4%
RSRZ outliers		0.2%

Worse　　　　　　　　　　　　　　　*Better*

▌ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 <=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	435	89%　　　　　　　　　　11%　　・
2	B	3	100%
2	C	3	67%　　　　　　33%
2	D	3	67%　　　　　　33%

2 Entry composition ⓘ

There are 9 unique types of molecules in this entry. The entry contains 4287 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total 3609	C 2359	N 580	O 659	S 11	0	5	0

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	3	Total 44	C 25	N 1	O 18	0	0	0
2	C	3	Total 44	C 25	N 1	O 18	0	0	0
2	D	3	Total 33	C 18	N 1	O 14	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Zn 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Ca 1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total 3	Mg 3	0	0

- Molecule 6 is 2-{2-[2-2-(METHOXY-ETHOXY)-ETHOXY]-ETHOXY}-ETHANOL (three-letter code: ETE) (formula: $C_9H_{20}O_5$).

ETE
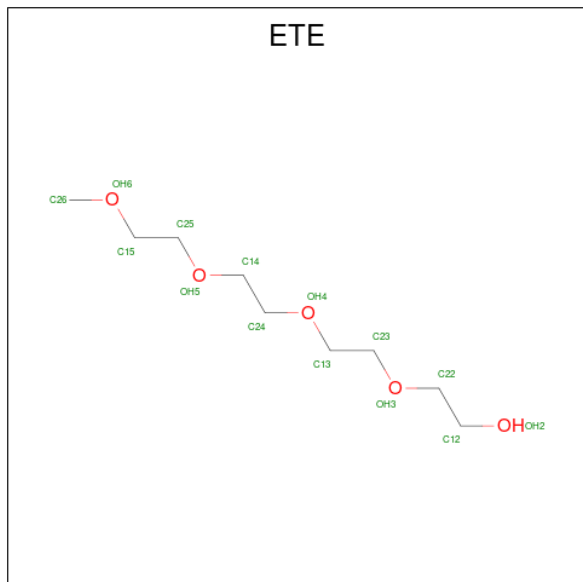
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total 14	C 9	O 5	0	0
6	A	1	Total 9	C 6	O 3	0	0

- Molecule 7 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).

TRS
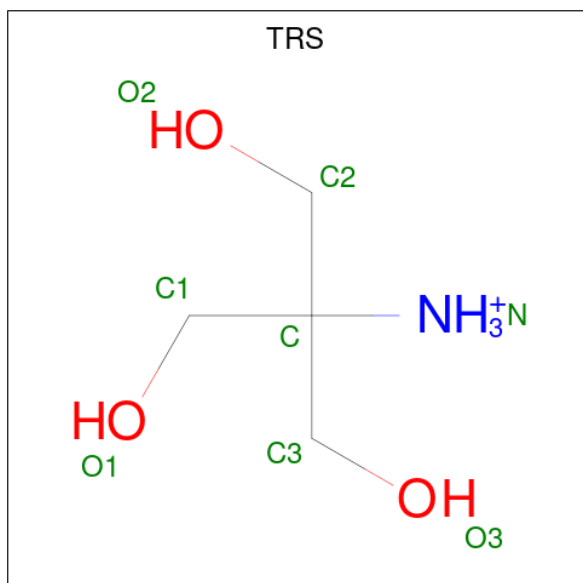
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total 8	C 4	N 1	O 3	0	0

- Molecule 8 is ETHANOL (three-letter code: EOH) (formula: C_2H_6O).
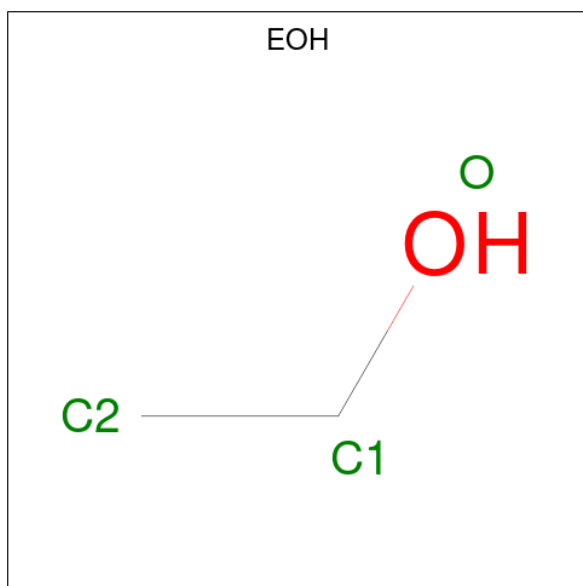
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total 3	C 2	O 1	0	0
8	A	1	Total 3	C 2	O 1	0	0
8	A	1	Total 3	C 2	O 1	0	0
8	A	1	Total 3	C 2	O 1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	509	Total 509	O 509	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: alpha amylase

Chain A:
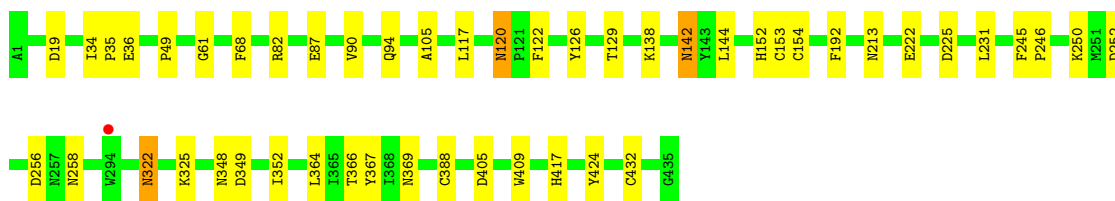
- Molecule 2: 4,6-dideoxy-4-{[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain B:

- Molecule 2: 4,6-dideoxy-4-{[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain C:

- Molecule 2: 4,6-dideoxy-4-{[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.48Å 76.51Å 136.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.60 68.23 – 1.52	Depositor EDS
% Data completeness (in resolution range)	93.3 (20.00-1.60) 89.8 (68.23-1.52)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.53 (at 1.52Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.159 , 0.172 0.156 , 0.150	Depositor DCC
R_{free} test set	1331 reflections (1.77%)	wwPDB-VP
Wilson B-factor (Å^2)	12.8	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 60.8	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	4287	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, AC1, ZN, GLC, EOH, TRS, ETE, CA

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	0.30	0/3749	0.58	0/5125

There are no bond length outliers.

There are no bond angle outliers.

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	3609	0	3309	37	1
2	B	44	0	30	4	1
2	C	44	0	30	0	0
2	D	33	0	21	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	3	0	0	0	0
6	A	23	0	31	1	0
7	A	8	0	12	2	0
8	A	12	0	24	1	0
9	A	509	0	0	3	0
All	All	4287	0	3457	42	1

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 6.

The worst 5 of 42 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:B:1:GLC:H62	2:B:2:GLC:C5	2.18	0.73
1:A:348[B]:ASN:HD22	1:A:348[B]:ASN:H	1.40	0.69
2:B:1:GLC:H62	2:B:2:GLC:O5	1.92	0.69
1:A:222[B]:GLU:OE1	7:A:447:TRS:H12	1.93	0.69
1:A:322:ASN:HD22	1:A:322:ASN:C	2.03	0.61

All (1) 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)
1:A:250:LYS:NZ	2:B:1:GLC:O2[4_455]	2.16	0.04

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	438/435 (101%)	422 (96%)	16 (4%)	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	366/361 (101%)	361 (99%)	5 (1%)	67 47

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	ASN
1	A	142	ASN
1	A	152	HIS
1	A	322	ASN
1	A	367	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	258	ASN
1	A	322	ASN
1	A	417	HIS
1	A	328	ASN
1	A	213	ASN

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 ⓘ

9 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	GLC	B	1	2	12,12,12	0.95	0	17,17,17	0.77	0
2	GLC	B	2	2	11,11,12	0.86	0	15,15,17	0.70	0
2	AC1	B	3	2	21,22,23	2.67	5 (23%)	22,32,34	1.99	5 (22%)
2	GLC	C	1	2	12,12,12	0.84	0	17,17,17	0.63	0
2	GLC	C	2	2	11,11,12	0.95	0	15,15,17	0.64	0
2	AC1	C	3	2	21,22,23	2.66	5 (23%)	22,32,34	1.50	3 (13%)
2	GLC	D	1	2	12,12,12	0.80	0	17,17,17	0.60	0
2	GLC	D	2	2	11,11,12	0.95	0	15,15,17	0.68	0
2	AC1	D	3	2	9,10,23	1.11	1 (11%)	14,14,34	0.77	1 (7%)

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	GLC	B	1	2	-	0/2/22/22	0/1/1/1
2	GLC	B	2	2	-	0/2/19/22	0/1/1/1
2	AC1	B	3	2	-	1/6/43/46	0/2/2/2
2	GLC	C	1	2	-	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	AC1	C	3	2	-	1/6/43/46	0/2/2/2
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1
2	AC1	D	3	2	-	-	0/1/1/2

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	AC1	C4A-C5B	6.93	1.57	1.51
2	B	3	AC1	C4A-C5B	6.91	1.57	1.51
2	B	3	AC1	C7B-C5B	6.47	1.41	1.32
2	C	3	AC1	C7B-C5B	6.27	1.41	1.32
2	B	3	AC1	C3B-C4A	4.24	1.59	1.53

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	AC1	O6B-C6B-C5B	6.47	127.98	112.50
2	C	3	AC1	C7B-C1B-N4A	-4.91	103.31	110.68
2	B	3	AC1	O4-C4A-C3B	-4.46	101.33	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	3	AC1	O3B-C3B-C2B	-2.54	104.48	110.35
2	B	3	AC1	O4-C4A-C5B	2.47	115.58	110.82

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	3	AC1	C2B-C1B-N4A-C4
2	C	3	AC1	C5-C4-N4A-C1B

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	GLC	4	0
2	B	1	GLC	4	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.

Oligosaccharide Chain B

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain D

Bond lengths

Bond angles

Torsions

Rings
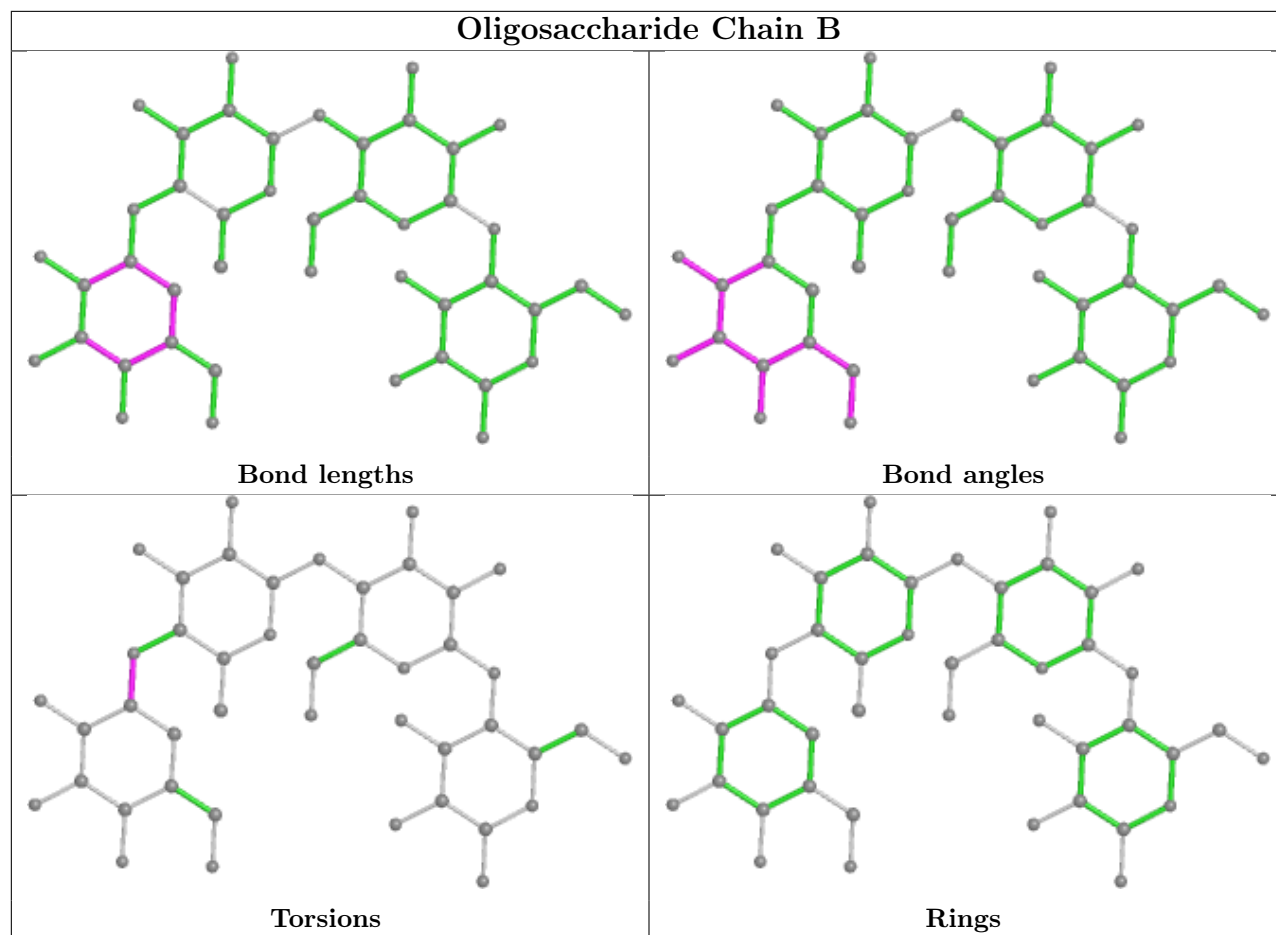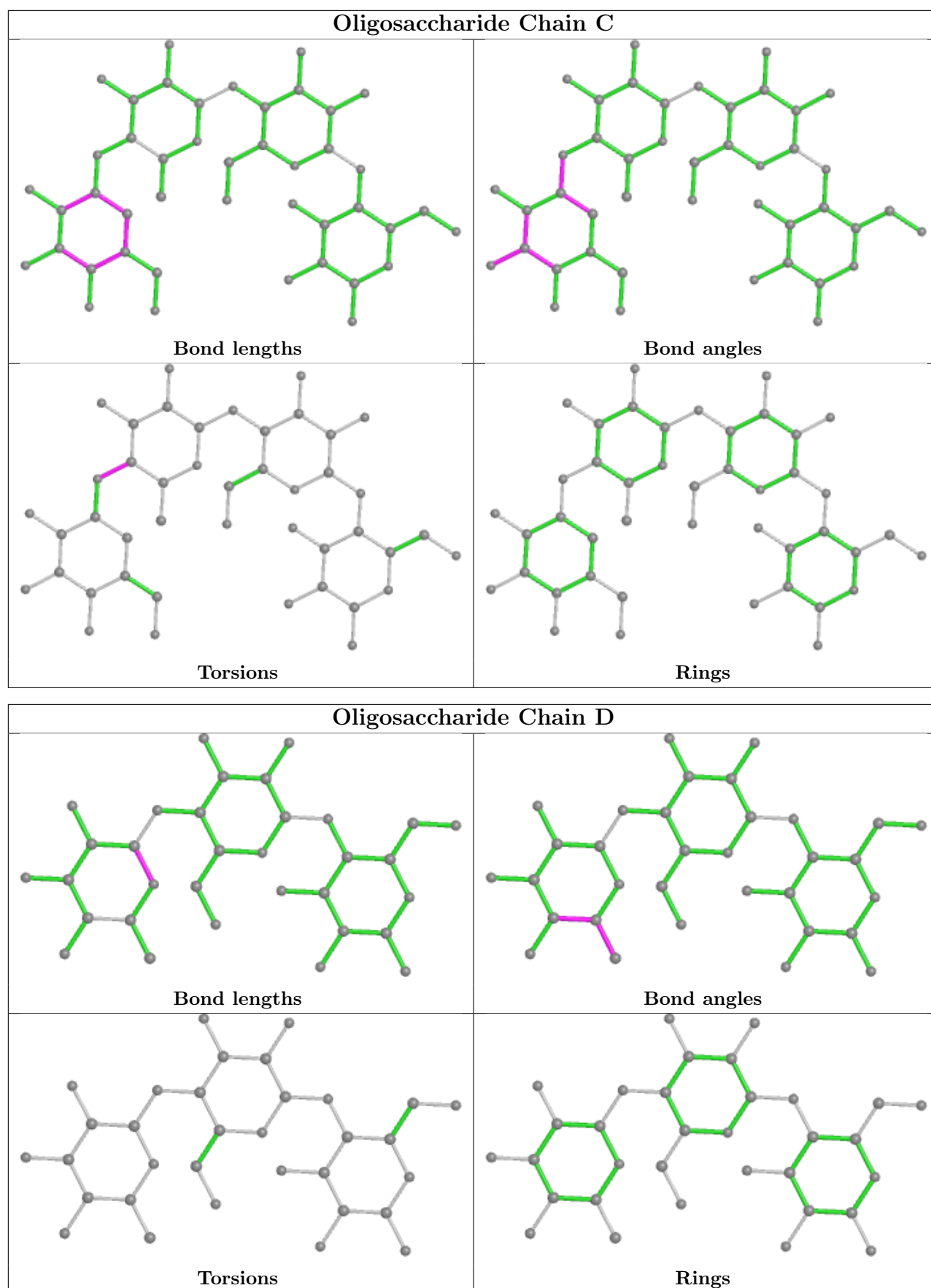
5.6   Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 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	#$|Z| > 2$
6	ETE	A	446	-	8,8,13	0.65	0	7,7,12	0.36	0
8	EOH	A	449	-	2,2,2	0.10	0	1,1,1	0.03	0
7	TRS	A	447	-	7,7,7	0.44	0	9,9,9	0.40	0
8	EOH	A	451	-	2,2,2	0.12	0	1,1,1	0.01	0
6	ETE	A	445	-	13,13,13	2.03	4 (30%)	12,12,12	0.88	0
8	EOH	A	448	-	2,2,2	0.12	0	1,1,1	0.01	0
8	EOH	A	450	-	2,2,2	0.13	0	1,1,1	0.01	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
6	ETE	A	446	-	-	4/6/6/11	-
7	TRS	A	447	-	-	0/9/9/9	-
6	ETE	A	445	-	-	2/11/11/11	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	445	ETE	OH4-C24	4.34	1.60	1.42
6	A	445	ETE	OH3-C23	3.42	1.56	1.42
6	A	445	ETE	OH6-C26	2.86	1.56	1.42
6	A	445	ETE	OH5-C25	2.17	1.51	1.42

There are no bond angle outliers.

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

WORLDWIDE
wwPDB
PROTEIN DATA BANK

Mol	Chain	Res	Type	Atoms
6	A	446	ETE	OH5-C14-C24-OH4
6	A	445	ETE	OH2-C12-C22-OH3
6	A	446	ETE	OH6-C15-C25-OH5
6	A	446	ETE	C15-C25-OH5-C14
6	A	445	ETE	OH6-C15-C25-OH5

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	446	ETE	1	0
7	A	447	TRS	2	0
8	A	450	EOH	1	0

5.7   Other polymers ⓘ

There are no such residues in this entry.

5.8   Polymer linkage issues ⓘ

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	435/435 (100%)	-0.50	1 (0%)	95	94	8, 13, 29, 38	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	294	TRP	3.8

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

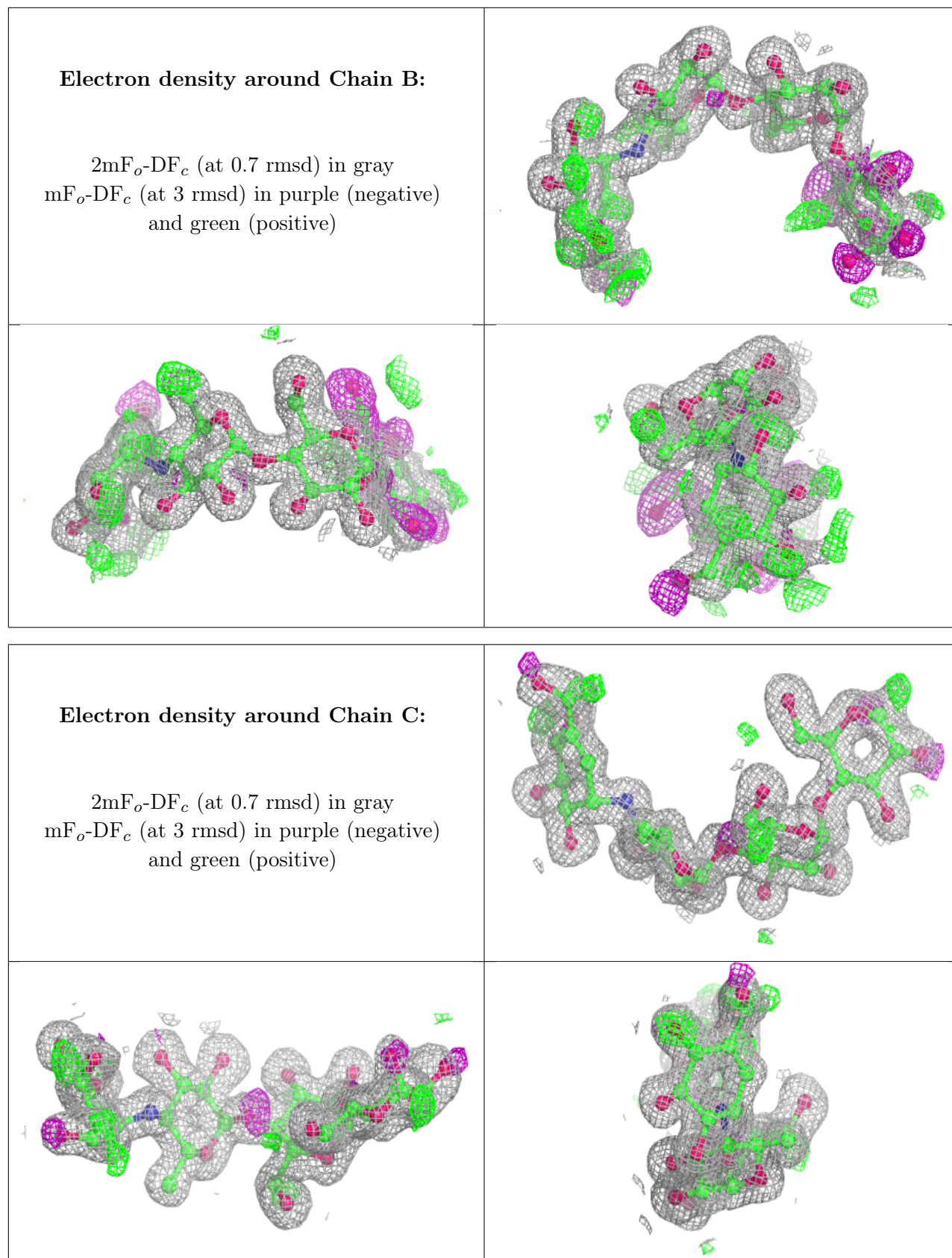There are no non-standard protein/DNA/RNA residues in this entry.
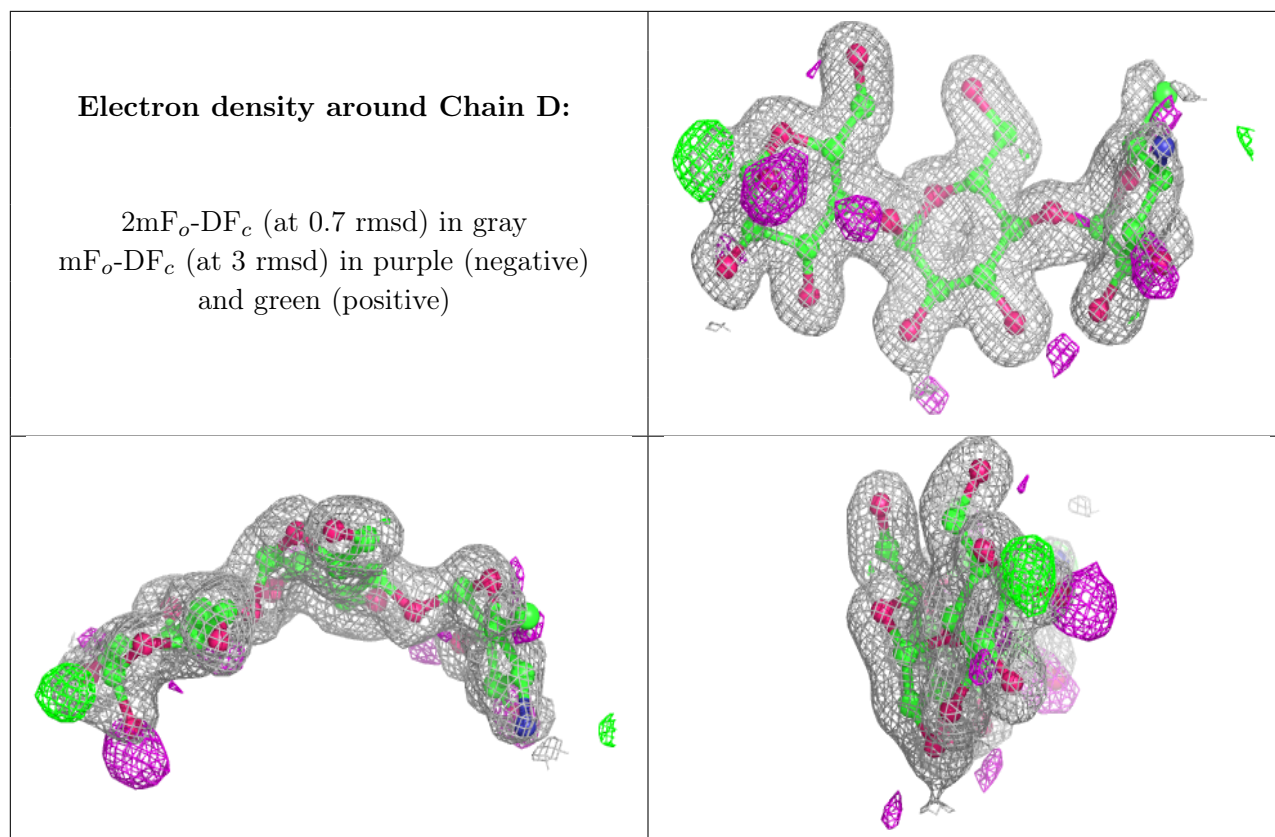
6.3 Carbohydrates (i)

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, 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	GLC	B	1	12/12	0.69	0.38	21,33,37,37	0
2	AC1	D	3	10/22	0.81	0.23	30,33,35,36	0
2	GLC	D	1	12/12	0.87	0.10	22,25,28,29	0
2	GLC	D	2	11/12	0.90	0.08	22,23,25,28	0
2	AC1	B	3	21/22	0.92	0.10	12,15,27,31	0
2	GLC	C	1	12/12	0.93	0.12	16,21,27,29	0
2	AC1	C	3	21/22	0.95	0.07	8,10,18,23	0
2	GLC	B	2	11/12	0.97	0.06	13,14,17,17	0
2	GLC	C	2	11/12	0.97	0.06	9,10,13,13	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain B:

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

Electron density around Chain C:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (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 ⓘ

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
6	ETE	A	445	14/14	0.72	0.20	27,28,31,35	0
6	ETE	A	446	9/14	0.73	0.26	33,34,35,35	0
8	EOH	A	450	3/3	0.81	0.11	28,28,29,30	0
8	EOH	A	451	3/3	0.86	0.19	31,31,32,33	0
7	TRS	A	447	8/8	0.87	0.14	19,25,27,28	0
8	EOH	A	449	3/3	0.97	0.08	16,16,19,20	0
5	MG	A	440	1/1	0.97	0.07	18,18,18,18	0
8	EOH	A	448	3/3	0.97	0.05	21,21,21,22	0
5	MG	A	439	1/1	0.99	0.04	11,11,11,11	0
4	CA	A	438	1/1	1.00	0.06	12,12,12,12	0
5	MG	A	441	1/1	1.00	0.07	12,12,12,12	0
3	ZN	A	437	1/1	1.00	0.03	12,12,12,12	0

6.5 Other polymers (i)

There are no such residues in this entry.