



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 7, 2024 – 09:57 pm GMT

PDB ID : 5N0A
Title : Crystal structure of A259C covalently linked dengue 2 virus envelope glycoprotein dimer in complex with the Fab fragment of the broadly neutralizing human antibody EDE2 A11
Authors : Vaney, M.C.; Rouvinski, A.; Guardado-Calvo, P.; Sharma, A.; Rey, F.A.
Deposited on : 2017-02-02
Resolution : 3.90 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

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>.

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

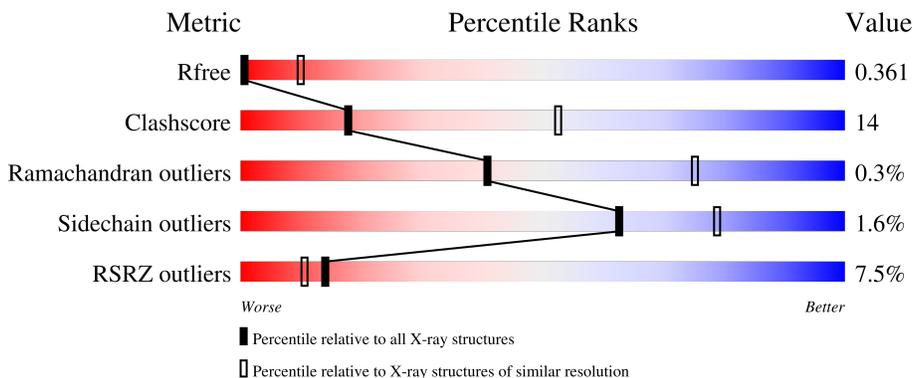
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 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.



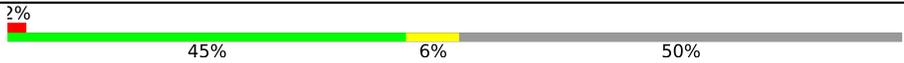
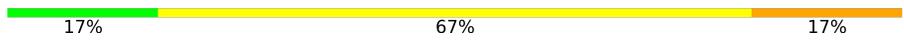
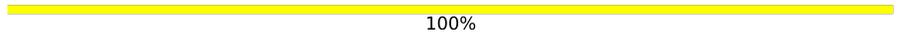
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1002 (4.14-3.66)
Clashscore	141614	1004 (4.12-3.68)
Ramachandran outliers	138981	1021 (4.14-3.66)
Sidechain outliers	138945	1014 (4.14-3.66)
RSRZ outliers	127900	1275 (4.20-3.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	430	
1	B	430	
2	H	283	
2	I	283	
3	L	218	

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Mol	Chain	Length	Quality of chain
3	M	218	
4	C	6	
4	D	6	

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
4	MAN	C	4	-	-	-	X
4	MAN	D	4	-	-	-	X
4	FUC	D	6	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9683 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	373	2914	1842	497	551	24	0	0	0
1	B	379	2943	1859	501	558	25	0	0	0

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	118	LYS	MET	conflict	UNP Q68Y26
A	259	CYS	ALA	engineered mutation	UNP Q68Y26
A	396	LEU	-	expression tag	UNP Q68Y26
A	397	VAL	-	expression tag	UNP Q68Y26
A	398	PRO	-	expression tag	UNP Q68Y26
A	399	ARG	-	expression tag	UNP Q68Y26
A	400	GLY	-	expression tag	UNP Q68Y26
A	401	SER	-	expression tag	UNP Q68Y26
A	402	SER	-	expression tag	UNP Q68Y26
A	403	ALA	-	expression tag	UNP Q68Y26
A	404	TRP	-	expression tag	UNP Q68Y26
A	405	SER	-	expression tag	UNP Q68Y26
A	406	HIS	-	expression tag	UNP Q68Y26
A	407	PRO	-	expression tag	UNP Q68Y26
A	408	GLN	-	expression tag	UNP Q68Y26
A	409	PHE	-	expression tag	UNP Q68Y26
A	410	GLU	-	expression tag	UNP Q68Y26
A	411	LYS	-	expression tag	UNP Q68Y26
A	412	GLY	-	expression tag	UNP Q68Y26
A	413	GLY	-	expression tag	UNP Q68Y26
A	414	SER	-	expression tag	UNP Q68Y26
A	415	GLY	-	expression tag	UNP Q68Y26
A	416	GLY	-	expression tag	UNP Q68Y26
A	417	GLY	-	expression tag	UNP Q68Y26
A	418	SER	-	expression tag	UNP Q68Y26

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Chain	Residue	Modelled	Actual	Comment	Reference
A	419	GLY	-	expression tag	UNP Q68Y26
A	420	GLY	-	expression tag	UNP Q68Y26
A	421	SER	-	expression tag	UNP Q68Y26
A	422	ALA	-	expression tag	UNP Q68Y26
A	423	TRP	-	expression tag	UNP Q68Y26
A	424	SER	-	expression tag	UNP Q68Y26
A	425	HIS	-	expression tag	UNP Q68Y26
A	426	PRO	-	expression tag	UNP Q68Y26
A	427	GLN	-	expression tag	UNP Q68Y26
A	428	PHE	-	expression tag	UNP Q68Y26
A	429	GLU	-	expression tag	UNP Q68Y26
A	430	LYS	-	expression tag	UNP Q68Y26
B	118	LYS	MET	conflict	UNP Q68Y26
B	259	CYS	ALA	engineered mutation	UNP Q68Y26
B	396	LEU	-	expression tag	UNP Q68Y26
B	397	VAL	-	expression tag	UNP Q68Y26
B	398	PRO	-	expression tag	UNP Q68Y26
B	399	ARG	-	expression tag	UNP Q68Y26
B	400	GLY	-	expression tag	UNP Q68Y26
B	401	SER	-	expression tag	UNP Q68Y26
B	402	SER	-	expression tag	UNP Q68Y26
B	403	ALA	-	expression tag	UNP Q68Y26
B	404	TRP	-	expression tag	UNP Q68Y26
B	405	SER	-	expression tag	UNP Q68Y26
B	406	HIS	-	expression tag	UNP Q68Y26
B	407	PRO	-	expression tag	UNP Q68Y26
B	408	GLN	-	expression tag	UNP Q68Y26
B	409	PHE	-	expression tag	UNP Q68Y26
B	410	GLU	-	expression tag	UNP Q68Y26
B	411	LYS	-	expression tag	UNP Q68Y26
B	412	GLY	-	expression tag	UNP Q68Y26
B	413	GLY	-	expression tag	UNP Q68Y26
B	414	SER	-	expression tag	UNP Q68Y26
B	415	GLY	-	expression tag	UNP Q68Y26
B	416	GLY	-	expression tag	UNP Q68Y26
B	417	GLY	-	expression tag	UNP Q68Y26
B	418	SER	-	expression tag	UNP Q68Y26
B	419	GLY	-	expression tag	UNP Q68Y26
B	420	GLY	-	expression tag	UNP Q68Y26
B	421	SER	-	expression tag	UNP Q68Y26
B	422	ALA	-	expression tag	UNP Q68Y26
B	423	TRP	-	expression tag	UNP Q68Y26

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Chain	Residue	Modelled	Actual	Comment	Reference
B	424	SER	-	expression tag	UNP Q68Y26
B	425	HIS	-	expression tag	UNP Q68Y26
B	426	PRO	-	expression tag	UNP Q68Y26
B	427	GLN	-	expression tag	UNP Q68Y26
B	428	PHE	-	expression tag	UNP Q68Y26
B	429	GLU	-	expression tag	UNP Q68Y26
B	430	LYS	-	expression tag	UNP Q68Y26

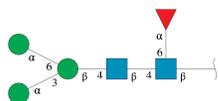
- Molecule 2 is a protein called BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11 HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	131	1032	652	178	197	5	0	0	0
2	I	131	1032	652	178	197	5	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	110	810	502	137	167	4	0	0	0
3	M	110	810	502	137	167	4	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.

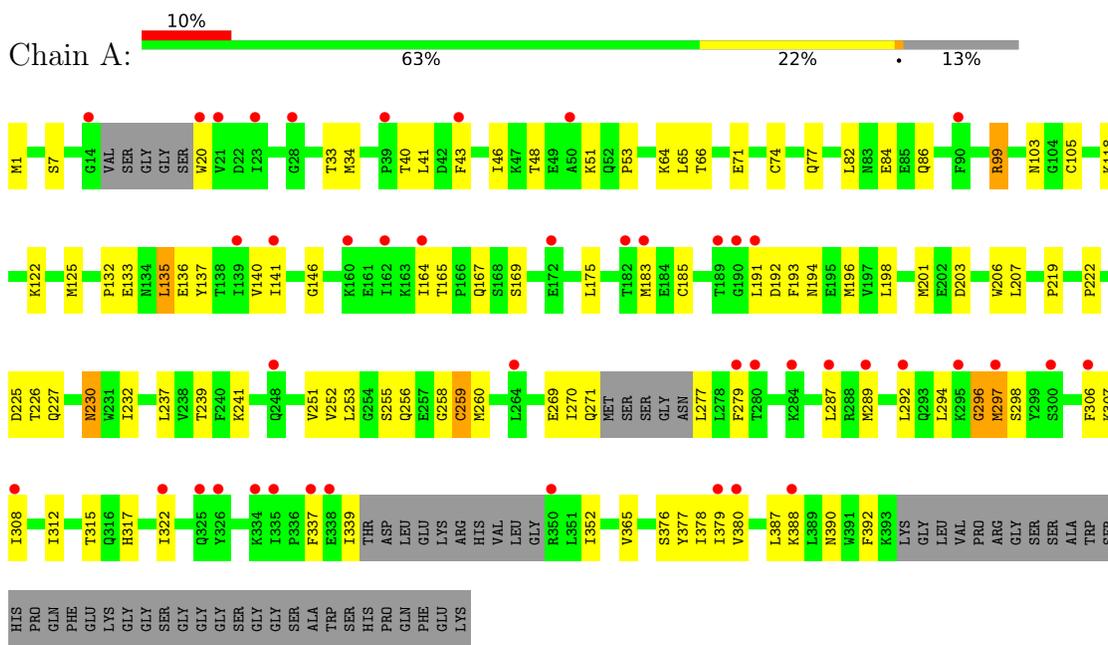


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	C	6	71	40	2	29	0	0	0
4	D	6	71	40	2	29	0	0	0

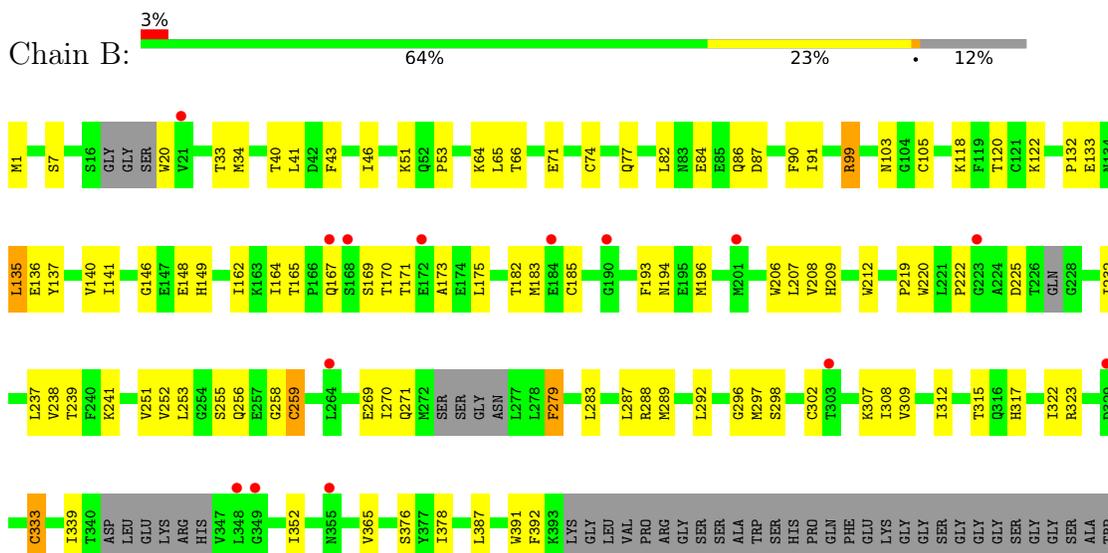
3 Residue-property plots [i](#)

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: Envelope Glycoprotein E

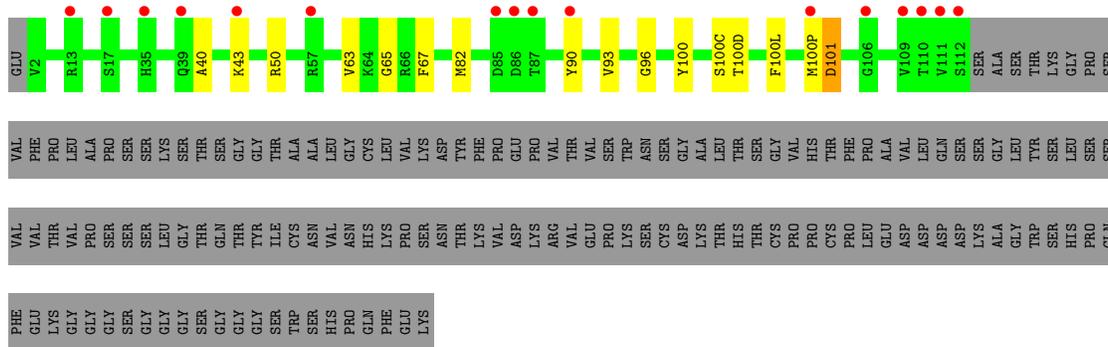
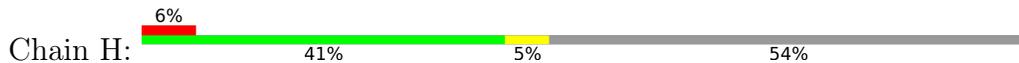


- Molecule 1: Envelope Glycoprotein E

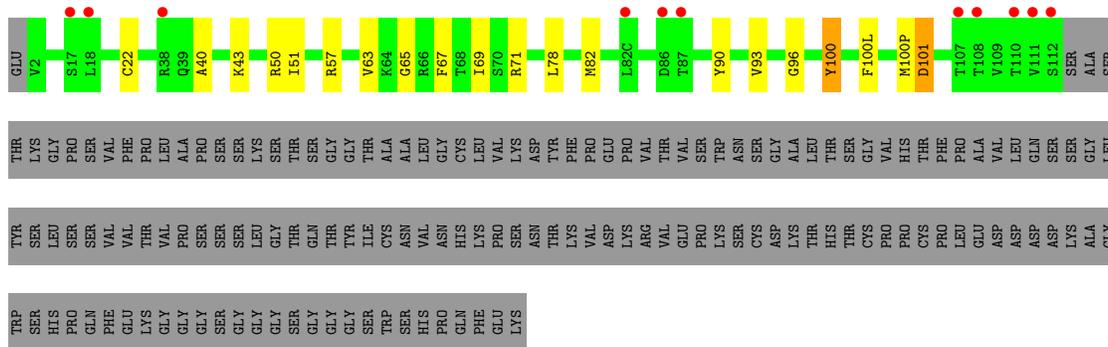
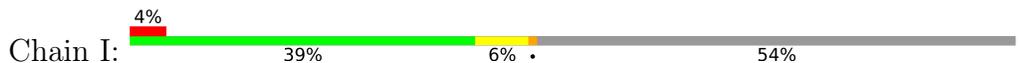


SER
HIS
PRO
GLN
PHE
GLU
LYS

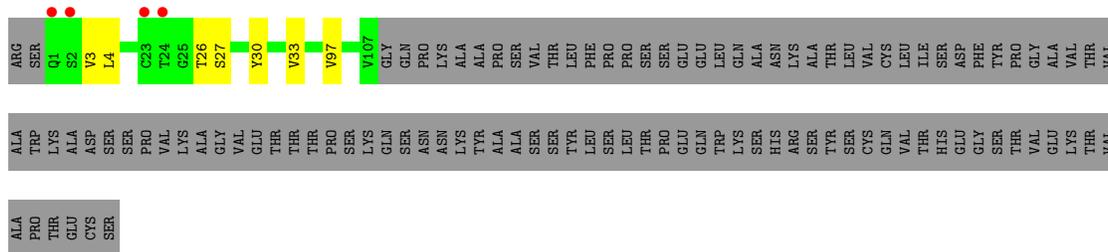
● Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11 HEAVY CHAIN



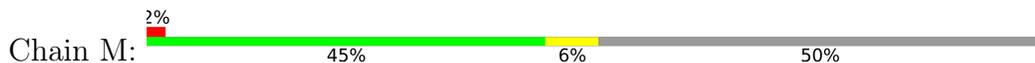
● Molecule 2: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11 HEAVY CHAIN



● Molecule 3: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11



● Molecule 3: BROADLY NEUTRALIZING HUMAN ANTIBODY EDE2 A11



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.77Å 182.30Å 208.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.90 49.40 – 3.90	Depositor EDS
% Data completeness (in resolution range)	90.7 (20.00-3.90) 90.7 (49.40-3.90)	Depositor EDS
R_{merge}	0.32	Depositor
R_{sym}	0.32	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 3.88Å)	Xtrriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.285 , 0.300 0.325 , 0.361	Depositor DCC
R_{free} test set	955 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	78.7	Xtrriage
Anisotropy	1.122	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 110.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	9683	wwPDB-VP
Average B, all atoms (Å ²)	143.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: FUC, MAN, BMA, NAG

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	0.37	0/2972	0.70	2/4014 (0.0%)
1	B	0.37	0/2998	0.68	0/4047
2	H	0.34	0/1060	0.66	1/1439 (0.1%)
2	I	0.35	0/1060	0.65	1/1439 (0.1%)
3	L	0.36	0/825	0.66	0/1122
3	M	0.36	0/825	0.64	0/1122
All	All	0.36	0/9740	0.67	4/13183 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	296	GLY	C-N-CA	5.36	135.09	121.70
2	I	101	ASP	CB-CG-OD2	5.23	123.01	118.30
2	H	101	ASP	CB-CG-OD2	5.23	123.00	118.30
1	A	203	ASP	CB-CG-OD1	5.14	122.93	118.30

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	2914	0	2907	98	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2943	0	2943	103	0
2	H	1032	0	963	12	0
2	I	1032	0	963	41	0
3	L	810	0	798	5	0
3	M	810	0	798	11	0
4	C	71	0	61	1	0
4	D	71	0	61	2	0
All	All	9683	0	9494	260	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 14.

The worst 5 of 260 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:B:46:ILE:HD12	1:B:140:VAL:CG2	1.28	1.56
1:B:46:ILE:CD1	1:B:140:VAL:CG2	1.85	1.51
2:I:50:ARG:NH1	2:I:100(L):PHE:HE2	1.28	1.31
1:B:46:ILE:CD1	1:B:140:VAL:HG21	1.52	1.24
2:I:67:PHE:CE1	2:I:82:MET:HB3	1.72	1.22

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	365/430 (85%)	346 (95%)	18 (5%)	1 (0%)	41 75
1	B	369/430 (86%)	351 (95%)	17 (5%)	1 (0%)	41 75
2	H	129/283 (46%)	124 (96%)	4 (3%)	1 (1%)	19 57
2	I	129/283 (46%)	124 (96%)	4 (3%)	1 (1%)	19 57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	L	108/218 (50%)	104 (96%)	4 (4%)	0	100	100
3	M	108/218 (50%)	104 (96%)	4 (4%)	0	100	100
All	All	1208/1862 (65%)	1153 (95%)	51 (4%)	4 (0%)	41	75

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	297	MET
1	B	297	MET
2	I	65	GLY
2	H	65	GLY

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	326/368 (89%)	319 (98%)	7 (2%)	53	73
1	B	330/368 (90%)	322 (98%)	8 (2%)	49	69
2	H	109/236 (46%)	108 (99%)	1 (1%)	78	87
2	I	109/236 (46%)	108 (99%)	1 (1%)	78	87
3	L	92/185 (50%)	92 (100%)	0	100	100
3	M	92/185 (50%)	92 (100%)	0	100	100
All	All	1058/1578 (67%)	1041 (98%)	17 (2%)	62	79

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	333	CYS
2	I	100	TYR
1	B	1	MET
1	B	86	GLN
1	B	99	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	386	GLN
3	L	37	GLN
3	M	37	GLN
1	A	271	GLN
1	B	134	ASN

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](#)

12 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	4,1	14,14,15	0.29	0	17,19,21	0.56	0
4	NAG	C	2	4	14,14,15	0.34	0	17,19,21	1.33	3 (17%)
4	BMA	C	3	4	11,11,12	0.40	0	15,15,17	1.12	1 (6%)
4	MAN	C	4	4	11,11,12	0.38	0	15,15,17	0.98	1 (6%)
4	MAN	C	5	4	11,11,12	0.36	0	15,15,17	0.89	1 (6%)
4	FUC	C	6	4	10,10,11	0.44	0	14,14,16	0.93	1 (7%)
4	NAG	D	1	4,1	14,14,15	0.30	0	17,19,21	0.58	0
4	NAG	D	2	4	14,14,15	0.35	0	17,19,21	1.31	3 (17%)
4	BMA	D	3	4	11,11,12	0.36	0	15,15,17	1.16	1 (6%)
4	MAN	D	4	4	11,11,12	0.39	0	15,15,17	0.98	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	D	5	4	11,11,12	0.37	0	15,15,17	0.84	1 (6%)
4	FUC	D	6	4	10,10,11	0.44	0	14,14,16	0.94	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	4,1	-	0/6/23/26	0/1/1/1
4	NAG	C	2	4	-	2/6/23/26	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
4	NAG	D	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
4	BMA	D	3	4	-	2/2/19/22	1/1/1/1
4	MAN	D	4	4	-	1/2/19/22	1/1/1/1
4	MAN	D	5	4	-	0/2/19/22	0/1/1/1
4	FUC	D	6	4	-	-	0/1/1/1

There are no bond length outliers.

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	3	BMA	C1-O5-C5	3.92	117.51	112.19
4	C	3	BMA	C1-O5-C5	3.69	117.19	112.19
4	D	4	MAN	C1-O5-C5	3.52	116.96	112.19
4	C	4	MAN	C1-O5-C5	3.50	116.94	112.19
4	D	2	NAG	C1-O5-C5	3.47	116.89	112.19

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	3	BMA	O5-C5-C6-O6
4	D	3	BMA	O5-C5-C6-O6
4	C	3	BMA	C4-C5-C6-O6
4	D	3	BMA	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	D	2	NAG	C4-C5-C6-O6

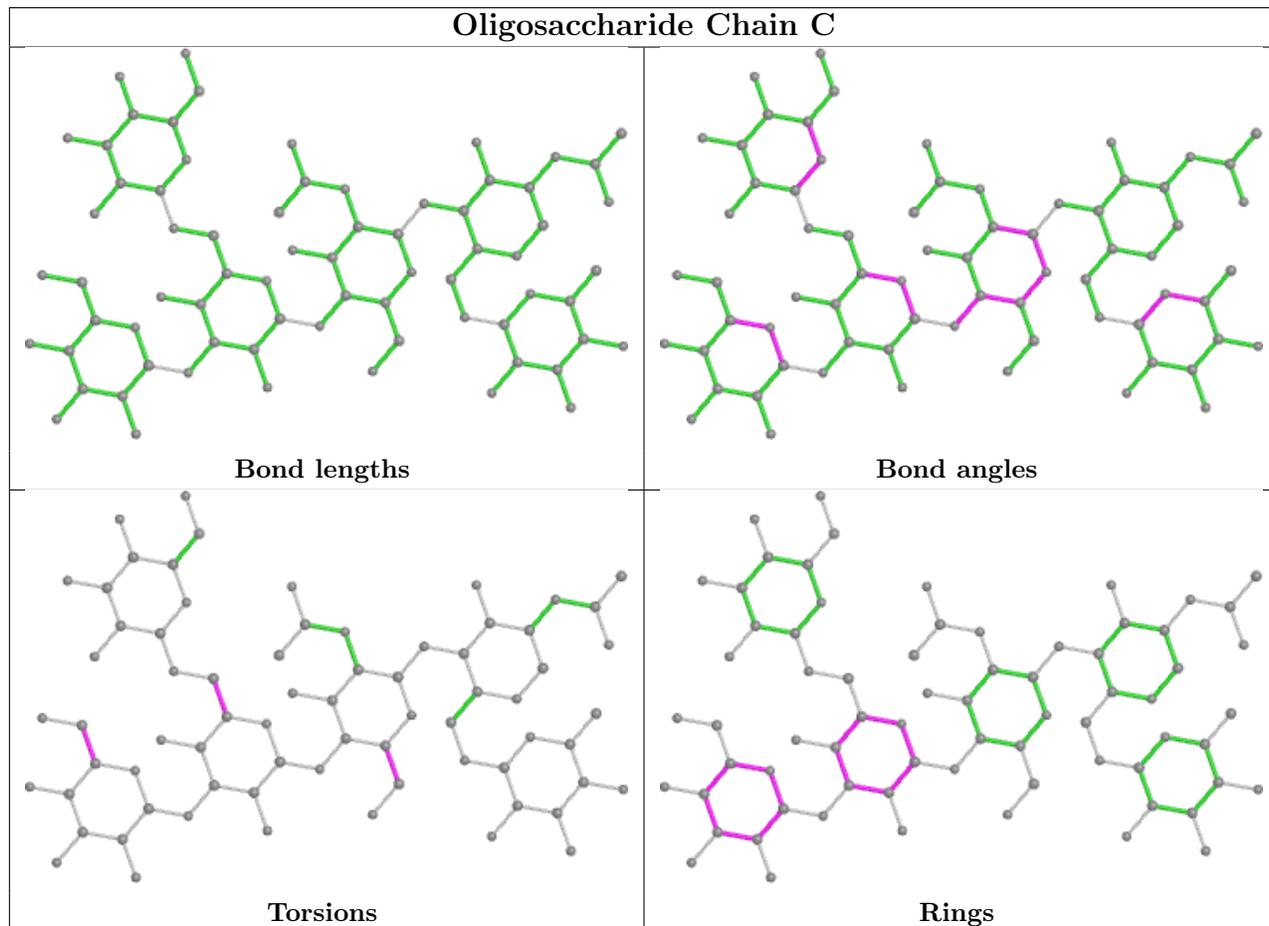
All (4) ring outliers are listed below:

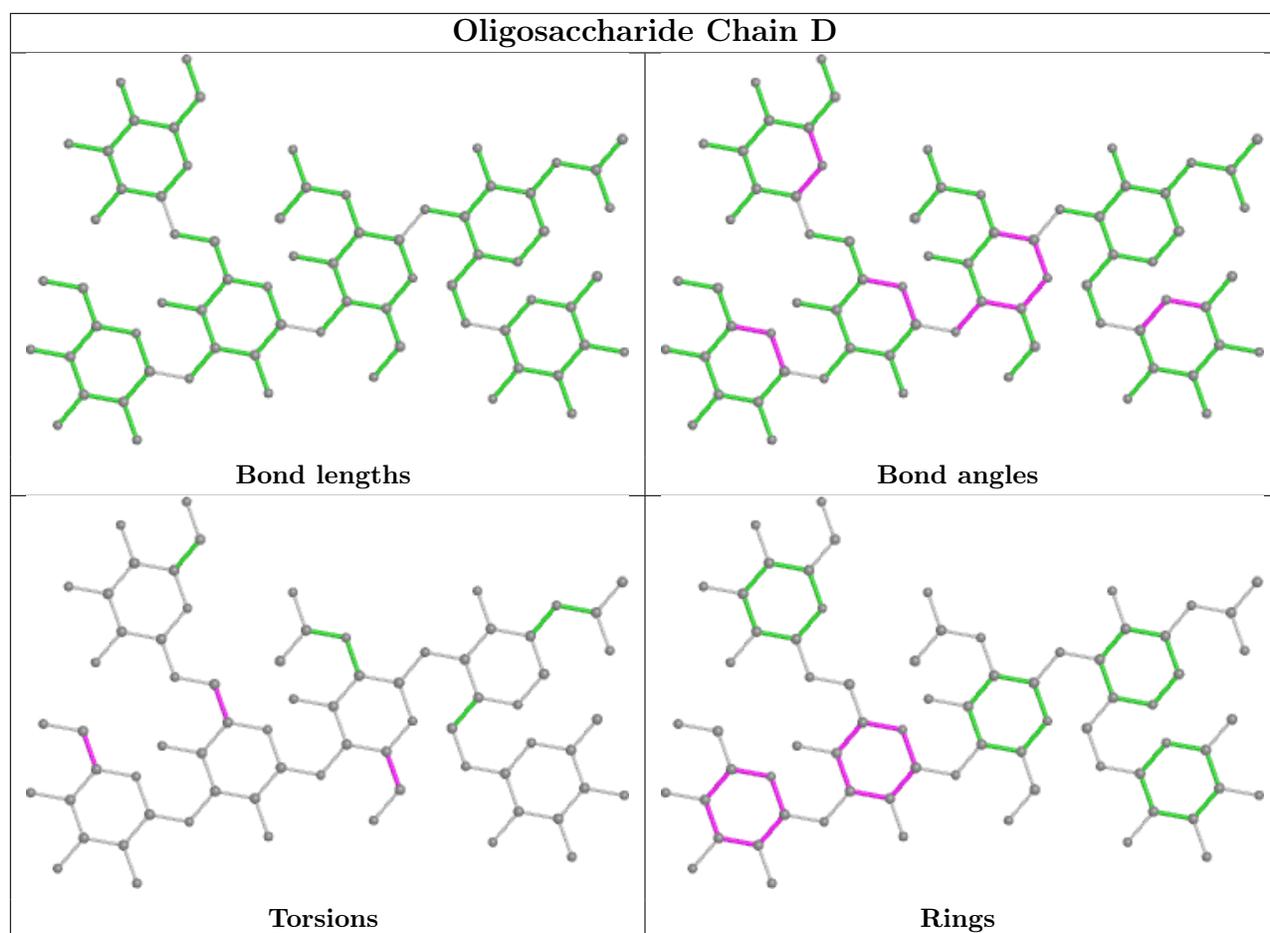
Mol	Chain	Res	Type	Atoms
4	C	3	BMA	C1-C2-C3-C4-C5-O5
4	D	3	BMA	C1-C2-C3-C4-C5-O5
4	C	4	MAN	C1-C2-C3-C4-C5-O5
4	D	4	MAN	C1-C2-C3-C4-C5-O5

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	2	NAG	1	0
4	D	1	NAG	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](#)

There are no ligands in this entry.

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	373/430 (86%)	0.89	44 (11%) 4 4	81, 143, 236, 287	0
1	B	379/430 (88%)	0.58	14 (3%) 41 32	68, 142, 233, 287	0
2	H	131/283 (46%)	0.80	16 (12%) 4 4	89, 133, 218, 295	0
2	I	131/283 (46%)	0.72	11 (8%) 11 9	91, 130, 189, 229	0
3	L	110/218 (50%)	0.30	4 (3%) 42 33	83, 121, 163, 204	0
3	M	110/218 (50%)	0.38	4 (3%) 42 33	83, 128, 166, 235	0
All	All	1234/1862 (66%)	0.67	93 (7%) 14 10	68, 138, 218, 295	0

The worst 5 of 93 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	297	MET	5.6
1	A	20	TRP	4.4
1	A	28	GLY	4.3
1	A	334	LYS	4.0
1	A	50	ALA	4.0

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

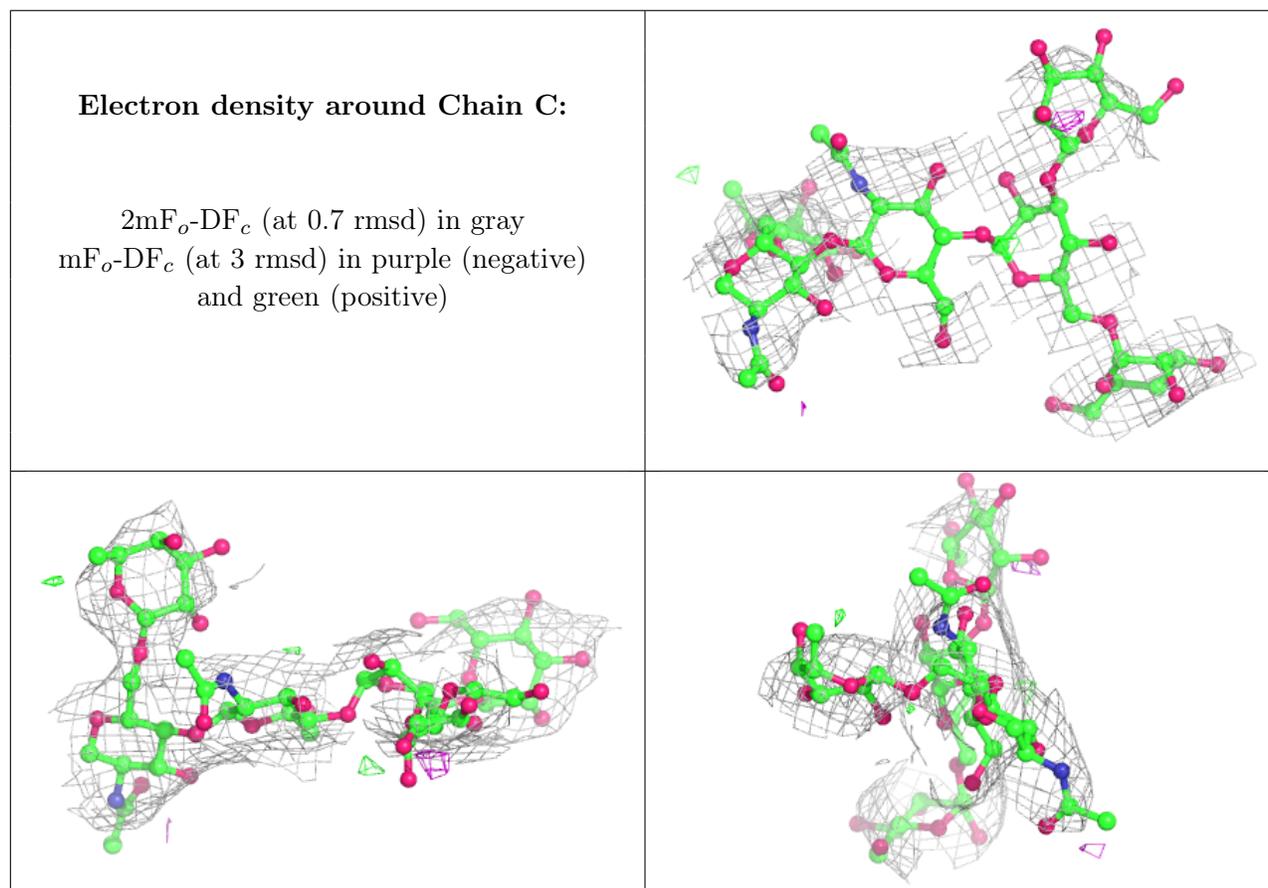
There are no non-standard protein/DNA/RNA residues in this entry.

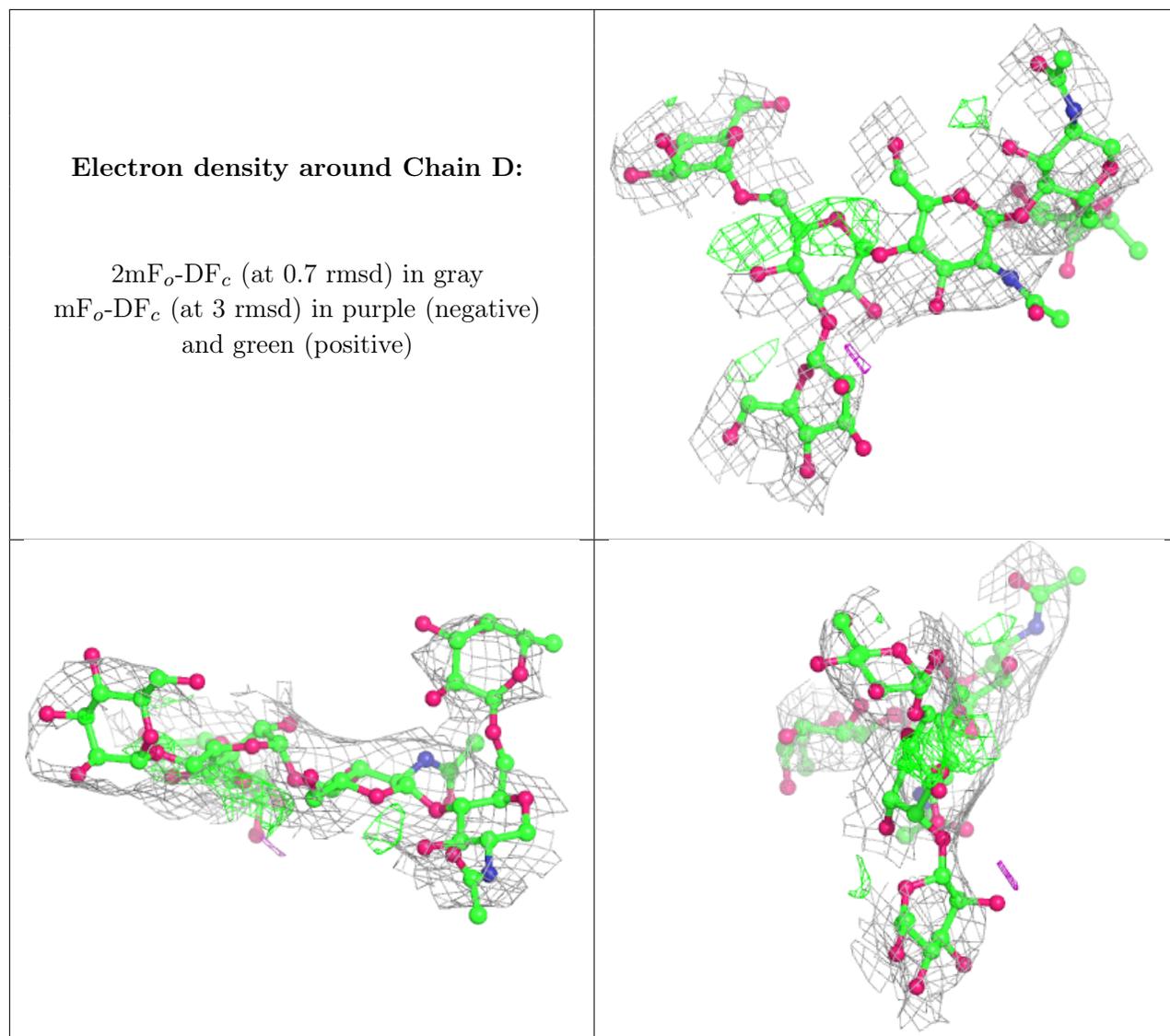
6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	MAN	D	4	11/12	0.72	0.43	172,184,188,190	0
4	FUC	D	6	10/11	0.72	0.53	179,185,188,189	0
4	MAN	C	4	11/12	0.79	0.42	162,183,190,199	0
4	FUC	C	6	10/11	0.80	0.29	171,175,183,190	0
4	BMA	D	3	11/12	0.80	0.29	97,109,131,155	0
4	MAN	C	5	11/12	0.84	0.36	109,116,124,136	0
4	MAN	D	5	11/12	0.86	0.30	104,125,132,133	0
4	NAG	D	1	14/15	0.86	0.29	87,101,171,184	0
4	NAG	D	2	14/15	0.88	0.26	83,99,149,159	0
4	NAG	C	1	14/15	0.90	0.27	96,118,153,167	0
4	NAG	C	2	14/15	0.93	0.25	95,113,144,146	0
4	BMA	C	3	11/12	0.96	0.26	81,104,118,137	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.





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.