wwPDB X-ray Structure Validation Summary Report (i)

Oct 1, 2024 – 04:59 am BST

PDB ID	:	8R49
Title	:	Plastidial phosphorylase Pho1 from Solanum tuberosum in complex with beta cyclodextrin
Authors	:	Koulas, S.M.; Leonidas, D.D.
Deposited on	:	2023-11-13
Resolution	:	2.80 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 (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 (1)) were used in the production of this report:

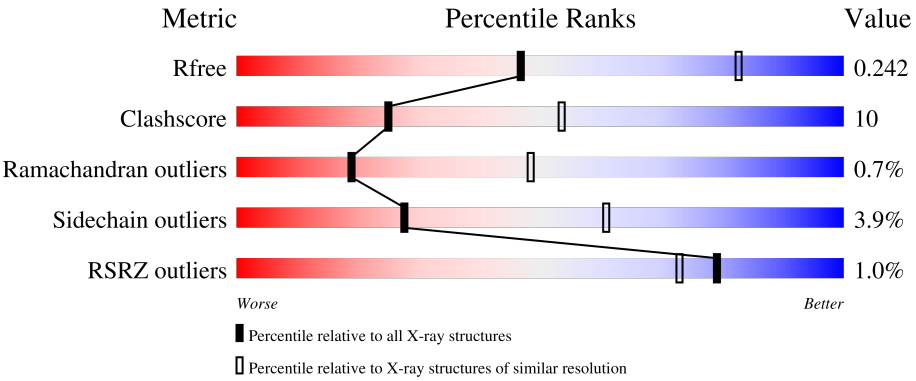
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.80 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}	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (2.80-2.80)

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	916	75% 14% • 9%
1	B	916	74% 15% • 9%
1	C	916	2% 69% 20% • 9%
2	D	7	29% 71%
2	E	7	57% 43%

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Mol	Chain	Length	Quality of chain
2	F	7	43% 57%

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
2	GLC	D	1	X	-	-	-
2	GLC	D	3	X	-	-	-
2	GLC	D	4	X	-	-	-
2	GLC	E	1	X	-	-	-
2	GLC	E	3	X	-	-	-
2	GLC	F	1	X	-	-	-
2	GLC	F	6	X	-	-	-

2 Entry composition (i)

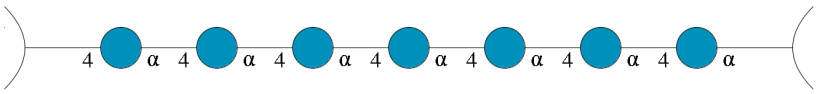
There are 4 unique types of molecules in this entry. The entry contains 20374 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alpha-1,4 glucan phosphorylase L-1 isozyme, chloroplastic/amyloplastic.

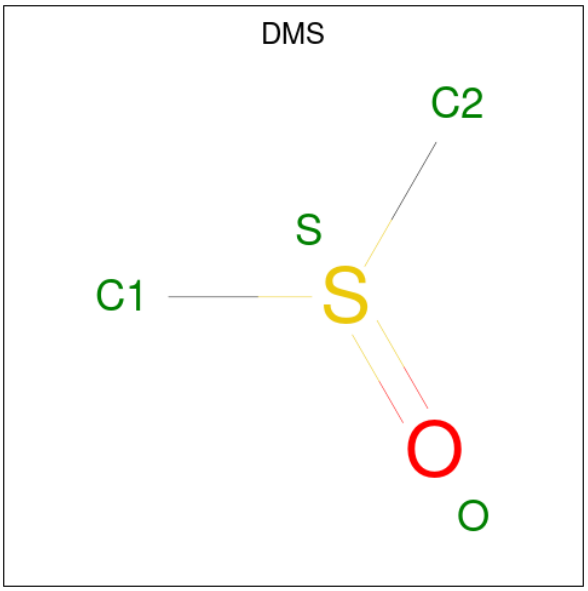
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	830	Total	C	N	O	P	S	0	0	0
			6670	4261	1119	1264	1	25			
1	B	830	Total	C	N	O	P	S	0	0	0
			6670	4261	1119	1264	1	25			
1	C	830	Total	C	N	O	P	S	0	0	0
			6670	4261	1119	1264	1	25			

- Molecule 2 is an oligosaccharide called Cycloheptakis-(1-4)-(alpha-D-glucopyranose).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	7	Total	C	O		0	0	0
			77	42	35				
2	E	7	Total	C	O		0	0	0
			77	42	35				
2	F	7	Total	C	O		0	0	0
			77	42	35				

- Molecule 3 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



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

- Molecule 4 is water.

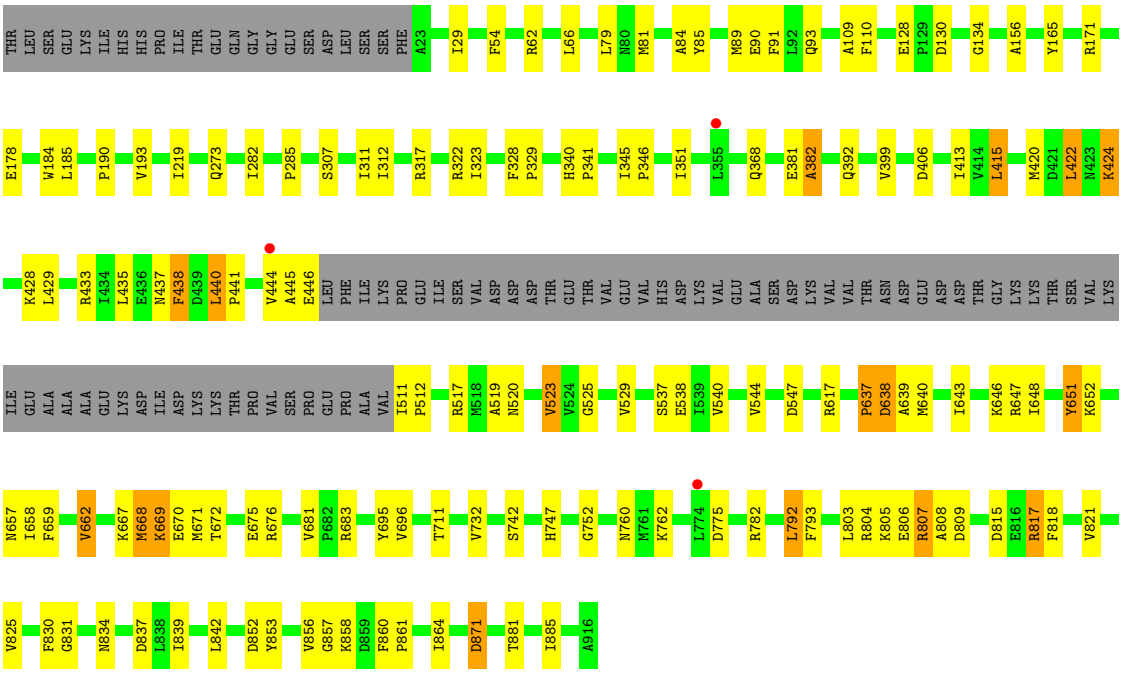
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	48	Total O 48 48	0	0
4	B	33	Total O 33 33	0	0
4	C	20	Total O 20 20	0	0

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.

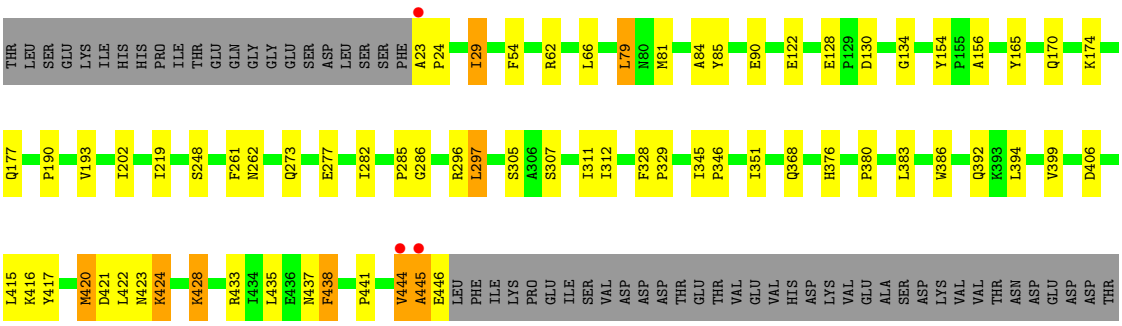
• Molecule 1: Alpha-1,4 glucan phosphorylase L-1 isozyme, chloroplastic/amyloplastic

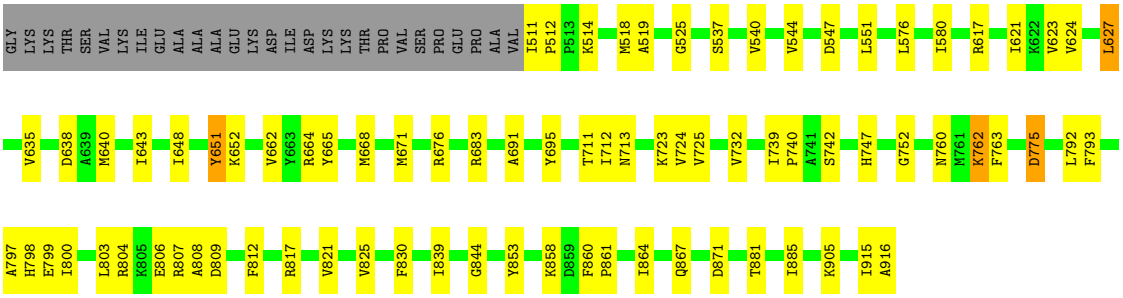
Chain A:



• Molecule 1: Alpha-1,4 glucan phosphorylase L-1 isozyme, chloroplastic/amyloplastic

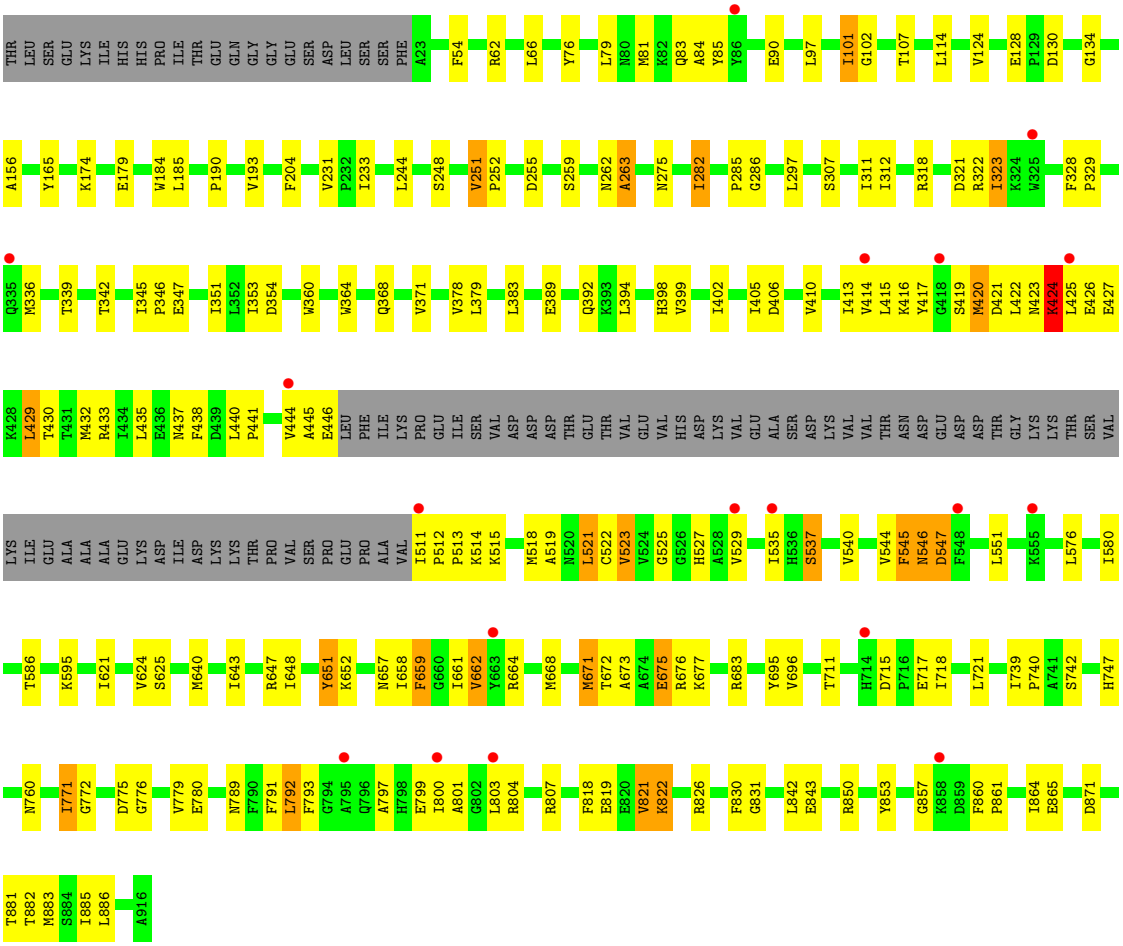
Chain B:





• Molecule 1: Alpha-1,4 glucan phosphorylase L-1 isozyme, chloroplastic/amyloplastic

Chain C:



• Molecule 2: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)

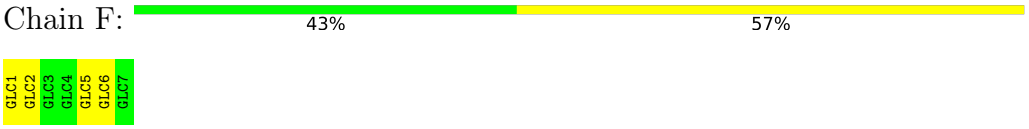
Chain D:

• Molecule 2: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)

Chain E:

• Molecule 2: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	218.63Å 136.73Å 123.16Å 90.00° 91.47° 90.00°	Depositor
Resolution (Å)	123.11 – 2.80 123.11 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.8 (123.11-2.80) 97.8 (123.11-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.66 (at 2.82Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.218 , 0.239 0.226 , 0.242	Depositor DCC
R_{free} test set	4496 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	77.1	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 62.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	20374	wwPDB-VP
Average B, all atoms (Å^2)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: LLP, GLC, DMS

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	0.31	0/6796	0.59	0/9195
1	B	0.31	0/6796	0.59	0/9195
1	C	0.31	0/6796	0.62	1/9195 (0.0%)
All	All	0.31	0/20388	0.60	1/27585 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	1
1	C	1	2
All	All	1	6

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	671	MET	CG-SD-CE	5.90	109.64	100.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	444	VAL	CA

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	317	ARG	Sidechain
1	A	817	ARG	Sidechain
1	A	831	GLY	Peptide
1	B	762	LLP	Mainchain
1	C	826	ARG	Sidechain

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	6670	0	6550	94	0
1	B	6670	0	6550	119	0
1	C	6670	0	6550	177	0
2	D	77	0	63	0	0
2	E	77	0	63	0	0
2	F	77	0	63	0	0
3	A	12	0	18	0	0
3	B	16	0	24	3	0
3	C	4	0	6	0	0
4	A	48	0	0	1	0
4	B	33	0	0	1	0
4	C	20	0	0	1	0
All	All	20374	0	19887	385	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 10.

The worst 5 of 385 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:C:342:THR:HG22	1:C:522:CYS:SG	1.91	1.09
1:A:668:MET:O	1:A:670:GLU:N	2.03	0.90
1:C:523:VAL:HG12	1:C:529:VAL:HG21	1.54	0.90
1:C:398:HIS:O	1:C:402:ILE:HD13	1.72	0.89
1:A:523:VAL:HG12	1:A:529:VAL:HG21	1.54	0.89

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	825/916 (90%)	771 (94%)	43 (5%)	11 (1%)	10	32
1	B	825/916 (90%)	778 (94%)	44 (5%)	3 (0%)	30	61
1	C	825/916 (90%)	775 (94%)	46 (6%)	4 (0%)	25	56
All	All	2475/2748 (90%)	2324 (94%)	133 (5%)	18 (1%)	19	48

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	445	ALA
1	A	637	PRO
1	A	638	ASP
1	A	668	MET
1	A	669	LYS

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	708/786 (90%)	687 (97%)	21 (3%)	36	70
1	B	708/786 (90%)	687 (97%)	21 (3%)	36	70
1	C	708/786 (90%)	667 (94%)	41 (6%)	17	45
All	All	2124/2358 (90%)	2041 (96%)	83 (4%)	27	61

5 of 83 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	430	THR
1	C	675	GLU
1	C	521	LEU
1	C	547	ASP
1	C	775	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	649	HIS
1	C	697	GLN
1	C	845	ASN
1	B	376	HIS
1	B	262	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)

3 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	LLP	A	762	1	23,24,25	0.54	0	25,32,34	0.60	0
1	LLP	C	762	1	23,24,25	0.49	0	25,32,34	0.55	0
1	LLP	B	762	1	23,24,25	0.47	0	25,32,34	0.67	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
1	LLP	A	762	1	-	5/16/17/19	0/1/1/1
1	LLP	C	762	1	-	3/16/17/19	0/1/1/1
1	LLP	B	762	1	-	4/16/17/19	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	762	LLP	C5'-OP4-P-OP1
1	A	762	LLP	C5'-OP4-P-OP2
1	A	762	LLP	C5'-OP4-P-OP3
1	B	762	LLP	C4-C5-C5'-OP4
1	B	762	LLP	C5'-OP4-P-OP2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	762	LLP	1	0
1	B	762	LLP	2	0

5.5 Carbohydrates (i)

21 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	GLC	D	1	2	11,11,12	0.46	0	15,15,17	0.83	0
2	GLC	D	2	2	11,11,12	0.50	0	15,15,17	0.72	0
2	GLC	D	3	2	11,11,12	0.61	0	15,15,17	0.97	1 (6%)
2	GLC	D	4	2	11,11,12	0.51	0	15,15,17	0.95	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	GLC	D	5	2	11,11,12	0.38	0	15,15,17	1.05	2 (13%)
2	GLC	D	6	2	11,11,12	0.36	0	15,15,17	0.99	1 (6%)
2	GLC	D	7	2	11,11,12	0.50	0	15,15,17	1.05	1 (6%)
2	GLC	E	1	2	11,11,12	0.48	0	15,15,17	0.81	1 (6%)
2	GLC	E	2	2	11,11,12	0.41	0	15,15,17	0.81	0
2	GLC	E	3	2	11,11,12	0.43	0	15,15,17	0.62	0
2	GLC	E	4	2	11,11,12	0.36	0	15,15,17	0.71	0
2	GLC	E	5	2	11,11,12	0.32	0	15,15,17	0.88	1 (6%)
2	GLC	E	6	2	11,11,12	0.43	0	15,15,17	0.76	0
2	GLC	E	7	2	11,11,12	0.46	0	15,15,17	1.21	2 (13%)
2	GLC	F	1	2	11,11,12	0.50	0	15,15,17	1.18	2 (13%)
2	GLC	F	2	2	11,11,12	0.54	0	15,15,17	0.91	1 (6%)
2	GLC	F	3	2	11,11,12	0.43	0	15,15,17	0.81	0
2	GLC	F	4	2	11,11,12	0.41	0	15,15,17	0.69	0
2	GLC	F	5	2	11,11,12	0.51	0	15,15,17	1.45	2 (13%)
2	GLC	F	6	2	11,11,12	0.43	0	15,15,17	1.50	2 (13%)
2	GLC	F	7	2	11,11,12	0.50	0	15,15,17	0.84	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	GLC	D	1	2	1/1/4/5	2/2/19/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1
2	GLC	D	3	2	1/1/4/5	0/2/19/22	0/1/1/1
2	GLC	D	4	2	1/1/4/5	2/2/19/22	0/1/1/1
2	GLC	D	5	2	-	1/2/19/22	0/1/1/1
2	GLC	D	6	2	-	1/2/19/22	0/1/1/1
2	GLC	D	7	2	-	2/2/19/22	0/1/1/1
2	GLC	E	1	2	1/1/4/5	2/2/19/22	0/1/1/1
2	GLC	E	2	2	-	2/2/19/22	0/1/1/1
2	GLC	E	3	2	1/1/4/5	2/2/19/22	0/1/1/1
2	GLC	E	4	2	-	1/2/19/22	0/1/1/1
2	GLC	E	5	2	-	0/2/19/22	0/1/1/1
2	GLC	E	6	2	-	0/2/19/22	0/1/1/1
2	GLC	E	7	2	-	0/2/19/22	0/1/1/1
2	GLC	F	1	2	1/1/4/5	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	F	2	2	-	1/2/19/22	0/1/1/1
2	GLC	F	3	2	-	0/2/19/22	0/1/1/1
2	GLC	F	4	2	-	1/2/19/22	0/1/1/1
2	GLC	F	5	2	-	0/2/19/22	0/1/1/1
2	GLC	F	6	2	1/1/4/5	0/2/19/22	0/1/1/1
2	GLC	F	7	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	5	GLC	C1-O5-C5	4.10	117.75	112.19
2	F	6	GLC	C1-O5-C5	3.95	117.55	112.19
2	F	6	GLC	C1-C2-C3	3.50	113.97	109.67
2	F	1	GLC	C1-C2-C3	3.42	113.87	109.67
2	F	2	GLC	C1-O5-C5	3.01	116.27	112.19

5 of 7 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	1	GLC	C1
2	D	3	GLC	C1
2	D	4	GLC	C1
2	E	1	GLC	C1
2	E	3	GLC	C1

5 of 19 torsion outliers are listed below:

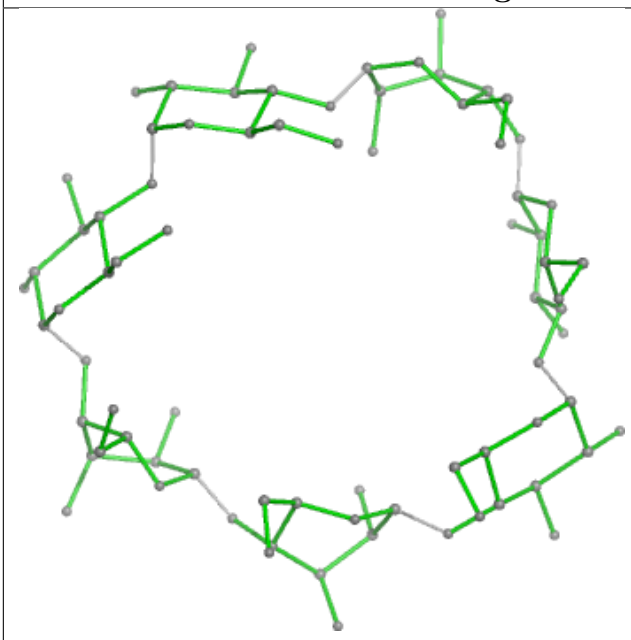
Mol	Chain	Res	Type	Atoms
2	D	1	GLC	O5-C5-C6-O6
2	E	1	GLC	O5-C5-C6-O6
2	D	4	GLC	O5-C5-C6-O6
2	E	1	GLC	C4-C5-C6-O6
2	F	7	GLC	O5-C5-C6-O6

There are no ring outliers.

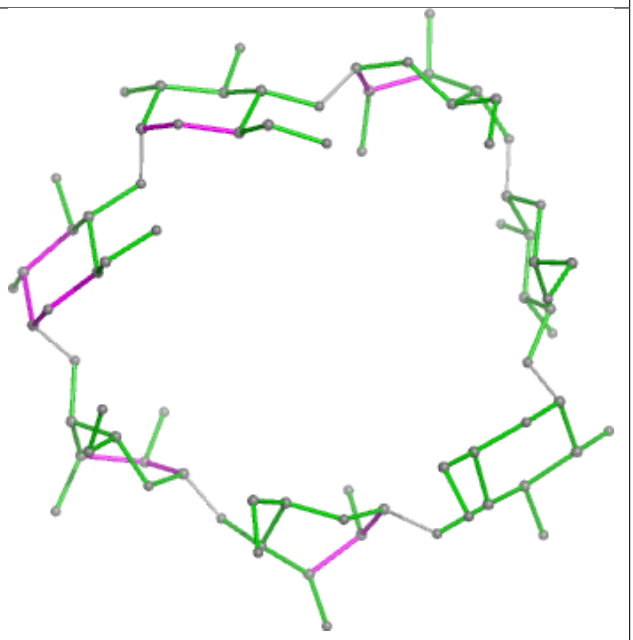
No monomer is involved in short contacts.

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

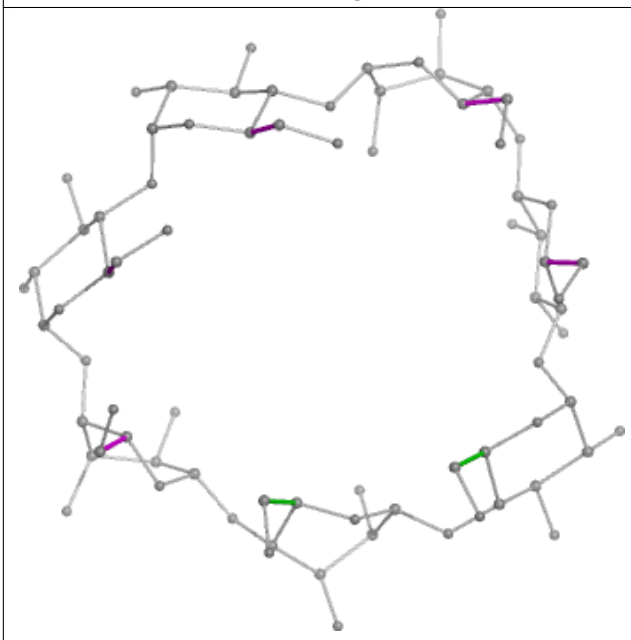
Oligosaccharide Chain D



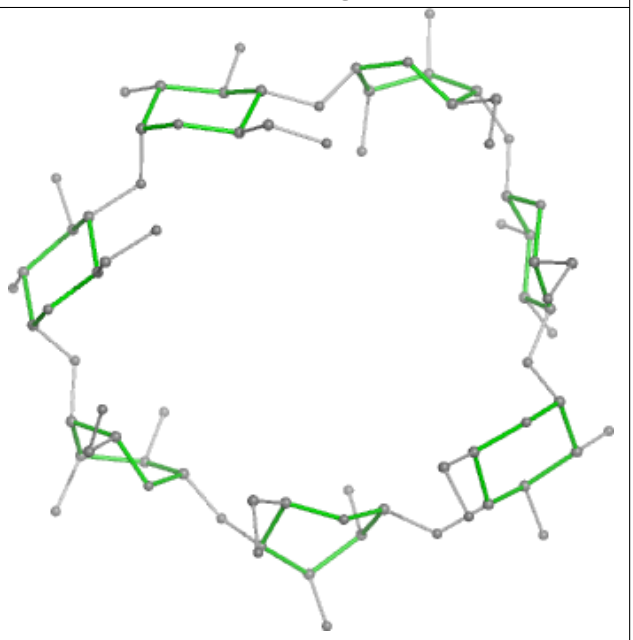
Bond lengths



Bond angles

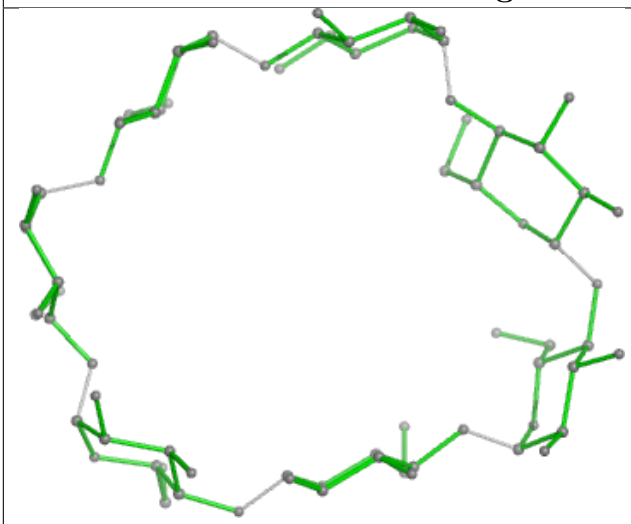


Torsions

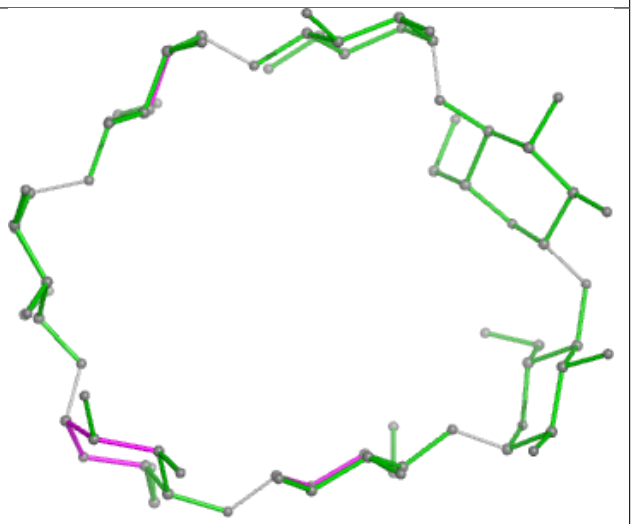


Rings

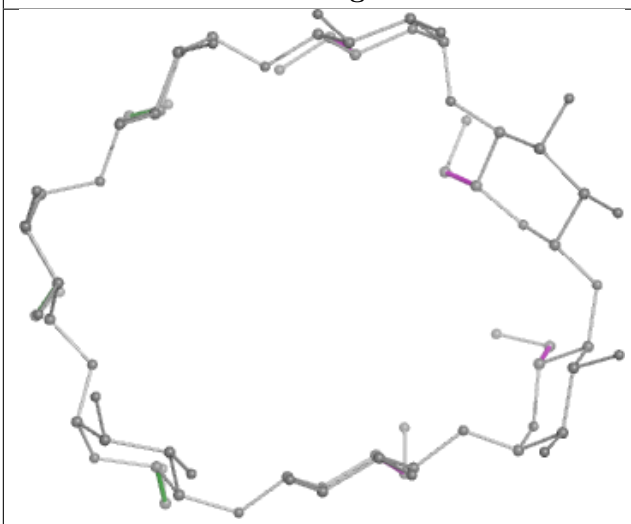
Oligosaccharide Chain E



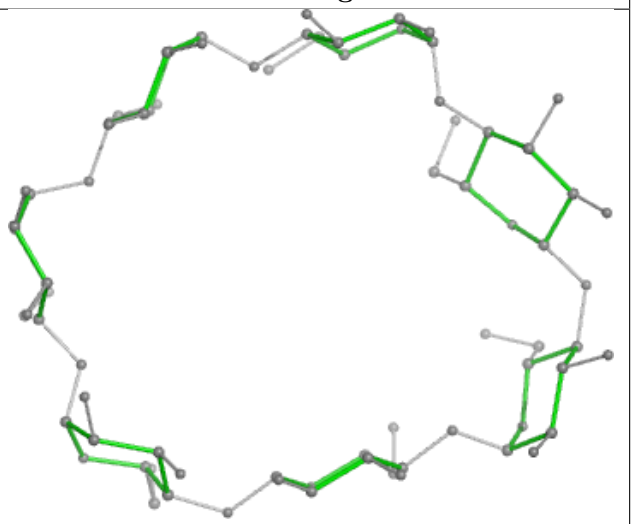
Bond lengths



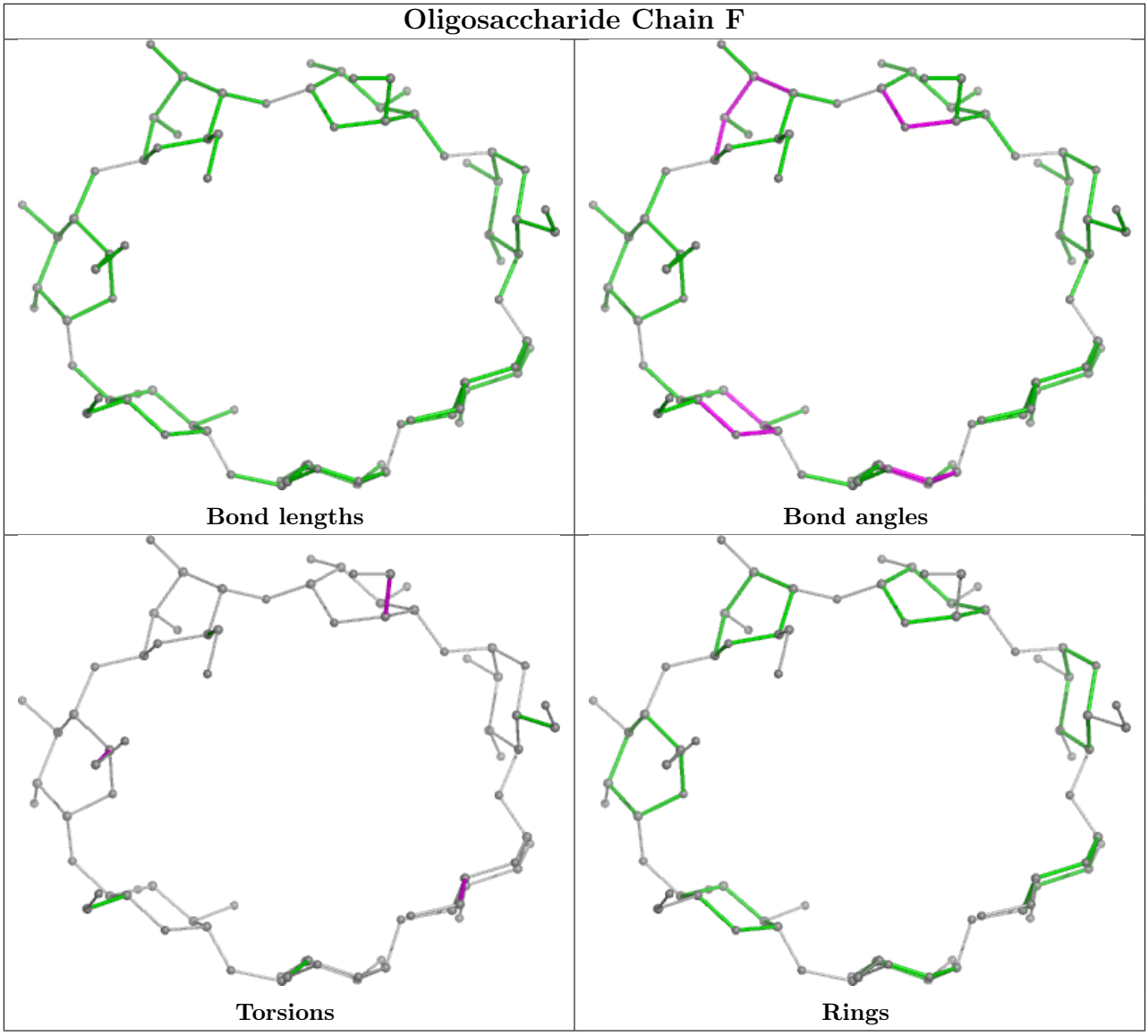
Bond angles



Torsions



Rings



5.6 Ligand geometry

8 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	$\#\|Z\| > 2$
3	DMS	A	1002	-	3,3,3	0.26	0	3,3,3	0.05	0
3	DMS	A	1003	-	3,3,3	0.29	0	3,3,3	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	DMS	A	1001	-	3,3,3	0.27	0	3,3,3	0.06	0
3	DMS	B	1002	-	3,3,3	0.24	0	3,3,3	0.11	0
3	DMS	B	1001	-	3,3,3	0.18	0	3,3,3	0.15	0
3	DMS	B	1003	-	3,3,3	0.26	0	3,3,3	0.05	0
3	DMS	B	1004	-	3,3,3	0.27	0	3,3,3	0.09	0
3	DMS	C	1001	-	3,3,3	0.29	0	3,3,3	0.07	0

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.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1001	DMS	3	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	829/916 (90%)	-0.04	3 (0%)	89	85	50, 78, 113, 143	0
1	B	829/916 (90%)	0.01	3 (0%)	89	85	56, 82, 117, 155	0
1	C	829/916 (90%)	0.33	18 (2%)	62	53	63, 98, 138, 174	0
All	All	2487/2748 (90%)	0.10	24 (0%)	79	73	50, 85, 128, 174	0

The worst 5 of 24 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	511	ILE	3.2
1	C	529	VAL	3.1
1	B	445	ALA	3.0
1	B	444	VAL	3.0
1	C	548	PHE	3.0

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

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
1	LLP	C	762	24/25	0.95	0.09	71,83,91,94	0
1	LLP	B	762	24/25	0.96	0.09	65,70,73,78	0
1	LLP	A	762	24/25	0.96	0.11	59,63,72,73	0

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,

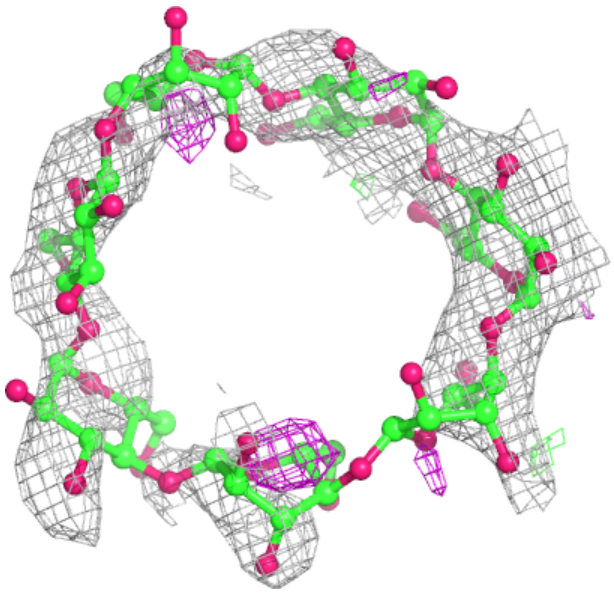
median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

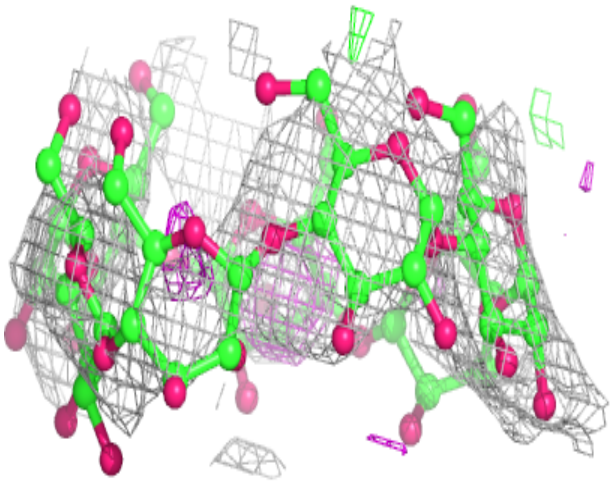
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GLC	F	1	11/12	0.39	0.13	174,190,197,197	0
2	GLC	F	6	11/12	0.40	0.12	170,178,181,182	0
2	GLC	E	7	11/12	0.41	0.13	157,167,175,180	0
2	GLC	F	2	11/12	0.44	0.13	180,194,198,199	0
2	GLC	E	1	11/12	0.48	0.12	167,182,185,186	0
2	GLC	E	2	11/12	0.55	0.13	160,169,177,178	0
2	GLC	D	3	11/12	0.55	0.13	164,178,182,185	0
2	GLC	F	3	11/12	0.56	0.13	183,190,195,197	0
2	GLC	D	4	11/12	0.56	0.15	155,167,177,177	0
2	GLC	F	7	11/12	0.57	0.10	180,184,188,193	0
2	GLC	E	6	11/12	0.60	0.10	148,158,164,166	0
2	GLC	F	5	11/12	0.60	0.12	166,171,175,178	0
2	GLC	D	1	11/12	0.60	0.10	170,177,181,182	0
2	GLC	D	7	11/12	0.61	0.12	166,172,176,178	0
2	GLC	D	2	11/12	0.61	0.10	173,181,183,187	0
2	GLC	E	5	11/12	0.68	0.11	139,150,153,156	0
2	GLC	D	6	11/12	0.73	0.11	145,154,160,164	0
2	GLC	E	3	11/12	0.73	0.12	157,165,170,171	0
2	GLC	D	5	11/12	0.75	0.13	147,156,161,161	0
2	GLC	E	4	11/12	0.75	0.11	140,145,150,151	0
2	GLC	F	4	11/12	0.77	0.11	166,171,174,178	0

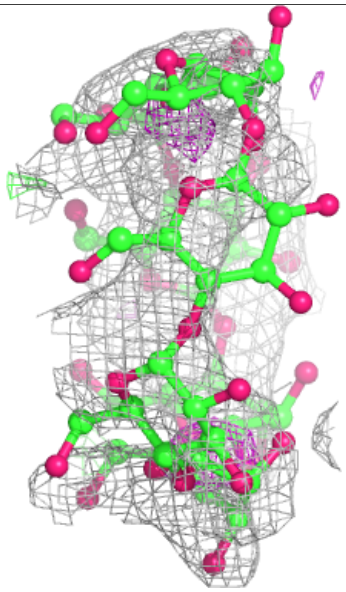
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain D:

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

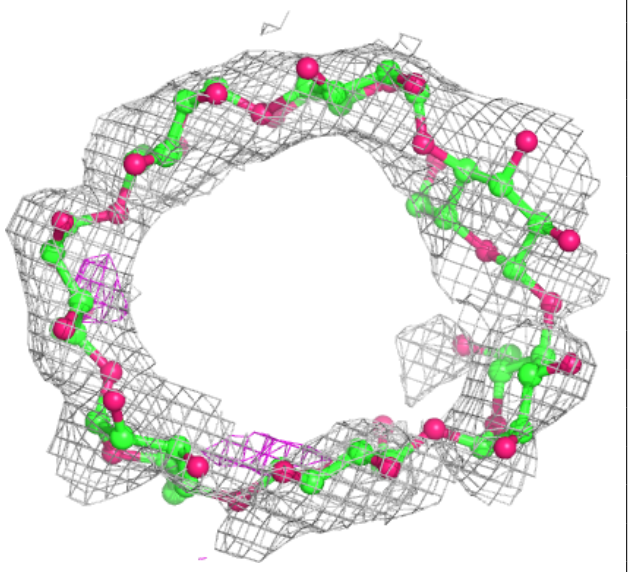


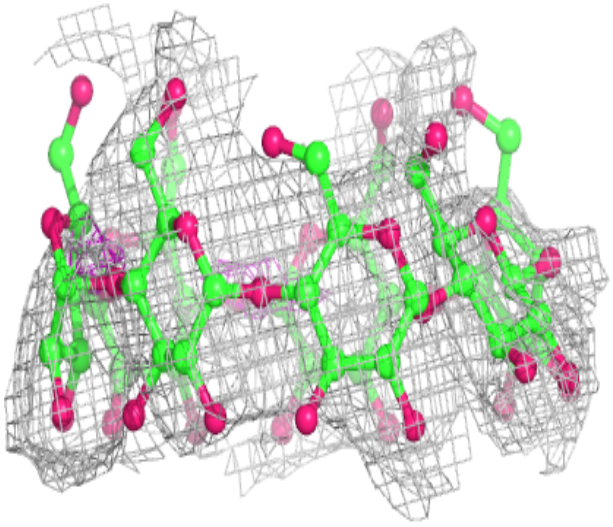


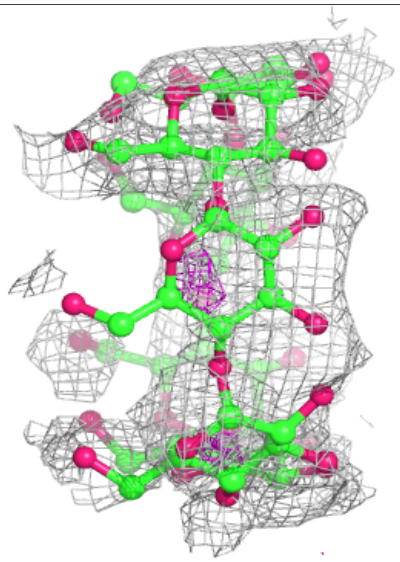


Electron density around Chain E:

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

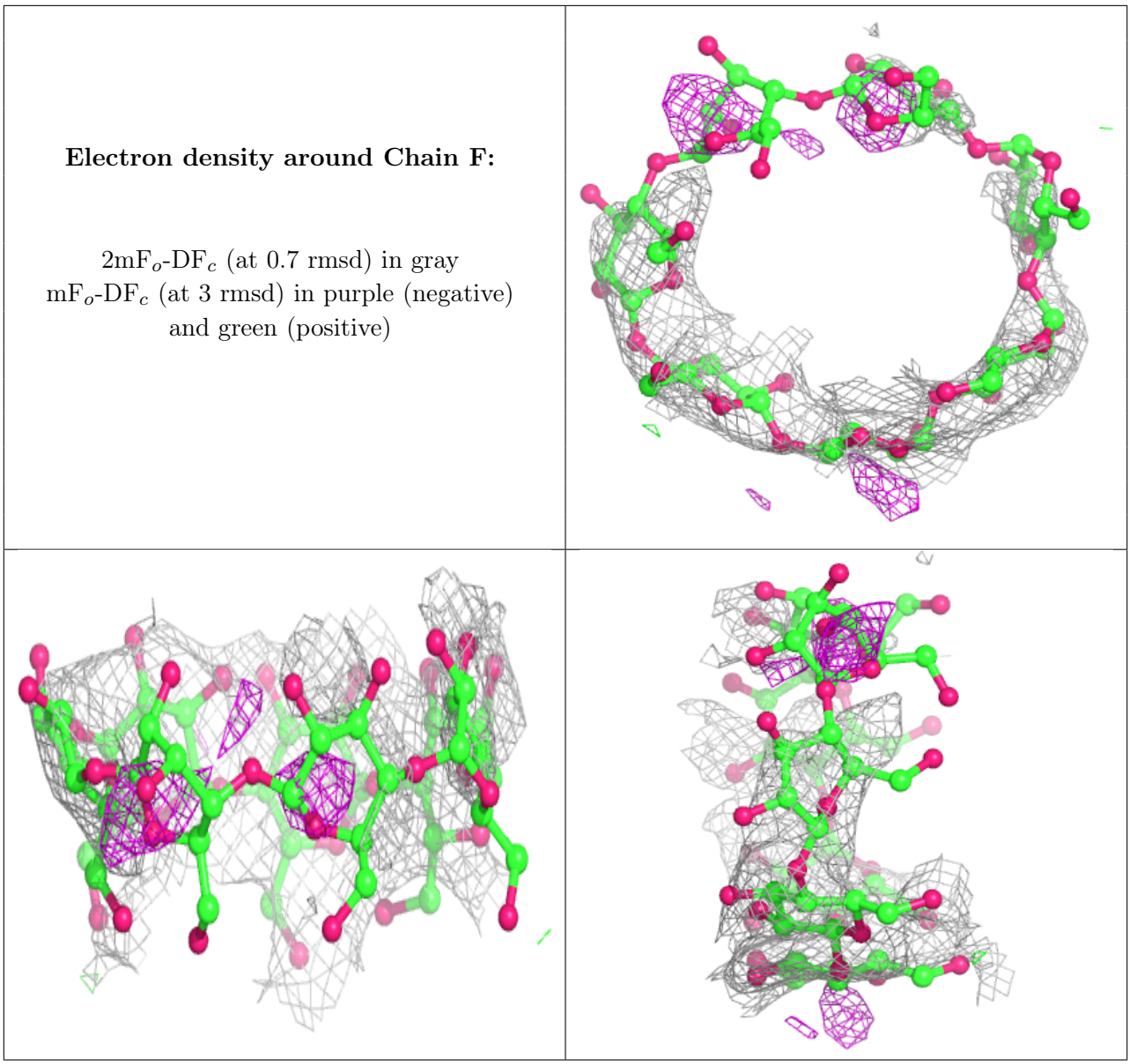






Electron density around Chain F:

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



6.4 Ligands

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	DMS	C	1001	4/4	0.71	0.27	97,106,107,109	0
3	DMS	A	1001	4/4	0.78	0.30	118,118,118,120	0
3	DMS	B	1003	4/4	0.82	0.22	118,120,121,122	0
3	DMS	A	1003	4/4	0.82	0.23	86,90,91,91	0
3	DMS	A	1002	4/4	0.83	0.20	125,125,130,134	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	DMS	B	1004	4/4	0.87	0.28	102,109,112,112	0
3	DMS	B	1002	4/4	0.87	0.24	121,121,121,124	0
3	DMS	B	1001	4/4	0.91	0.21	93,98,99,107	0

6.5 Other polymers (i)

There are no such residues in this entry.