Full wwPDB X-ray Structure Validation Report (i)

Jan 4, 2024 – 02:33 pm GMT

PDB ID	:	5A6Z
Title	:	Structure of the LecB lectin from Pseudomonas aeruginosa strain PA14 in complex with lewis a
Authors	:	Sommer, R.; Wagner, S.; Varrot, A.; Khaledi, A.; Haussler, S.; Imberty, A.; Titz, A.
Deposited on	:	2015-07-02
Resolution	:	1.50 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 (i) 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 (i)) were used in the production of this report:

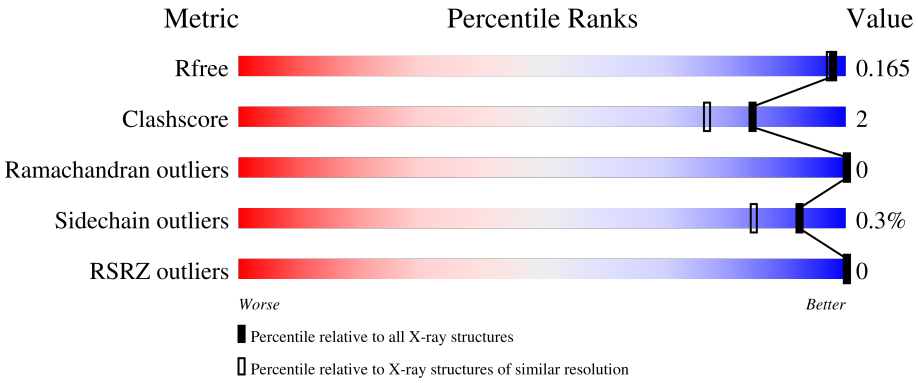
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.50 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	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 <=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	114	95% 5%
1	B	114	97% .
1	C	114	95% 5%
1	D	114	97% .
2	E	4	100%

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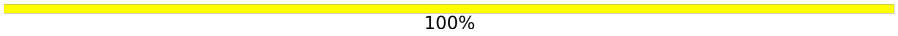
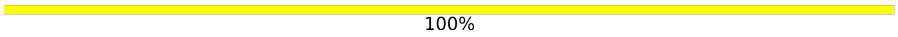
Mol	Chain	Length	Quality of chain
2	H	4	
3	F	3	
4	G	4	

2 Entry composition

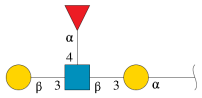
There are 8 unique types of molecules in this entry. The entry contains 4453 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 LECB.

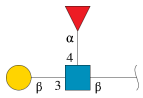
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	114	Total	C	N	O	0	5	0
			852	526	144	182			
1	B	114	Total	C	N	O	0	4	0
			843	523	142	178			
1	C	114	Total	C	N	O	0	3	0
			844	520	145	179			
1	D	114	Total	C	N	O	0	2	0
			839	519	144	176			

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



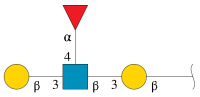
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	4	Total	C	N	O	0	0	0
			45	25	1	19			
2	H	4	Total	C	N	O	0	0	0
			47	26	1	20			

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



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

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

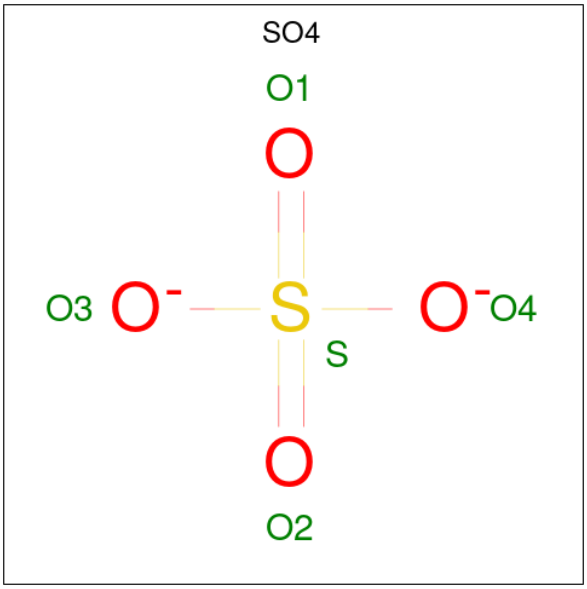


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	4	Total	C	N	O	0	0	0
			47	26	1	20			

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Ca	0	0
			2	2		
5	B	2	Total	Ca	0	0
			2	2		
5	C	2	Total	Ca	0	0
			2	2		
5	D	2	Total	Ca	0	0
			2	2		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		

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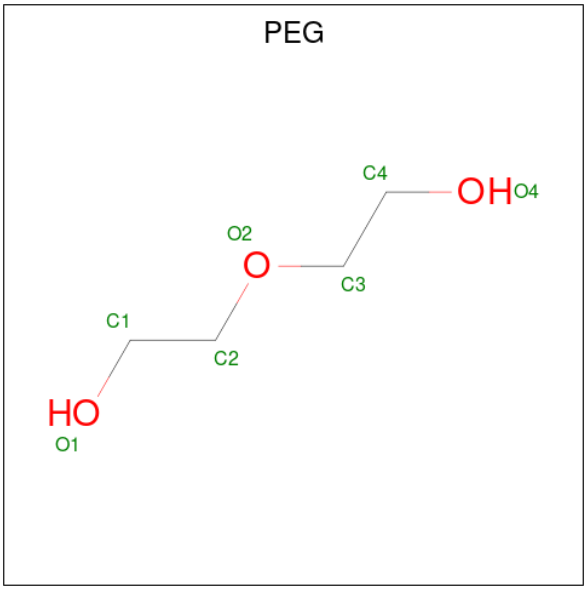
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total O S 5 4 1	0	0
6	C	1	Total O S 5 4 1	0	0

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	D	1	Total C O 14 8 6	0	1

- Molecule 8 is water.

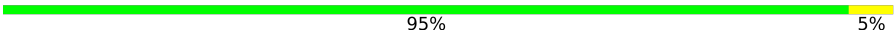
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	221	Total O 221 221	0	0
8	B	219	Total O 219 219	0	0
8	C	232	Total O 232 232	0	0
8	D	191	Total O 191 191	0	0

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.

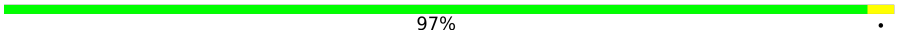
• Molecule 1: LECB

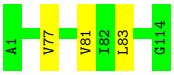
Chain A:



• Molecule 1: LECB

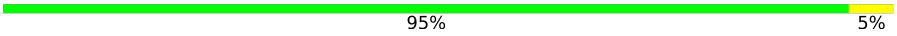
Chain B:





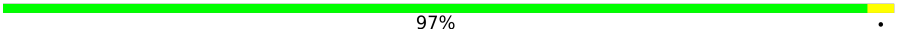
• Molecule 1: LECB

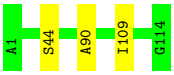
Chain C:



• Molecule 1: LECB

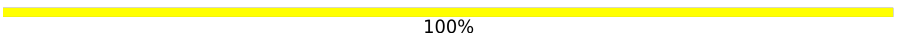
Chain D:





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

Chain E:

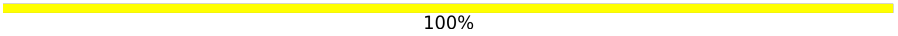


• Molecule 2: beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-3)-alpha-D-galactopyranose

Chain H:

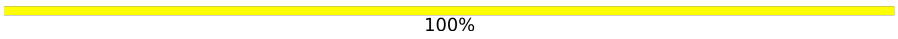
• Molecule 3: beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:



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

Chain G:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	52.98Å 63.77Å 64.51Å 90.00° 91.25° 90.00°	Depositor
Resolution (Å)	27.82 – 1.50 27.32 – 1.50	Depositor EDS
% Data completeness (in resolution range)	96.4 (27.82-1.50) 96.4 (27.32-1.50)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.98 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.131 , 0.155 0.141 , 0.165	Depositor DCC
R_{free} test set	3384 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	10.1	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 45.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -h,l,k 0.018 for -h,-l,-k 0.028 for h,-k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	4453	wwPDB-VP
Average B, all atoms (Å^2)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.68% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: PEG, GAL, FUC, GLA, CA, NAG, SO4

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	0.73	0/861	0.79	0/1178
1	B	0.74	0/855	0.79	0/1169
1	C	0.76	0/853	0.79	0/1166
1	D	0.73	0/848	0.79	0/1158
All	All	0.74	0/3417	0.79	0/4671

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	852	0	822	5	0
1	B	843	0	824	4	0
1	C	844	0	819	5	0
1	D	839	0	822	2	0
2	E	45	0	35	0	0
2	H	47	0	39	0	0
3	F	36	0	30	0	0
4	G	47	0	39	0	0
5	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	2	0	0	0	0
5	C	2	0	0	0	0
5	D	2	0	0	0	0
6	A	5	0	0	0	0
6	B	5	0	0	0	0
6	C	5	0	0	0	0
7	D	14	0	20	3	0
8	A	221	0	0	2	0
8	B	219	0	0	2	0
8	C	232	0	0	3	0
8	D	191	0	0	2	0
All	All	4453	0	3450	16	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 (16) 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:81[B]:VAL:HG21	1:B:81[B]:VAL:HG11	1.35	1.08
7:D:210[B]:PEG:H42	8:D:2106:HOH:O	1.63	0.95
7:D:210[A]:PEG:H41	8:D:2102:HOH:O	1.70	0.91
1:C:86:GLU:OE2	8:C:2193:HOH:O	2.05	0.74
1:C:29[B]:ASN:ND2	8:C:2087:HOH:O	1.62	0.67
1:D:44:SER:HA	7:D:210[A]:PEG:H42	1.79	0.65
8:A:2179:HOH:O	1:C:83:LEU:HD22	2.03	0.59
1:A:86:GLU:HG3	1:A:86:GLU:O	2.07	0.54
1:B:77[A]:VAL:HG12	8:B:2014:HOH:O	2.10	0.51
8:A:2178:HOH:O	1:B:83:LEU:HD22	2.11	0.49
1:C:29[B]:ASN:ND2	8:C:2074:HOH:O	2.47	0.47
1:A:83:LEU:HD22	8:B:2173:HOH:O	2.16	0.46
1:A:91:LEU:CD1	1:B:81[B]:VAL:HG21	2.47	0.45
1:C:90:ALA:HB3	1:C:109:ILE:HB	1.99	0.44
1:D:90:ALA:HB3	1:D:109:ILE:HB	2.01	0.42
1:A:90:ALA:HB3	1:A:109:ILE:HB	2.02	0.41

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	117/114 (103%)	114 (97%)	3 (3%)	0	100	100
1	B	116/114 (102%)	112 (97%)	4 (3%)	0	100	100
1	C	115/114 (101%)	111 (96%)	4 (4%)	0	100	100
1	D	114/114 (100%)	111 (97%)	3 (3%)	0	100	100
All	All	462/456 (101%)	448 (97%)	14 (3%)	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	95/91 (104%)	95 (100%)	0	100	100
1	B	94/91 (103%)	94 (100%)	0	100	100
1	C	94/91 (103%)	92 (98%)	2 (2%)	53	23
1	D	93/91 (102%)	93 (100%)	0	100	100
All	All	376/364 (103%)	374 (100%)	2 (0%)	92	78

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	52[A]	SER
1	C	52[B]	SER

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

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

15 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	GLA	E	1	2	10,10,12	1.04	1 (10%)	14,14,17	1.73	2 (14%)
2	NAG	E	2	2	14,14,15	0.78	0	17,19,21	1.14	1 (5%)
2	GAL	E	3	2	11,11,12	0.48	0	15,15,17	1.03	1 (6%)
2	FUC	E	4	2,5	10,10,11	1.17	1 (10%)	14,14,16	1.39	1 (7%)
3	NAG	F	1	3	15,15,15	0.61	0	21,21,21	1.00	1 (4%)
3	GAL	F	2	3	11,11,12	0.42	0	15,15,17	0.98	1 (6%)
3	FUC	F	3	3,5	10,10,11	0.99	1 (10%)	14,14,16	1.24	1 (7%)
4	GAL	G	1	4	12,12,12	0.84	0	17,17,17	1.22	3 (17%)
4	NAG	G	2	4	14,14,15	1.00	1 (7%)	17,19,21	1.33	1 (5%)
4	GAL	G	3	4	11,11,12	0.62	0	15,15,17	1.16	2 (13%)
4	FUC	G	4	4,5	10,10,11	0.90	0	14,14,16	1.27	2 (14%)
2	GLA	H	1	2	12,12,12	0.53	0	17,17,17	1.28	2 (11%)
2	NAG	H	2	2	14,14,15	0.75	0	17,19,21	1.02	0
2	GAL	H	3	2	11,11,12	0.68	0	15,15,17	0.90	0
2	FUC	H	4	2,5	10,10,11	1.28	2 (20%)	14,14,16	1.33	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
2	GLA	E	1	2	-	-	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	GAL	E	3	2	-	0/2/19/22	0/1/1/1
2	FUC	E	4	2,5	-	-	0/1/1/1
3	NAG	F	1	3	-	0/6/26/26	0/1/1/1
3	GAL	F	2	3	-	0/2/19/22	0/1/1/1
3	FUC	F	3	3,5	-	-	0/1/1/1
4	GAL	G	1	4	-	0/2/22/22	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	GAL	G	3	4	-	0/2/19/22	0/1/1/1
4	FUC	G	4	4,5	-	-	0/1/1/1
2	GLA	H	1	2	-	0/2/22/22	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
2	GAL	H	3	2	-	0/2/19/22	0/1/1/1
2	FUC	H	4	2,5	-	-	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	2	NAG	O5-C1	-3.50	1.38	1.43
2	E	4	FUC	O5-C1	-2.78	1.39	1.43
2	H	4	FUC	O4-C4	2.50	1.48	1.43
2	E	1	GLA	O3-C3	2.26	1.48	1.43
3	F	3	FUC	O5-C1	-2.24	1.40	1.43
2	H	4	FUC	O5-C1	-2.12	1.40	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	E	1	GLA	O3-C3-C4	5.11	119.78	109.99
2	H	4	FUC	O2-C2-C1	3.69	116.69	109.15
2	E	2	NAG	O5-C5-C6	-2.84	102.75	107.20
4	G	1	GAL	O3-C3-C2	-2.81	103.86	110.35
4	G	1	GAL	O3-C3-C4	2.62	116.41	110.35
4	G	4	FUC	O4-C4-C3	-2.62	104.30	110.35
2	E	3	GAL	C1-C2-C3	2.58	112.84	109.67
2	E	1	GLA	O4-C4-C3	2.45	115.05	110.14
4	G	3	GAL	C1-O5-C5	2.33	115.36	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	F	3	FUC	O5-C1-C2	-2.21	107.36	110.77
3	F	1	NAG	O1-C1-C2	2.19	113.77	109.22
4	G	3	GAL	C1-C2-C3	2.12	112.27	109.67
2	H	1	GLA	O5-C5-C6	2.09	111.63	106.44
2	H	1	GLA	O4-C4-C5	2.08	114.47	109.30
3	F	2	GAL	C1-C2-C3	2.07	112.21	109.67
4	G	4	FUC	O2-C2-C1	2.05	113.36	109.15
4	G	1	GAL	C1-O5-C5	-2.05	109.80	113.66
2	E	4	FUC	C1-C2-C3	2.02	112.14	109.67
4	G	2	NAG	C2-N2-C7	-2.01	120.04	122.90

There are no chirality outliers.

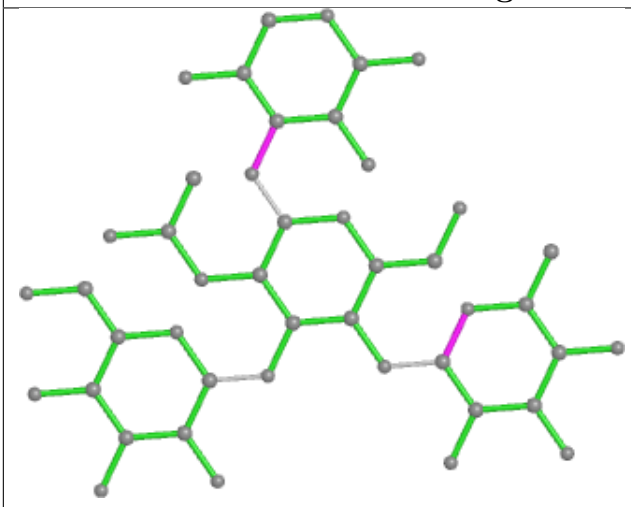
There are no torsion outliers.

There are no ring outliers.

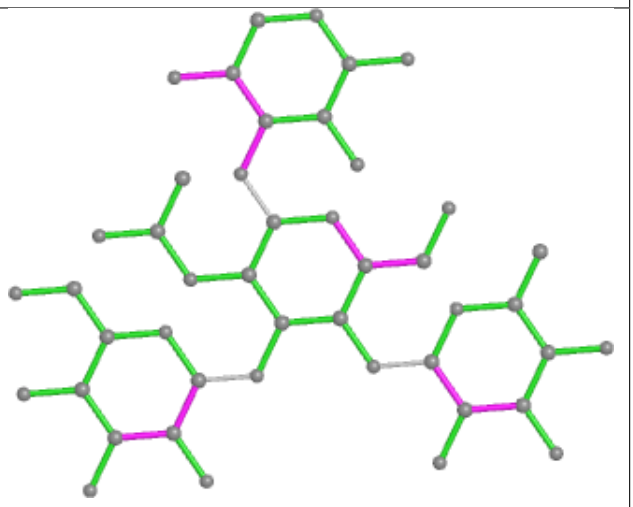
No monomer is involved in short contacts.

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

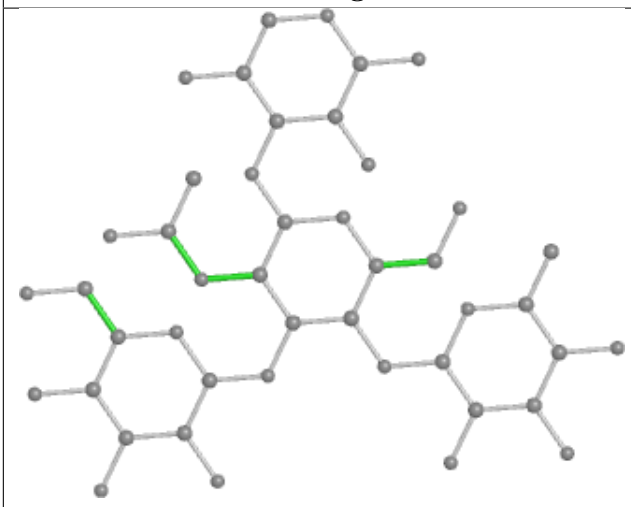
Oligosaccharide Chain E



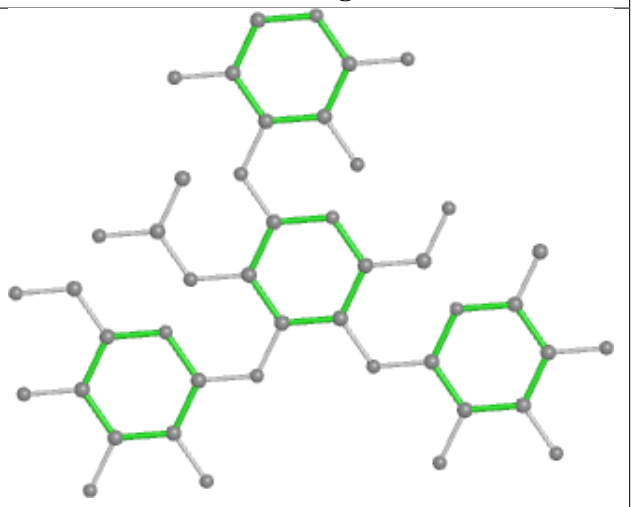
Bond lengths



Bond angles

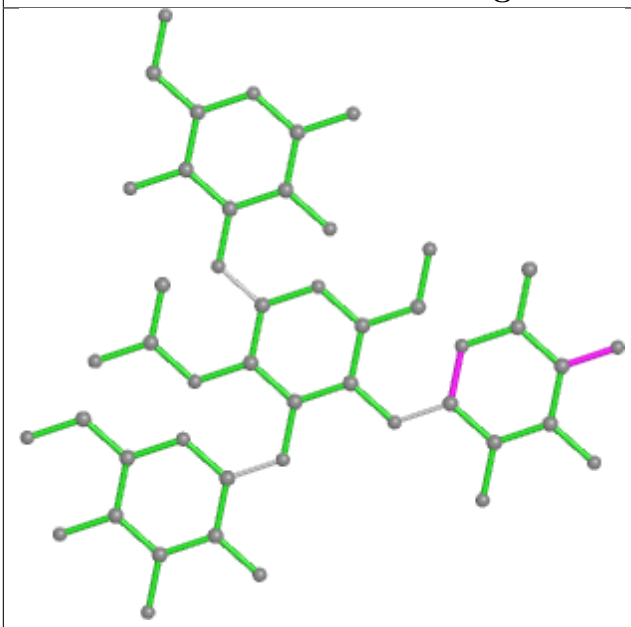


Torsions

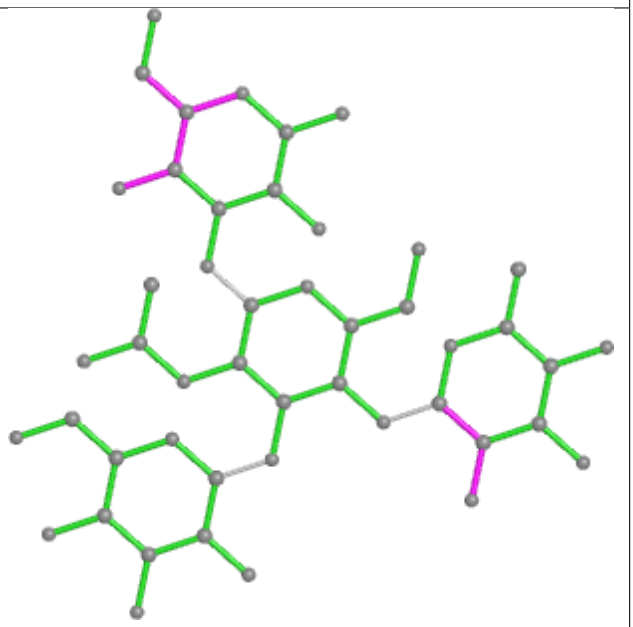


Rings

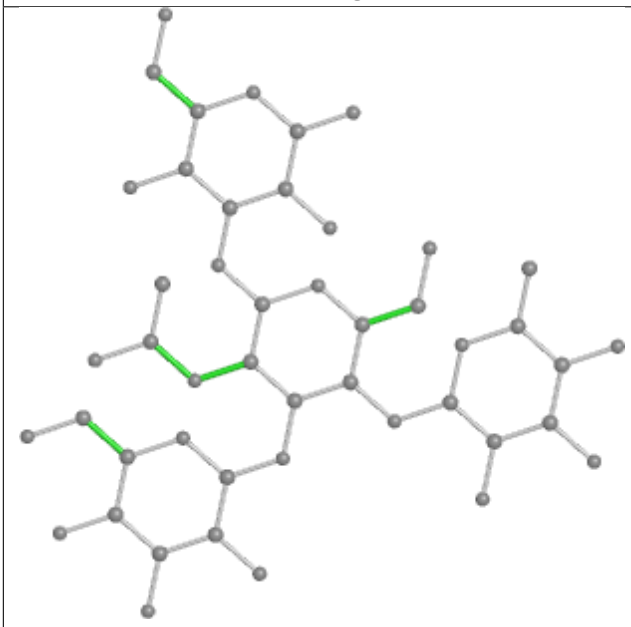
Oligosaccharide Chain H



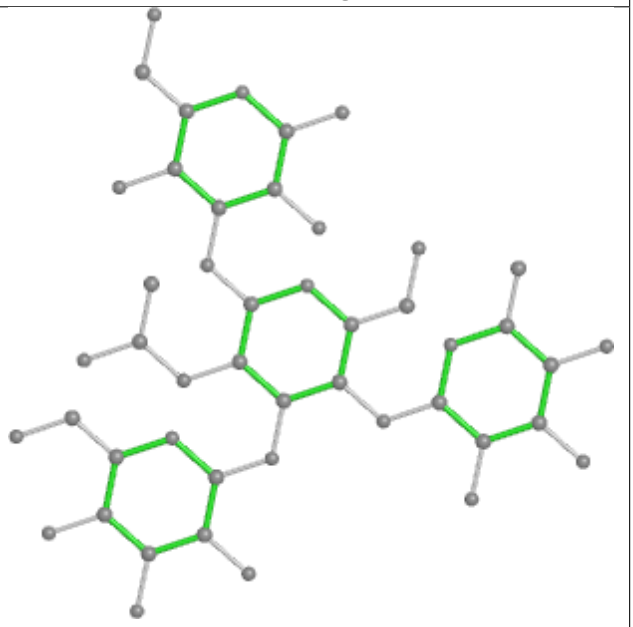
Bond lengths



Bond angles

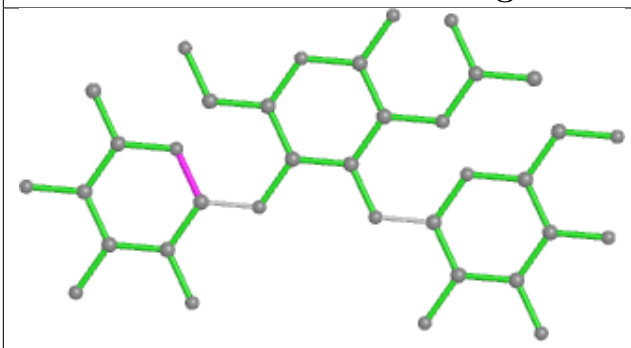


Torsions

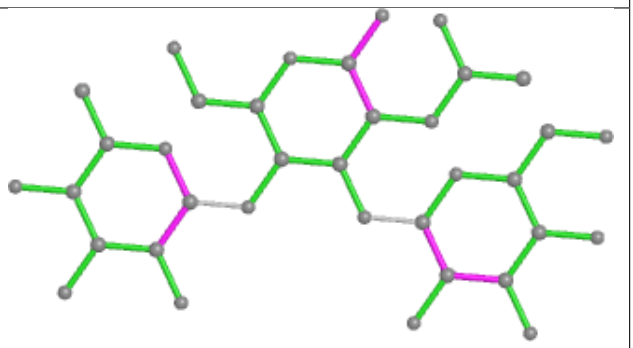


Rings

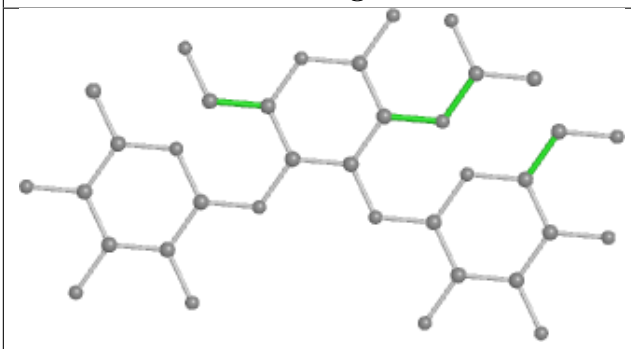
Oligosaccharide Chain F



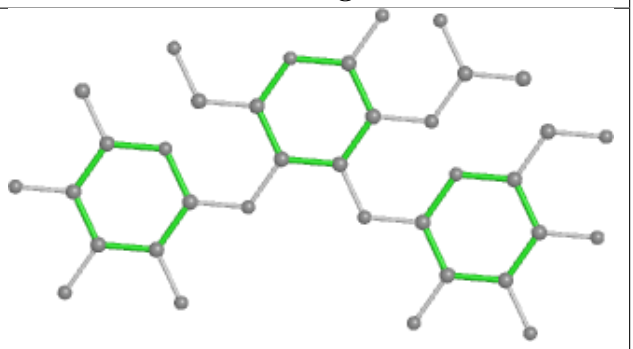
Bond lengths



Bond angles

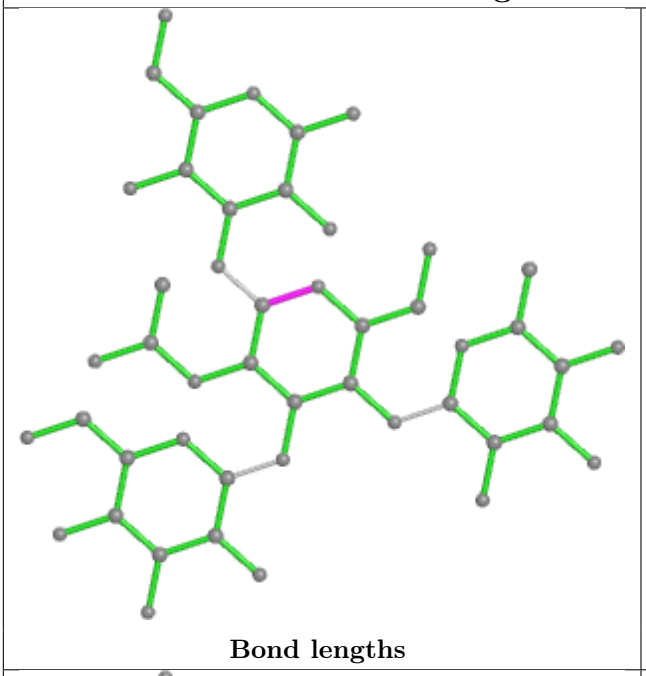


Torsions

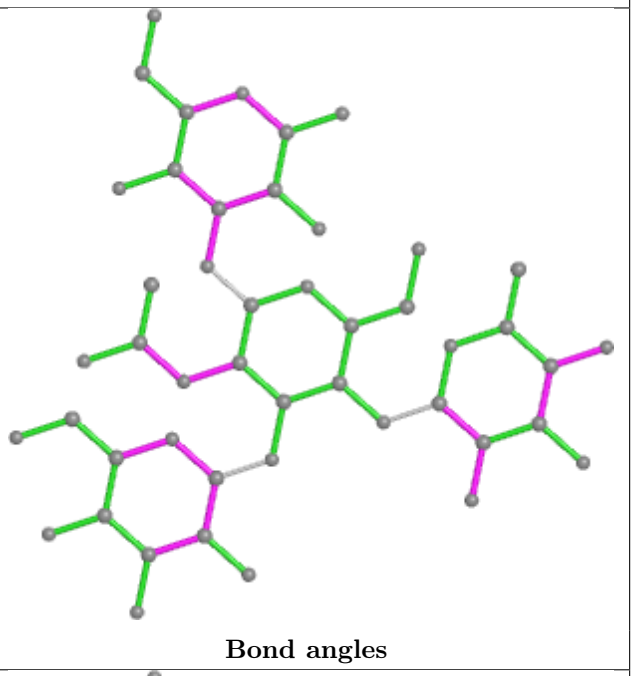


Rings

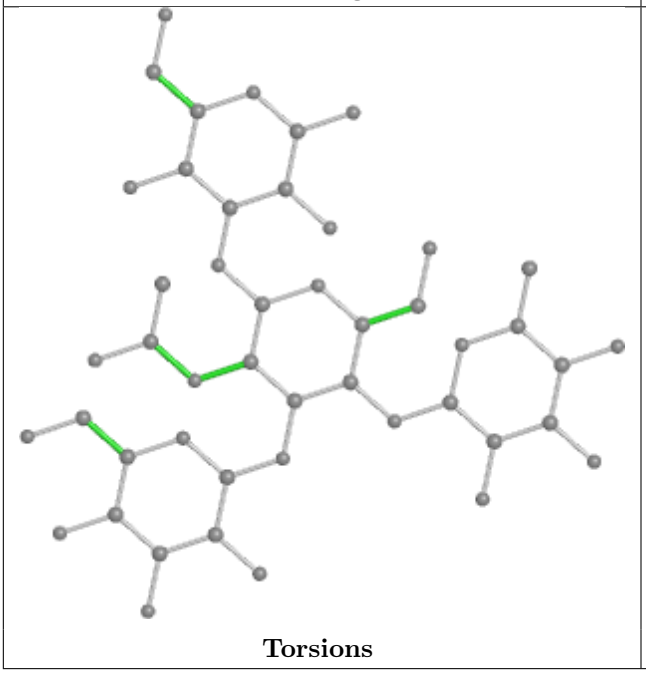
Oligosaccharide Chain G



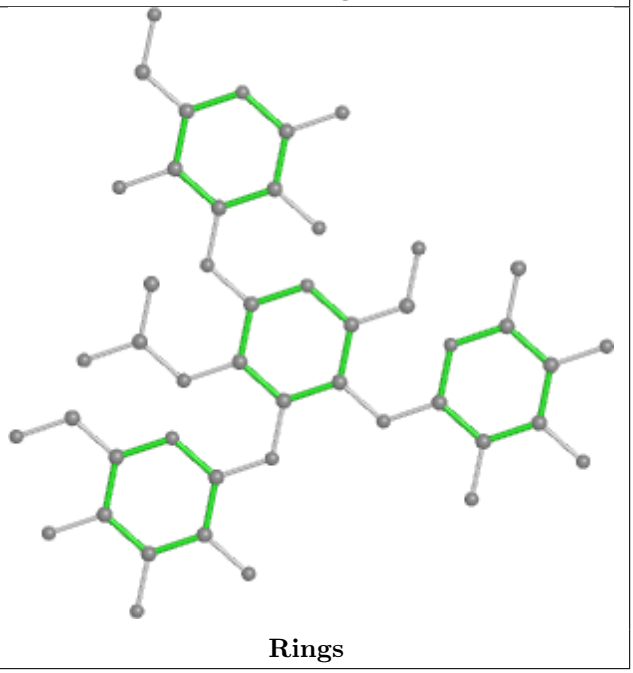
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

Of 13 ligands modelled in this entry, 8 are monoatomic - leaving 5 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	$\#\|Z\|>2$
6	SO4	A	211	-	4,4,4	0.54	0	6,6,6	0.85	0
7	PEG	D	210[A]	-	6,6,6	0.27	0	5,5,5	1.25	1 (20%)
6	SO4	B	211	-	4,4,4	0.57	0	6,6,6	0.18	0
6	SO4	C	211	-	4,4,4	0.83	0	6,6,6	0.61	0
7	PEG	D	210[B]	-	6,6,6	0.27	0	5,5,5	1.22	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
7	PEG	D	210[B]	-	-	4/4/4/4	-
7	PEG	D	210[A]	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	D	210[A]	PEG	C3-O2-C2	-2.24	103.57	113.29

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	D	210[A]	PEG	O2-C3-C4-O4
7	D	210[A]	PEG	O1-C1-C2-O2
7	D	210[B]	PEG	O1-C1-C2-O2
7	D	210[B]	PEG	O2-C3-C4-O4
7	D	210[B]	PEG	C4-C3-O2-C2
7	D	210[B]	PEG	C1-C2-O2-C3

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	210[A]	PEG	2	0
7	D	210[B]	PEG	1	0

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	114/114 (100%)	-0.26	0	100	100	6, 10, 16, 23	0
1	B	114/114 (100%)	-0.23	0	100	100	7, 10, 15, 21	3 (2%)
1	C	114/114 (100%)	-0.28	0	100	100	5, 8, 13, 22	2 (1%)
1	D	114/114 (100%)	-0.33	0	100	100	6, 8, 12, 14	2 (1%)
All	All	456/456 (100%)	-0.28	0	100	100	5, 9, 15, 23	7 (1%)

There are no RSRZ outliers to report.

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	GLA	E	1	10/12	0.78	0.21	25,29,30,30	8
2	GLA	H	1	12/12	0.84	0.21	21,30,34,40	1
4	GAL	G	1	12/12	0.86	0.13	17,21,26,32	0
3	NAG	F	1	15/15	0.92	0.21	11,17,31,31	1
2	NAG	E	2	14/15	0.92	0.15	12,18,28,29	0
4	NAG	G	2	14/15	0.93	0.11	9,12,26,26	0
2	GAL	E	3	11/12	0.94	0.09	17,19,23,24	0
2	NAG	H	2	14/15	0.95	0.13	9,13,22,24	0
3	GAL	F	2	11/12	0.95	0.13	16,19,22,23	0

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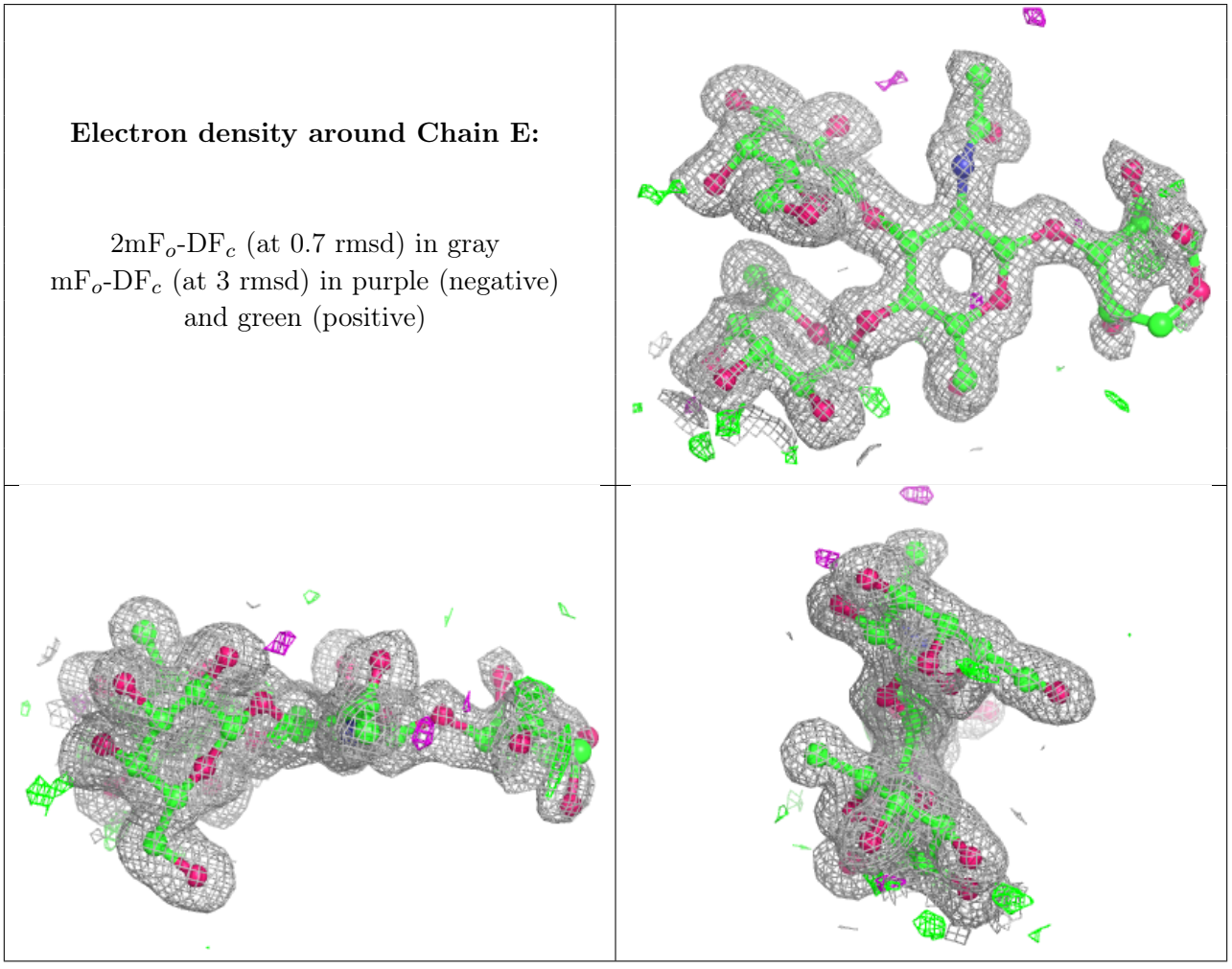
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	FUC	F	3	10/11	0.96	0.09	7,8,8,9	0
2	FUC	E	4	10/11	0.96	0.09	10,11,11,11	0
2	GAL	H	3	11/12	0.96	0.10	14,16,18,20	0
4	FUC	G	4	10/11	0.96	0.10	6,7,7,8	0
4	GAL	G	3	11/12	0.97	0.10	13,15,17,19	0
2	FUC	H	4	10/11	0.98	0.06	7,8,8,8	0

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.

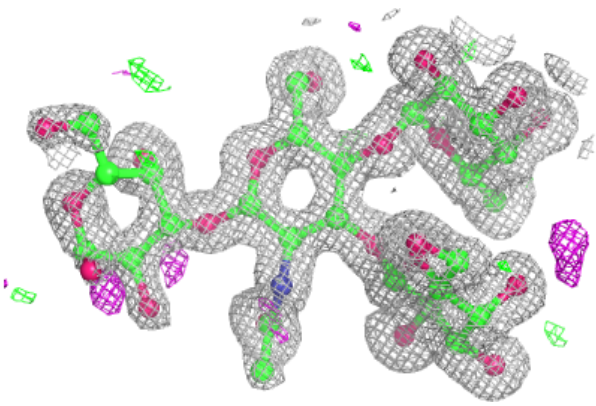
Electron density around Chain E:

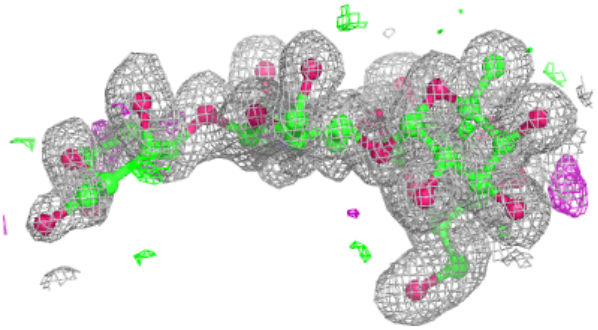
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

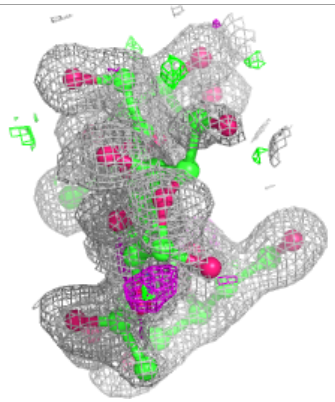


Electron density around Chain H:

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

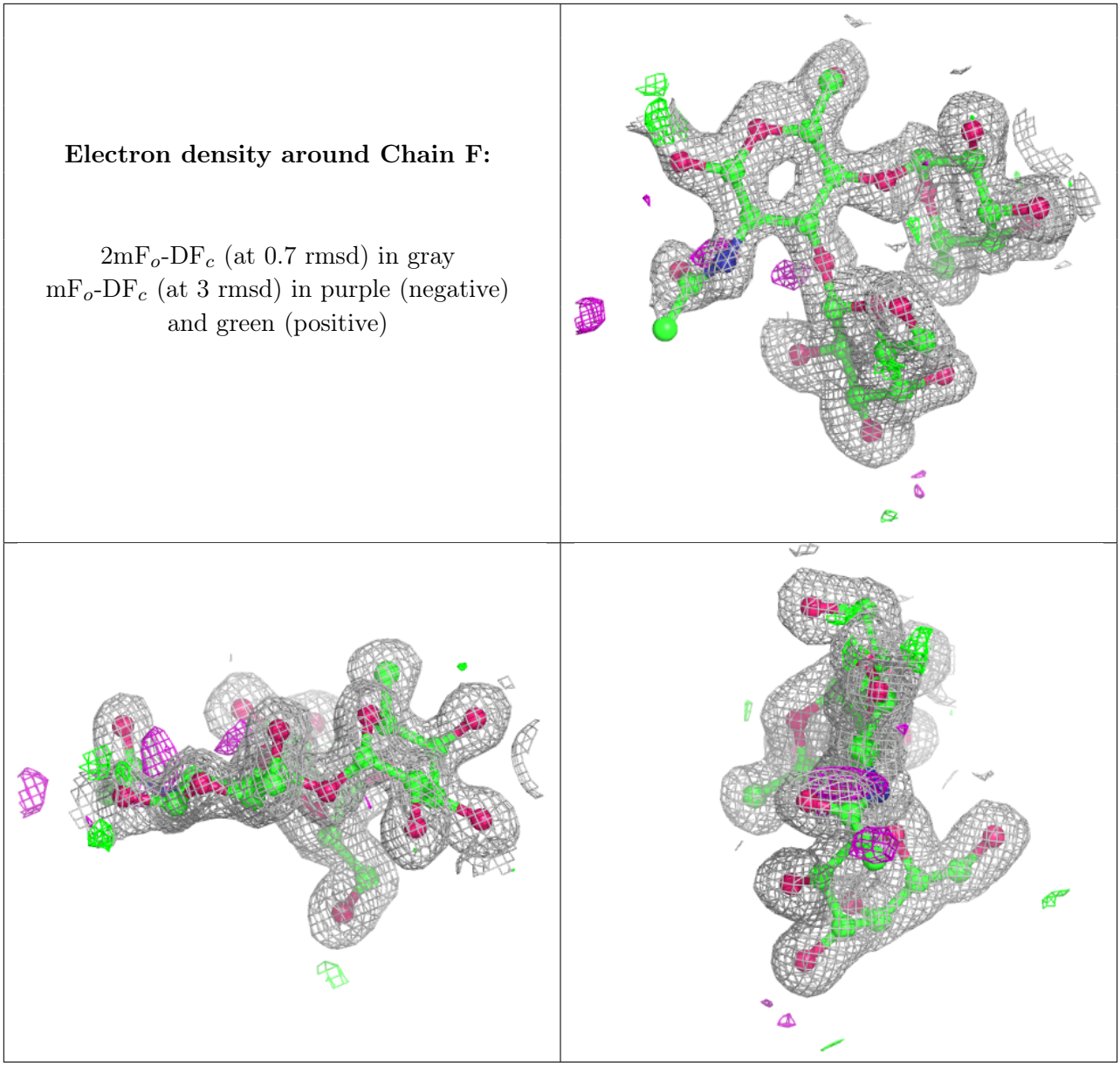


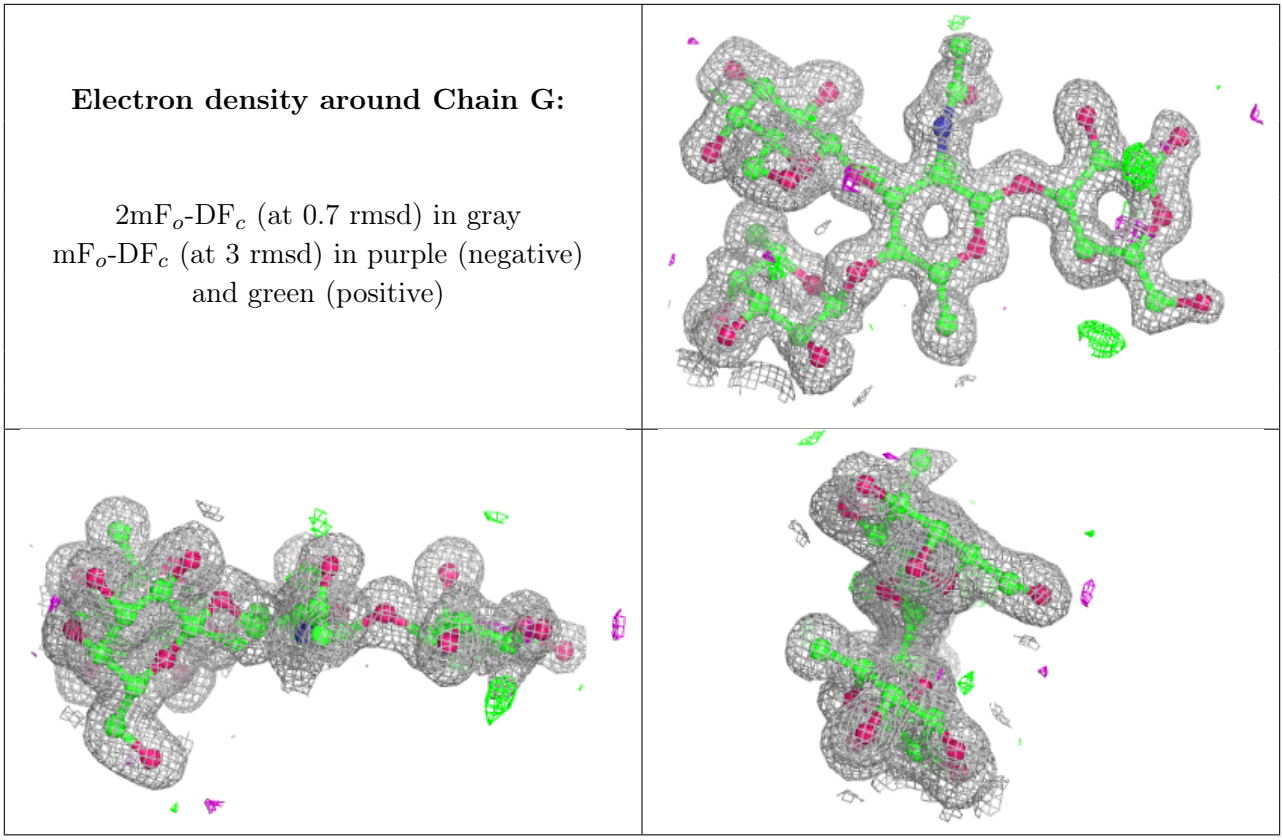




Electron density around Chain F:

$2mF_o$-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^{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
6	SO4	C	211	5/5	0.86	0.26	23,24,28,28	5
6	SO4	A	211	5/5	0.94	0.20	23,29,33,34	0
7	PEG	D	210[A]	7/7	0.94	0.13	18,21,28,31	7
7	PEG	D	210[B]	7/7	0.94	0.13	17,20,28,30	7
6	SO4	B	211	5/5	0.97	0.25	34,35,39,41	0
5	CA	A	200	1/1	0.99	0.04	9,9,9,9	0
5	CA	A	199	1/1	0.99	0.04	10,10,10,10	0
5	CA	D	200	1/1	1.00	0.04	7,7,7,7	0
5	CA	B	199	1/1	1.00	0.03	8,8,8,8	0
5	CA	B	200	1/1	1.00	0.04	7,7,7,7	0
5	CA	C	199	1/1	1.00	0.04	6,6,6,6	0
5	CA	C	200	1/1	1.00	0.04	6,6,6,6	0
5	CA	D	199	1/1	1.00	0.04	7,7,7,7	0

6.5 Other polymers (i)

There are no such residues in this entry.