



# wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 29, 2026 – 04:08 am BST

PDB ID : 9I4L / pdb\_00009i4l  
Title : Galectin-1 in Complex with Methyl 2'-fluoro-N-acetyllactosaminide  
Authors : Pachl, P.  
Deposited on : 2025-01-25  
Resolution : 1.43 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	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
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

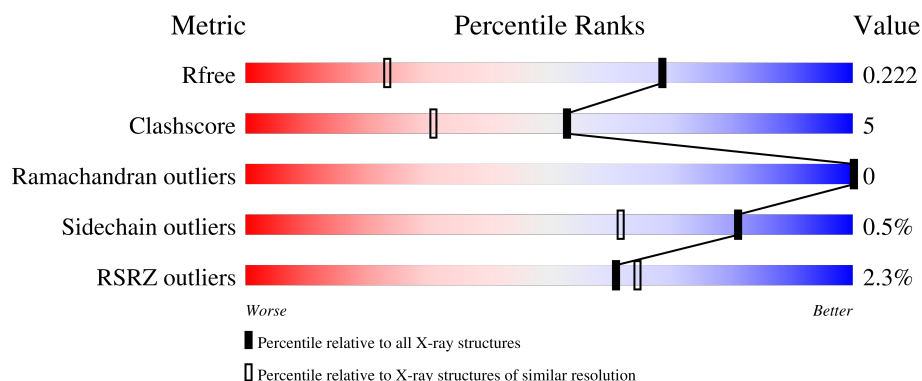
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.43 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	4041 (1.44-1.40)
Clashscore	190562	4154 (1.44-1.40)
Ramachandran outliers	187476	4083 (1.44-1.40)
Sidechain outliers	187428	4082 (1.44-1.40)
RSRZ outliers	180081	4039 (1.44-1.40)

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	136	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>..</div> </div> </div>
1	B	136	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>..</div> </div> </div>
2	C	2	<div> <div>50%</div> <div>50%</div> </div>
2	D	2	<div> <div>100%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4544 atoms, of which 2107 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 Galectin-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	133	Total	C	H	N	O	S	23	2	0
			2054	656	1015	177	195	11			
1	B	133	Total	C	H	N	O	S	24	5	0
			2081	664	1028	179	200	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP P09382
B	-1	SER	-	expression tag	UNP P09382

- Molecule 2 is an oligosaccharide called 2-deoxy-2-fluoro-beta-D-galactopyranose-(1-4)-methyl 2-acetamido-2-deoxy-beta-D-glucopyranoside.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	2	Total	C	F	H	N	O	8	0	0
			53	15	1	26	1	10			
2	D	2	Total	C	F	H	N	O	8	0	0
			53	15	1	26	1	10			

- Molecule 3 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
3	A	1	Total	C	H	O	2	0
			10	2	6	2		
3	B	1	Total	C	H	O	2	0
			10	2	6	2		


- Molecule 4 is water.

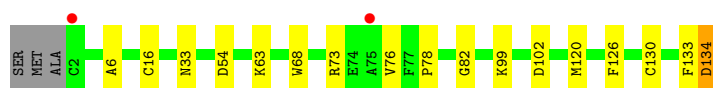
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	134	Total	O	0	2
			135	135		
4	B	146	Total	O	0	4
			148	148		

### 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: Galectin-1

Chain A: 



- Molecule 1: Galectin-1

Chain B: 



- Molecule 2: 2-deoxy-2-fluoro-beta-D-galactopyranose-(1-4)-methyl 2-acetamido-2-deoxy-beta-D-glucopyranoside

Chain C: 



- Molecule 2: 2-deoxy-2-fluoro-beta-D-galactopyranose-(1-4)-methyl 2-acetamido-2-deoxy-beta-D-glucopyranoside

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.25Å 58.27Å 111.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.18 – 1.43 41.18 – 1.43	Depositor EDS
% Data completeness (in resolution range)	99.3 (41.18-1.43) 99.3 (41.18-1.43)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.99 (at 1.43Å)	Xtriage
Refinement program	REFMAC 5.8.0430 (refmacat 0.4.88)	Depositor
R, $R_{free}$	0.165 , 0.220 0.166 , 0.222	Depositor DCC
$R_{free}$ test set	2718 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.6	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 49.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	4544	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.94% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, CME, 2FG, MAG

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	0.76	0/1034	1.19	2/1395 (0.1%)
1	B	0.76	0/1056	1.15	4/1425 (0.3%)
All	All	0.76	0/2090	1.17	6/2820 (0.2%)

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	B	0	2

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	PHE	CA-CB-CG	6.46	120.25	113.80
1	B	26	ASP	CA-CB-CG	6.06	118.66	112.60
1	B	39	ASN	CA-CB-CG	-5.91	106.69	112.60
1	B	105	GLU	N-CA-CB	-5.86	100.62	111.52
1	A	102	ASP	CB-CA-C	5.63	120.66	111.26

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	18[A]	ARG	Sidechain
1	B	18[B]	ARG	Sidechain

## 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	1039	1015	1007	9	0
1	B	1053	1028	1020	10	0
2	C	27	26	25	2	0
2	D	27	26	25	0	0
3	A	4	6	6	1	0
3	B	4	6	6	2	0
4	A	135	0	0	0	0
4	B	148	0	0	2	0
All	All	2437	2107	2089	21	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 5.

The worst 5 of 21 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:C:2:2FG:F2	2:C:2:2FG:C2	1.56	1.43
1:A:6:ALA:HB3	1:A:120[A]:MET:HG2	1.85	0.57
1:A:82:GLY:C	3:A:201:EDO:H12	2.32	0.53
1:A:133:PHE:O	1:A:134:ASP:C	2.51	0.53
1:A:76:VAL:HG12	1:A:78:PRO:HD3	1.90	0.52

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	130/136 (96%)	126 (97%)	4 (3%)	0	100	100
1	B	132/136 (97%)	130 (98%)	2 (2%)	0	100	100
All	All	262/272 (96%)	256 (98%)	6 (2%)	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	107/108 (99%)	106 (99%)	1 (1%)	70	43
1	B	110/108 (102%)	110 (100%)	0	100	100
All	All	217/216 (100%)	216 (100%)	1 (0%)	81	62

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

Mol	Chain	Res	Type
1	A	134	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	56	ASN
1	B	52	HIS
1	B	56	ASN
1	B	80	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

6 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	CME	A	88	1	8,9,10	0.86	0	5,9,11	0.83	0
1	CME	B	16	1	8,9,10	0.48	0	5,9,11	0.63	0
1	CME	B	130	1	8,9,10	0.56	0	5,9,11	0.95	0
1	CME	A	16	1	8,9,10	0.41	0	5,9,11	0.50	0
1	CME	A	130	1	8,9,10	1.03	0	5,9,11	1.74	2 (40%)
1	CME	B	88	1	8,9,10	0.78	0	5,9,11	0.80	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
1	CME	A	88	1	-	0/5/8/10	-
1	CME	B	16	1	-	1/5/8/10	-
1	CME	B	130	1	-	0/5/8/10	-
1	CME	A	16	1	-	0/5/8/10	-
1	CME	A	130	1	-	2/5/8/10	-
1	CME	B	88	1	-	0/5/8/10	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	CME	CZ-CE-SD	-2.68	104.05	113.37
1	A	130	CME	CB-SG-SD	2.03	109.07	103.82

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	16	CME	SD-CE-CZ-OH
1	A	130	CME	CA-CB-SG-SD
1	A	130	CME	CE-SD-SG-CB

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	16	CME	1	0

## 5.5 Carbohydrates [i](#)

4 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	MAG	C	1	2	16,16,16	1.55	4 (25%)	22,22,22	1.11	3 (13%)
2	2FG	C	2	2	11,11,12	3.09	5 (45%)	10,15,17	2.20	3 (30%)
2	MAG	D	1	2	16,16,16	2.26	7 (43%)	22,22,22	1.66	5 (22%)
2	2FG	D	2	2	11,11,12	2.53	4 (36%)	10,15,17	1.41	1 (10%)

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	MAG	C	1	2	-	0/8/28/28	0/1/1/1
2	2FG	C	2	2	-	0/2/19/22	0/1/1/1
2	MAG	D	1	2	-	2/8/28/28	0/1/1/1
2	2FG	D	2	2	-	0/2/19/22	0/1/1/1

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	2FG	F2-C2	6.31	1.56	1.40
2	D	2	2FG	C2-C3	4.37	1.57	1.51
2	D	2	2FG	O4-C4	4.36	1.53	1.43
2	C	2	2FG	O4-C4	4.31	1.53	1.43
2	D	1	MAG	C2-N2	4.28	1.52	1.45

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	2FG	O5-C5-C6	5.42	115.70	107.20
2	D	1	MAG	C4-C3-C2	3.75	115.84	110.34
2	D	1	MAG	C1-O5-C5	-3.73	106.37	113.69
2	C	1	MAG	O5-C1-O1	-3.15	103.67	110.97
2	D	2	2FG	C6-C5-C4	-2.50	107.14	113.00

There are no chirality outliers.

All (2) torsion outliers are listed below:

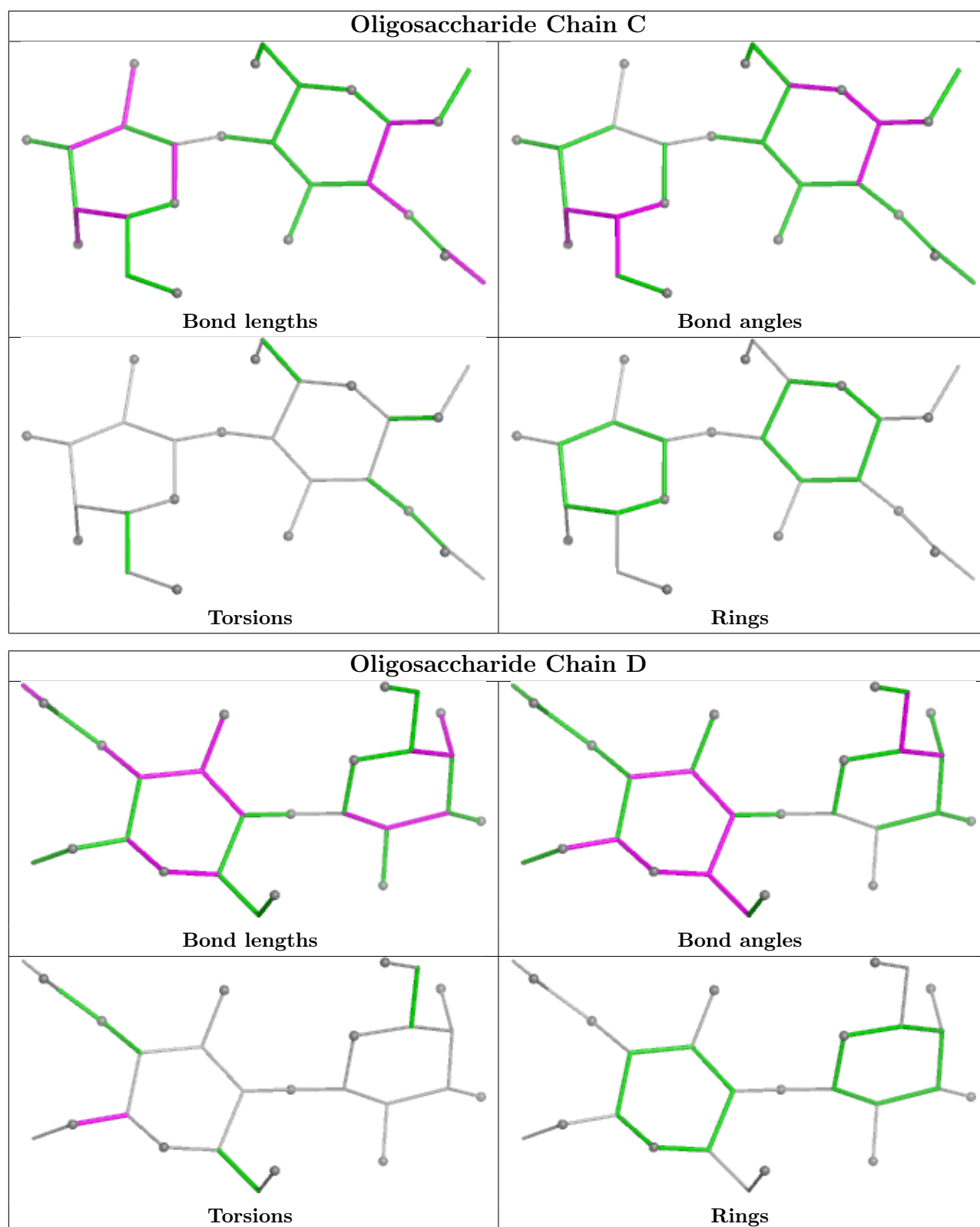
Mol	Chain	Res	Type	Atoms
2	D	1	MAG	C2-C1-O1-CM
2	D	1	MAG	O5-C1-O1-CM

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	2FG	2	0

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



## 5.6 Ligand geometry [i](#)

2 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	$\# Z  > 2$
3	EDO	B	201	-	3,3,3	0.49	0	2,2,2	0.38	0
3	EDO	A	201	-	3,3,3	0.66	0	2,2,2	0.61	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
3	EDO	B	201	-	-	0/1/1/1	-
3	EDO	A	201	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	201	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	201	EDO	2	0
3	A	201	EDO	1	0

## 5.7 Other polymers [i](#)

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, 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	130/136 (95%)	-0.36	2 (1%) 72 76	14, 27, 47, 62	2 (1%)
1	B	130/136 (95%)	-0.30	4 (3%) 51 54	10, 25, 44, 56	5 (3%)
All	All	260/272 (95%)	-0.33	6 (2%) 61 64	10, 26, 47, 62	7 (2%)

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	CYS	2.3
1	B	106	PHE	2.2
1	B	124	GLY	2.2
1	B	134[A]	ASP	2.1
1	B	123[A]	ASP	2.1

### 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	CME	A	130	10/11	0.97	0.07	17,23,42,45	1
1	CME	A	16	10/11	0.98	0.06	26,29,43,47	1
1	CME	B	16	10/11	0.98	0.06	21,24,47,50	1
1	CME	B	88	10/11	0.98	0.06	21,25,35,36	1
1	CME	A	88	10/11	0.99	0.05	21,26,41,44	1
1	CME	B	130	10/11	0.99	0.04	17,20,30,32	1

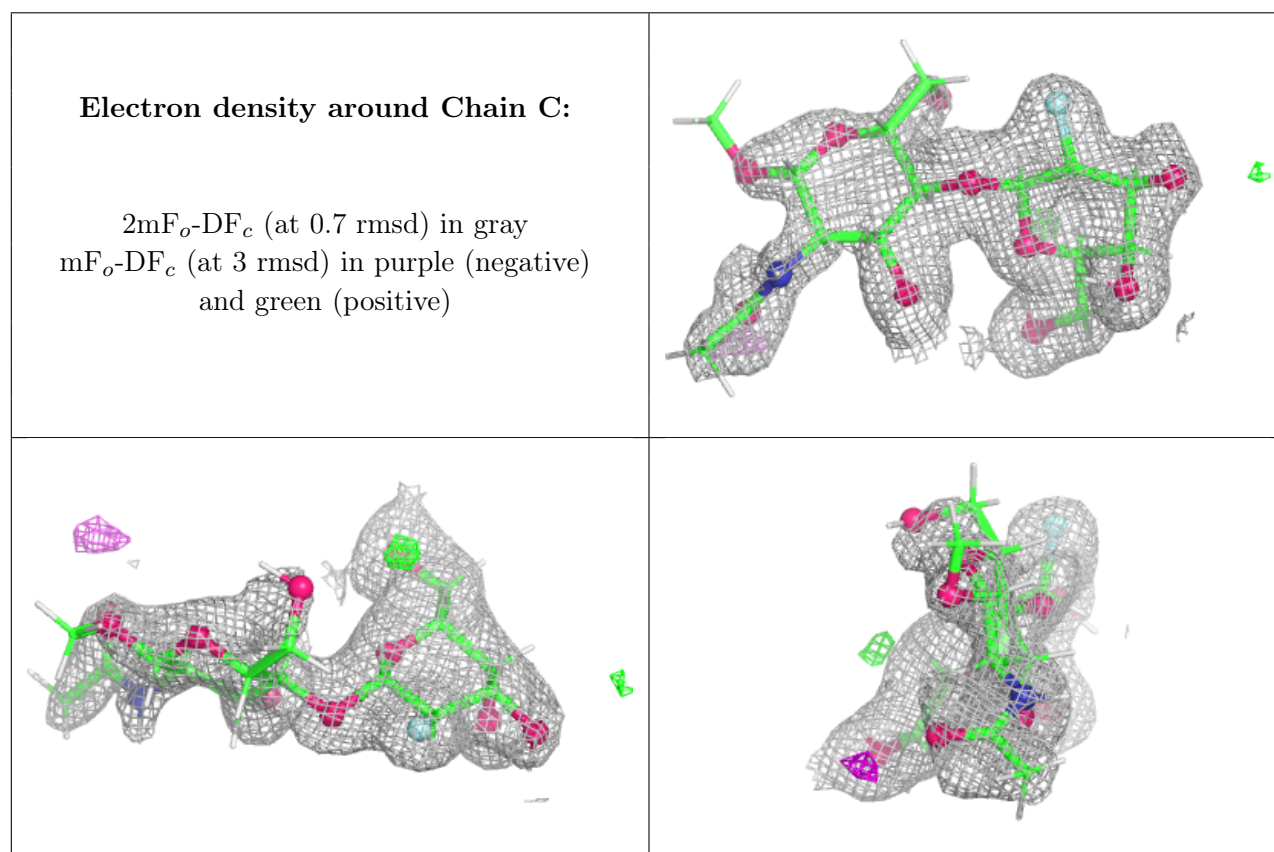


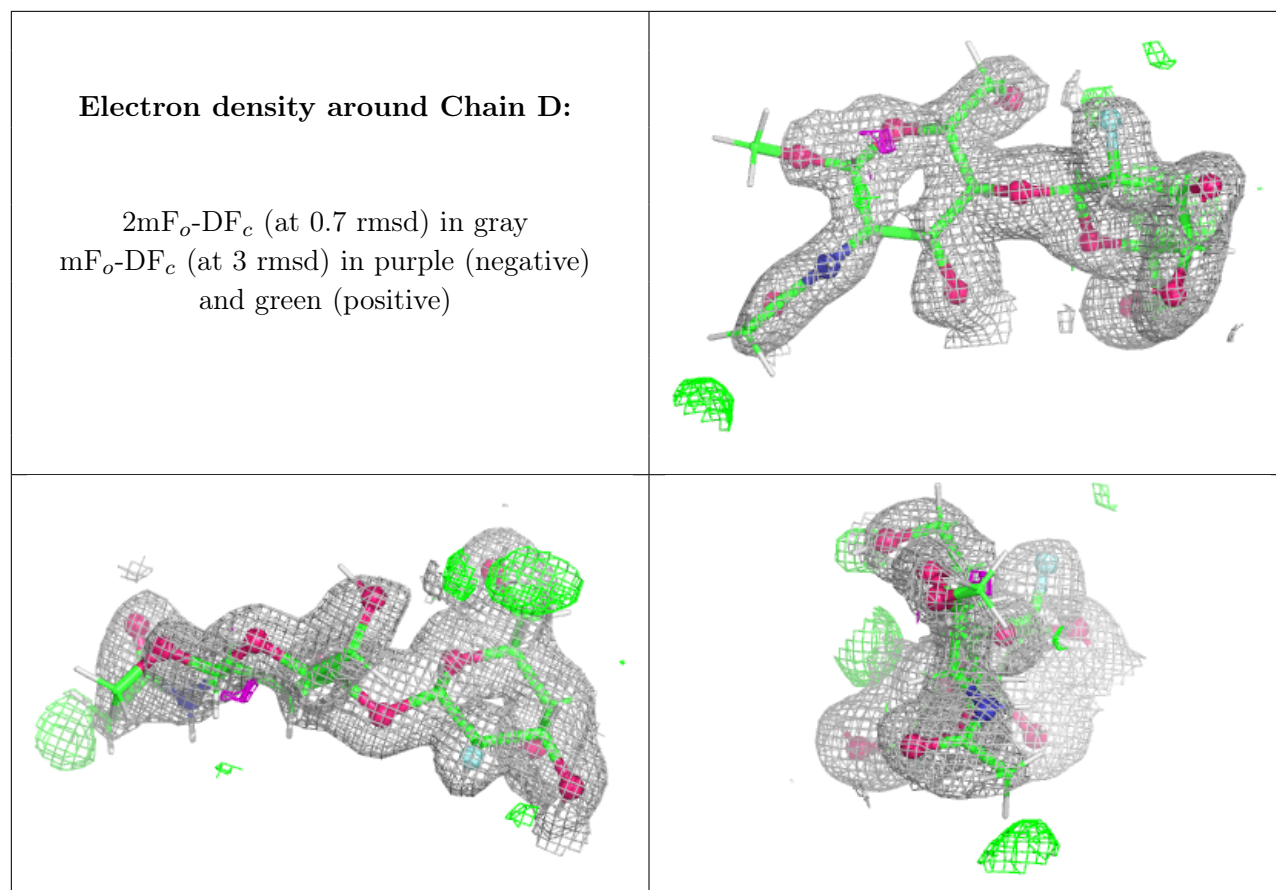
### 6.3 Carbohydrates [i](#)

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( $\text{\AA}^2$ )	Q<0.9
2	MAG	C	1	16/16	0.86	0.12	39,67,75,76	5
2	MAG	D	1	16/16	0.87	0.11	39,54,65,67	5
2	2FG	D	2	11/12	0.93	0.08	32,41,47,55	3
2	2FG	C	2	11/12	0.94	0.07	33,41,53,55	3

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.





## 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
3	EDO	A	201	4/4	0.85	0.11	20,35,39,43	2
3	EDO	B	201	4/4	0.92	0.08	20,34,37,40	2

## 6.5 Other polymers [i](#)

There are no such residues in this entry.