WORLDWIDE wwPDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 28, 2024 – 10:59 pm BST

PDB ID	:	6EUF
Title	:	The GH43, Beta 1,3 Galactosidase, BT0265
Authors	:	Cartmell, A.; Gilbert, H.J.
Deposited on	:	2017-10-30
Resolution	:	2.20 Å(reported)

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

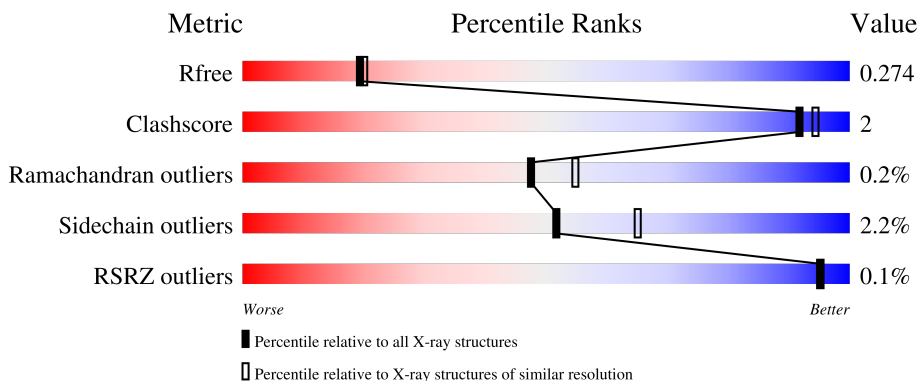|		
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.20 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.
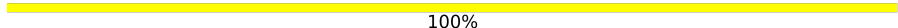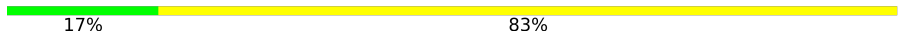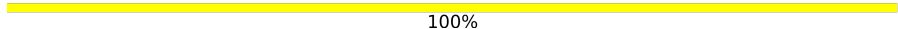
Metric	Percentile Ranks	Value
Rfree		0.274
Clashscore		2
Ramachandran outliers		0.2%
Sidechain outliers		2.2%
RSRZ outliers		0.1%

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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	475	93% ... 7% •
1	B	475	94% ... 5% •
1	C	475	91% ... 8% •
1	D	475	91% ... 8% •
2	E	5	100%

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Mol	Chain	Length	Quality of chain
2	H	5	100%
3	F	6	17% 83%
3	G	6	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
3	RAM	F	4	-	-	-	X

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 15689 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-glucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	0	0
			3726	2382	608	717	19			
1	B	472	Total	C	N	O	S	0	1	0
			3744	2390	613	722	19			
1	C	472	Total	C	N	O	S	0	0	0
			3726	2382	608	717	19			
1	D	472	Total	C	N	O	S	0	2	0
			3723	2379	608	717	19			

- Molecule 2 is an oligosaccharide called alpha-L-arabinofuranose-(1-3)-[alpha-L-arabinofuranose-(1-4)][beta-D-glucopyranuronic acid-(1-6)]beta-D-galactopyranose-(1-6)-beta-D-galactopyranose.
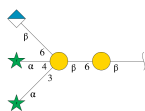
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	5	Total	C	O	0	0	0
			53	28	25			
2	H	5	Total	C	O	0	0	0
			53	28	25			

- Molecule 3 is an oligosaccharide called alpha-L-rhamnopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-6)-[alpha-L-arabinofuranose-(1-3)][alpha-L-arabinofuranose-(1-4)]beta-D-galactopyranose-(1-6)-beta-D-galactopyranose.
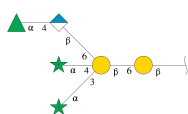
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	F	6	Total 63	C 34	O 29	0	0	0
3	G	6	Total 63	C 34	O 29	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	146	Total 146	O 146	0	0
4	B	166	Total 166	O 166	0	0
4	C	124	Total 124	O 124	0	0
4	D	102	Total 102	O 102	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.
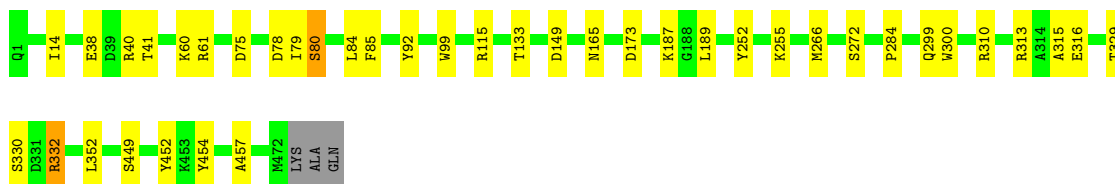
- Molecule 1: Beta-glucanase

Chain A:
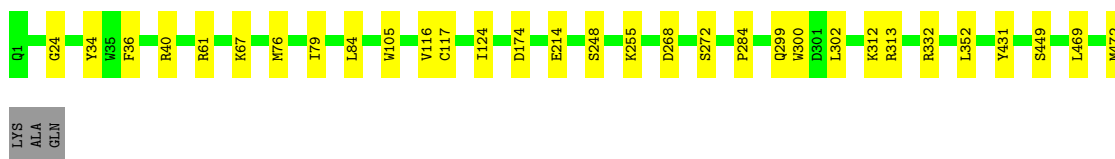
- Molecule 1: Beta-glucanase

Chain B:

- Molecule 1: Beta-glucanase

Chain C:

- Molecule 1: Beta-glucanase

Chain D:
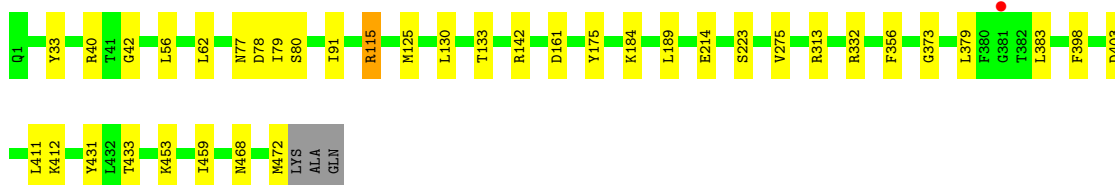
• Molecule 2: alpha-L-arabinofuranose-(1-3)-[alpha-L-arabinofuranose-(1-4)][beta-D-glucopyranur onic acid-(1-6)]beta-D-galactopyranose-(1-6)-beta-D-galactopyranose

Chain E: 100%

GAL1 GAL2 AHR3 AHR4 BDP5

• Molecule 2: alpha-L-arabinofuranose-(1-3)-[alpha-L-arabinofuranose-(1-4)][beta-D-glucopyranur onic acid-(1-6)]beta-D-galactopyranose-(1-6)-beta-D-galactopyranose

Chain H: 100%

GAL1 GAL2 AHR3 AHR4 BDP5

• Molecule 3: alpha-L-rhamnopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-6)-[alpha-L-arabin ofuranose-(1-3)][alpha-L-arabinofuranose-(1-4)]beta-D-galactopyranose-(1-6)-beta-D-galactopyran ose

Chain F: 17%   83%

GAL1 GAL2 BDP3 RAM4 AHR5 AHR6

• Molecule 3: alpha-L-rhamnopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-6)-[alpha-L-arabin ofuranose-(1-3)][alpha-L-arabinofuranose-(1-4)]beta-D-galactopyranose-(1-6)-beta-D-galactopyran ose

Chain G: 100%

GAL1 GAL2 BDP3 RAM4 AHR5 AHR6

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	106.00Å 119.70Å 182.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.04 – 2.20 100.04 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (100.04-2.20) 99.9 (100.04-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	2.86 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.208 , 0.271 0.215 , 0.274	Depositor DCC
R_{free} test set	5837 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å^2)	34.2	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 36.1	EDS
L-test for twinning [2]	$<|L|> = 0.50$, $<L^2> = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	15689	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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: BDP, RAM, GAL, AHR

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	0.86	0/3825	0.89	6/5201 (0.1%)
1	B	0.86	0/3843	0.91	10/5223 (0.2%)
1	C	0.83	0/3825	0.87	5/5201 (0.1%)
1	D	0.80	0/3828	0.88	8/5206 (0.2%)
All	All	0.84	0/15321	0.89	29/20831 (0.1%)

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	40	ARG	NE-CZ-NH2	-8.83	115.89	120.30
1	D	40	ARG	NE-CZ-NH1	8.72	124.66	120.30
1	B	40	ARG	NE-CZ-NH1	8.36	124.48	120.30
1	B	40	ARG	NE-CZ-NH2	-7.34	116.63	120.30
1	A	40	ARG	NE-CZ-NH1	7.18	123.89	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3726	0	3441	8	0
1	B	3744	0	3478	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3726	0	3441	15	0
1	D	3723	0	3434	12	0
2	E	53	0	25	0	0
2	H	53	0	24	0	0
3	F	63	0	34	0	0
3	G	63	0	34	0	0
4	A	146	0	0	0	0
4	B	166	0	0	1	0
4	C	124	0	0	0	0
4	D	102	0	0	0	0
All	All	15689	0	13911	45	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 45 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:C:38:GLU:OE2	1:C:40:ARG:NH1	2.25	0.69
1:C:79:ILE:HG22	1:C:79:ILE:O	2.00	0.61
1:B:322:ILE:HD12	1:B:370:THR:HG21	1.84	0.59
1:A:79:ILE:HD11	1:A:117:CYS:SG	2.43	0.59
1:A:312:LYS:HE2	1:A:449:SER:OG	2.04	0.58

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	470/475 (99%)	448 (95%)	22 (5%)	0	100	100
1	B	471/475 (99%)	449 (95%)	21 (4%)	1 (0%)	47	55

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	470/475 (99%)	451 (96%)	18 (4%)	1 (0%)	47	55
1	D	472/475 (99%)	450 (95%)	20 (4%)	2 (0%)	34	37
All	All	1883/1900 (99%)	1798 (96%)	81 (4%)	4 (0%)	47	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	173	ASP
1	B	437	LYS
1	D	373	GLY
1	D	412	LYS

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	382/417 (92%)	372 (97%)	10 (3%)	46	58
1	B	389/417 (93%)	385 (99%)	4 (1%)	76	86
1	C	382/417 (92%)	371 (97%)	11 (3%)	42	54
1	D	382/417 (92%)	373 (98%)	9 (2%)	49	62
All	All	1535/1668 (92%)	1501 (98%)	34 (2%)	52	65

5 of 34 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	214	GLU
1	D	223	SER
1	D	431	TYR
1	B	420	SER
1	B	299	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	339	ASN
1	B	467	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 ⓘ

22 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	GAL	E	1	2	12,12,12	0.63	0	17,17,17	1.12	2 (11%)
2	GAL	E	2	2	11,11,12	0.64	0	15,15,17	1.55	1 (6%)
2	AHR	E	3	2	9,9,10	0.66	0	10,12,14	1.98	2 (20%)
2	AHR	E	4	2	9,9,10	0.50	0	10,12,14	1.44	1 (10%)
2	BDP	E	5	2	12,12,13	1.05	0	14,17,19	1.27	1 (7%)
3	GAL	F	1	3	12,12,12	0.47	0	17,17,17	1.32	2 (11%)
3	GAL	F	2	3	11,11,12	0.34	0	15,15,17	1.68	6 (40%)
3	BDP	F	3	3	12,12,13	0.77	0	14,17,19	1.81	4 (28%)
3	RAM	F	4	3	10,10,11	0.65	0	14,14,16	1.04	1 (7%)
3	AHR	F	5	3	9,9,10	0.46	0	10,12,14	1.75	2 (20%)
3	AHR	F	6	3	9,9,10	0.63	0	10,12,14	0.98	0
3	GAL	G	1	3	12,12,12	0.69	0	17,17,17	1.74	5 (29%)
3	GAL	G	2	3	11,11,12	0.81	0	15,15,17	1.29	2 (13%)
3	BDP	G	3	3	12,12,13	0.95	0	14,17,19	2.12	3 (21%)

Mol	Type	Chain	Res	Link	Bond lengths				Bond angles		
					Counts	RMSZ	#|Z| > 2		Counts	RMSZ	#|Z| > 2
3	RAM	G	4	3	10,10,11	0.56	0		14,14,16	1.30	1 (7%)
3	AHR	G	5	3	9,9,10	0.41	0		10,12,14	1.33	2 (20%)
3	AHR	G	6	3	9,9,10	0.77	0		10,12,14	1.19	1 (10%)
2	GAL	H	1	2	12,12,12	0.90	0		17,17,17	1.94	5 (29%)
2	GAL	H	2	2	11,11,12	0.36	0		15,15,17	1.30	1 (6%)
2	AHR	H	3	2	9,9,10	0.60	0		10,12,14	1.19	1 (10%)
2	AHR	H	4	2	9,9,10	0.61	0		10,12,14	1.73	2 (20%)
2	BDP	H	5	2	12,12,13	0.99	1 (8%)		14,17,19	1.00	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GAL	E	1	2	-	0/2/22/22	0/1/1/1
2	GAL	E	2	2	-	2/2/19/22	0/1/1/1
2	AHR	E	3	2	-	0/2/15/18	0/1/1/1
2	AHR	E	4	2	-	0/2/15/18	0/1/1/1
2	BDP	E	5	2	-	0/4/21/24	0/1/1/1
3	GAL	F	1	3	-	0/2/22/22	0/1/1/1
3	GAL	F	2	3	-	2/2/19/22	0/1/1/1
3	BDP	F	3	3	-	0/4/21/24	0/1/1/1
3	RAM	F	4	3	-	-	0/1/1/1
3	AHR	F	5	3	-	2/2/15/18	0/1/1/1
3	AHR	F	6	3	-	2/2/15/18	0/1/1/1
3	GAL	G	1	3	-	0/2/22/22	0/1/1/1
3	GAL	G	2	3	-	1/2/19/22	0/1/1/1
3	BDP	G	3	3	-	2/4/21/24	0/1/1/1
3	RAM	G	4	3	-	-	0/1/1/1
3	AHR	G	5	3	-	2/2/15/18	0/1/1/1
3	AHR	G	6	3	-	2/2/15/18	0/1/1/1
2	GAL	H	1	2	-	0/2/22/22	0/1/1/1
2	GAL	H	2	2	-	2/2/19/22	0/1/1/1
2	AHR	H	3	2	-	2/2/15/18	0/1/1/1
2	AHR	H	4	2	-	0/2/15/18	0/1/1/1
2	BDP	H	5	2	-	0/4/21/24	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	5	BDP	O6B-C6	-2.04	1.23	1.30

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	G	3	BDP	O4-C4-C5	5.15	121.29	109.74
2	E	3	AHR	O4-C4-C3	4.60	108.78	104.70
3	F	5	AHR	O4-C4-C3	-4.59	100.64	104.70
2	H	1	GAL	C1-O5-C5	-4.36	105.43	113.66
2	E	2	GAL	O5-C5-C6	4.18	113.75	107.20

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

| Mol	Chain	Res	Type	Atoms
3	G	5	AHR	O4-C4-C5-O5
3	G	6	AHR	O4-C4-C5-O5
3	F	6	AHR	O4-C4-C5-O5
3	G	6	AHR	C3-C4-C5-O5
3	F	5	AHR	C3-C4-C5-O5

There are no ring outliers.

No monomer is involved in short contacts.

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 E

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain H

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain F

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain G

Bond lengths

Bond angles

Torsions

Rings
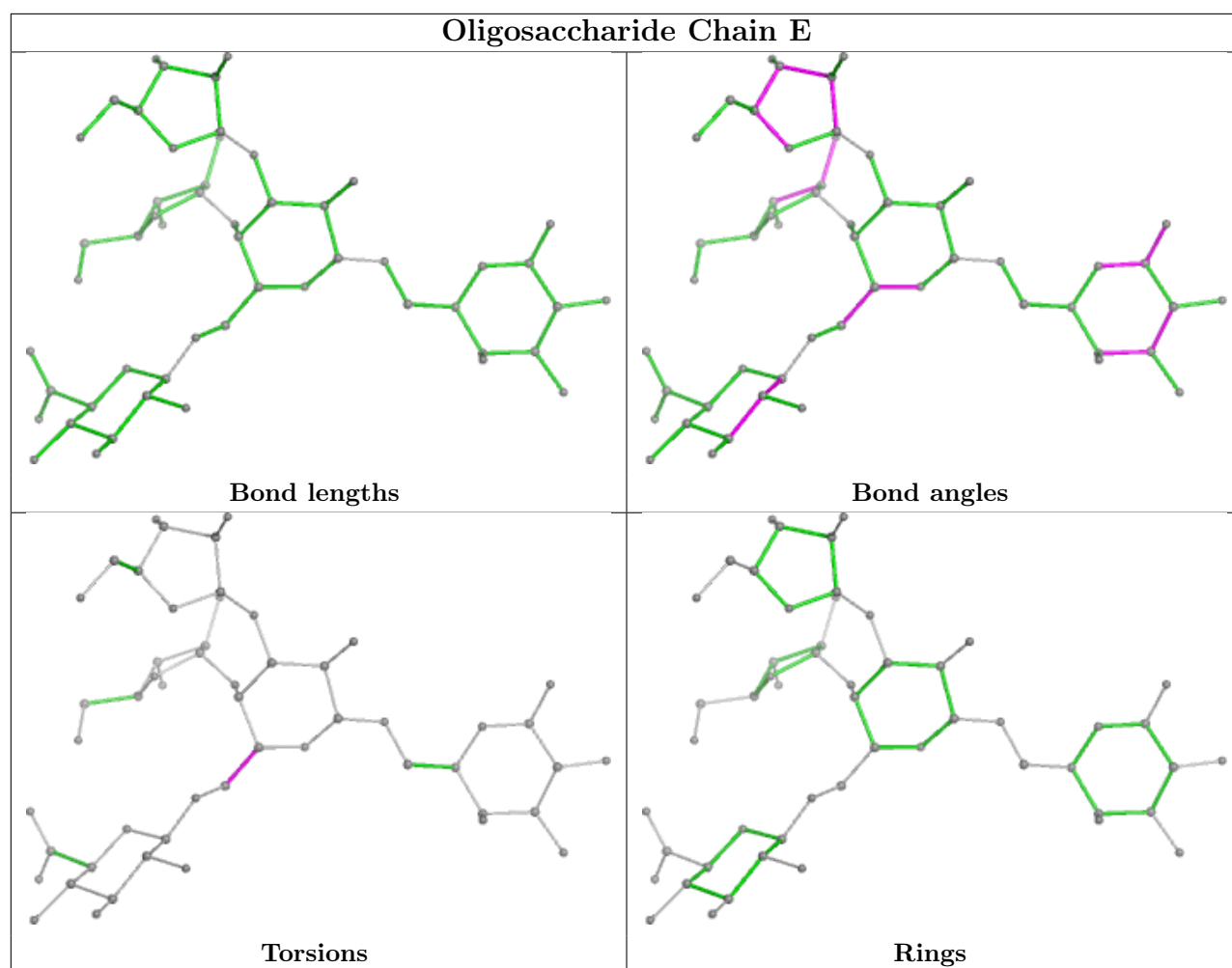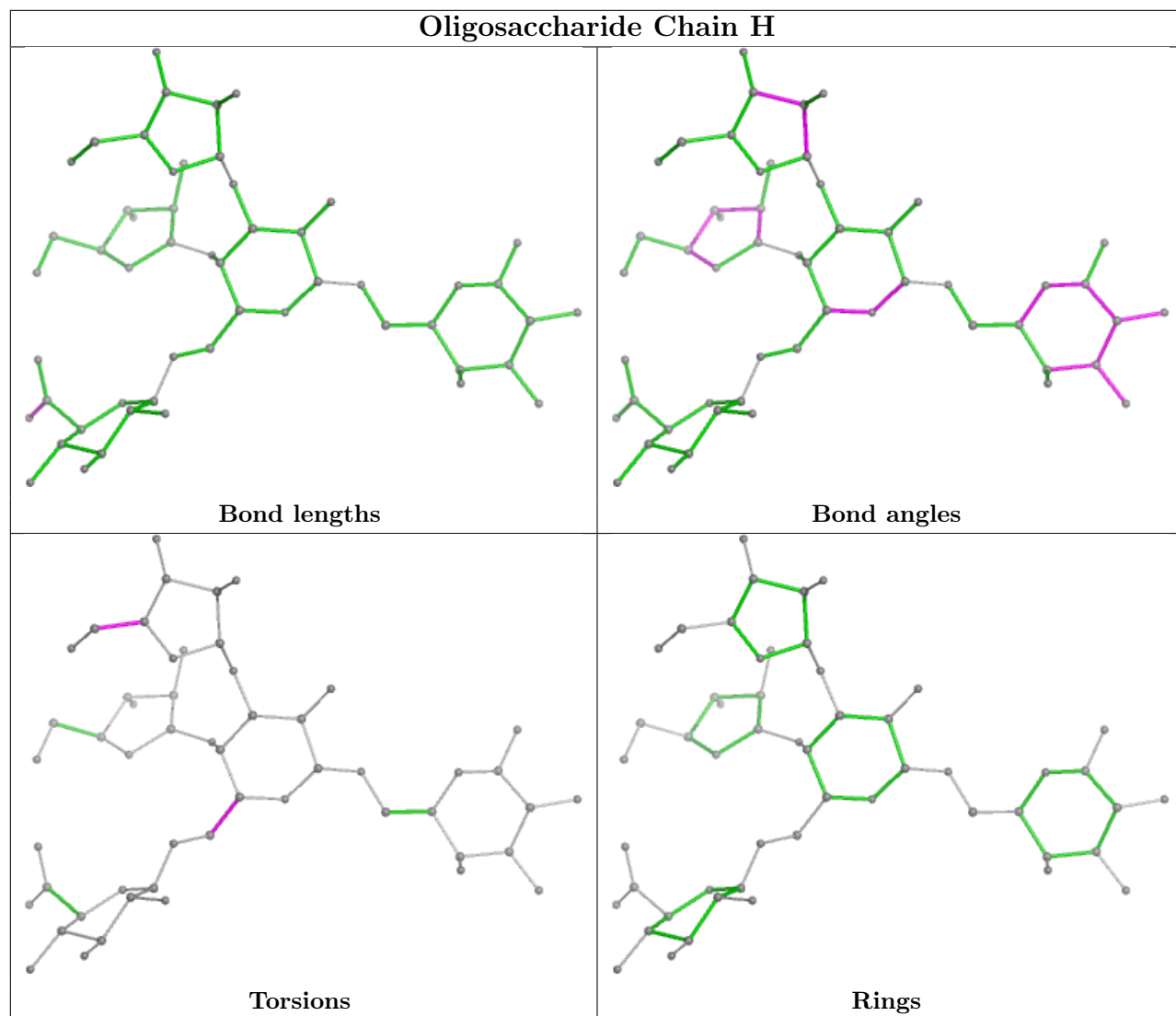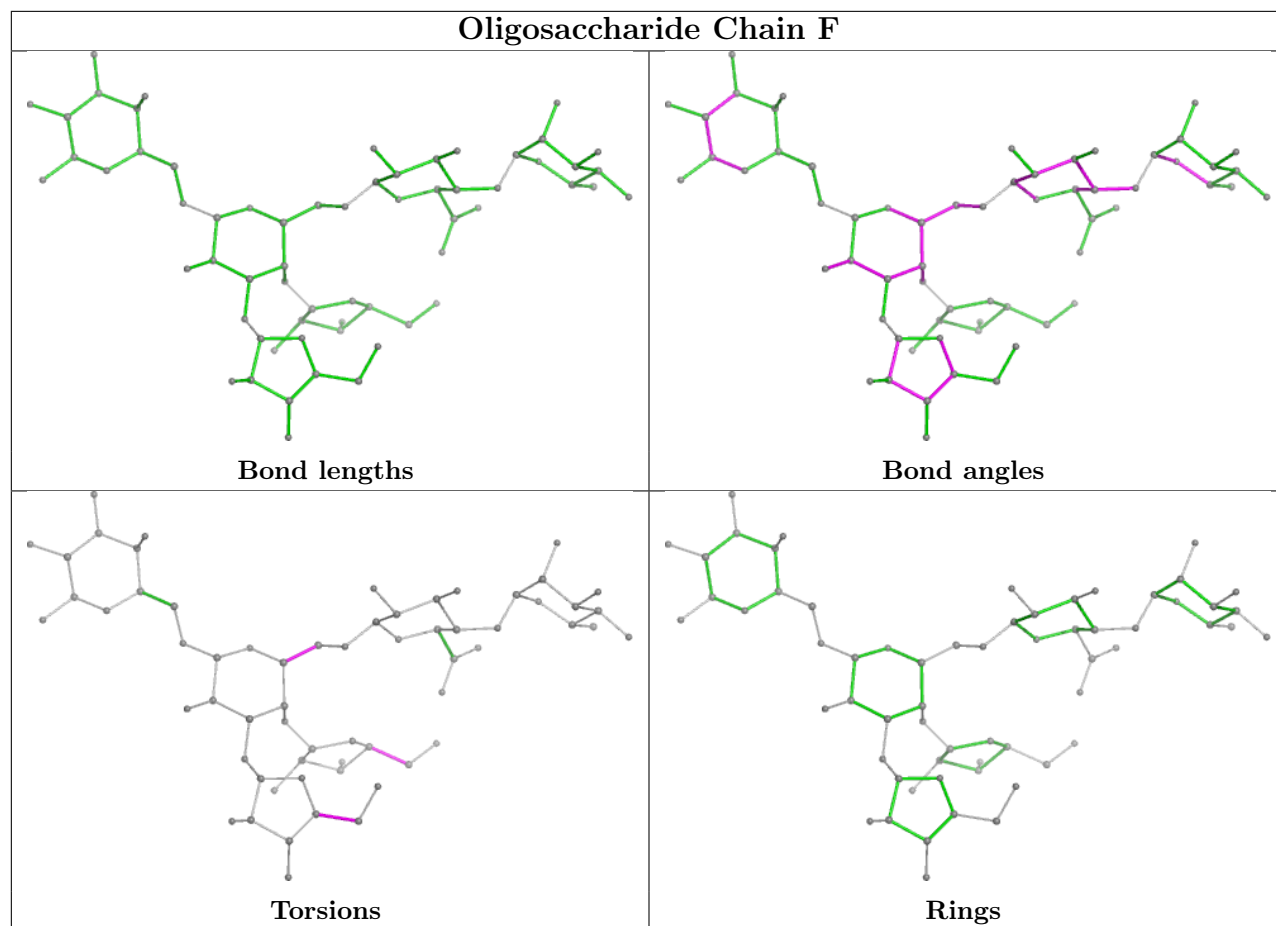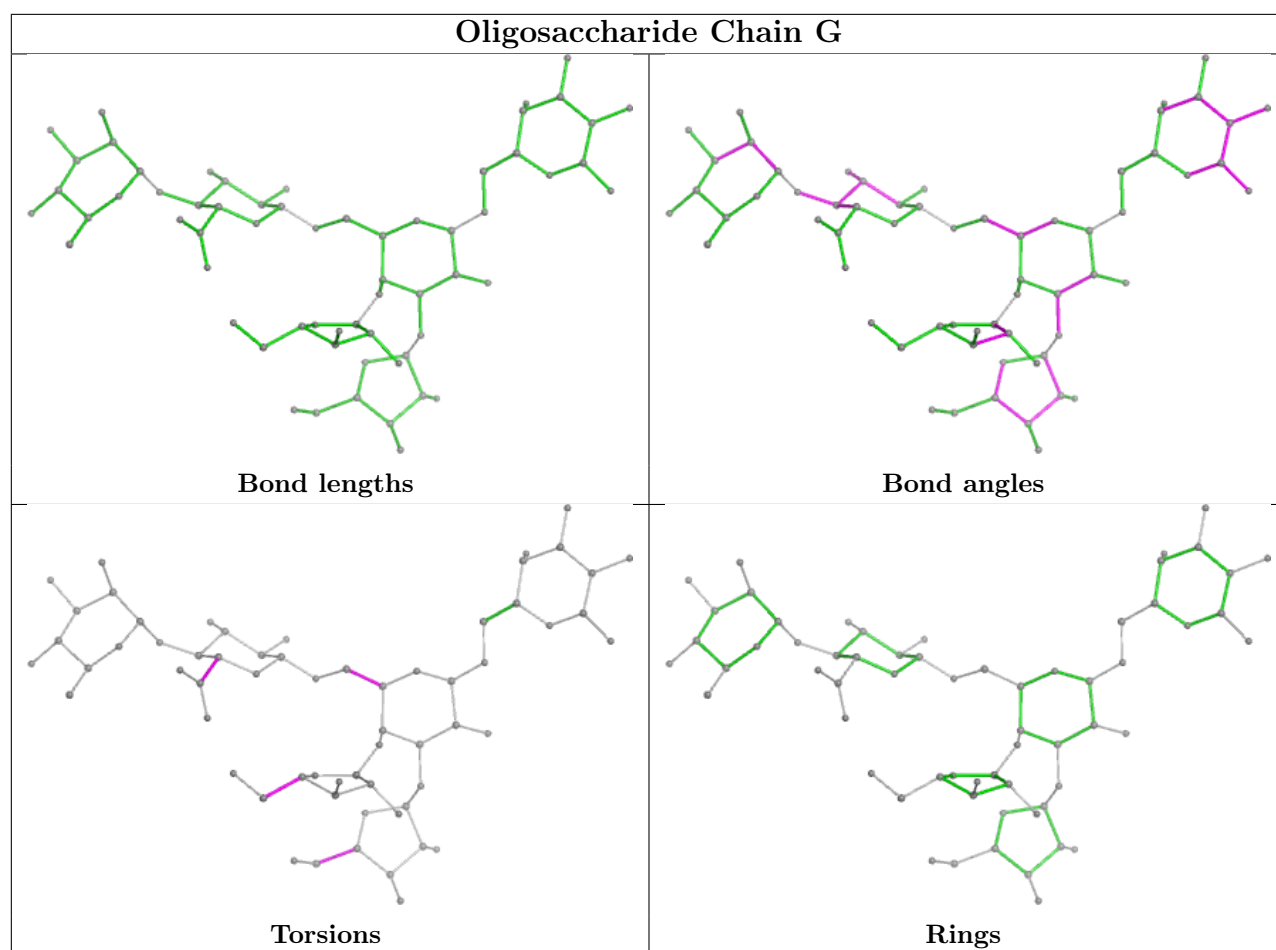
5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	A	472/475 (99%)	-0.75	0	100	100	22, 33, 50, 76	0
1	B	472/475 (99%)	-0.75	0	100	100	21, 34, 49, 70	0
1	C	472/475 (99%)	-0.63	0	100	100	26, 39, 59, 77	0
1	D	472/475 (99%)	-0.48	1 (0%)	95	94	22, 41, 68, 96	0
All	All	1888/1900 (99%)	-0.65	1 (0%)	95	95	21, 36, 60, 96	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	381	GLY	2.1

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

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

6.3 Carbohydrates ⓘ

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.
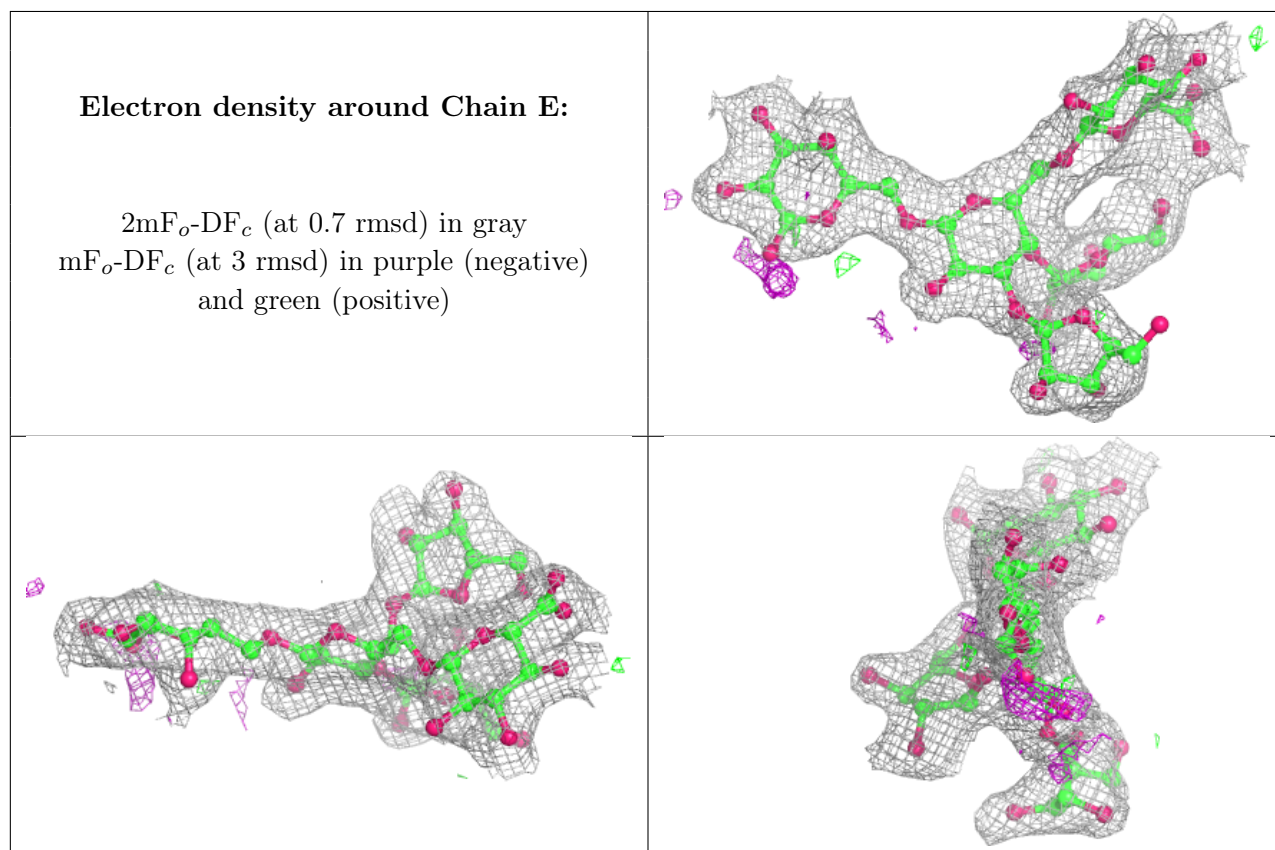
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å²)	Q<0.9
3	RAM	G	4	10/11	0.67	0.38	81,106,116,121	0
3	AHR	F	6	9/10	0.73	0.30	61,74,88,91	0
2	AHR	H	4	9/10	0.75	0.16	72,78,85,88	0
3	RAM	F	4	10/11	0.76	0.47	95,103,115,116	0
2	AHR	E	3	9/10	0.78	0.16	71,75,81,81	0
3	AHR	G	6	9/10	0.80	0.14	75,82,93,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	AHR	E	4	9/10	0.81	0.15	59,65,70,70	0
3	AHR	G	5	9/10	0.83	0.12	74,77,83,84	0
3	AHR	F	5	9/10	0.85	0.14	69,71,77,84	0
3	BDP	G	3	12/13	0.88	0.11	58,69,73,83	0
2	BDP	E	5	12/13	0.88	0.12	48,61,63,64	0
2	AHR	H	3	9/10	0.89	0.15	81,85,90,91	0
2	BDP	H	5	12/13	0.91	0.11	49,54,56,59	0
2	GAL	H	2	11/12	0.91	0.10	34,40,47,51	0
3	BDP	F	3	12/13	0.92	0.12	61,65,73,88	0
3	GAL	G	1	12/12	0.92	0.12	31,38,43,51	0
2	GAL	E	2	11/12	0.93	0.07	35,42,48,48	0
3	GAL	F	2	11/12	0.94	0.08	35,39,46,52	0
3	GAL	G	2	11/12	0.94	0.07	36,47,56,57	0
2	GAL	E	1	12/12	0.95	0.10	22,28,31,41	0
3	GAL	F	1	12/12	0.96	0.10	29,32,37,43	0
2	GAL	H	1	12/12	0.96	0.09	24,27,30,31	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:

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

Electron density around Chain H:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
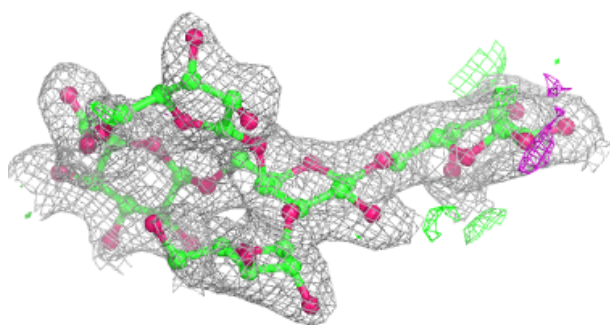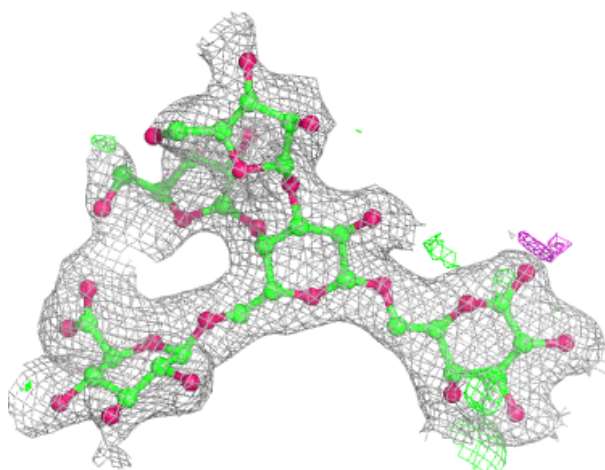
Electron density around Chain F:

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

Electron density around Chain G:

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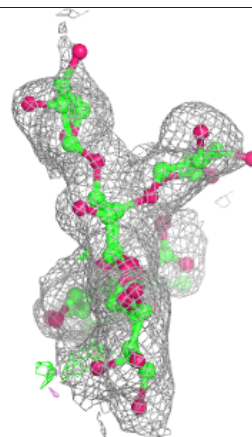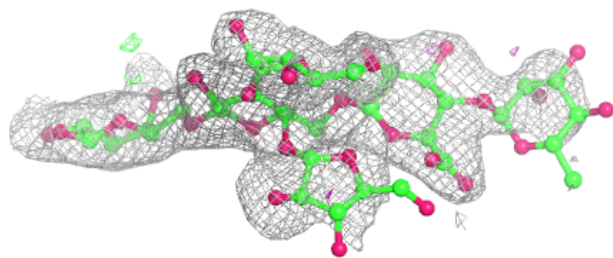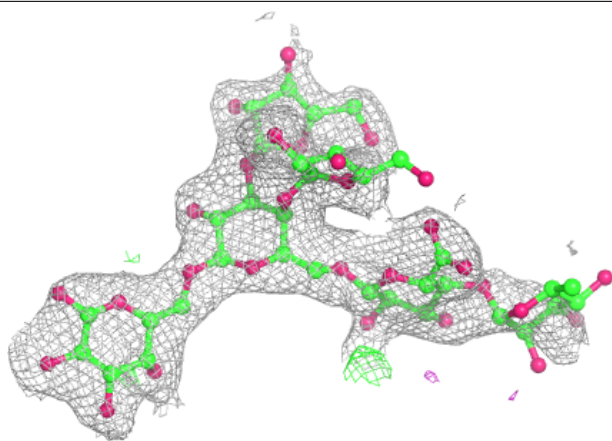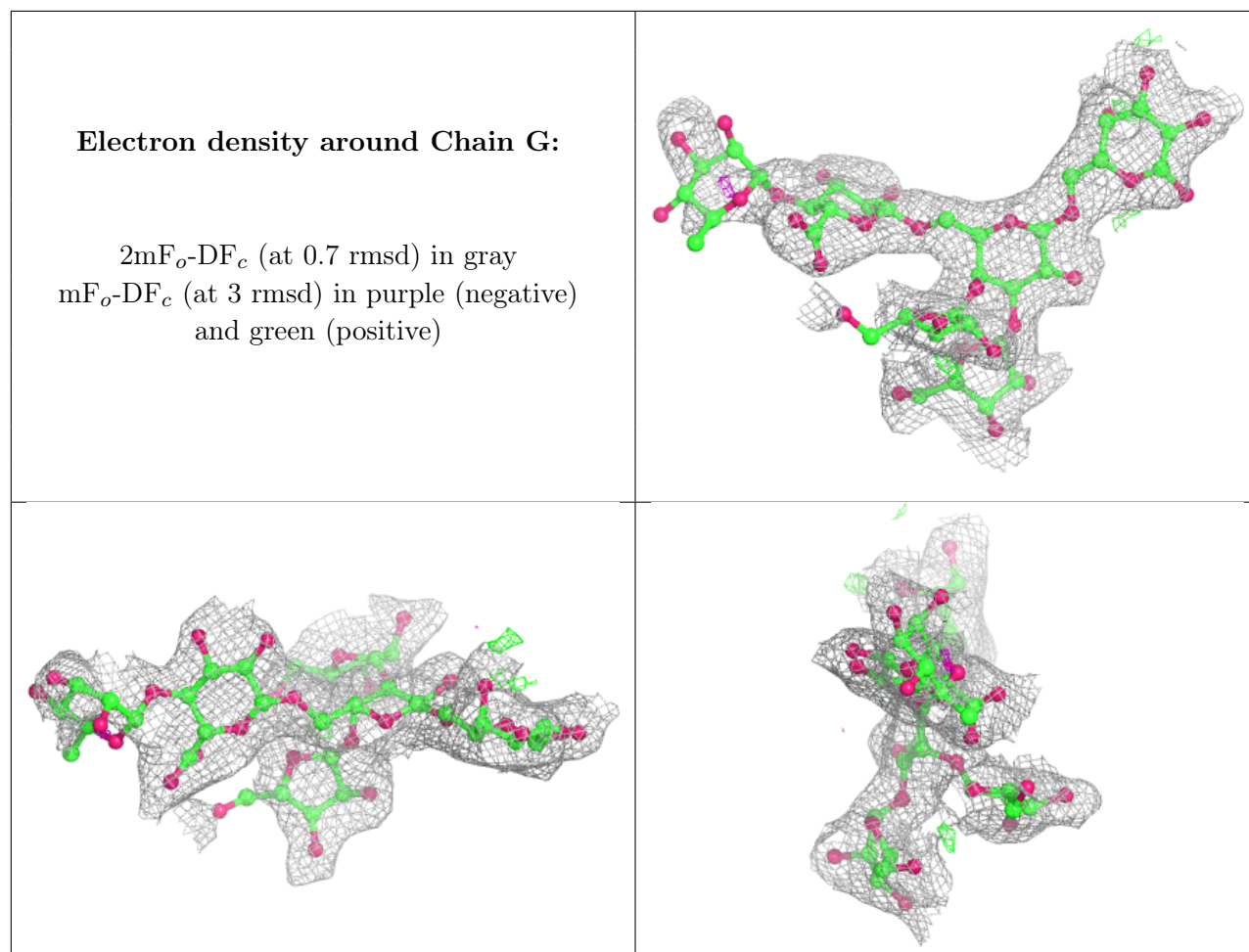
6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.