



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 7, 2023 – 06:45 am GMT

PDB ID : 1HHC  
Title : Crystal structure of Decaplanin - space group P21, second form  
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Deposited on : 2000-12-22  
Resolution : 1.13 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)

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>.

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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  
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

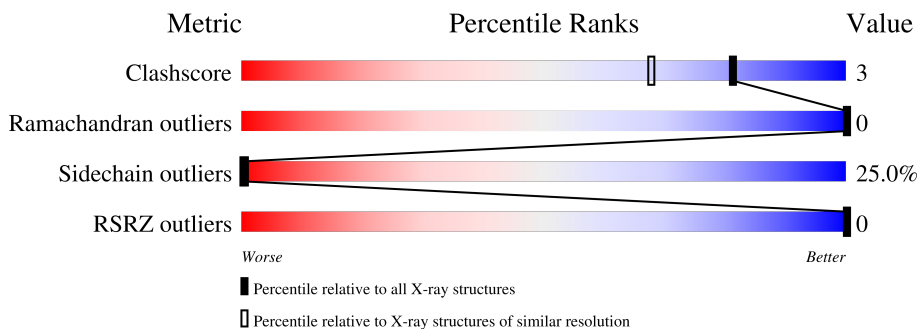
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.13 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))
Clashscore	141614	1205 (1.14-1.10)
Ramachandran outliers	138981	1168 (1.14-1.10)
Sidechain outliers	138945	1165 (1.14-1.10)
RSRZ outliers	127900	1146 (1.14-1.10)

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	7	
1	B	7	
1	C	7	
1	D	7	
2	E	2	
2	F	2	
2	G	2	

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Mol	Chain	Length	Quality of chain
2	H	2	 100%

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 629 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 DECAPLANIN.

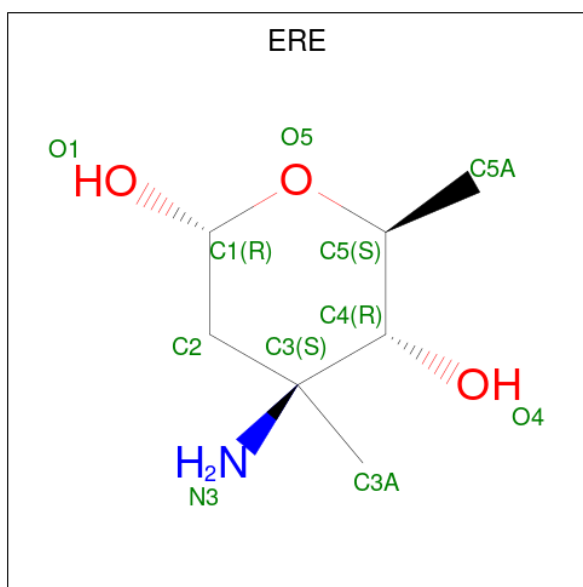
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	Cl	N	O			
1	A	7	Total 79	C 53	Cl 1	N 8	O 17	0	0	0
1	B	7	Total 76	C 50	Cl 1	N 8	O 17	0	0	0
1	C	7	Total 79	C 53	Cl 1	N 8	O 17	0	0	0
1	D	7	Total 79	C 53	Cl 1	N 8	O 17	0	0	0

- Molecule 2 is an oligosaccharide called alpha-L-rhamnopyranose-(1-2)-beta-D-glucopyranoside.



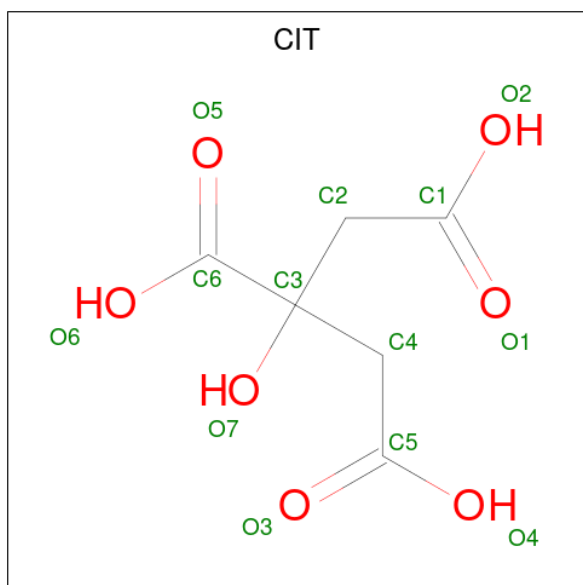
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	E	2	Total 21	C 12	O 9	0	0	0
2	F	2	Total 21	C 12	O 9	0	0	0
2	G	2	Total 21	C 12	O 9	0	0	0
2	H	2	Total 21	C 12	O 9	0	0	0

- Molecule 3 is 4-epi-vancosamine (three-letter code: ERE) (formula: C<sub>7</sub>H<sub>15</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	Total	C	N	O	0	0
			10	7	1	2		
3	B	1	Total	C	N	O	0	0
			10	7	1	2		
3	C	1	Total	C	N	O	0	0
			10	7	1	2		
3	D	1	Total	C	N	O	0	0
			10	7	1	2		

- Molecule 4 is CITRIC ACID (three-letter code: CIT) (formula: C<sub>6</sub>H<sub>8</sub>O<sub>7</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 13 6 7	0	0
4	C	1	Total C O 3 1 2	0	0
4	D	1	Total C O 13 6 7	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	41	Total O 41 41	0	0
5	B	42	Total O 42 42	0	0
5	C	38	Total O 38 38	0	0
5	D	42	Total O 42 42	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: DECAPLANIN

Chain A: 



- Molecule 1: DECAPLANIN

Chain B: 



- Molecule 1: DECAPLANIN

Chain C: 

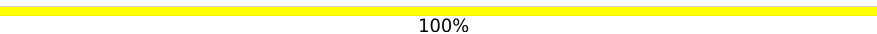


- Molecule 1: DECAPLANIN

Chain D: 



- Molecule 2: alpha-L-rhamnopyranose-(1-2)-beta-D-glucopyranose

Chain E: 



- Molecule 2: alpha-L-rhamnopyranose-(1-2)-beta-D-glucopyranose

Chain F: 




- Molecule 2: alpha-L-rhamnopyranose-(1-2)-beta-D-glucopyranose

Chain G:  100%

BGC1  
RAM2

- Molecule 2: alpha-L-rhamnopyranose-(1-2)-beta-D-glucopyranose

Chain H:  100%

BGC1  
RAM2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	28.95Å 31.91Å 34.73Å 90.00° 109.87° 90.00°	Depositor
Resolution (Å)	18.08 – 1.13 18.10 – 1.13	Depositor EDS
% Data completeness (in resolution range)	96.8 (18.08-1.13) 89.8 (18.10-1.13)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.75 (at 1.13Å)	Xtrriage
Refinement program	SHELXL-97	Depositor
R, $R_{free}$	0.224 , 0.264 0.205 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	5.1	Xtrriage
Anisotropy	0.325	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 82.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	629	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	8.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.88% 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: GHP, OMZ, 3FG, OMX, CIT, MLU, ERE, RAM, BGC

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	0.70	0/7	1.21	0/8
1	B	0.98	0/7	2.41	0/8
1	C	0.70	0/7	0.87	0/8
1	D	0.39	0/7	1.16	0/8
All	All	0.72	0/28	1.53	0/32

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 [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	79	0	45	1	0
1	B	76	0	38	0	0
1	C	79	0	45	0	0
1	D	79	0	46	0	0
2	E	21	0	18	0	0
2	F	21	0	18	0	0
2	G	21	0	18	0	0
2	H	21	0	18	0	0
3	A	10	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	10	0	13	0	0
3	C	10	0	13	0	0
3	D	10	0	13	0	0
4	A	13	0	5	2	0
4	C	3	0	0	0	0
4	D	13	0	5	0	0
5	A	41	0	0	0	0
5	B	42	0	0	0	0
5	C	38	0	0	0	0
5	D	42	0	0	0	0
All	All	629	0	308	2	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 3.

All (2) 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)
4:A:1002:CIT:H42	4:A:1002:CIT:O1	2.03	0.59
1:A:1:MLU:HD12	4:A:1002:CIT:C6	2.49	0.43

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	1/7 (14%)	1 (100%)	0	0	100	100
1	B	1/7 (14%)	0	1 (100%)	0	100	100
1	C	1/7 (14%)	0	1 (100%)	0	100	100
1	D	1/7 (14%)	1 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	4/28 (14%)	2 (50%)	2 (50%)	0	100	100

There are no Ramachandran outliers to report.

### 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	1/1 (100%)	1 (100%)	0	100	100
1	B	1/1 (100%)	0	1 (100%)	0	0
1	C	1/1 (100%)	1 (100%)	0	100	100
1	D	1/1 (100%)	1 (100%)	0	100	100
All	All	4/4 (100%)	3 (75%)	1 (25%)	0	0

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	3	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

24 non-standard protein/DNA/RNA residues 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
1	MLU	B	1	1	4,5,9	1.30	1 (25%)	1,5,11	1.12	0
1	MLU	C	1	1	7,8,9	0.68	0	6,9,11	0.54	0
1	GHP	B	5	1	10,11,12	1.37	1 (10%)	11,14,16	1.62	2 (18%)
1	OMX	B	6	1,3	11,13,14	0.92	0	15,17,19	2.46	5 (33%)
1	MLU	D	1	1	7,8,9	0.99	1 (14%)	6,9,11	1.52	1 (16%)
1	3FG	C	7	1	12,13,13	0.98	1 (8%)	14,18,18	2.02	5 (35%)
1	OMX	C	6	1,3	11,13,14	0.94	1 (9%)	15,17,19	1.27	2 (13%)
1	3FG	D	7	1	12,13,13	1.36	2 (16%)	14,18,18	1.93	3 (21%)
1	MLU	A	1	1	7,8,9	1.11	1 (14%)	6,9,11	0.80	0
1	GHP	A	5	1	10,11,12	0.68	0	11,14,16	1.70	2 (18%)
1	GHP	D	4	1,2	10,11,12	1.26	1 (10%)	11,14,16	1.83	3 (27%)
1	GHP	C	5	1	10,11,12	0.96	1 (10%)	11,14,16	2.40	4 (36%)
1	3FG	A	7	1	12,13,13	1.26	1 (8%)	14,18,18	1.67	5 (35%)
1	OMX	A	6	1,3	11,13,14	1.11	1 (9%)	15,17,19	1.74	5 (33%)
1	OMZ	B	2	1	12,14,15	0.74	0	17,19,21	1.00	0
1	OMZ	C	2	1	12,14,15	1.20	1 (8%)	17,19,21	1.63	5 (29%)
1	GHP	B	4	1,2	10,11,12	1.31	1 (10%)	11,14,16	1.79	3 (27%)
1	OMX	D	6	1,3	11,13,14	1.04	2 (18%)	15,17,19	2.62	8 (53%)
1	GHP	C	4	1,2	10,11,12	1.14	0	11,14,16	0.88	0
1	GHP	D	5	1	10,11,12	0.91	0	11,14,16	2.17	4 (36%)
1	GHP	A	4	1,2	10,11,12	1.23	1 (10%)	11,14,16	1.92	5 (45%)
1	OMZ	D	2	1	12,14,15	1.40	2 (16%)	17,19,21	1.52	3 (17%)
1	OMZ	A	2	1	12,14,15	1.05	0	17,19,21	2.36	4 (23%)
1	3FG	B	7	1	12,13,13	1.15	0	14,18,18	1.38	3 (21%)

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
1	MLU	B	1	1	-	0/1/4/10	-
1	MLU	C	1	1	-	0/5/8/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	GHP	B	5	1	-	0/4/6/8	0/1/1/1
1	OMX	B	6	1,3	-	1/9/10/12	0/1/1/1
1	MLU	D	1	1	-	0/5/8/10	-
1	3FG	C	7	1	-	1/8/8/8	0/1/1/1
1	OMX	C	6	1,3	-	1/9/10/12	0/1/1/1
1	3FG	D	7	1	-	2/8/8/8	0/1/1/1
1	MLU	A	1	1	-	0/5/8/10	-
1	GHP	A	5	1	-	0/4/6/8	0/1/1/1
1	GHP	D	4	1,2	-	1/4/6/8	0/1/1/1
1	GHP	C	5	1	-	0/4/6/8	0/1/1/1
1	3FG	A	7	1	-	2/8/8/8	0/1/1/1
1	OMX	A	6	1,3	-	1/9/10/12	0/1/1/1
1	OMZ	B	2	1	-	1/9/10/12	0/1/1/1
1	OMZ	C	2	1	-	1/9/10/12	0/1/1/1
1	GHP	B	4	1,2	-	3/4/6/8	0/1/1/1
1	OMX	D	6	1,3	-	1/9/10/12	0/1/1/1
1	GHP	C	4	1,2	-	1/4/6/8	0/1/1/1
1	GHP	D	5	1	-	0/4/6/8	0/1/1/1
1	GHP	A	4	1,2	-	2/4/6/8	0/1/1/1
1	OMZ	D	2	1	-	0/9/10/12	0/1/1/1
1	OMZ	A	2	1	-	1/9/10/12	0/1/1/1
1	3FG	B	7	1	-	2/8/8/8	0/1/1/1

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	4	GHP	C6-C1	-3.54	1.33	1.39
1	B	5	GHP	C1-CA	3.22	1.56	1.52
1	D	4	GHP	C6-C1	-3.11	1.34	1.39
1	D	2	OMZ	CG-CB	-2.95	1.47	1.51
1	C	2	OMZ	CZ-CE1	-2.81	1.36	1.39

The worst 5 of 72 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	OMZ	CD1-CE1-CZ	-7.82	116.38	120.91
1	D	6	OMX	CE1-CD1-CG	6.06	127.30	121.20
1	B	6	OMX	CD1-CE1-CZ	-5.78	113.54	119.88
1	D	5	GHP	C3-C2-C1	5.10	126.33	121.20
1	B	6	OMX	CE1-CD1-CG	4.96	126.19	121.20

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	6	OMX	O-C-CA-CB
1	B	6	OMX	O-C-CA-CB
1	D	6	OMX	O-C-CA-CB
1	A	7	3FG	C-CA-CB-CG1
1	A	7	3FG	C-CA-CB-CG2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	MLU	1	0

## 5.5 Carbohydrates [i](#)

8 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	E	1	1,2	11,11,12	1.83	2 (18%)	15,15,17	1.41	3 (20%)
2	RAM	E	2	2	10,10,11	2.32	3 (30%)	14,14,16	2.12	5 (35%)
2	BGC	F	1	1,2	11,11,12	1.57	2 (18%)	15,15,17	1.84	4 (26%)
2	RAM	F	2	2	10,10,11	3.44	4 (40%)	14,14,16	1.59	2 (14%)
2	BGC	G	1	1,2	11,11,12	1.42	2 (18%)	15,15,17	1.15	1 (6%)
2	RAM	G	2	2	10,10,11	2.73	1 (10%)	14,14,16	1.59	1 (7%)
2	BGC	H	1	1,2	11,11,12	1.47	1 (9%)	15,15,17	1.41	2 (13%)
2	RAM	H	2	2	10,10,11	2.62	1 (10%)	14,14,16	1.91	6 (42%)

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	E	1	1,2	-	0/2/19/22	0/1/1/1
2	RAM	E	2	2	-	-	0/1/1/1
2	BGC	F	1	1,2	-	0/2/19/22	0/1/1/1
2	RAM	F	2	2	-	-	0/1/1/1
2	BGC	G	1	1,2	-	0/2/19/22	0/1/1/1
2	RAM	G	2	2	-	-	0/1/1/1
2	BGC	H	1	1,2	-	0/2/19/22	0/1/1/1
2	RAM	H	2	2	-	-	0/1/1/1

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	RAM	C2-C3	-8.86	1.39	1.52
2	G	2	RAM	C2-C3	-8.00	1.40	1.52
2	H	2	RAM	C2-C3	-7.52	1.41	1.52
2	E	2	RAM	C2-C3	-5.73	1.44	1.52
2	F	2	RAM	O5-C1	4.79	1.51	1.43

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	RAM	C1-C2-C3	5.24	116.11	109.67
2	E	2	RAM	O3-C3-C2	-3.95	102.42	109.99
2	F	2	RAM	C2-C3-C4	3.76	117.40	110.89
2	E	2	RAM	C3-C4-C5	-3.52	104.30	109.77
2	E	2	RAM	O5-C5-C4	3.48	115.77	109.52

There are no chirality outliers.

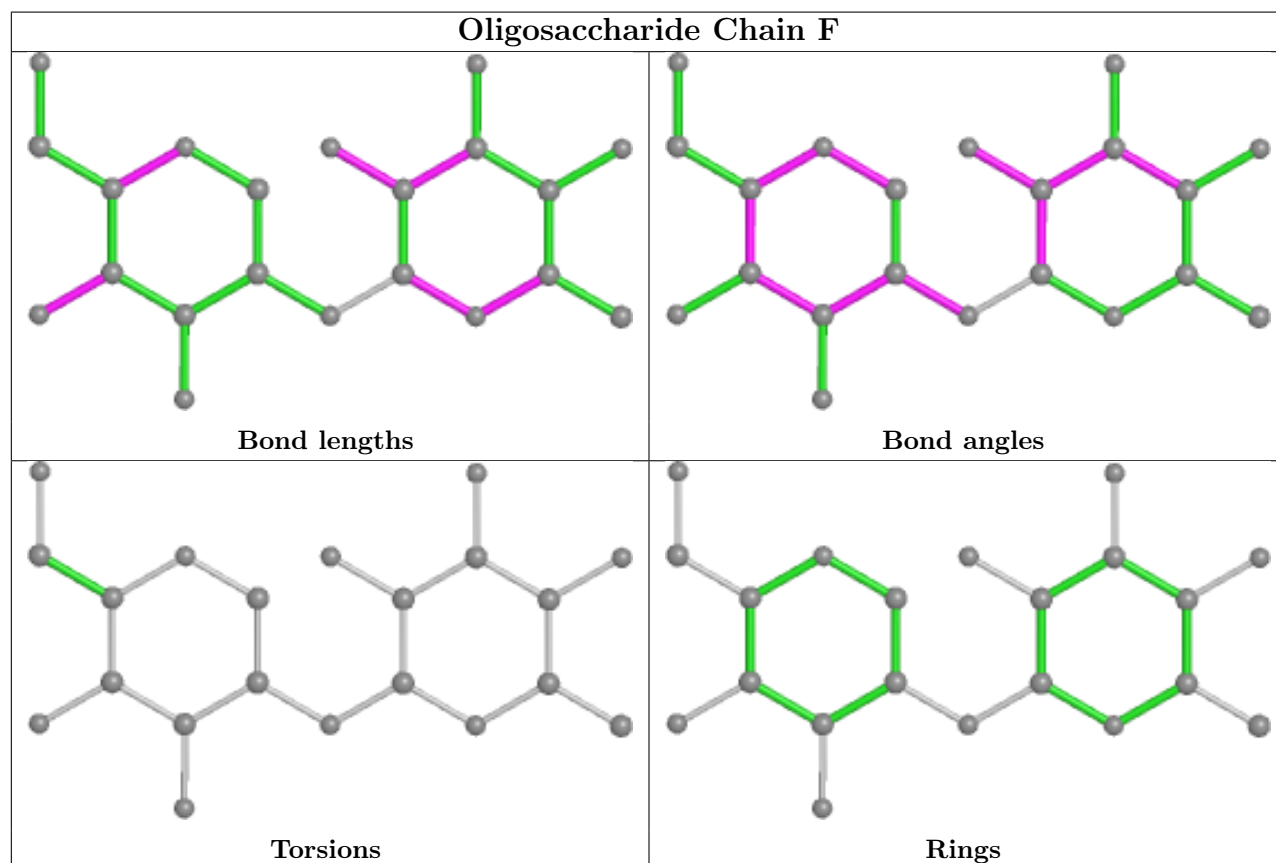
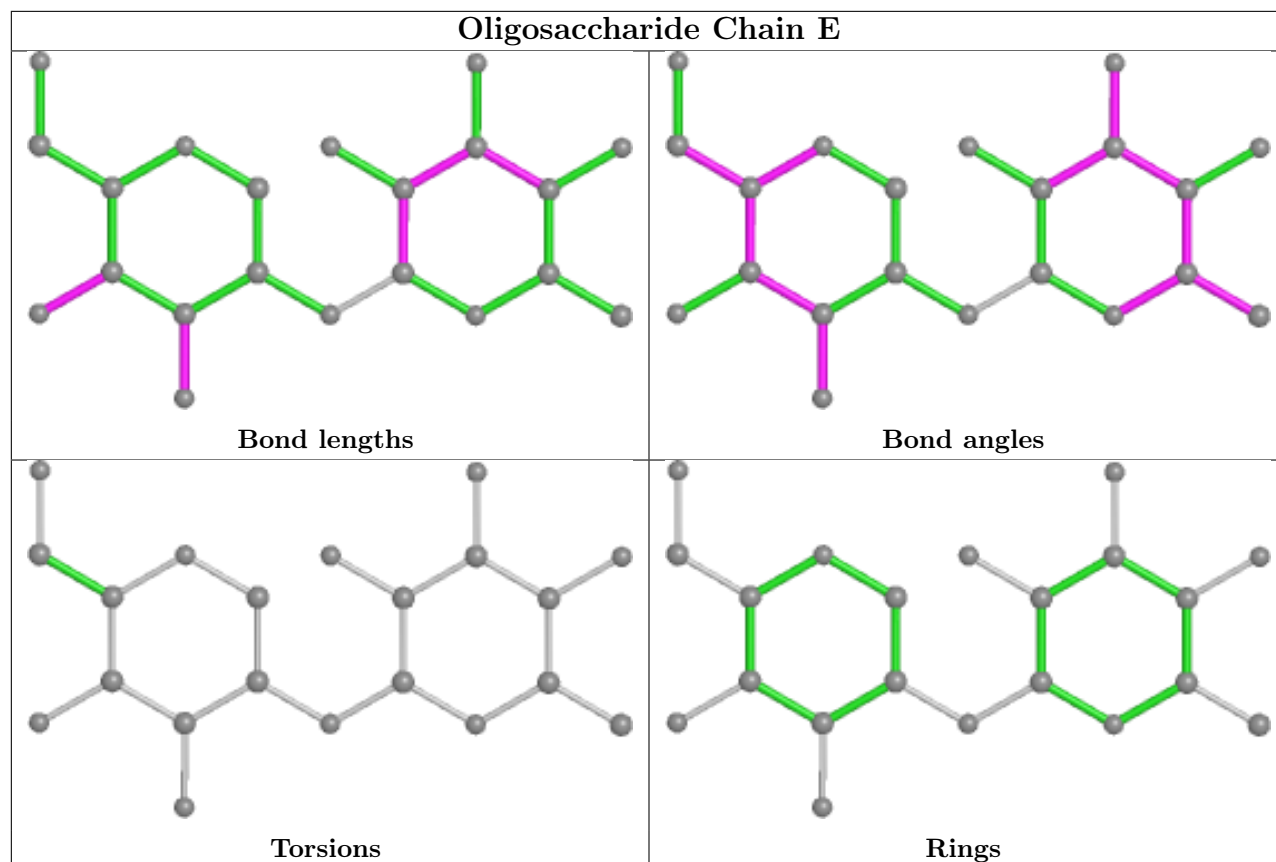
There are no torsion outliers.

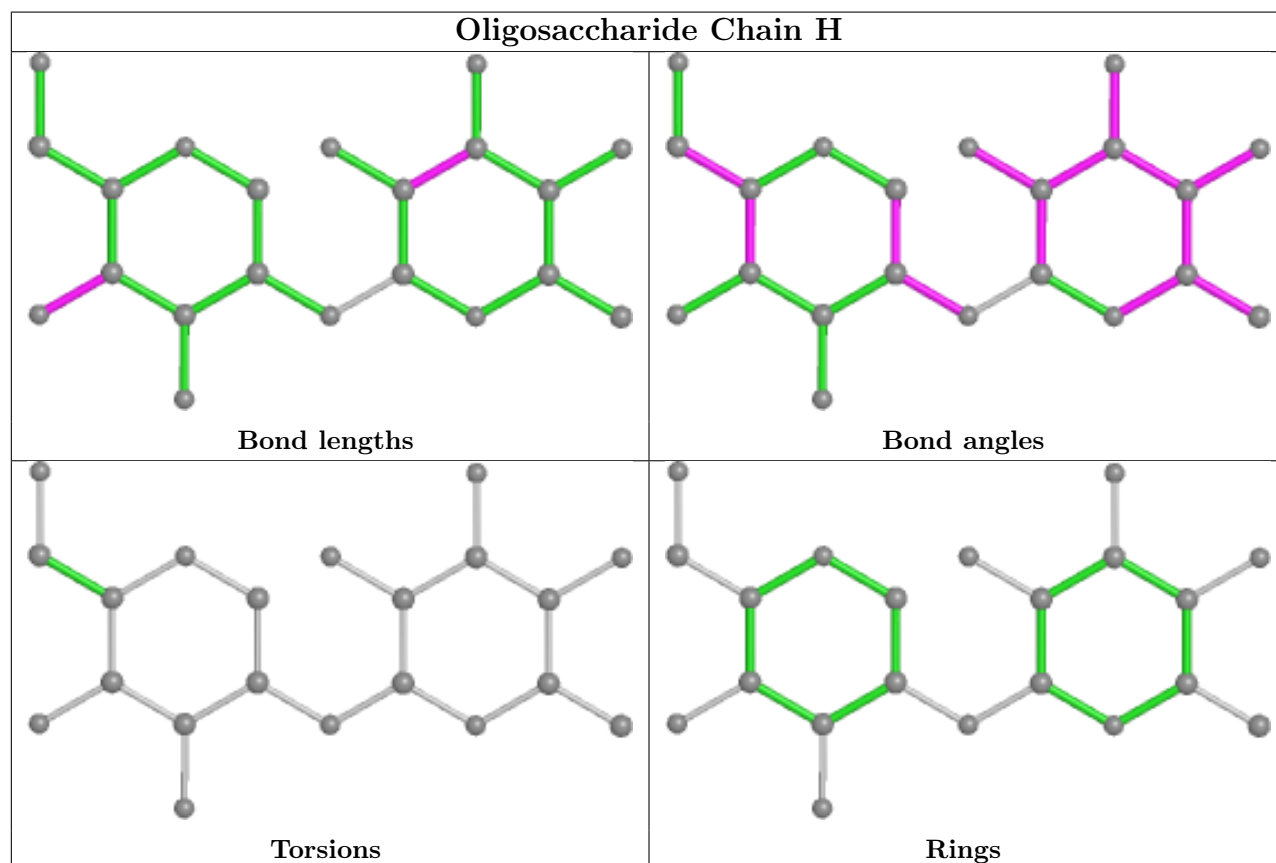
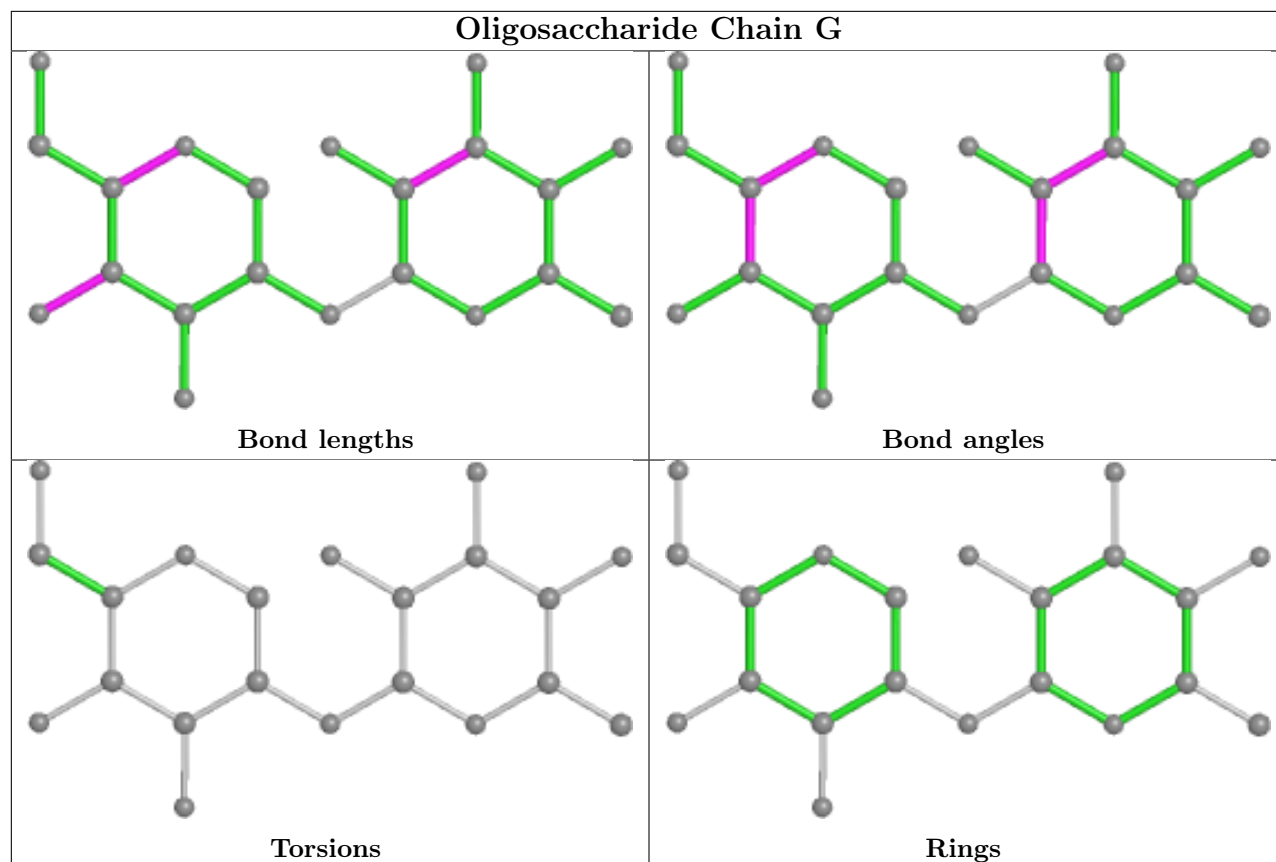
There are no ring outliers.

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.







## 5.6 Ligand geometry [i](#)

7 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
4	CIT	A	1002	-	12,12,12	1.54	3 (25%)	17,17,17	1.86	5 (29%)
3	ERE	B	8	1	6,10,11	0.96	0	6,15,17	0.98	1 (16%)
4	CIT	C	1003	-	2,2,12	0.55	0	1,1,17	0.34	0
3	ERE	A	8	1	6,10,11	0.90	0	6,15,17	1.54	1 (16%)
4	CIT	D	1001	-	12,12,12	1.31	2 (16%)	17,17,17	1.80	4 (23%)
3	ERE	C	8	1	6,10,11	0.92	0	6,15,17	1.60	2 (33%)
3	ERE	D	8	1	6,10,11	1.24	1 (16%)	6,15,17	1.24	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
4	CIT	A	1002	-	-	1/16/16/16	-
3	ERE	D	8	1	-	-	0/1/1/1
3	ERE	B	8	1	-	-	0/1/1/1
3	ERE	A	8	1	-	-	0/1/1/1
4	CIT	D	1001	-	-	2/16/16/16	-
3	ERE	C	8	1	-	-	0/1/1/1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	CIT	O4-C5	-2.79	1.21	1.30
4	D	1001	CIT	O4-C5	-2.48	1.22	1.30
4	D	1001	CIT	C2-C3	2.42	1.56	1.53
4	A	1002	CIT	C3-C6	-2.40	1.50	1.53
3	D	8	ERE	O4-C4	2.31	1.47	1.42

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

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
4	D	1001	CIT	O6-C6-C3	3.92	119.86	113.05
4	A	1002	CIT	O4-C5-C4	3.34	125.08	114.35
3	A	8	ERE	C5A-C5-C4	-3.09	107.30	112.57
4	A	1002	CIT	O4-C5-O3	-3.03	115.76	123.30
4	D	1001	CIT	C2-C3-C6	-2.95	103.77	110.11

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	1001	CIT	C3-C4-C5-O3
4	D	1001	CIT	C3-C4-C5-O4
4	A	1002	CIT	O7-C3-C4-C5

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	CIT	2	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	1/7 (14%)	0.62	0 100 100	6, 6, 6, 6	0
1	B	1/7 (14%)	0.69	0 100 100	6, 6, 6, 6	0
1	C	1/7 (14%)	0.09	0 100 100	5, 5, 5, 5	0
1	D	1/7 (14%)	0.60	0 100 100	6, 6, 6, 6	0
All	All	4/28 (14%)	0.50	0 100 100	5, 6, 6, 6	0

There are no RSRZ outliers to report.

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	MLU	B	1	6/10	0.88	0.16	9,10,11,12	0
1	MLU	A	1	9/10	0.94	0.12	6,7,9,11	0
1	MLU	D	1	9/10	0.94	0.11	6,7,8,9	0
1	MLU	C	1	9/10	0.96	0.10	4,5,7,8	0
1	OMX	A	6	13/14	0.96	0.10	5,5,6,6	0
1	GHP	A	4	11/12	0.96	0.11	3,4,4,4	0
1	GHP	B	4	11/12	0.96	0.10	3,4,4,5	0
1	GHP	D	4	11/12	0.96	0.11	3,3,4,4	0
1	GHP	C	5	11/12	0.96	0.11	2,3,4,6	0
1	3FG	A	7	13/13	0.96	0.10	7,8,10,11	0
1	GHP	A	5	11/12	0.97	0.08	4,5,6,7	0
1	GHP	B	5	11/12	0.97	0.11	2,3,4,4	0
1	OMZ	B	2	14/15	0.97	0.11	5,6,8,8	0
1	GHP	D	5	11/12	0.97	0.09	2,3,3,6	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	OMX	B	6	13/14	0.97	0.10	2,3,4,4	0
1	3FG	D	7	13/13	0.97	0.09	3,3,6,6	0
1	OMX	D	6	13/14	0.98	0.08	2,3,3,4	0
1	OMZ	A	2	14/15	0.98	0.09	4,5,5,6	0
1	GHP	C	4	11/12	0.98	0.09	2,2,3,3	0
1	OMX	C	6	13/14	0.98	0.08	2,3,3,4	0
1	3FG	B	7	13/13	0.98	0.09	2,2,3,4	0
1	3FG	C	7	13/13	0.98	0.07	4,5,7,8	0
1	OMZ	D	2	14/15	0.98	0.10	4,5,6,7	0
1	OMZ	C	2	14/15	0.99	0.08	3,4,5,6	0

### 6.3 Carbohydrates [i](#)

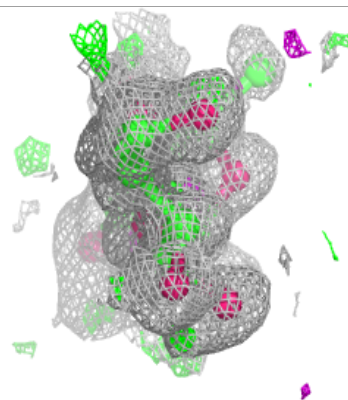
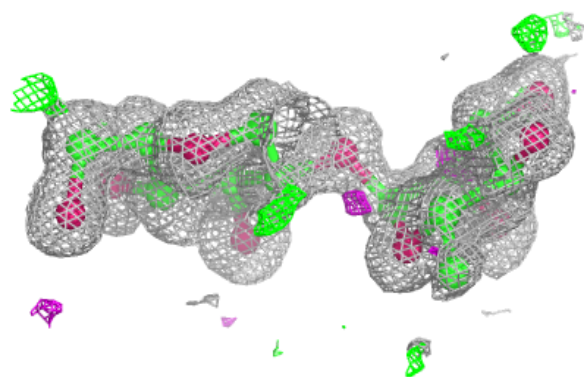
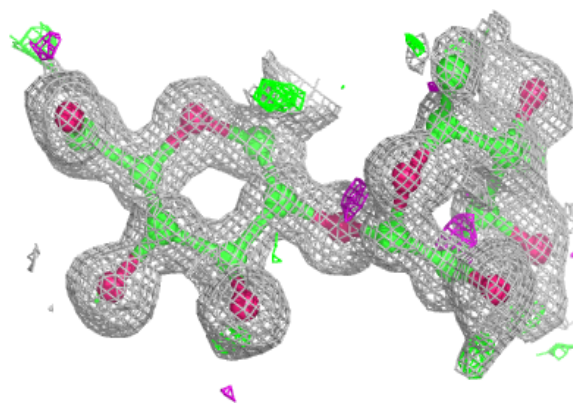
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<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	RAM	E	2	10/11	0.91	0.16	7,9,12,13	0
2	BGC	F	1	11/12	0.94	0.13	5,7,8,9	0
2	RAM	H	2	10/11	0.94	0.12	6,7,9,9	0
2	RAM	G	2	10/11	0.95	0.10	5,6,9,9	0
2	BGC	H	1	11/12	0.95	0.12	5,5,7,7	0
2	RAM	F	2	10/11	0.95	0.13	6,7,9,10	0
2	BGC	G	1	11/12	0.96	0.10	3,4,4,6	0
2	BGC	E	1	11/12	0.97	0.10	4,5,6,6	0

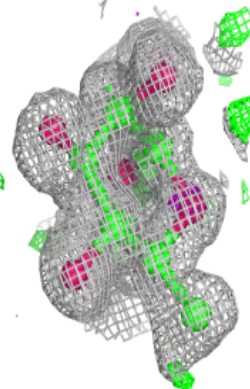
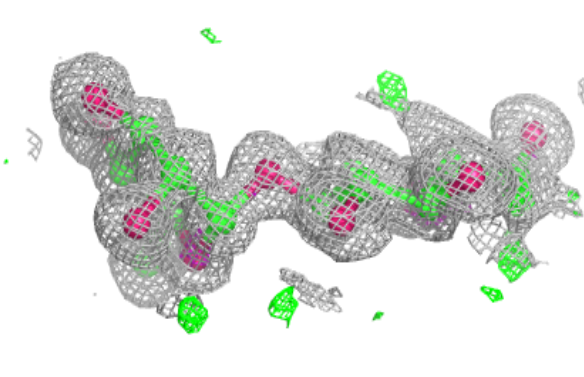
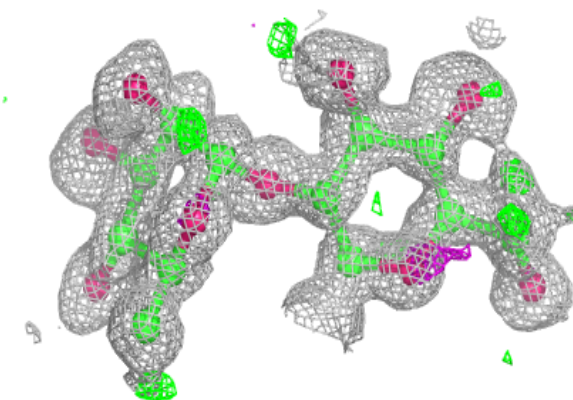
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

**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 F:**

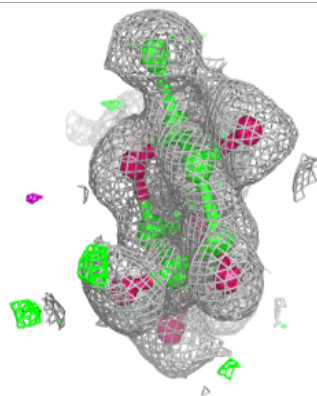
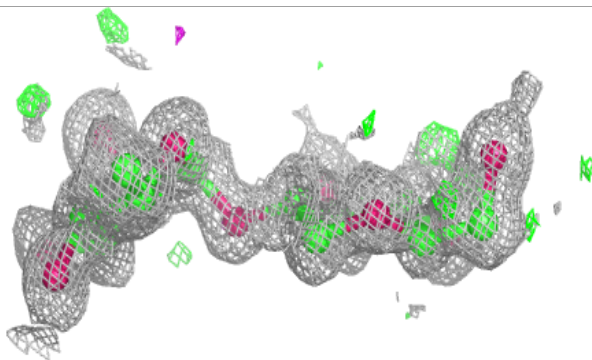
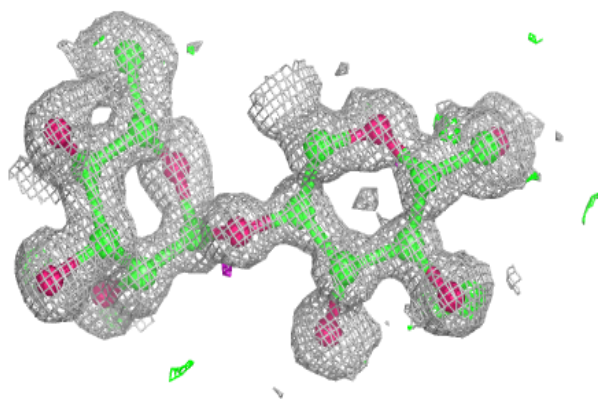
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



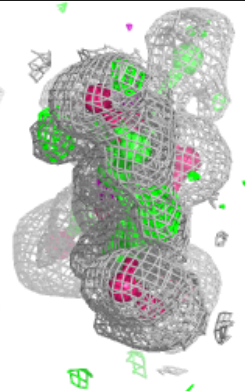
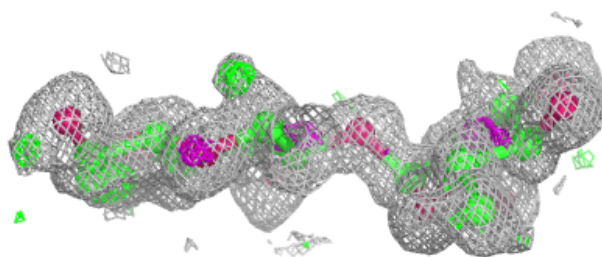
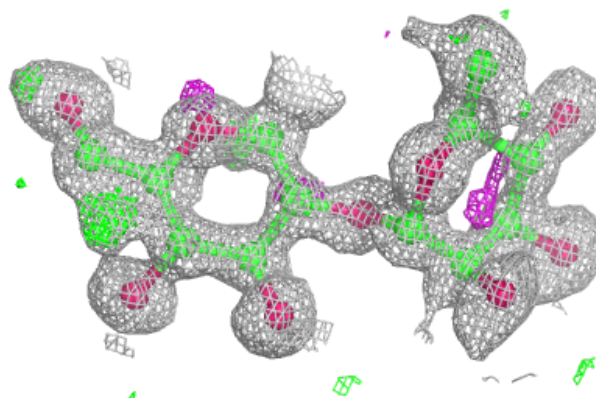


**Electron density around Chain G:**

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

**Electron density around Chain H:**

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





## 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( $\text{\AA}^2$ )	Q<0.9
4	CIT	A	1002	13/13	0.89	0.23	2,5,7,7	13
3	ERE	C	8	10/11	0.95	0.11	4,5,5,7	0
3	ERE	A	8	10/11	0.95	0.10	6,7,9,10	0
3	ERE	D	8	10/11	0.96	0.10	3,4,5,7	0
4	CIT	C	1003	3/13	0.97	0.07	8,8,8,8	0
3	ERE	B	8	10/11	0.98	0.08	3,4,5,6	0
4	CIT	D	1001	13/13	0.98	0.08	4,4,5,5	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.