Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 04:32 PM BST

PDB ID	:	4WZL
Title	:	Crystal structure of P domain from norovirus strain Saga4 in complex with HBGA type Lea (triglycan)
Authors	:	Singh, B.K.; Hansman, G.S.
Deposited on	:	2014-11-20
Resolution	:	1.70 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ⓘ) were used in the production of this report:

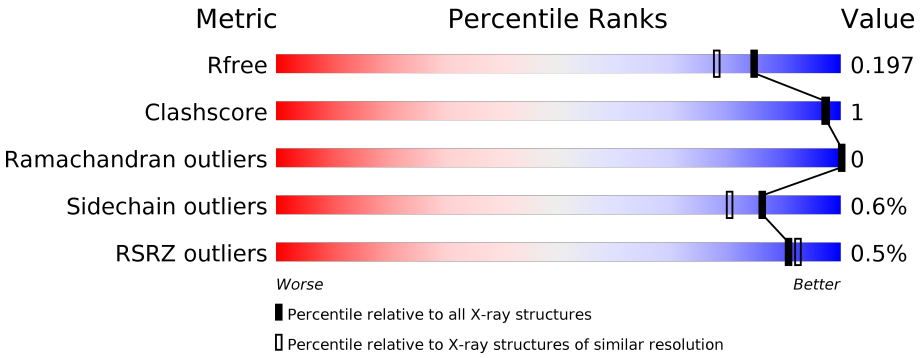
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 on 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 <=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	308	% 98% .
1	B	308	98% .
2	C	3	100%
2	D	3	100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9983 atoms, of which 4478 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.

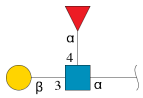
- Molecule 1 is a protein called VP1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	308	Total	C	H	N	O	S	0	2	0
			4643	1534	2235	409	455	10			
1	B	308	Total	C	H	N	O	S	0	2	0
			4647	1531	2243	408	455	10			

There are 4 discrepancies between the modelled and reference sequences:

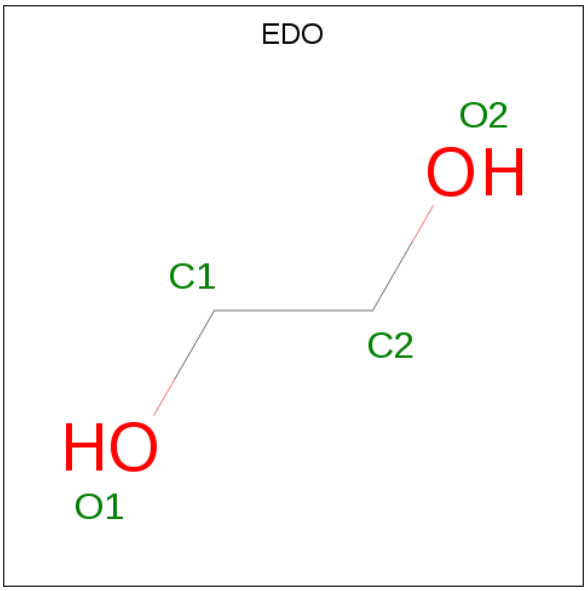
Chain	Residue	Modelled	Actual	Comment	Reference
A	223	GLY	-	expression tag	UNP B5BTR7
A	224	SER	-	expression tag	UNP B5BTR7
B	223	GLY	-	expression tag	UNP B5BTR7
B	224	SER	-	expression tag	UNP B5BTR7

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



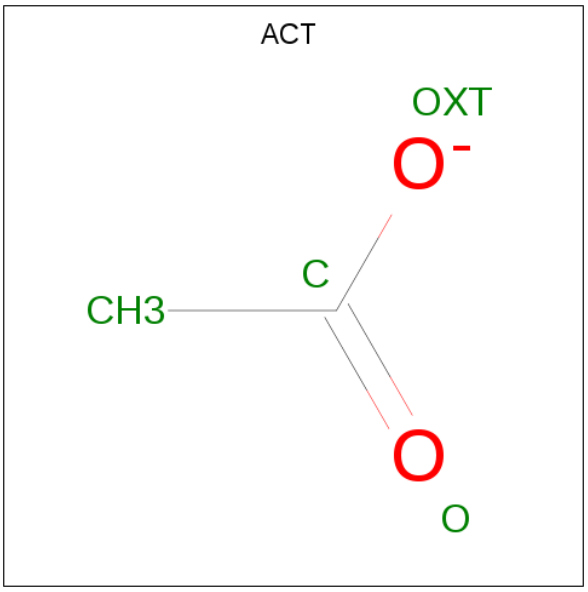
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			36	20	1	15			
2	D	3	Total	C	N	O	0	0	0
			36	20	1	15			

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 5 is water.

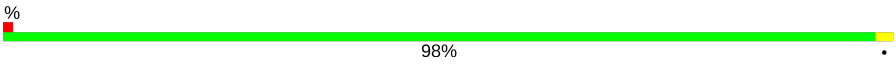
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	284	Total O 285 285	0	1
5	B	283	Total O 284 284	0	1

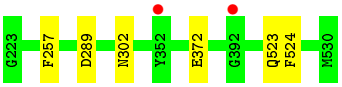
3 Residue-property plots (i)

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: VP1

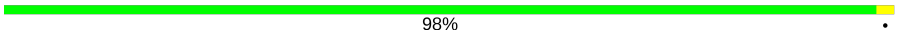
Chain A:

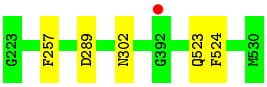




• Molecule 1: VP1

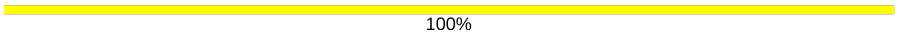
Chain B:





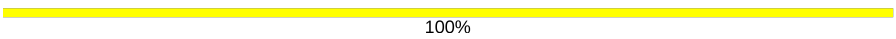
• Molecule 2: beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-alpha-D-glucopyranose

Chain C:



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

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	96.87Å 58.83Å 124.44Å 90.00° 119.80° 90.00°	Depositor
Resolution (Å)	42.47 – 1.70 49.59 – 1.57	Depositor EDS
% Data completeness (in resolution range)	98.0 (42.47-1.70) 97.6 (49.59-1.57)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.36 (at 1.57Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.170 , 0.196 0.171 , 0.197	Depositor DCC
R_{free} test set	4126 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	15.6	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.42 , 50.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.23$	Xtriage
Estimated twinning fraction	0.027 for 1/2*h-3/2*k,-1/2*h-1/2*k,-h+k-l 0.040 for 1/2*h+3/2*k,1/2*h-1/2*k,-h-k-l	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	9983	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 96.60 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.5630e-10. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, GAL, EDO, NDG, ACT

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	0.28	0/2487	0.51	0/3402
1	B	0.28	0/2483	0.52	0/3398
All	All	0.28	0/4970	0.51	0/6800

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	2408	2235	2321	3	0
1	B	2404	2243	2310	3	0
2	C	36	0	30	1	0
2	D	36	0	30	1	0
3	A	16	0	24	0	0
3	B	20	0	30	0	0
4	A	8	0	6	0	0
4	B	8	0	6	0	0
5	A	285	0	0	2	0
5	B	284	0	0	2	0
All	All	5505	4478	4757	8	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 1.

All (8) 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:302:ASN:ND2	5:A:948:HOH:O	2.07	0.88
1:B:302:ASN:ND2	5:B:944:HOH:O	2.08	0.85
1:B:257:PHE:O	5:B:780:HOH:O	2.18	0.54
1:A:257:PHE:O	5:A:783:HOH:O	2.19	0.51
2:C:2:GAL:H2	2:C:3:FUC:C6	2.42	0.49
2:D:2:GAL:H2	2:D:3:FUC:C6	2.44	0.48
1:A:523:GLN:HG2	1:A:524:PHE:CD1	2.55	0.42
1:B:523:GLN:HG2	1:B:524:PHE:CD1	2.55	0.42

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	308/308 (100%)	301 (98%)	7 (2%)	0	100	100
1	B	308/308 (100%)	301 (98%)	7 (2%)	0	100	100
All	All	616/616 (100%)	602 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	269/267 (101%)	267 (99%)	2 (1%)	84 77
1	B	268/267 (100%)	267 (100%)	1 (0%)	91 87
All	All	537/534 (101%)	534 (99%)	3 (1%)	86 80

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	289	ASP
1	A	372	GLU
1	B	289	ASP

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

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

6 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	NDG	C	1	2	15,15,15	0.67	0	21,21,21	1.09	1 (4%)
2	GAL	C	2	2	11,11,12	0.53	0	15,15,17	0.89	0
2	FUC	C	3	2	10,10,11	0.30	0	14,14,16	0.63	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	NDG	D	1	2	15,15,15	0.67	0	21,21,21	1.08	1 (4%)
2	GAL	D	2	2	11,11,12	0.52	0	15,15,17	0.89	0
2	FUC	D	3	2	10,10,11	0.30	0	14,14,16	0.62	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
2	NDG	C	1	2	-	2/6/26/26	0/1/1/1
2	GAL	C	2	2	-	1/2/19/22	0/1/1/1
2	FUC	C	3	2	-	-	0/1/1/1
2	NDG	D	1	2	-	2/6/26/26	0/1/1/1
2	GAL	D	2	2	-	1/2/19/22	0/1/1/1
2	FUC	D	3	2	-	-	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	NDG	C1-C2-N2	3.82	115.15	110.73
2	D	1	NDG	C1-C2-N2	3.76	115.09	110.73

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	NDG	O5-C5-C6-O6
2	C	1	NDG	O5-C5-C6-O6
2	C	1	NDG	C4-C5-C6-O6
2	D	1	NDG	C4-C5-C6-O6
2	D	2	GAL	O5-C5-C6-O6
2	C	2	GAL	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	GAL	1	0

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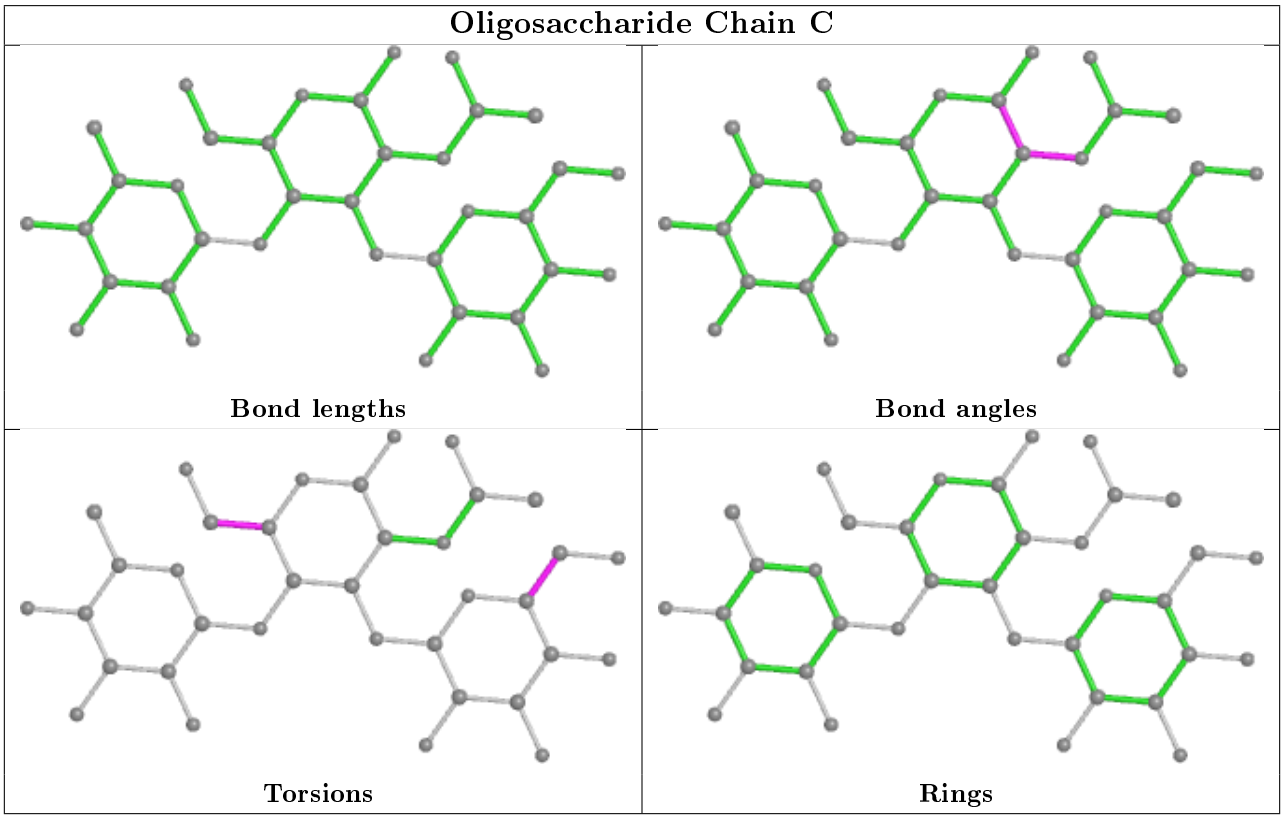
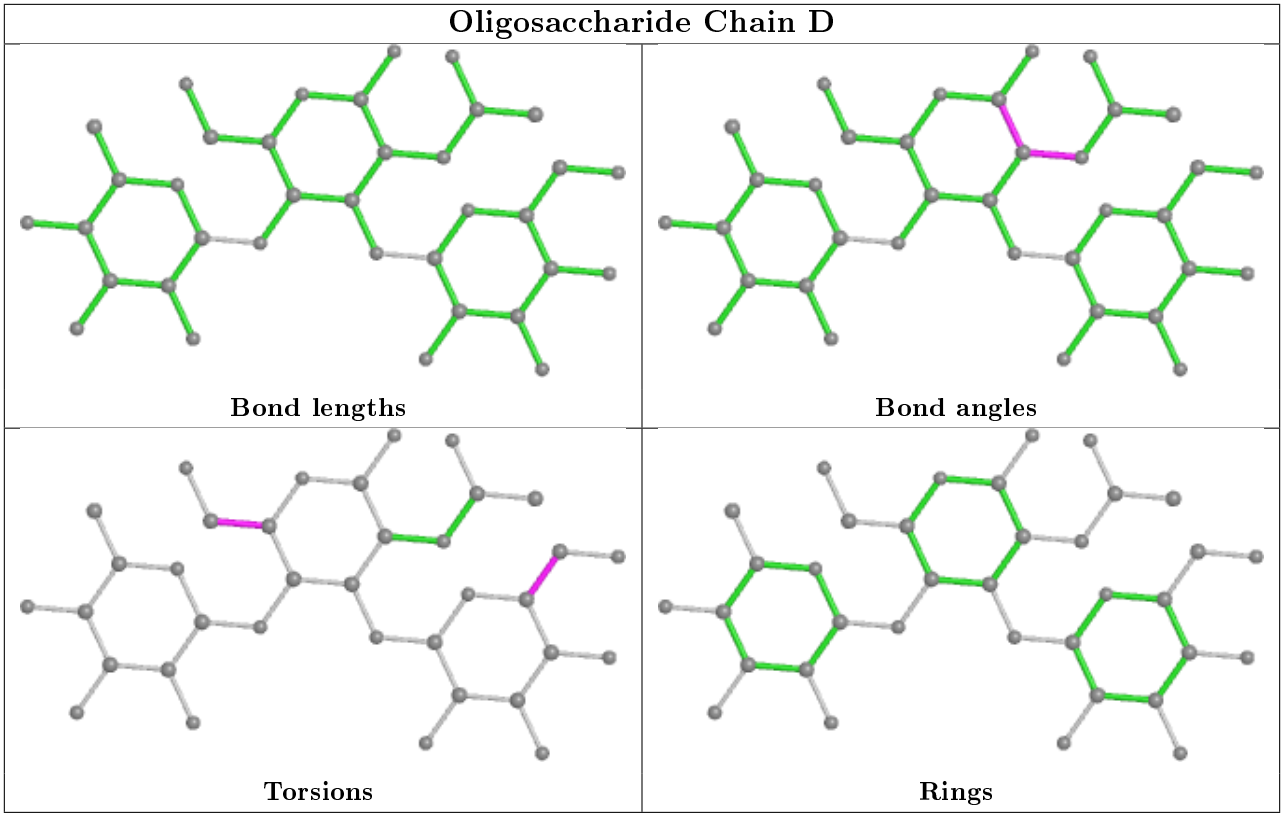
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	3	FUC	1	0
2	D	3	FUC	1	0
2	D	2	GAL	1	0

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)

13 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	ACT	B	606	-	1,3,3	1.71	0	0,3,3	0.00	-
3	EDO	A	603	-	3,3,3	0.53	0	2,2,2	0.15	0
3	EDO	A	602	-	3,3,3	0.43	0	2,2,2	0.46	0
3	EDO	B	602	-	3,3,3	0.53	0	2,2,2	0.16	0
4	ACT	B	607	-	1,3,3	0.97	0	0,3,3	0.00	-
4	ACT	A	605	-	1,3,3	1.65	0	0,3,3	0.00	-
3	EDO	A	601	-	3,3,3	0.47	0	2,2,2	0.43	0
3	EDO	B	603	-	3,3,3	0.44	0	2,2,2	0.46	0
3	EDO	B	604	-	3,3,3	0.48	0	2,2,2	0.40	0
3	EDO	A	604	-	3,3,3	0.47	0	2,2,2	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	EDO	B	605	-	3,3,3	0.45	0	2,2,2	0.39	0
4	ACT	A	606	-	1,3,3	1.42	0	0,3,3	0.00	-
3	EDO	B	601	-	3,3,3	0.46	0	2,2,2	0.44	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	A	603	-	-	0/1/1/1	-
3	EDO	A	602	-	-	0/1/1/1	-
3	EDO	B	602	-	-	0/1/1/1	-
3	EDO	A	601	-	-	0/1/1/1	-
3	EDO	B	603	-	-	0/1/1/1	-
3	EDO	B	604	-	-	0/1/1/1	-
3	EDO	A	604	-	-	0/1/1/1	-
3	EDO	B	605	-	-	0/1/1/1	-
3	EDO	B	601	-	-	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

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^{th} 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	308/308 (100%)	-0.38	2 (0%)	89	91	7, 17, 32, 51	0
1	B	308/308 (100%)	-0.57	1 (0%)	94	94	5, 14, 31, 47	0
All	All	616/616 (100%)	-0.48	3 (0%)	91	92	5, 16, 32, 51	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	392	GLY	3.6
1	A	392	GLY	2.6
1	A	352[A]	TYR	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

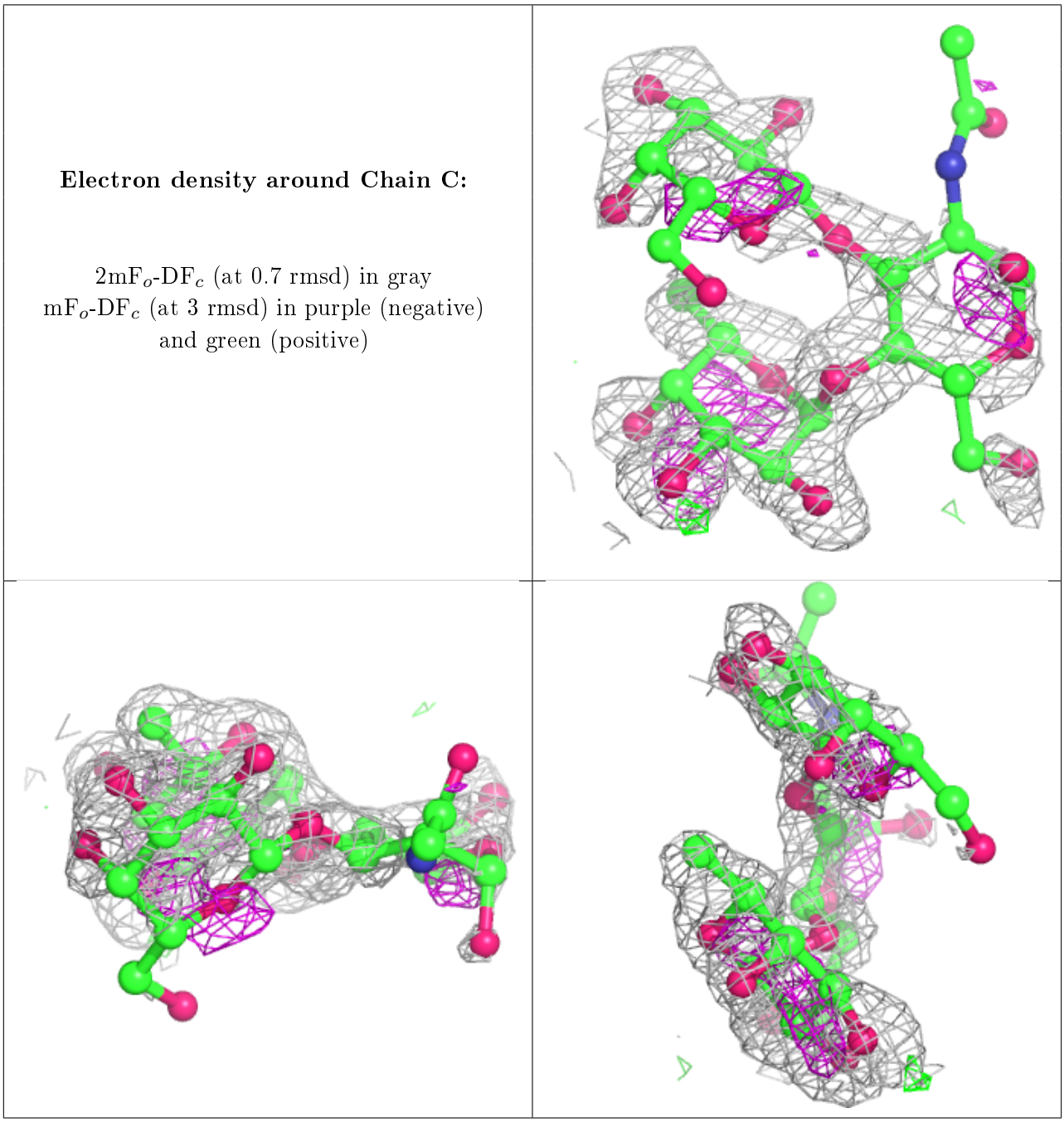
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GAL	D	2	11/12	0.67	0.29	50,52,54,54	0
2	GAL	C	2	11/12	0.68	0.30	50,51,54,54	0
2	NDG	C	1	15/15	0.74	0.35	36,53,59,63	0
2	NDG	D	1	15/15	0.74	0.36	36,53,60,64	0
2	FUC	D	3	10/11	0.75	0.22	22,28,33,37	0
2	FUC	C	3	10/11	0.77	0.21	21,28,32,37	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.

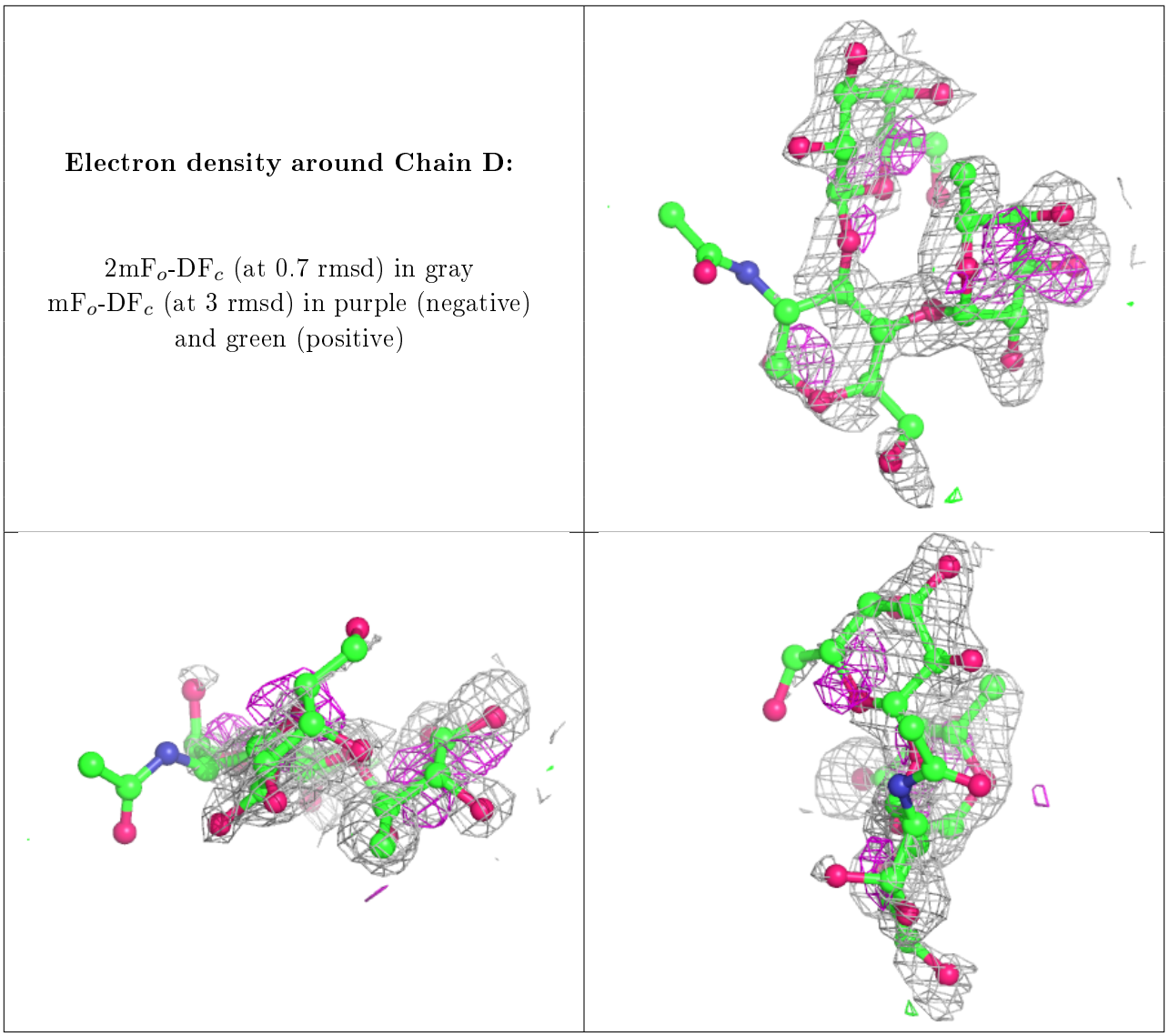
Electron density around Chain C:

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



Electron density around Chain D:

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



6.4 Ligands

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^{th} 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^2)	Q<0.9
3	EDO	B	605	4/4	0.70	0.29	38,41,41,44	0
4	ACT	B	607	4/4	0.87	0.14	25,29,33,45	0
4	ACT	A	605	4/4	0.91	0.09	18,20,22,25	0
3	EDO	A	603	4/4	0.91	0.10	13,16,21,23	0
4	ACT	A	606	4/4	0.91	0.15	25,32,35,36	0
3	EDO	B	602	4/4	0.92	0.09	12,16,21,23	0
3	EDO	A	604	4/4	0.92	0.17	17,20,25,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	B	604	4/4	0.93	0.18	17,20,26,35	0
4	ACT	B	606	4/4	0.93	0.09	16,20,21,26	0
3	EDO	A	602	4/4	0.95	0.08	20,20,22,23	0
3	EDO	A	601	4/4	0.95	0.13	16,21,25,30	0
3	EDO	B	603	4/4	0.96	0.08	19,20,21,23	0
3	EDO	B	601	4/4	0.97	0.13	16,20,25,29	0

6.5 Other polymers

There are no such residues in this entry.