



# wwPDB X-ray Structure Validation Summary Report

Mar 8, 2026 – 03:17 PM UTC

PDB ID : 4UTB / pdb\_00004utb  
Title : Crystal structure of dengue 2 virus envelope glycoprotein in complex with the Fab fragment of the broadly neutralizing human antibody EDE2 A11  
Authors : Rouvinski, A.; Guardado-Calvo, P.; Barba-Spaeth, G.; Duquerroy, S.; Vaney, M.C.; Rey, F.A.  
Deposited on : 2014-07-18  
Resolution : 3.85 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 : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

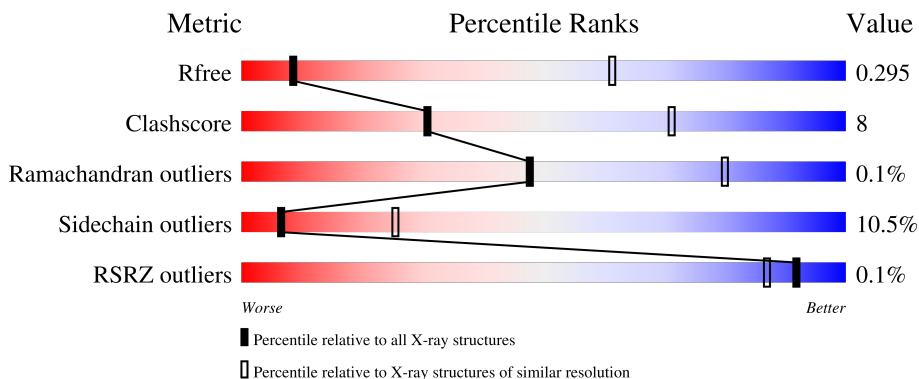
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.85 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.




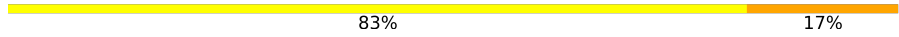


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1165 (4.02-3.70)
Clashscore	190562	1207 (4.02-3.70)
Ramachandran outliers	187476	1149 (4.02-3.70)
Sidechain outliers	187428	1142 (4.02-3.70)
RSRZ outliers	180081	1164 (4.02-3.70)

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	422	
1	B	422	
2	H	283	
2	I	283	
3	L	218	

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Mol	Chain	Length	Quality of chain
3	M	218	 83% 10% . .
4	C	6	 83% 17%
4	E	6	 17% 50% 33%
5	D	2	 50% 50%

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 12714 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 ENVELOPE GLYCOPROTEIN E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	391	3045	1919	524	578	24	0	0	0
1	B	390	3036	1914	522	576	24	0	0	0

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1392	LEU	-	expression tag	UNP Q68Y26
A	1393	ARG	-	expression tag	UNP Q68Y26
A	1394	PRO	-	expression tag	UNP Q68Y26
A	1395	LEU	-	expression tag	UNP Q68Y26
A	1396	GLU	-	expression tag	UNP Q68Y26
A	1397	SER	-	expression tag	UNP Q68Y26
A	1398	ARG	-	expression tag	UNP Q68Y26
A	1399	GLY	-	expression tag	UNP Q68Y26
A	1400	PRO	-	expression tag	UNP Q68Y26
A	1401	PHE	-	expression tag	UNP Q68Y26
A	1402	GLU	-	expression tag	UNP Q68Y26
A	1403	GLY	-	expression tag	UNP Q68Y26
A	1404	LYS	-	expression tag	UNP Q68Y26
A	1405	PRO	-	expression tag	UNP Q68Y26
A	1406	ILE	-	expression tag	UNP Q68Y26
A	1407	PRO	-	expression tag	UNP Q68Y26
A	1408	ASN	-	expression tag	UNP Q68Y26
A	1409	PRO	-	expression tag	UNP Q68Y26
A	1410	LEU	-	expression tag	UNP Q68Y26
A	1411	LEU	-	expression tag	UNP Q68Y26
A	1412	GLY	-	expression tag	UNP Q68Y26
A	1413	LEU	-	expression tag	UNP Q68Y26
A	1414	ASP	-	expression tag	UNP Q68Y26
A	1415	SER	-	expression tag	UNP Q68Y26
A	1416	THR	-	expression tag	UNP Q68Y26

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1417	ARG	-	expression tag	UNP Q68Y26
A	1418	THR	-	expression tag	UNP Q68Y26
A	1419	GLY	-	expression tag	UNP Q68Y26
A	1420	HIS	-	expression tag	UNP Q68Y26
A	1421	HIS	-	expression tag	UNP Q68Y26
A	1422	HIS	-	expression tag	UNP Q68Y26
A	118	LYS	MET	conflict	UNP Q68Y26
B	1392	LEU	-	expression tag	UNP Q68Y26
B	1393	ARG	-	expression tag	UNP Q68Y26
B	1394	PRO	-	expression tag	UNP Q68Y26
B	1395	LEU	-	expression tag	UNP Q68Y26
B	1396	GLU	-	expression tag	UNP Q68Y26
B	1397	SER	-	expression tag	UNP Q68Y26
B	1398	ARG	-	expression tag	UNP Q68Y26
B	1399	GLY	-	expression tag	UNP Q68Y26
B	1400	PRO	-	expression tag	UNP Q68Y26
B	1401	PHE	-	expression tag	UNP Q68Y26
B	1402	GLU	-	expression tag	UNP Q68Y26
B	1403	GLY	-	expression tag	UNP Q68Y26
B	1404	LYS	-	expression tag	UNP Q68Y26
B	1405	PRO	-	expression tag	UNP Q68Y26
B	1406	ILE	-	expression tag	UNP Q68Y26
B	1407	PRO	-	expression tag	UNP Q68Y26
B	1408	ASN	-	expression tag	UNP Q68Y26
B	1409	PRO	-	expression tag	UNP Q68Y26
B	1410	LEU	-	expression tag	UNP Q68Y26
B	1411	LEU	-	expression tag	UNP Q68Y26
B	1412	GLY	-	expression tag	UNP Q68Y26
B	1413	LEU	-	expression tag	UNP Q68Y26
B	1414	ASP	-	expression tag	UNP Q68Y26
B	1415	SER	-	expression tag	UNP Q68Y26
B	1416	THR	-	expression tag	UNP Q68Y26
B	1417	ARG	-	expression tag	UNP Q68Y26
B	1418	THR	-	expression tag	UNP Q68Y26
B	1419	GLY	-	expression tag	UNP Q68Y26
B	1420	HIS	-	expression tag	UNP Q68Y26
B	1421	HIS	-	expression tag	UNP Q68Y26
B	1422	HIS	-	expression tag	UNP Q68Y26
B	118	LYS	MET	conflict	UNP Q68Y26

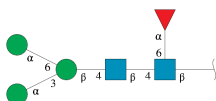
- Molecule 2 is a protein called BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	225	Total 1718	C 1088	N 291	O 332	S 7	0	0	0
2	I	203	Total 1573	C 999	N 267	O 300	S 7	0	0	0

- Molecule 3 is a protein called BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	213	Total 1582	C 987	N 265	O 324	S 6	0	0	0
3	M	210	Total 1563	C 975	N 262	O 320	S 6	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	C	6	Total 71	C 40	N 2	O 29	0	0	0
4	E	6	Total 71	C 40	N 2	O 29	0	0	0

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	D	2	Total 28	C 16	N 2	O 10	0	0	0

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
6	B	1	14	8	1	5	0	0

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
7	H	1	5	4	1	0	0
7	I	1	5	4	1	0	0

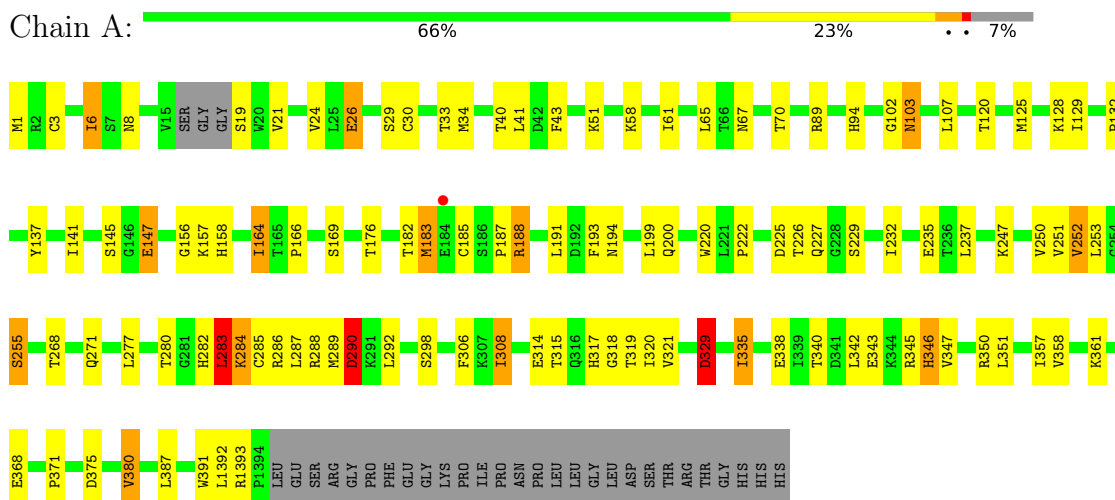
- Molecule 8 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
8	I	2	Total O 2 2	0	0
8	M	1	Total O 1 1	0	0

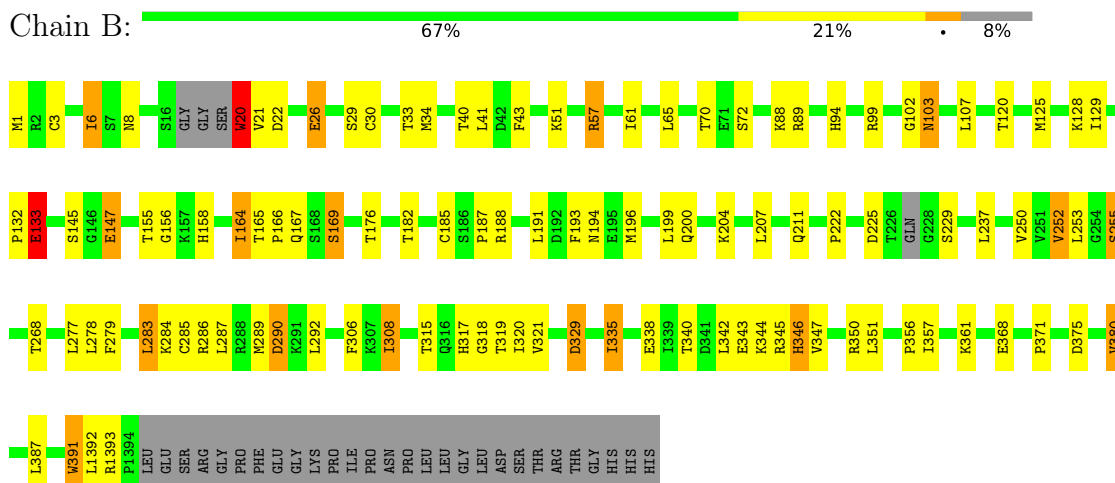
### 3 Residue-property plots

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: ENVELOPE GLYCOPROTEIN E



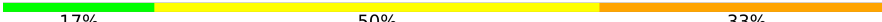
- Molecule 1: ENVELOPE GLYCOPROTEIN E



- Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11





Chain E:  17% 50% 33%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50% 50%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.78Å 181.85Å 204.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.85 20.00 – 3.85	Depositor EDS
% Data completeness (in resolution range)	97.9 (20.00-3.85) 97.0 (20.00-3.85)	Depositor EDS
$R_{merge}$	0.27	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.56 (at 3.86Å)	Xtrriage
Refinement program	BUSTER 2.11.4	Depositor
R, $R_{free}$	0.230 , 0.257 0.278 , 0.295	Depositor DCC
$R_{free}$ test set	1074 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	109.8	Xtrriage
Anisotropy	0.530	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 117.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	12714	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	129.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.23% 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: MAN, SO4, BMA, NAG, FUC

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	0.75	1/3106 (0.0%)	1.28	17/4199 (0.4%)
1	B	0.72	0/3096	1.22	12/4184 (0.3%)
2	H	0.58	0/1764	1.05	6/2406 (0.2%)
2	I	0.62	0/1614	1.09	4/2195 (0.2%)
3	L	0.64	0/1618	1.01	1/2209 (0.0%)
3	M	0.67	0/1598	1.07	4/2180 (0.2%)
All	All	0.68	1/12796 (0.0%)	1.15	44/17373 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	183	MET	SD-CE	-6.76	1.62	1.79

The worst 5 of 44 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	67	ASN	CA-CB-CG	15.62	128.22	112.60
3	M	171	ASN	CA-CB-CG	10.96	123.56	112.60
1	A	103	ASN	N-CA-CB	-7.14	101.40	110.90
1	A	194	ASN	CA-CB-CG	7.13	119.73	112.60
1	B	133	GLU	CB-CG-CD	7.11	124.68	112.60

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	3045	0	3044	68	0
1	B	3036	0	3035	67	0
2	H	1718	0	1644	25	0
2	I	1573	0	1498	35	0
3	L	1582	0	1553	12	0
3	M	1563	0	1534	23	0
4	C	71	0	61	2	0
4	E	71	0	61	2	0
5	D	28	0	25	0	0
6	B	14	0	13	0	0
7	H	5	0	0	0	0
7	I	5	0	0	0	0
8	I	2	0	0	0	0
8	M	1	0	0	0	0
All	All	12714	0	12468	210	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 8.

The worst 5 of 210 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:A:335:ILE:CD1	1:A:358:VAL:HG23	1.64	1.25
1:A:335:ILE:HD11	1:A:358:VAL:CG2	1.70	1.19
1:B:57:ARG:HG3	1:B:57:ARG:HH11	1.17	1.02
1:B:34:MET:HG2	1:B:40:THR:HG22	1.49	0.94
1:A:371:PRO:HG2	1:A:391:TRP:CZ3	2.04	0.92

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	387/422 (92%)	370 (96%)	16 (4%)	1 (0%)	36	69
1	B	384/422 (91%)	367 (96%)	16 (4%)	1 (0%)	36	69
2	H	221/283 (78%)	211 (96%)	10 (4%)	0	100	100
2	I	193/283 (68%)	185 (96%)	8 (4%)	0	100	100
3	L	211/218 (97%)	205 (97%)	6 (3%)	0	100	100
3	M	208/218 (95%)	203 (98%)	5 (2%)	0	100	100
All	All	1604/1846 (87%)	1541 (96%)	61 (4%)	2 (0%)	48	80

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	147	GLU
1	B	147	GLU

### 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	341/366 (93%)	293 (86%)	48 (14%)	3	17
1	B	340/366 (93%)	289 (85%)	51 (15%)	3	16
2	H	190/236 (80%)	176 (93%)	14 (7%)	13	39
2	I	174/236 (74%)	163 (94%)	11 (6%)	16	43
3	L	180/185 (97%)	170 (94%)	10 (6%)	19	45
3	M	178/185 (96%)	165 (93%)	13 (7%)	13	39
All	All	1403/1574 (89%)	1256 (90%)	147 (10%)	6	25

5 of 147 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	I	53	ASP
3	M	171	ASN

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Mol	Chain	Res	Type
2	I	111	VAL
3	L	136	LEU
1	B	6	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
2	I	76	ASN
3	M	171	ASN
2	I	164	HIS
2	H	76	ASN
2	I	39	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

14 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	$\# Z  > 2$
4	NAG	C	1	1,4	14,14,15	0.30	0	17,19,21	0.52	0
4	NAG	C	2	4	14,14,15	0.31	0	17,19,21	1.69	3 (17%)
4	BMA	C	3	4	11,11,12	0.37	0	15,15,17	1.11	1 (6%)
4	MAN	C	4	4	11,11,12	0.40	0	15,15,17	1.12	1 (6%)
4	MAN	C	5	4	11,11,12	0.30	0	15,15,17	0.73	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	FUC	C	6	4	10,10,11	0.47	0	14,14,16	1.06	1 (7%)
5	NAG	D	1	1,5	14,14,15	0.27	0	17,19,21	1.09	2 (11%)
5	NAG	D	2	5	14,14,15	0.37	0	17,19,21	0.74	0
4	NAG	E	1	1,4	14,14,15	0.32	0	17,19,21	0.58	0
4	NAG	E	2	4	14,14,15	0.29	0	17,19,21	1.72	3 (17%)
4	BMA	E	3	4	11,11,12	0.41	0	15,15,17	1.17	1 (6%)
4	MAN	E	4	4	11,11,12	0.40	0	15,15,17	1.06	1 (6%)
4	MAN	E	5	4	11,11,12	0.27	0	15,15,17	0.83	1 (6%)
4	FUC	E	6	4	10,10,11	0.47	0	14,14,16	1.05	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
4	NAG	C	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	C	2	4	-	0/6/23/26	0/1/1/1
4	BMA	C	3	4	-	2/2/19/22	1/1/1/1
4	MAN	C	4	4	-	1/2/19/22	1/1/1/1
4	MAN	C	5	4	-	0/2/19/22	0/1/1/1
4	FUC	C	6	4	-	-	0/1/1/1
5	NAG	D	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	D	2	5	-	0/6/23/26	0/1/1/1
4	NAG	E	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	1/1/1/1
4	MAN	E	4	4	-	1/2/19/22	1/1/1/1
4	MAN	E	5	4	-	0/2/19/22	0/1/1/1
4	FUC	E	6	4	-	-	0/1/1/1

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2	NAG	C1-O5-C5	4.68	118.46	112.19
4	E	2	NAG	C1-O5-C5	4.36	118.03	112.19
4	E	2	NAG	O5-C1-C2	4.21	117.80	111.29
4	C	4	MAN	C1-O5-C5	3.99	117.54	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	4	MAN	C1-O5-C5	3.87	117.38	112.19

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	1	NAG	O5-C5-C6-O6
4	C	3	BMA	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
4	C	3	BMA	C4-C5-C6-O6
4	E	3	BMA	C4-C5-C6-O6

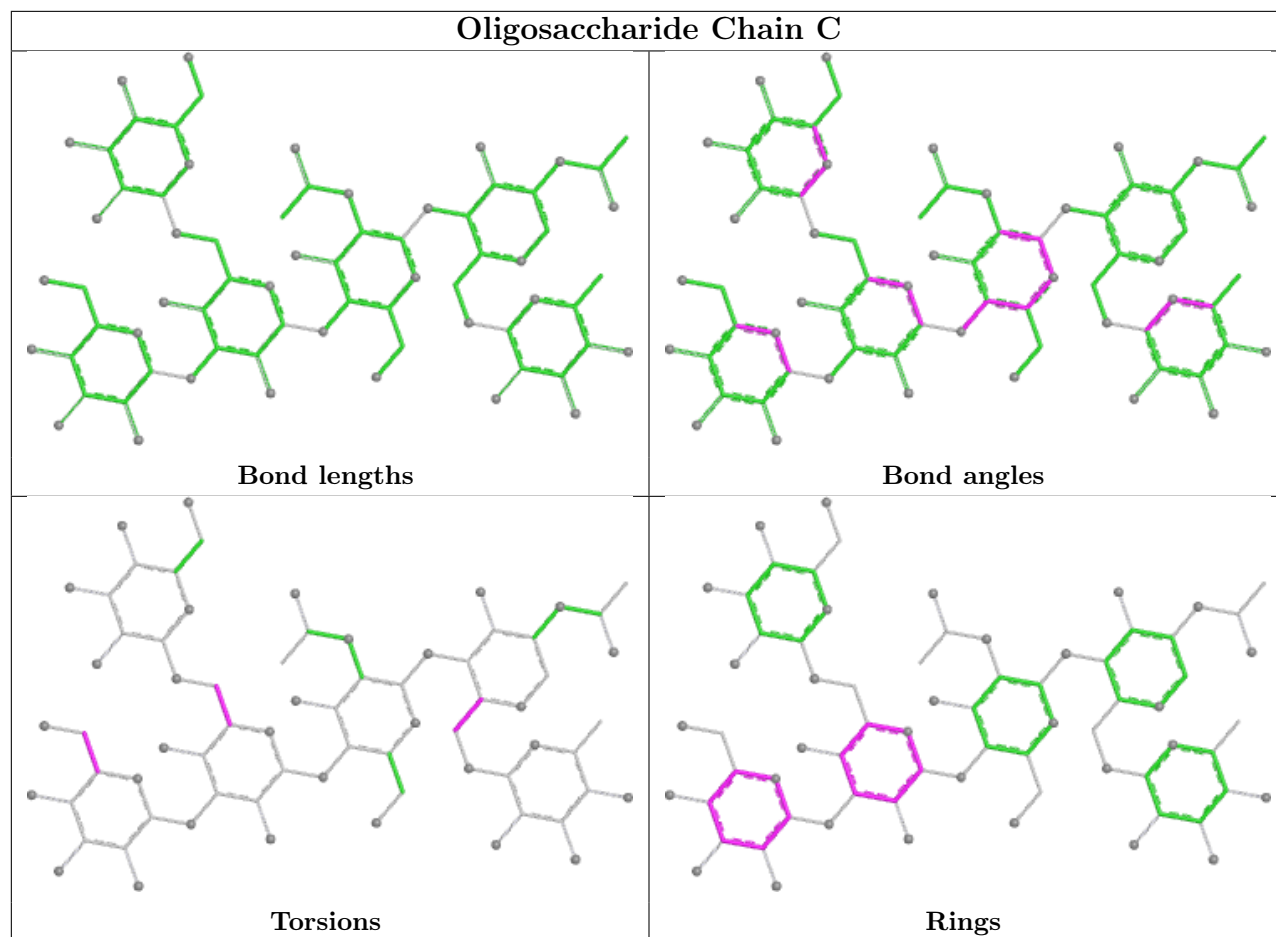
All (4) ring outliers are listed below:

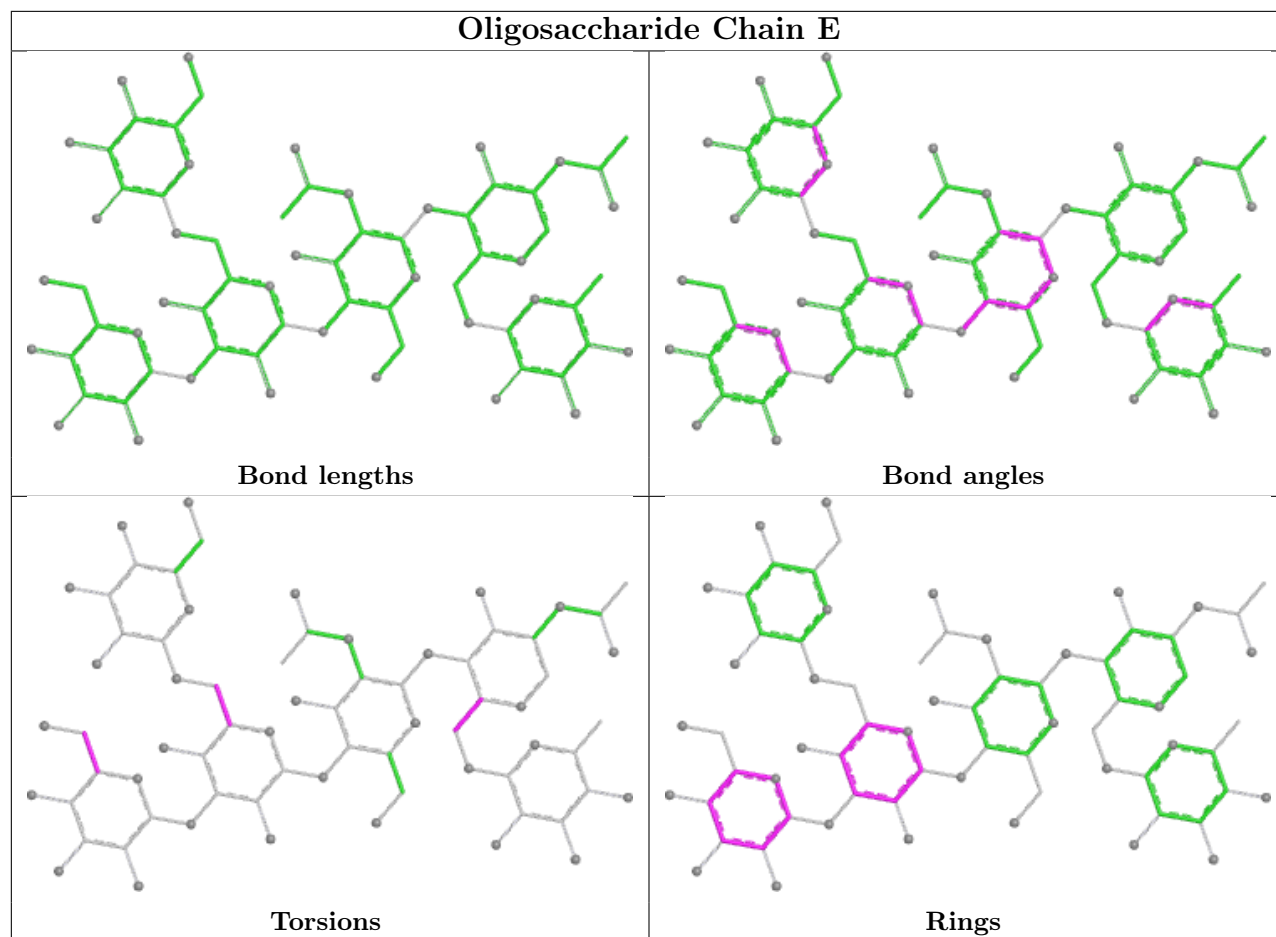
Mol	Chain	Res	Type	Atoms
4	C	3	BMA	C1-C2-C3-C4-C5-O5
4	E	3	BMA	C1-C2-C3-C4-C5-O5
4	E	4	MAN	C1-C2-C3-C4-C5-O5
4	C	4	MAN	C1-C2-C3-C4-C5-O5

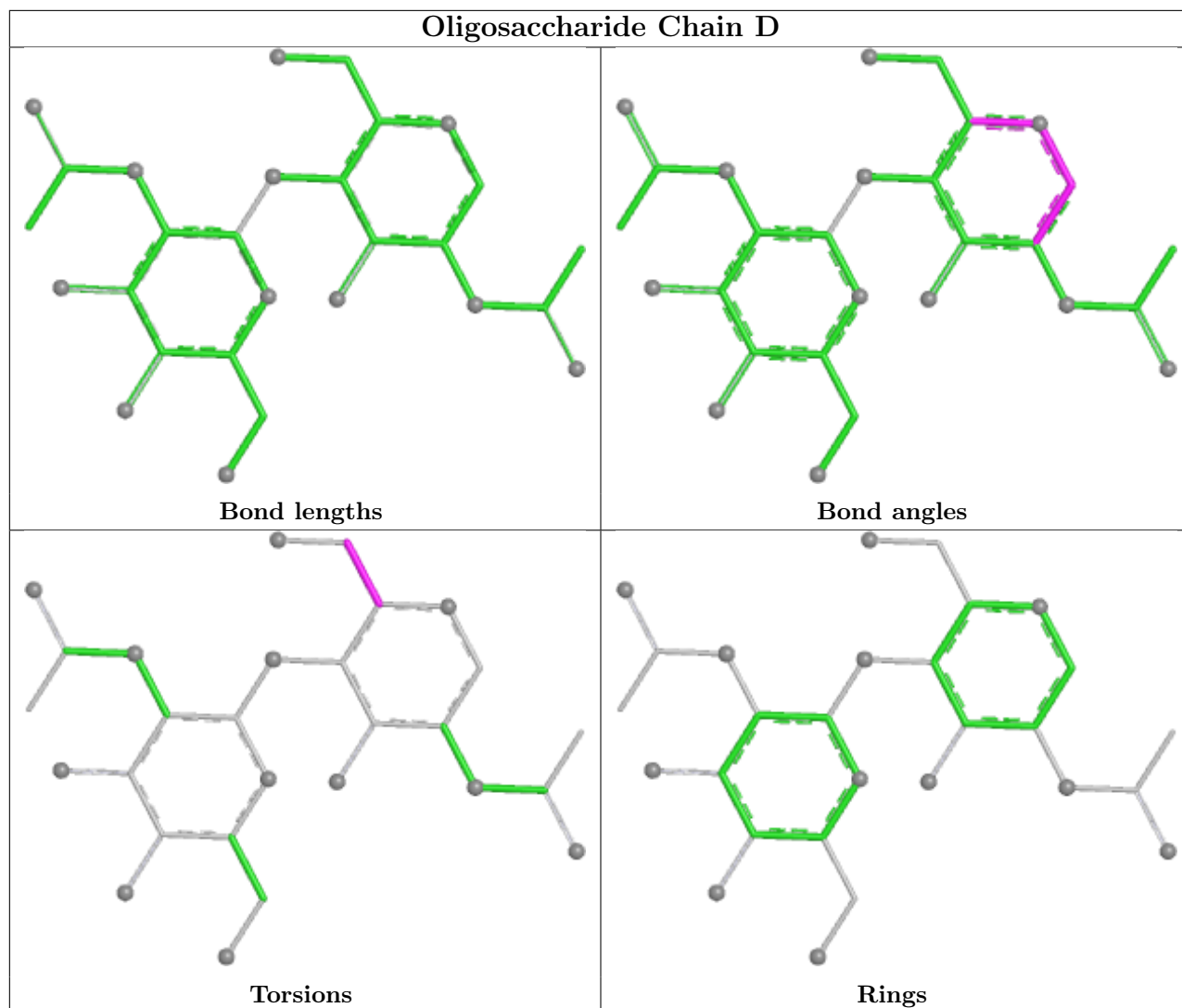
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1	NAG	1	0
4	E	2	NAG	1	0
4	C	2	NAG	1	0
4	E	3	BMA	1	0

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







## 5.6 Ligand geometry [i](#)

3 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	$\# Z  > 2$
7	SO4	H	581	-	4,4,4	0.33	0	6,6,6	0.19	0
7	SO4	I	581	-	4,4,4	0.34	0	6,6,6	0.15	0
6	NAG	B	567	1	14,14,15	0.39	0	17,19,21	0.66	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	NAG	B	567	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	391/422 (92%)	-0.05	1 (0%) 90 77	100, 125, 158, 180	0
1	B	390/422 (92%)	-0.16	0 100 100	94, 125, 158, 177	0
2	H	225/283 (79%)	-0.20	0 100 100	94, 126, 146, 162	0
2	I	203/283 (71%)	-0.13	0 100 100	91, 122, 187, 206	0
3	L	213/218 (97%)	-0.26	0 100 100	90, 126, 160, 184	0
3	M	210/218 (96%)	-0.19	0 100 100	92, 128, 178, 193	0
All	All	1632/1846 (88%)	-0.15	1 (0%) 92 87	90, 125, 172, 206	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	184	GLU	2.1

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

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

### 6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

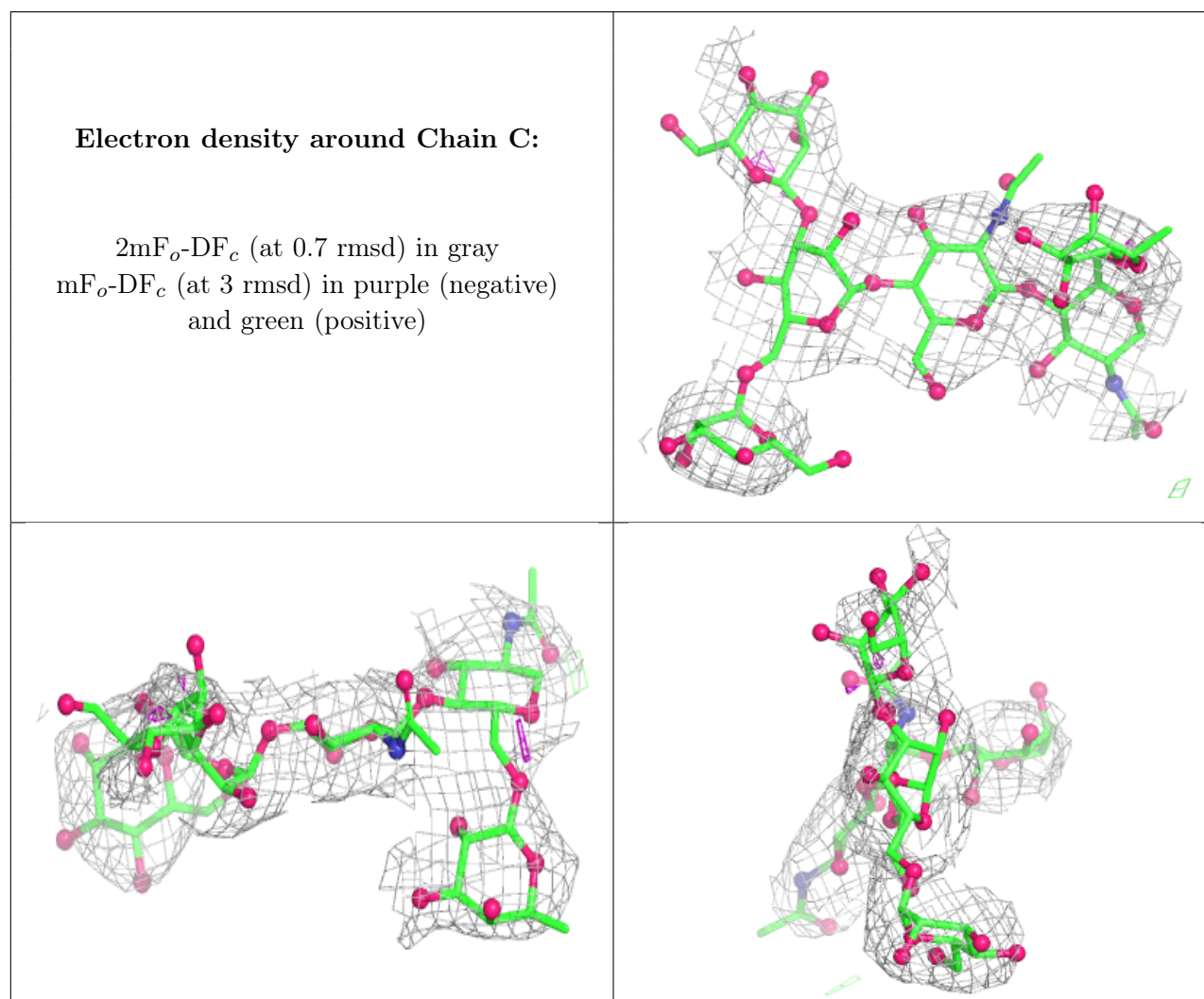
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	NAG	C	1	14/15	-	-	110,112,124,128	0
4	NAG	C	2	14/15	-	-	117,121,130,131	0
4	BMA	C	3	11/12	-	-	125,127,135,144	0

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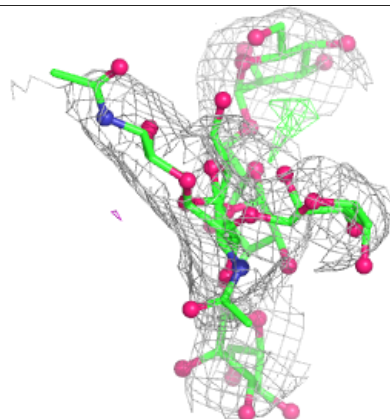
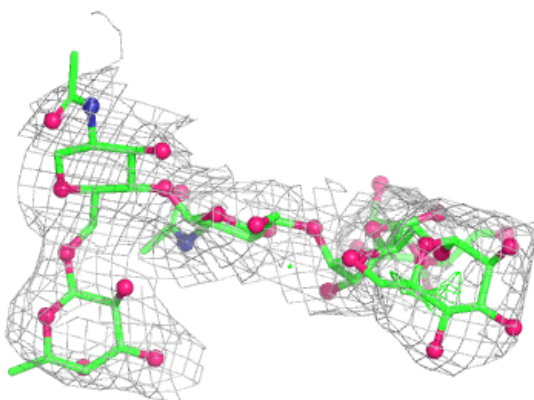
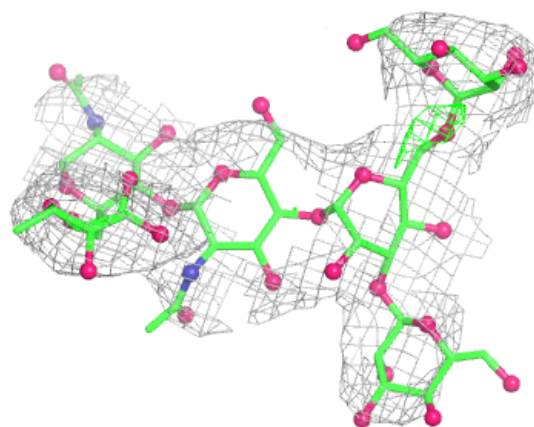
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	MAN	C	4	11/12	-	-	151,155,160,161	0
4	MAN	C	5	11/12	-	-	125,127,130,131	0
4	FUC	C	6	10/11	-	-	130,133,135,137	0
4	NAG	E	1	14/15	-	-	102,108,118,124	0
4	NAG	E	2	14/15	-	-	112,115,118,125	0
4	BMA	E	3	11/12	-	-	132,133,143,151	0
4	MAN	E	4	11/12	-	-	157,161,163,163	0
4	MAN	E	5	11/12	-	-	133,136,137,138	0
4	FUC	E	6	10/11	-	-	129,133,135,137	0
5	NAG	D	1	14/15	-	-	169,172,173,174	0
5	NAG	D	2	14/15	-	-	169,174,177,177	0

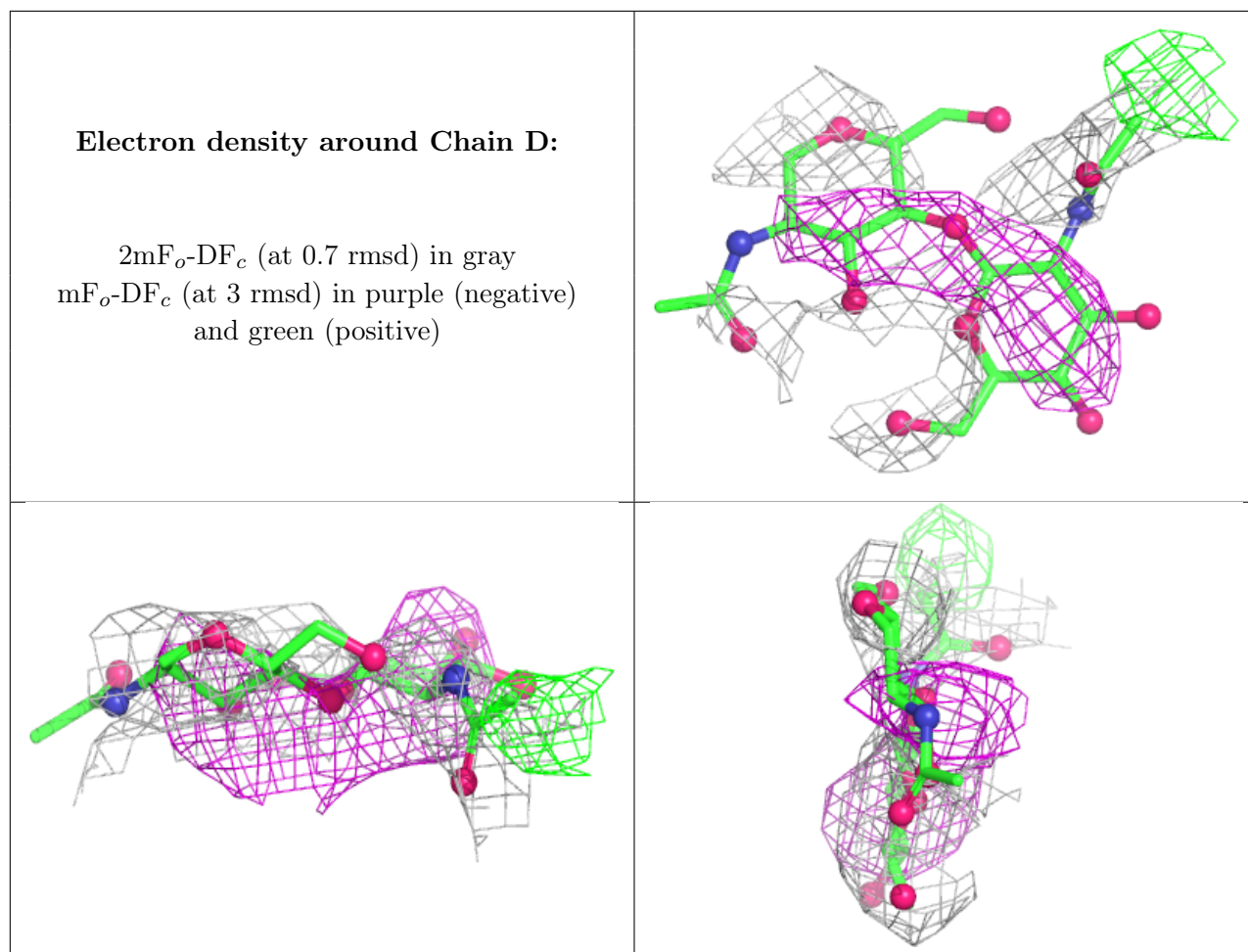
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



**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)





## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
6	NAG	B	567	14/15	0.61	0.10	154,156,158,158	0
7	SO4	I	581	5/5	0.76	0.08	169,169,170,170	0
7	SO4	H	581	5/5	0.83	0.09	144,144,145,146	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.