



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2021 – 04:03 PM EDT

PDB ID : 7LO4  
Title : SARS-CoV-2 spike receptor-binding domain with a G485R mutation in complex with human ACE2  
Authors : Weekley, C.M.; Parker, M.W.  
Deposited on : 2021-02-09  
Resolution : 2.46 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.17.1  
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) : 2.17.1

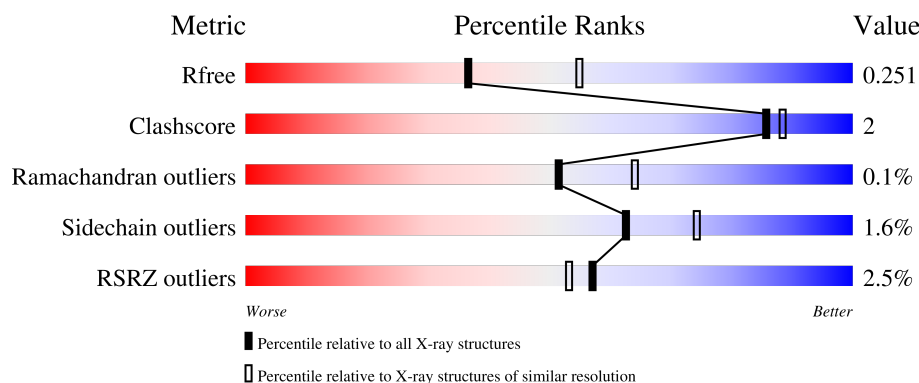
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


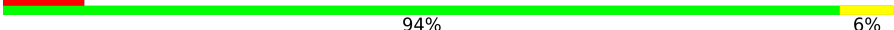
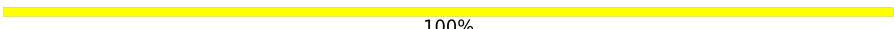
The reported resolution of this entry is 2.46 Å.

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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	596	
2	B	200	
3	C	3	

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6692 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 Processed angiotensin-converting enzyme 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4862	3111	805	917	29			

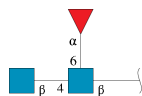
- Molecule 2 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	200	Total	C	N	O	S	0	0	0
			1591	1017	268	298	8			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	485	ARG	GLY	engineered mutation	UNP P0DTC2
B	531	GLU	-	expression tag	UNP P0DTC2
B	532	ASN	-	expression tag	UNP P0DTC2

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



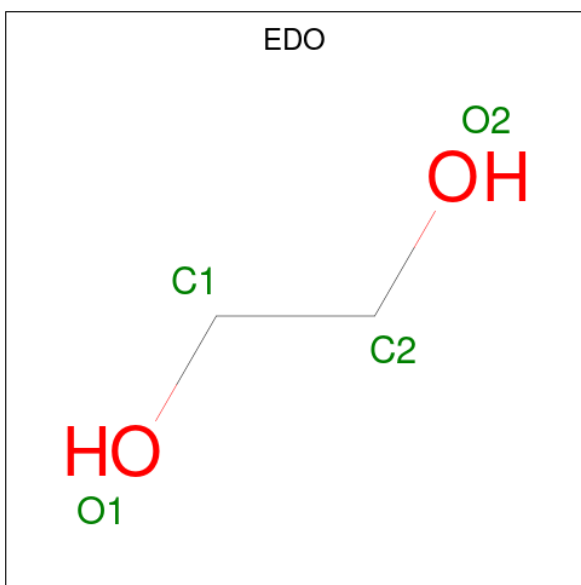
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	3	Total	C	N	O		0	0	0
			38	22	2	14				

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Zn 1 1	0	0

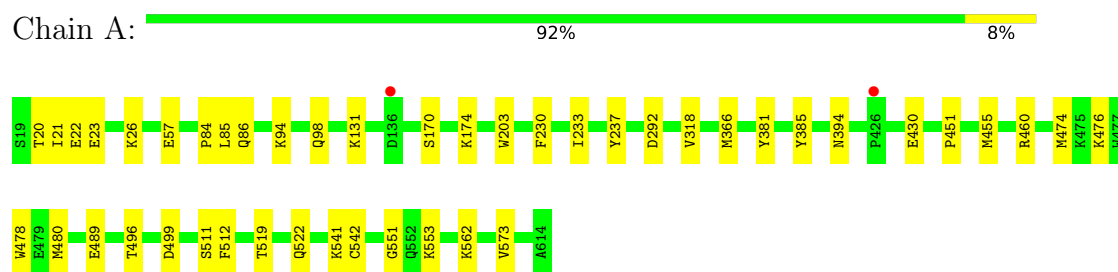
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	97	Total O 97 97	0	0
7	B	35	Total O 35 35	0	0

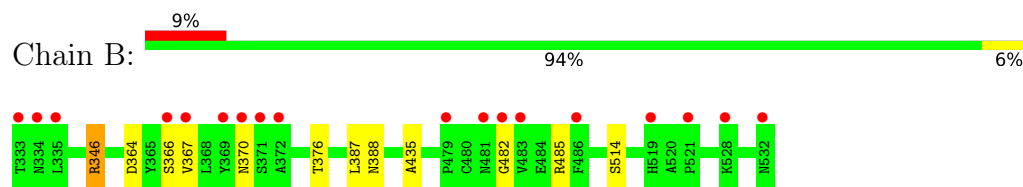
### 3 Residue-property plots

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: Processed angiotensin-converting enzyme 2



- Molecule 2: Spike protein S1



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.61Å 112.18Å 145.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.79 – 2.46 48.79 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.79-2.46) 99.8 (48.79-2.46)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.05 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.193 , 0.251 0.193 , 0.251	Depositor DCC
$R_{free}$ test set	1968 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.0	Xtriage
Anisotropy	0.162	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 37.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6692	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: FUC, NAG, EDO, ZN

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	0.44	0/4999	0.55	0/6792
2	B	0.45	0/1635	0.57	0/2223
All	All	0.44	0/6634	0.55	0/9015

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

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	4862	0	4635	21	0
2	B	1591	0	1512	7	0
3	C	38	0	34	0	0
4	A	42	0	39	0	0
4	B	14	0	13	0	0
5	A	12	0	18	0	0
6	A	1	0	0	0	0
7	A	97	0	0	0	0
7	B	35	0	0	1	0
All	All	6692	0	6251	28	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including



hydrogen atoms). The all-atom clashscore for this structure is 2.

All (28) 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:B:366:SER:O	2:B:370:ASN:ND2	2.18	0.75
2:B:364:ASP:O	2:B:367:VAL:HG22	1.95	0.65
1:A:86:GLN:OE1	1:A:86:GLN:N	2.28	0.62
1:A:85:LEU:HD22	1:A:94:LYS:HG3	1.82	0.60
1:A:230:PHE:O	1:A:233:ILE:HG22	2.08	0.53
1:A:21:ILE:HG21	1:A:84:PRO:HD2	1.91	0.52
1:A:22:GLU:HG2	1:A:26:LYS:HE3	1.91	0.51
2:B:366:SER:HB3	2:B:388:ASN:HD21	1.77	0.49
1:A:20:THR:OG1	1:A:23:GLU:HG2	2.14	0.47
1:A:474:MET:HE1	1:A:499:ASP:HB2	1.96	0.46
1:A:476:LYS:O	1:A:480:MET:HG3	2.16	0.46
1:A:553:LYS:HE2	1:A:573:VAL:O	2.16	0.45
2:B:387:LEU:HD23	2:B:387:LEU:HA	1.84	0.45
1:A:174:LYS:HG2	1:A:496:THR:O	2.17	0.45
1:A:292:ASP:HA	1:A:366:MET:SD	2.57	0.45
1:A:460:ARG:NH2	1:A:512:PHE:HB2	2.32	0.45
1:A:394:ASN:HB3	1:A:562:LYS:HD3	2.00	0.44
1:A:23:GLU:HG2	1:A:23:GLU:H	1.60	0.43
1:A:430:GLU:OE1	1:A:541:LYS:HD3	2.18	0.43
1:A:478:TRP:CD2	1:A:489:GLU:HB3	2.53	0.43
1:A:519:THR:O	1:A:522:GLN:HG2	2.18	0.43
1:A:237:TYR:CE1	1:A:451:PRO:HG2	2.54	0.43
1:A:318:VAL:O	1:A:551:GLY:HA3	2.19	0.43
2:B:376:THR:HB	2:B:435:ALA:HB3	2.01	0.42
2:B:346:ARG:HA	7:B:707:HOH:O	2.18	0.42
2:B:370:ASN:OD1	2:B:370:ASN:O	2.38	0.41
1:A:430:GLU:H	1:A:430:GLU:HG3	1.64	0.41
1:A:203:TRP:CD2	1:A:511:SER:HB3	2.57	0.40

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	594/596 (100%)	577 (97%)	17 (3%)	0	100	100
2	B	198/200 (99%)	186 (94%)	11 (6%)	1 (0%)	29	34
All	All	792/796 (100%)	763 (96%)	28 (4%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	482	GLY

### 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	526/526 (100%)	518 (98%)	8 (2%)	65	76
2	B	174/174 (100%)	171 (98%)	3 (2%)	60	73
All	All	700/700 (100%)	689 (98%)	11 (2%)	62	74

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	GLU
1	A	98	GLN
1	A	131	LYS
1	A	170	SER
1	A	381	TYR
1	A	385	TYR
1	A	455	MET
1	A	542	CYS
2	B	346	ARG
2	B	485	ARG
2	B	514	SER

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

Mol	Chain	Res	Type
1	A	42	GLN
2	B	370	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

3 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	$\# Z  > 2$
3	NAG	C	1	3,1	14,14,15	0.31	0	17,19,21	0.67	1 (5%)
3	NAG	C	2	3	14,14,15	0.75	1 (7%)	17,19,21	0.58	0
3	FUC	C	3	3	10,10,11	1.94	3 (30%)	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
3	NAG	C	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
3	FUC	C	3	3	-	-	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	3	FUC	C2-C3	4.43	1.59	1.52
3	C	2	NAG	C1-C2	2.35	1.55	1.52
3	C	3	FUC	C1-C2	2.32	1.57	1.52
3	C	3	FUC	C4-C3	2.22	1.58	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	NAG	C1-O5-C5	2.18	115.15	112.19
3	C	3	FUC	C1-C2-C3	2.18	112.35	109.67

There are no chirality outliers.

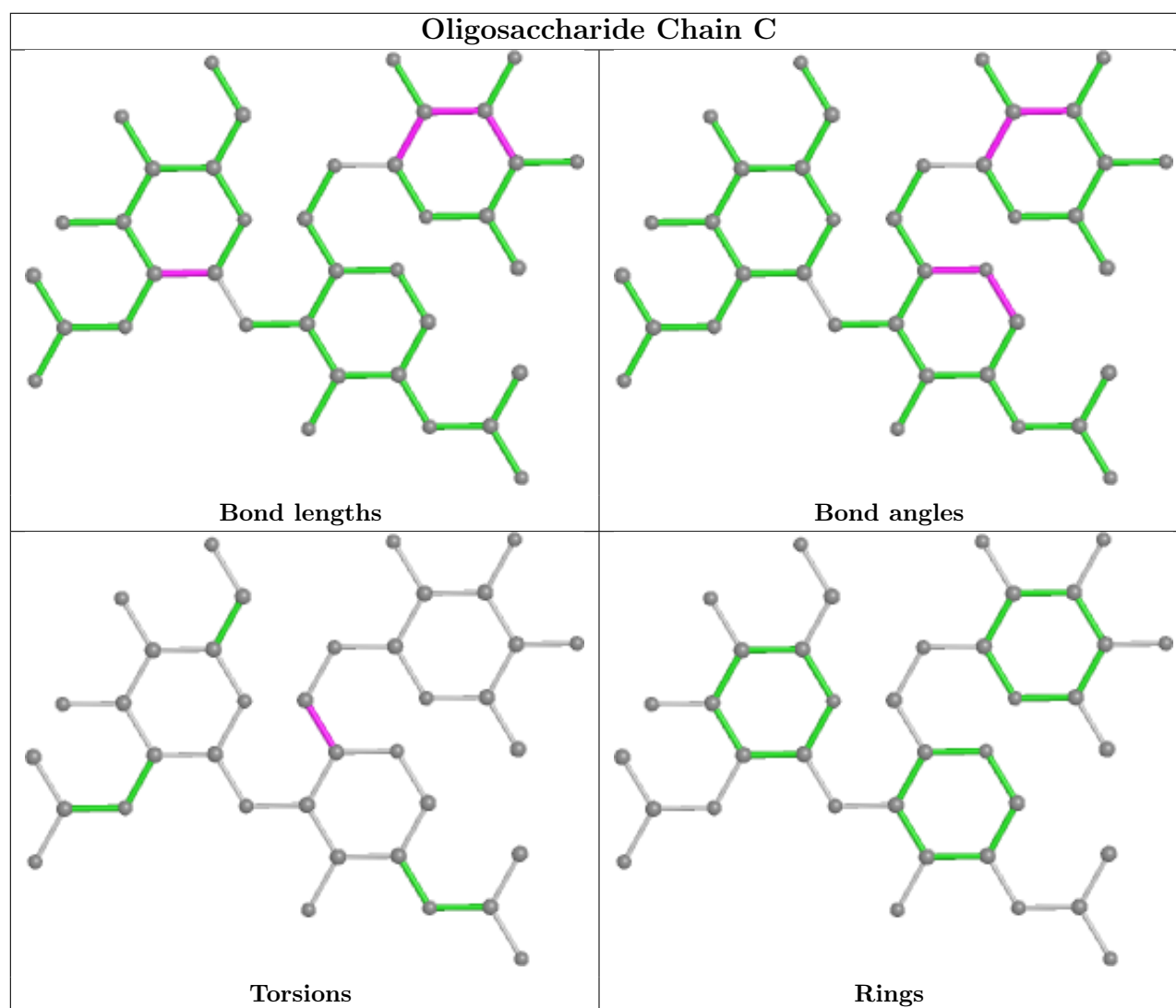
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1	NAG	O5-C5-C6-O6

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](#)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 for Mogul analysis.

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	A	702	1	14,14,15	0.51	0	17,19,21	0.53	0
4	NAG	A	703	1	14,14,15	1.26	2 (14%)	17,19,21	0.91	1 (5%)
4	NAG	B	601	2	14,14,15	1.16	1 (7%)	17,19,21	0.91	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	EDO	A	706	-	3,3,3	0.52	0	2,2,2	0.44	0
4	NAG	A	701	1	14,14,15	0.35	0	17,19,21	0.53	0
5	EDO	A	704	-	3,3,3	0.43	0	2,2,2	0.58	0
5	EDO	A	705	-	3,3,3	0.50	0	2,2,2	0.26	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	NAG	A	702	1	-	2/6/23/26	0/1/1/1
4	NAG	A	703	1	-	2/6/23/26	0/1/1/1
4	NAG	B	601	2	-	2/6/23/26	0/1/1/1
5	EDO	A	706	-	-	0/1/1/1	-
4	NAG	A	701	1	-	2/6/23/26	0/1/1/1
5	EDO	A	704	-	-	1/1/1/1	-
5	EDO	A	705	-	-	0/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	NAG	O5-C1	4.05	1.50	1.43
4	A	703	NAG	O5-C1	4.04	1.50	1.43
4	A	703	NAG	C1-C2	2.22	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	703	NAG	C1-O5-C5	3.04	116.32	112.19
4	B	601	NAG	C1-O5-C5	2.83	116.03	112.19

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	601	NAG	O5-C5-C6-O6
4	A	701	NAG	O5-C5-C6-O6
4	B	601	NAG	C4-C5-C6-O6
4	A	702	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	A	703	NAG	O5-C5-C6-O6
4	A	701	NAG	C4-C5-C6-O6
4	A	702	NAG	C4-C5-C6-O6
4	A	703	NAG	C4-C5-C6-O6
5	A	704	EDO	O1-C1-C2-O2

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

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	596/596 (100%)	-0.25	2 (0%) 94 94	24, 35, 53, 85	0
2	B	200/200 (100%)	0.20	18 (9%) 9 6	23, 38, 69, 82	0
All	All	796/796 (100%)	-0.13	20 (2%) 57 53	23, 35, 63, 85	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	369	TYR	4.9
1	A	136	ASP	4.1
2	B	370	ASN	4.0
2	B	333	THR	3.9
2	B	532	ASN	3.7
2	B	483	VAL	3.6
2	B	334	ASN	3.5
2	B	486	PHE	3.0
2	B	521	PRO	3.0
2	B	367	VAL	2.9
2	B	366	SER	2.8
2	B	519	HIS	2.8
2	B	371	SER	2.8
2	B	335	LEU	2.5
2	B	528	LYS	2.5
2	B	372	ALA	2.5
2	B	482	GLY	2.1
2	B	481	ASN	2.1
1	A	426	PRO	2.0
2	B	479	PRO	2.0



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

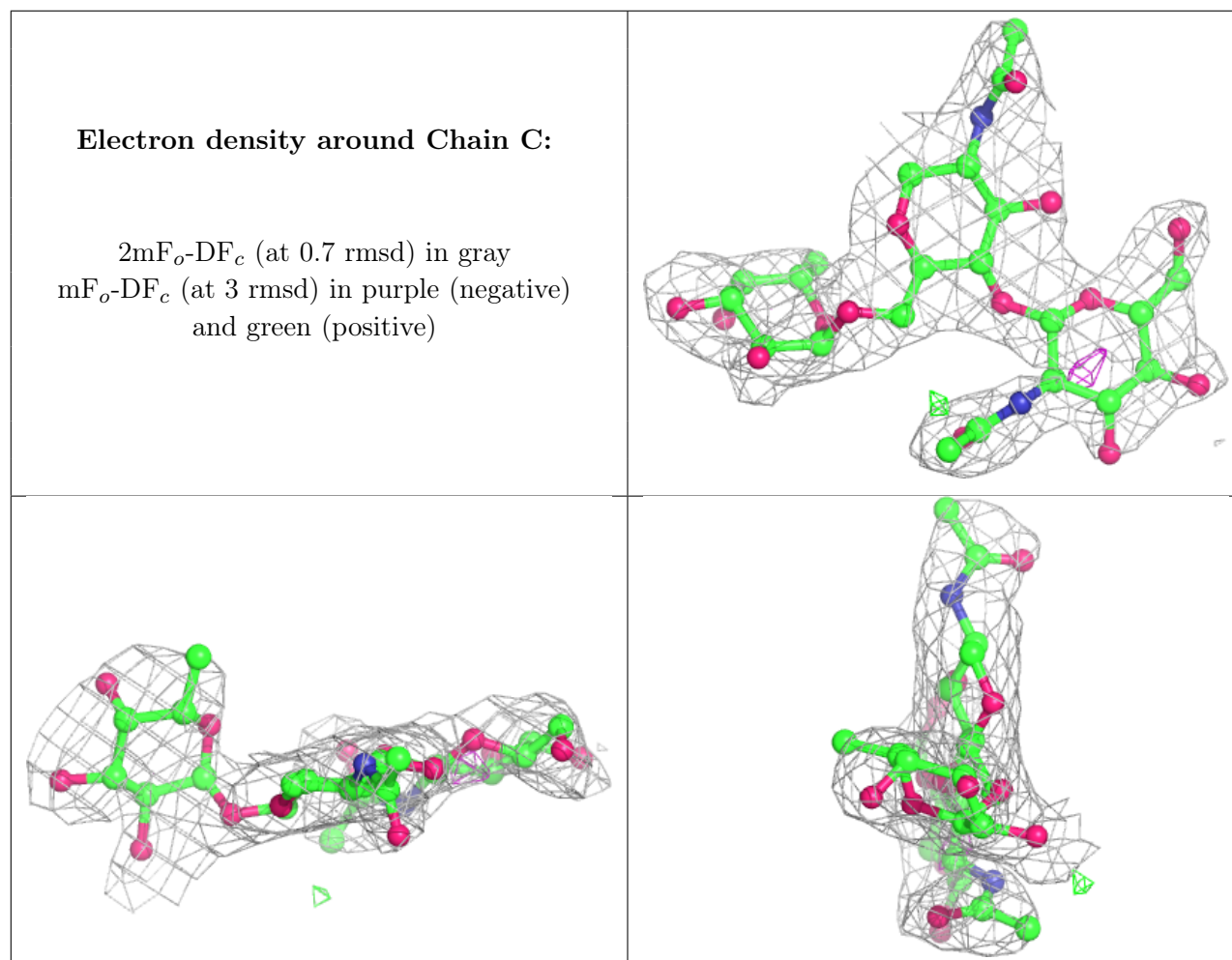
There are no non-standard protein/DNA/RNA residues in this entry.

## 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, 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	NAG	C	2	14/15	0.68	0.38	65,72,77,79	0
3	FUC	C	3	10/11	0.76	0.31	58,67,71,73	0
3	NAG	C	1	14/15	0.93	0.22	57,62,66,72	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](#)

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	A	703	14/15	0.86	0.16	44,50,57,59	0
4	NAG	A	702	14/15	0.87	0.19	50,56,60,64	0
5	EDO	A	706	4/4	0.87	0.21	38,44,45,45	0
4	NAG	B	601	14/15	0.91	0.18	39,54,61,68	0
4	NAG	A	701	14/15	0.91	0.15	51,57,67,79	0
5	EDO	A	705	4/4	0.93	0.20	31,36,38,39	0
5	EDO	A	704	4/4	0.94	0.22	32,35,42,46	0
6	ZN	A	707	1/1	0.98	0.17	27,27,27,27	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.