



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 03:35 AM BST

PDB ID : 4BLB
Title : Crystal structure of a human Suppressor of fused (SUFU)-GLI1p complex
Authors : Cherry, A.L.; Finta, C.; Karlstrom, M.; De Sanctis, D.; Toftgard, R.; Jovine, L.
Deposited on : 2013-05-02
Resolution : 2.80 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

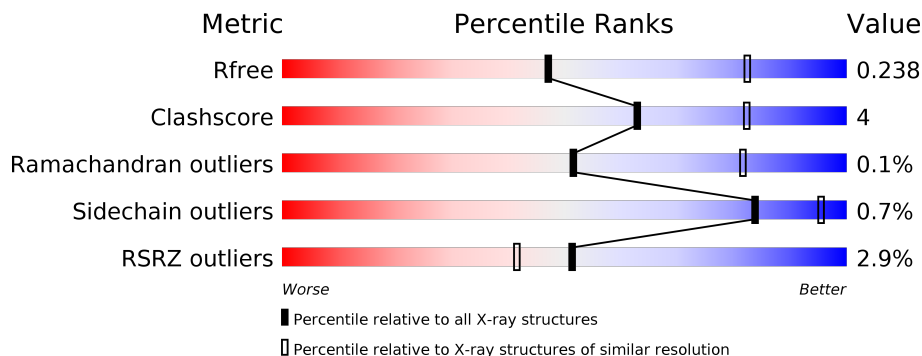
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




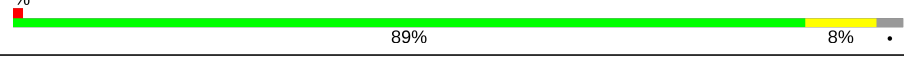


The reported resolution of this entry is 2.80 Å.

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.







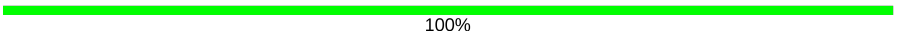

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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 $\leq 5\%$. 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	753	
1	B	753	
1	C	753	
1	D	753	
2	E	17	
2	F	17	

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Mol	Chain	Length	Quality of chain
2	G	17	
2	H	17	
3	I	2	
3	J	2	
3	K	2	
3	L	2	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23141 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 MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	734	5757	3681	967	1090	19	0	0	0
1	B	717	5621	3597	942	1064	18	0	0	0
1	C	725	5687	3640	957	1072	18	0	0	0
1	D	730	5727	3666	962	1081	18	0	0	0

There are 128 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP P0AEX9
A	3	THR	ILE	engineered mutation	UNP P0AEX9
A	360	ALA	GLU	engineered mutation	UNP P0AEX9
A	363	ALA	LYS	engineered mutation	UNP P0AEX9
A	364	ALA	ASP	engineered mutation	UNP P0AEX9
A	368	ASN	ARG	engineered mutation	UNP P0AEX9
A	369	ALA	-	linker	UNP P0AEX9
A	370	ALA	-	linker	UNP P0AEX9
A	371	ALA	-	linker	UNP P0AEX9
A	216	HIS	ALA	engineered mutation	UNP P0AEX9
A	220	HIS	LYS	engineered mutation	UNP P0AEX9
A	401	ASP	TRP	engineered mutation	UNP Q9UMX1
A	402	SER	LEU	engineered mutation	UNP Q9UMX1
A	403	PHE	GLY	engineered mutation	UNP Q9UMX1
A	619	PRO	-	linker	UNP Q9UMX1
A	620	SER	-	linker	UNP Q9UMX1
A	621	ARG	-	linker	UNP Q9UMX1
A	622	GLY	-	linker	UNP Q9UMX1
A	625	PRO	-	linker	UNP Q9UMX1
A	718	ALA	PRO	engineered mutation	UNP Q9UMX1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLU	deletion	UNP Q9UMX1
A	?	-	GLU	deletion	UNP Q9UMX1
A	?	-	PHE	deletion	UNP Q9UMX1
A	719	ALA	LYS	engineered mutation	UNP Q9UMX1
A	746	VAL	-	expression tag	UNP Q9UMX1
A	747	GLU	-	expression tag	UNP Q9UMX1
A	748	HIS	-	expression tag	UNP Q9UMX1
A	749	HIS	-	expression tag	UNP Q9UMX1
A	750	HIS	-	expression tag	UNP Q9UMX1
A	751	HIS	-	expression tag	UNP Q9UMX1
A	752	HIS	-	expression tag	UNP Q9UMX1
A	753	HIS	-	expression tag	UNP Q9UMX1
B	1	MET	-	expression tag	UNP P0AEX9
B	3	THR	ILE	engineered mutation	UNP P0AEX9
B	360	ALA	GLU	engineered mutation	UNP P0AEX9
B	363	ALA	LYS	engineered mutation	UNP P0AEX9
B	364	ALA	ASP	engineered mutation	UNP P0AEX9
B	368	ASN	ARG	engineered mutation	UNP P0AEX9
B	369	ALA	-	linker	UNP P0AEX9
B	370	ALA	-	linker	UNP P0AEX9
B	371	ALA	-	linker	UNP P0AEX9
B	216	HIS	ALA	engineered mutation	UNP P0AEX9
B	220	HIS	LYS	engineered mutation	UNP P0AEX9
B	401	ASP	TRP	engineered mutation	UNP Q9UMX1
B	402	SER	LEU	engineered mutation	UNP Q9UMX1
B	403	PHE	GLY	engineered mutation	UNP Q9UMX1
B	619	PRO	-	linker	UNP Q9UMX1
B	620	SER	-	linker	UNP Q9UMX1
B	621	ARG	-	linker	UNP Q9UMX1
B	622	GLY	-	linker	UNP Q9UMX1
B	625	PRO	-	linker	UNP Q9UMX1
B	718	ALA	PRO	engineered mutation	UNP Q9UMX1
B	?	-	GLU	deletion	UNP Q9UMX1
B	?	-	GLU	deletion	UNP Q9UMX1
B	?	-	PHE	deletion	UNP Q9UMX1
B	719	ALA	LYS	engineered mutation	UNP Q9UMX1
B	746	VAL	-	expression tag	UNP Q9UMX1
B	747	GLU	-	expression tag	UNP Q9UMX1
B	748	HIS	-	expression tag	UNP Q9UMX1
B	749	HIS	-	expression tag	UNP Q9UMX1
B	750	HIS	-	expression tag	UNP Q9UMX1
B	751	HIS	-	expression tag	UNP Q9UMX1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	752	HIS	-	expression tag	UNP Q9UMX1
B	753	HIS	-	expression tag	UNP Q9UMX1
C	1	MET	-	expression tag	UNP P0AEX9
C	3	THR	ILE	engineered mutation	UNP P0AEX9
C	360	ALA	GLU	engineered mutation	UNP P0AEX9
C	363	ALA	LYS	engineered mutation	UNP P0AEX9
C	364	ALA	ASP	engineered mutation	UNP P0AEX9
C	368	ASN	ARG	engineered mutation	UNP P0AEX9
C	369	ALA	-	linker	UNP P0AEX9
C	370	ALA	-	linker	UNP P0AEX9
C	371	ALA	-	linker	UNP P0AEX9
C	216	HIS	ALA	engineered mutation	UNP P0AEX9
C	220	HIS	LYS	engineered mutation	UNP P0AEX9
C	401	ASP	TRP	engineered mutation	UNP Q9UMX1
C	402	SER	LEU	engineered mutation	UNP Q9UMX1
C	403	PHE	GLY	engineered mutation	UNP Q9UMX1
C	619	PRO	-	linker	UNP Q9UMX1
C	620	SER	-	linker	UNP Q9UMX1
C	621	ARG	-	linker	UNP Q9UMX1
C	622	GLY	-	linker	UNP Q9UMX1
C	625	PRO	-	linker	UNP Q9UMX1
C	718	ALA	PRO	engineered mutation	UNP Q9UMX1
C	?	-	GLU	deletion	UNP Q9UMX1
C	?	-	GLU	deletion	UNP Q9UMX1
C	?	-	PHE	deletion	UNP Q9UMX1
C	719	ALA	LYS	engineered mutation	UNP Q9UMX1
C	746	VAL	-	expression tag	UNP Q9UMX1
C	747	GLU	-	expression tag	UNP Q9UMX1
C	748	HIS	-	expression tag	UNP Q9UMX1
C	749	HIS	-	expression tag	UNP Q9UMX1
C	750	HIS	-	expression tag	UNP Q9UMX1
C	751	HIS	-	expression tag	UNP Q9UMX1
C	752	HIS	-	expression tag	UNP Q9UMX1
C	753	HIS	-	expression tag	UNP Q9UMX1
D	1	MET	-	expression tag	UNP P0AEX9
D	3	THR	ILE	engineered mutation	UNP P0AEX9
D	360	ALA	GLU	engineered mutation	UNP P0AEX9
D	363	ALA	LYS	engineered mutation	UNP P0AEX9
D	364	ALA	ASP	engineered mutation	UNP P0AEX9
D	368	ASN	ARG	engineered mutation	UNP P0AEX9
D	369	ALA	-	linker	UNP P0AEX9
D	370	ALA	-	linker	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	371	ALA	-	linker	UNP P0AEX9
D	216	HIS	ALA	engineered mutation	UNP P0AEX9
D	220	HIS	LYS	engineered mutation	UNP P0AEX9
D	401	ASP	TRP	engineered mutation	UNP Q9UMX1
D	402	SER	LEU	engineered mutation	UNP Q9UMX1
D	403	PHE	GLY	engineered mutation	UNP Q9UMX1
D	619	PRO	-	linker	UNP Q9UMX1
D	620	SER	-	linker	UNP Q9UMX1
D	621	ARG	-	linker	UNP Q9UMX1
D	622	GLY	-	linker	UNP Q9UMX1
D	625	PRO	-	linker	UNP Q9UMX1
D	718	ALA	PRO	engineered mutation	UNP Q9UMX1
D	?	-	GLU	deletion	UNP Q9UMX1
D	?	-	GLU	deletion	UNP Q9UMX1
D	?	-	PHE	deletion	UNP Q9UMX1
D	719	ALA	LYS	engineered mutation	UNP Q9UMX1
D	746	VAL	-	expression tag	UNP Q9UMX1
D	747	GLU	-	expression tag	UNP Q9UMX1
D	748	HIS	-	expression tag	UNP Q9UMX1
D	749	HIS	-	expression tag	UNP Q9UMX1
D	750	HIS	-	expression tag	UNP Q9UMX1
D	751	HIS	-	expression tag	UNP Q9UMX1
D	752	HIS	-	expression tag	UNP Q9UMX1
D	753	HIS	-	expression tag	UNP Q9UMX1

- Molecule 2 is a protein called ZINC FINGER PROTEIN GLI1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	9	Total	C	N	O	0	0	0
			65	41	11	13			
2	F	9	Total	C	N	O	0	0	0
			65	41	11	13			
2	G	8	Total	C	N	O	0	0	0
			58	37	10	11			
2	H	9	Total	C	N	O	0	0	0
			65	41	11	13			

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	I	2	Total	C	O	0	0	0
			23	12	11			
3	J	2	Total	C	O	0	0	0
			23	12	11			
3	K	2	Total	C	O	0	0	0
			23	12	11			
3	L	2	Total	C	O	0	0	0
			23	12	11			

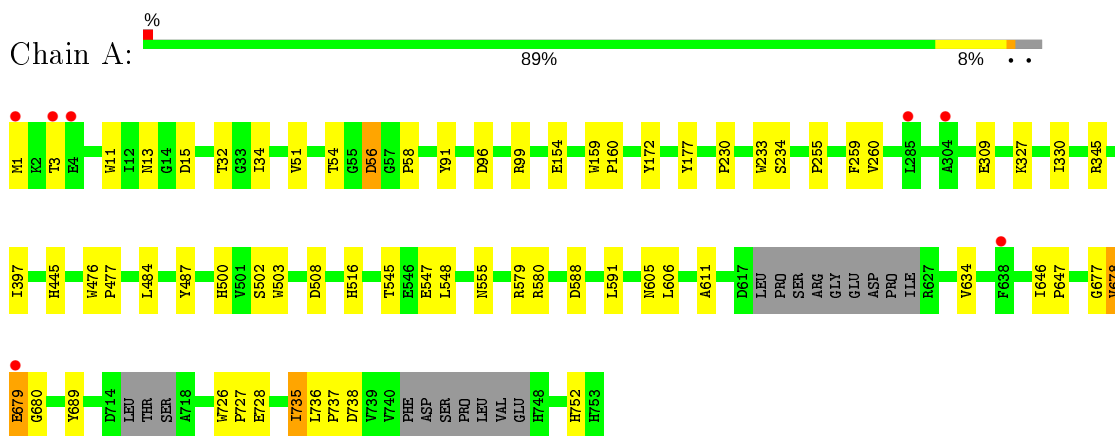
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	A	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		

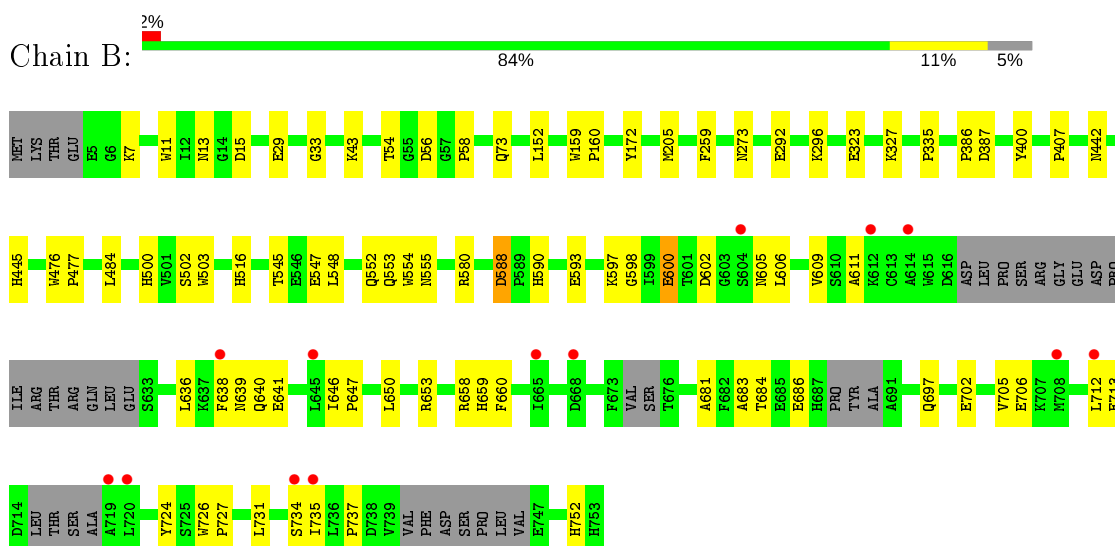
3 Residue-property plots [i](#)

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.

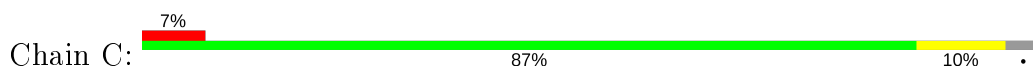
- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG

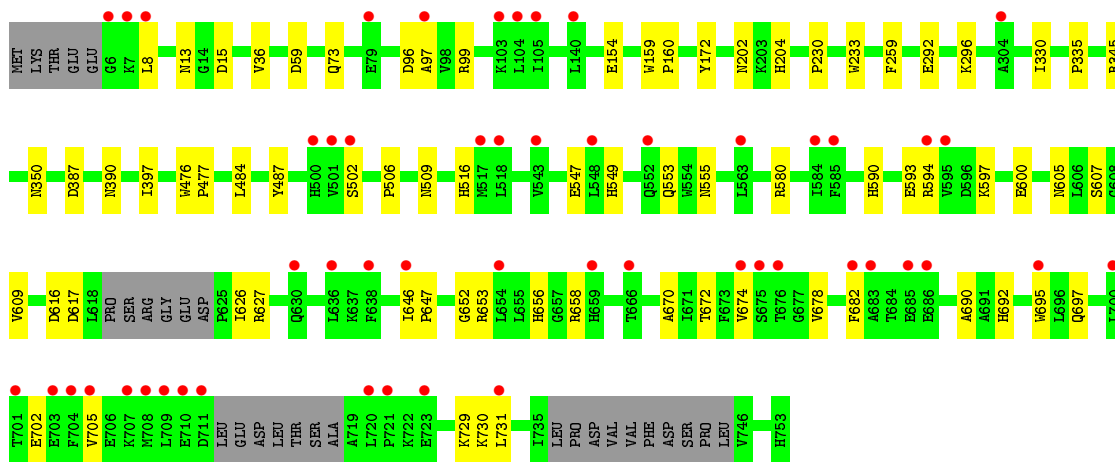


- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG

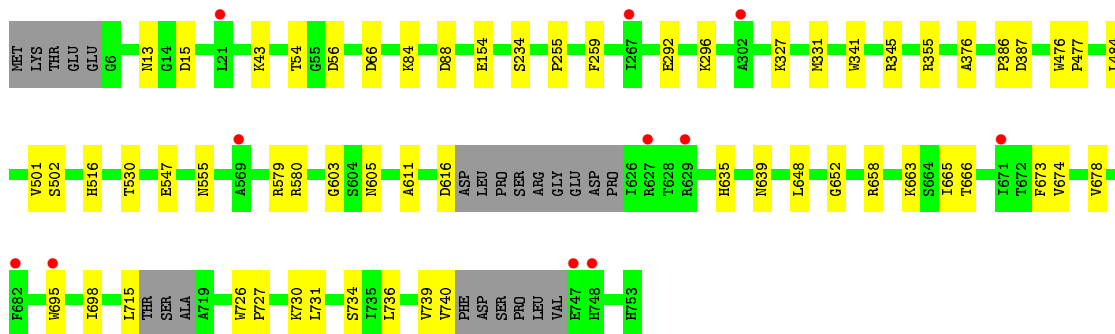
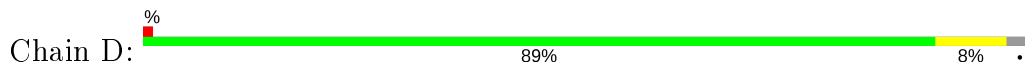


- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG

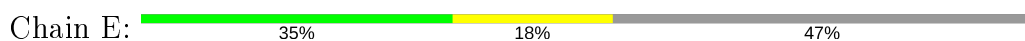




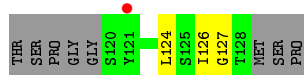
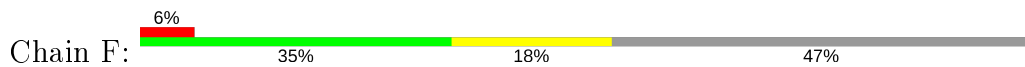
● Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG



● Molecule 2: ZINC FINGER PROTEIN GLI1



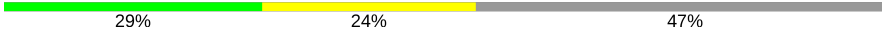
● Molecule 2: ZINC FINGER PROTEIN GLI1



● Molecule 2: ZINC FINGER PROTEIN GLI1



- Molecule 2: ZINC FINGER PROTEIN GLI1

Chain H:  29% 24% 47%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain I:  100%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain J:  100%




- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K:  100%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain L:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	116.30Å 137.60Å 116.54Å 90.00° 105.49° 90.00°	Depositor
Resolution (Å)	19.98 – 2.80 62.73 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.9 (19.98-2.80) 97.9 (62.73-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.81Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.197 , 0.234 0.203 , 0.238	Depositor DCC
R_{free} test set	1966 reflections (2.31%)	wwPDB-VP
Wilson B-factor (Å ²)	80.0	Xtrriage
Anisotropy	0.369	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.019 for l,-k,h	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23141	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, GLC

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	0.35	0/5913	0.51	1/8040 (0.0%)
1	B	0.38	2/5773 (0.0%)	0.51	2/7847 (0.0%)
1	C	0.36	0/5843	0.50	0/7945
1	D	0.34	0/5883	0.50	0/8001
2	E	0.36	0/66	0.63	0/88
2	F	0.40	0/66	0.59	0/88
2	G	0.32	0/59	0.49	0/78
2	H	0.46	0/66	0.75	0/88
All	All	0.36	2/23669 (0.0%)	0.51	3/32175 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	600	GLU	CG-CD	5.22	1.59	1.51
1	B	600	GLU	CD-OE1	5.18	1.31	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	588	ASP	CB-CG-OD1	-5.50	113.35	118.30
1	B	600	GLU	OE1-CD-OE2	5.14	129.47	123.30
1	A	56	ASP	CB-CG-OD1	-5.10	113.71	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5757	0	5587	38	0
1	B	5621	0	5442	63	0
1	C	5687	0	5522	55	0
1	D	5727	0	5562	35	0
2	E	65	0	60	4	0
2	F	65	0	60	7	0
2	G	58	0	53	7	0
2	H	65	0	60	4	0
3	I	23	0	21	0	0
3	J	23	0	21	0	0
3	K	23	0	21	0	0
3	L	23	0	21	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
All	All	23141	0	22430	195	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 4.

All (195) 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:652:GLY:O	1:C:658:ARG:NH2	1.97	0.97
1:B:7:LYS:O	1:B:273:ASN:ND2	2.01	0.93
1:B:681:ALA:HB1	1:B:697:GLN:HE22	1.36	0.88
1:B:636:LEU:HB2	1:B:731:LEU:HD13	1.66	0.77
1:B:681:ALA:HB1	1:B:697:GLN:NE2	2.02	0.74
1:C:593:GLU:HG2	1:C:597:LYS:HE3	1.72	0.72
1:B:683:ALA:HB3	1:B:697:GLN:NE2	2.05	0.72
1:B:43:LYS:NZ	1:B:386:PRO:O	2.28	0.66
1:B:597:LYS:HA	1:B:600:GLU:HG2	1.76	0.65
1:B:639:ASN:OD1	1:B:640:GLN:N	2.28	0.65
1:D:663:LYS:HE2	1:D:695:TRP:CH2	2.31	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:GLU:OE2	1:A:580:ARG:NH2	2.30	0.65
1:A:54:THR:OG1	1:A:56:ASP:CG	2.36	0.65
1:A:555:ASN:HB3	1:A:605:ASN:HB2	1.79	0.63
1:A:154:GLU:OE1	1:A:345:ARG:NH1	2.32	0.63
1:C:653:ARG:HA	1:C:658:ARG:NH2	2.14	0.62
1:D:13:ASN:ND2	1:D:15:ASP:OD1	2.32	0.62
1:A:678:VAL:O	1:A:680:GLY:N	2.33	0.61
1:C:597:LYS:HA	1:C:600:GLU:OE2	2.01	0.61
1:B:638:PHE:CD2	1:B:731:LEU:HD11	2.35	0.61
1:A:51:VAL:HB	1:A:56:ASP:OD1	2.02	0.60
1:B:54:THR:OG1	1:B:56:ASP:CG	2.42	0.58
1:C:13:ASN:ND2	1:C:15:ASP:OD1	2.36	0.58
1:B:609:VAL:HG21	2:F:124:LEU:HD11	1.85	0.57
2:H:127:GLY:O	2:H:128:THR:OG1	2.15	0.57
1:A:484:LEU:HD23	2:E:126:ILE:HD11	1.86	0.57
1:B:593:GLU:HG2	1:B:597:LYS:HE3	1.86	0.56
1:D:616:ASP:O	1:D:635:HIS:N	2.38	0.56
1:C:547:GLU:OE2	1:C:580:ARG:NH2	2.39	0.56
1:B:638:PHE:HD2	1:B:731:LEU:HD11	1.70	0.55
1:B:683:ALA:HB3	1:B:697:GLN:HE21	1.70	0.55
1:D:43:LYS:NZ	1:D:386:PRO:O	2.39	0.55
1:C:96:ASP:OD1	1:C:99:ARG:NH1	2.40	0.54
1:B:73:GLN:OE1	1:B:335:PRO:HB3	2.07	0.54
1:B:54:THR:OG1	1:B:56:ASP:OD1	2.26	0.54
1:C:609:VAL:HG21	2:G:124:LEU:HD11	1.90	0.54
1:B:555:ASN:HB3	1:B:605:ASN:HB2	1.88	0.54
1:B:552:GLN:O	1:B:606:LEU:HD13	2.07	0.53
1:B:554:TRP:CE2	1:B:598:GLY:HA3	2.42	0.53
1:A:54:THR:OG1	1:A:56:ASP:OD1	2.26	0.53
1:C:656:HIS:CB	1:C:658:ARG:NH1	2.72	0.53
1:B:547:GLU:OE2	1:B:580:ARG:NH2	2.42	0.53
1:B:597:LYS:HA	1:B:600:GLU:CG	2.39	0.53
1:C:154:GLU:OE1	1:C:345:ARG:NH1	2.42	0.53
1:C:674:VAL:HG13	1:C:678:VAL:HG21	1.89	0.53
1:C:8:LEU:HD23	1:C:36:VAL:HG22	1.90	0.53
1:C:670:ALA:HB3	1:C:695:TRP:CE3	2.43	0.53
1:B:724:TYR:HB3	1:B:726:TRP:CH2	2.44	0.52
1:D:555:ASN:HB3	1:D:605:ASN:HB2	1.91	0.52
1:C:484:LEU:HD23	2:G:126:ILE:CD1	2.39	0.52
1:B:13:ASN:ND2	1:B:15:ASP:OD1	2.42	0.52
1:C:653:ARG:HG2	1:C:658:ARG:HE	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:GLU:HG2	1:B:327:LYS:NZ	2.24	0.52
1:A:91:TYR:OH	1:A:309:GLU:OE1	2.28	0.51
1:C:506:PRO:HG2	1:C:509:ASN:HA	1.93	0.51
1:D:292:GLU:HG2	1:D:296:LYS:HE3	1.92	0.51
1:B:602:ASP:O	1:B:658:ARG:NH2	2.43	0.51
1:C:590:HIS:CE1	1:C:593:GLU:OE1	2.64	0.51
1:A:736:LEU:N	1:A:737:PRO:HD3	2.26	0.51
1:C:502:SER:HA	1:C:516:HIS:HD2	1.77	0.50
1:C:502:SER:OG	2:G:125:SER:N	2.40	0.50
1:C:682:PHE:CE2	1:C:692:HIS:NE2	2.79	0.50
1:C:555:ASN:HB3	1:C:605:ASN:HB2	1.93	0.49
1:A:445:HIS:HA	1:A:752:HIS:HB3	1.94	0.49
1:D:255:PRO:HB3	1:D:327:LYS:HD3	1.94	0.49
1:C:616:ASP:OD1	1:C:617:ASP:N	2.41	0.49
1:C:656:HIS:CG	1:C:658:ARG:NH1	2.81	0.49
1:B:502:SER:HA	1:B:516:HIS:HD2	1.78	0.48
1:D:674:VAL:CG1	1:D:678:VAL:HG21	2.43	0.48
1:B:484:LEU:HD23	2:F:126:ILE:HD11	1.95	0.48
1:B:735:ILE:HG23	1:B:735:ILE:O	2.13	0.48
1:B:684:THR:OG1	1:B:686:GLU:HB2	2.13	0.48
1:B:683:ALA:CB	1:B:697:GLN:NE2	2.77	0.48
1:C:682:PHE:CE1	1:C:690:ALA:HB3	2.49	0.48
1:C:670:ALA:CB	1:C:695:TRP:CE3	2.96	0.48
1:D:739:VAL:HG23	1:D:740:VAL:HG23	1.96	0.47
1:B:606:LEU:HD12	1:B:606:LEU:N	2.28	0.47
1:A:646:ILE:HB	1:A:647:PRO:HD3	1.96	0.47
1:B:588:ASP:OD1	1:B:590:HIS:CD2	2.68	0.47
1:C:73:GLN:OE1	1:C:335:PRO:CG	2.63	0.47
1:C:626:ILE:HG22	1:C:627:ARG:N	2.30	0.47
1:C:292:GLU:HG2	1:C:296:LYS:HE3	1.96	0.47
1:B:29:GLU:O	1:B:33:GLY:N	2.41	0.47
1:C:230:PRO:HA	1:C:233:TRP:CE2	2.49	0.47
1:A:13:ASN:ND2	1:A:15:ASP:OD1	2.47	0.46
1:B:611:ALA:HB2	2:F:124:LEU:HD12	1.97	0.46
1:D:484:LEU:HD23	2:H:126:ILE:HD11	1.97	0.46
1:A:11:TRP:CE2	1:A:58:PRO:HG3	2.50	0.46
1:C:702:GLU:HA	1:C:705:VAL:HB	1.97	0.46
1:D:84:LYS:NZ	1:D:88:ASP:OD1	2.46	0.46
1:C:292:GLU:O	1:C:296:LYS:HG3	2.16	0.46
1:C:730:LYS:C	1:C:731:LEU:HD12	2.36	0.46
1:D:547:GLU:OE2	1:D:580:ARG:NH2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:653:ARG:NE	1:B:659:HIS:O	2.40	0.46
1:C:672:THR:HB	1:C:697:GLN:HG2	1.96	0.46
1:B:152:LEU:HD11	1:B:205:MET:HE3	1.98	0.46
1:B:323:GLU:CG	1:B:327:LYS:NZ	2.79	0.46
1:A:735:ILE:HG22	1:A:735:ILE:O	2.16	0.45
1:C:656:HIS:CG	1:C:658:ARG:HH12	2.35	0.45
1:D:476:TRP:N	1:D:477:PRO:CD	2.80	0.45
1:B:597:LYS:CA	1:B:600:GLU:HG2	2.46	0.45
1:C:476:TRP:N	1:C:477:PRO:CD	2.80	0.45
1:C:594:ARG:HA	1:C:597:LYS:HB2	1.98	0.45
1:D:674:VAL:HG13	1:D:678:VAL:HG21	1.97	0.45
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.52	0.45
1:C:387:ASP:N	1:C:387:ASP:OD1	2.50	0.45
1:B:641:GLU:HG2	2:F:124:LEU:HD13	1.99	0.45
1:A:96:ASP:OD1	1:A:99:ARG:NH1	2.49	0.45
1:A:159:TRP:N	1:A:160:PRO:CD	2.80	0.45
1:C:626:ILE:HG22	1:C:627:ARG:HG3	1.98	0.45
1:D:730:LYS:O	1:D:731:LEU:HD12	2.17	0.45
1:B:683:ALA:H	1:B:697:GLN:NE2	2.13	0.44
1:D:736:LEU:O	1:D:739:VAL:HG22	2.18	0.44
1:B:650:LEU:HB3	1:B:705:VAL:HG13	1.98	0.44
1:B:503:TRP:CH2	2:F:127:GLY:HA3	2.52	0.44
1:B:712:LEU:HD12	1:B:713:GLU:N	2.32	0.44
1:C:159:TRP:N	1:C:160:PRO:CD	2.81	0.44
1:D:726:TRP:HB3	1:D:727:PRO:HD2	1.99	0.44
1:D:639:ASN:HA	1:D:734:SER:O	2.17	0.44
1:B:387:ASP:OD1	1:B:387:ASP:N	2.51	0.44
1:C:590:HIS:O	1:C:593:GLU:HB3	2.17	0.44
1:C:653:ARG:HA	1:C:658:ARG:HH21	1.81	0.44
1:D:501:VAL:HG11	2:H:126:ILE:HD11	1.99	0.43
1:D:54:THR:OG1	1:D:56:ASP:OD1	2.34	0.43
1:A:260:VAL:HB	1:A:330:ILE:HA	2.00	0.43
1:A:611:ALA:HB2	2:E:124:LEU:HD12	1.99	0.43
1:C:729:LYS:HG2	1:C:731:LEU:CD1	2.48	0.43
1:D:663:LYS:NZ	1:D:695:TRP:HH2	2.16	0.43
1:D:66:ASP:CB	1:D:331:MET:HE2	2.48	0.43
1:C:97:ALA:HB2	1:C:330:ILE:HD11	2.01	0.43
1:D:387:ASP:N	1:D:387:ASP:OD1	2.49	0.43
1:A:555:ASN:HB3	1:A:605:ASN:CB	2.48	0.43
1:B:588:ASP:OD1	1:B:588:ASP:C	2.57	0.43
1:B:11:TRP:CD2	1:B:58:PRO:HG3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:TRP:N	1:A:477:PRO:CD	2.82	0.43
1:D:673:PHE:CE2	1:D:698:ILE:HD11	2.53	0.43
1:A:502:SER:HA	1:A:516:HIS:HD2	1.84	0.43
1:C:202:ASN:HB2	1:C:204:HIS:CD2	2.53	0.43
1:C:607:SER:N	2:G:120:SER:O	2.49	0.43
1:D:611:ALA:HB2	2:H:124:LEU:HD12	2.00	0.43
1:D:331:MET:HE1	1:D:341:TRP:HE1	1.84	0.43
1:C:484:LEU:HD23	2:G:126:ILE:HD11	2.01	0.42
1:D:603:GLY:HA3	1:D:658:ARG:NH2	2.34	0.42
1:B:476:TRP:N	1:B:477:PRO:CD	2.82	0.42
1:A:255:PRO:HB3	1:A:327:LYS:HD3	2.01	0.42
1:B:11:TRP:CE2	1:B:58:PRO:HG3	2.54	0.42
1:B:653:ARG:HD3	1:B:660:PHE:HB2	2.00	0.42
1:D:665:ILE:HG13	1:D:666:THR:HG23	2.01	0.42
1:A:677:GLY:O	1:A:679:GLU:N	2.52	0.42
1:A:503:TRP:H	1:A:516:HIS:HD2	1.67	0.42
1:B:646:ILE:HB	1:B:647:PRO:HD3	2.01	0.42
1:B:292:GLU:HG2	1:B:296:LYS:HE3	2.02	0.42
1:C:476:TRP:CG	1:C:477:PRO:HD3	2.54	0.42
1:A:516:HIS:CE1	1:A:545:THR:HG22	2.55	0.42
1:A:736:LEU:N	1:A:737:PRO:CD	2.83	0.42
1:A:32:THR:HG22	1:A:34:ILE:HD13	2.02	0.42
1:D:502:SER:HA	1:D:516:HIS:HD2	1.84	0.42
1:B:516:HIS:CE1	1:B:545:THR:HG22	2.55	0.42
1:B:500:HIS:HD2	1:B:548:LEU:HD11	1.84	0.42
1:C:350:ASN:OD1	1:C:390:ASN:ND2	2.45	0.41
2:E:123:HIS:HB3	2:E:126:ILE:HG21	2.02	0.41
1:D:154:GLU:OE1	1:D:345:ARG:NH1	2.52	0.41
1:C:487:TYR:CE1	2:G:126:ILE:HG22	2.55	0.41
1:B:445:HIS:HA	1:B:752:HIS:HB3	2.02	0.41
1:C:397:ILE:H	1:C:397:ILE:HD12	1.84	0.41
1:D:665:ILE:HG13	1:D:666:THR:N	2.35	0.41
1:B:641:GLU:OE2	2:F:124:LEU:HD22	2.21	0.41
1:A:588:ASP:HB3	1:A:591:LEU:CD1	2.50	0.41
1:B:554:TRP:NE1	1:B:598:GLY:HA3	2.35	0.41
1:D:579:ARG:NE	1:D:579:ARG:HA	2.35	0.41
1:A:397:ILE:HD12	1:A:397:ILE:H	1.86	0.41
1:D:331:MET:HE3	1:D:341:TRP:HZ2	1.86	0.41
1:A:634:VAL:HG21	1:A:689:TYR:HB3	2.02	0.41
1:B:442:ASN:HA	1:B:445:HIS:O	2.20	0.41
1:B:726:TRP:HB3	1:B:727:PRO:HD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:GLN:OE1	1:B:335:PRO:CB	2.68	0.41
2:F:126:ILE:HG13	2:F:127:GLY:N	2.36	0.41
1:A:476:TRP:CZ2	1:A:508:ASP:HB3	2.56	0.41
1:A:487:TYR:CD1	2:E:126:ILE:HB	2.56	0.41
1:A:500:HIS:HE1	1:A:606:LEU:HD21	1.85	0.41
1:B:400:TYR:HB2	1:B:407:PRO:HA	2.03	0.41
1:A:516:HIS:CD2	1:A:548:LEU:HD22	2.57	0.40
1:B:734:SER:HB2	1:B:737:PRO:HG3	2.03	0.40
1:C:8:LEU:HD12	1:C:59:ASP:HB2	2.04	0.40
1:C:646:ILE:HB	1:C:647:PRO:HD3	2.04	0.40
1:C:656:HIS:HB2	1:C:658:ARG:NH1	2.36	0.40
1:B:159:TRP:N	1:B:160:PRO:CD	2.84	0.40
1:A:172:TYR:HD2	1:A:177:TYR:CZ	2.38	0.40
1:C:506:PRO:O	1:C:509:ASN:N	2.48	0.40
1:C:549:HIS:O	1:C:553:GLN:HG2	2.21	0.40
1:D:376:ALA:CB	1:D:530:THR:HG22	2.51	0.40
2:G:126:ILE:HG13	2:G:127:GLY:N	2.37	0.40
1:A:726:TRP:HB3	1:A:727:PRO:HD2	2.02	0.40
1:B:702:GLU:OE2	1:B:706:GLU:HG3	2.21	0.40
1:D:648:LEU:O	1:D:652:GLY:N	2.53	0.40

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	726/753 (96%)	704 (97%)	18 (2%)	4 (1%)	25	56
1	B	705/753 (94%)	687 (97%)	18 (3%)	0	100	100
1	C	717/753 (95%)	700 (98%)	17 (2%)	0	100	100
1	D	722/753 (96%)	706 (98%)	16 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	7/17 (41%)	7 (100%)	0	0	100	100
2	F	7/17 (41%)	7 (100%)	0	0	100	100
2	G	6/17 (35%)	6 (100%)	0	0	100	100
2	H	7/17 (41%)	7 (100%)	0	0	100	100
All	All	2897/3080 (94%)	2824 (98%)	69 (2%)	4 (0%)	51	81

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	679	GLU
1	A	678	VAL
1	A	738	ASP
1	A	735	ILE

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	608/626 (97%)	602 (99%)	6 (1%)	76	93
1	B	593/626 (95%)	590 (100%)	3 (0%)	88	96
1	C	600/626 (96%)	598 (100%)	2 (0%)	92	98
1	D	605/626 (97%)	601 (99%)	4 (1%)	84	95
2	E	7/13 (54%)	7 (100%)	0	100	100
2	F	7/13 (54%)	7 (100%)	0	100	100
2	G	6/13 (46%)	5 (83%)	1 (17%)	2	6
2	H	7/13 (54%)	7 (100%)	0	100	100
All	All	2433/2556 (95%)	2417 (99%)	16 (1%)	84	95

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	3	THR
1	A	234	SER
1	A	259	PHE
1	A	579	ARG
1	A	728	GLU
1	B	172	TYR
1	B	259	PHE
1	B	553	GLN
1	C	172	TYR
1	C	259	PHE
1	D	234	SER
1	D	259	PHE
1	D	355	ARG
1	D	715	LEU
2	G	124	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	500	HIS
1	A	516	HIS
1	B	516	HIS
1	B	590	HIS
1	B	697	GLN
1	C	516	HIS
1	D	500	HIS
1	D	516	HIS

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](#)

8 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	GLC	I	1	3	12,12,12	0.64	0	17,17,17	0.58	0
3	GLC	I	2	3	11,11,12	0.72	0	15,15,17	0.80	0
3	GLC	J	1	3	12,12,12	0.63	0	17,17,17	0.53	0
3	GLC	J	2	3	11,11,12	0.70	0	15,15,17	0.60	0
3	GLC	K	1	3	12,12,12	0.56	0	17,17,17	0.74	0
3	GLC	K	2	3	11,11,12	0.78	0	15,15,17	0.78	0
3	GLC	L	1	3	12,12,12	0.60	0	17,17,17	0.85	0
3	GLC	L	2	3	11,11,12	0.60	0	15,15,17	0.88	1 (6%)

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	GLC	I	1	3	-	0/2/22/22	0/1/1/1
3	GLC	I	2	3	-	0/2/19/22	0/1/1/1
3	GLC	J	1	3	-	2/2/22/22	0/1/1/1
3	GLC	J	2	3	-	0/2/19/22	0/1/1/1
3	GLC	K	1	3	-	1/2/22/22	0/1/1/1
3	GLC	K	2	3	-	0/2/19/22	0/1/1/1
3	GLC	L	1	3	-	2/2/22/22	0/1/1/1
3	GLC	L	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	2	GLC	C6-C5-C4	-2.08	108.13	113.00

There are no chirality outliers.

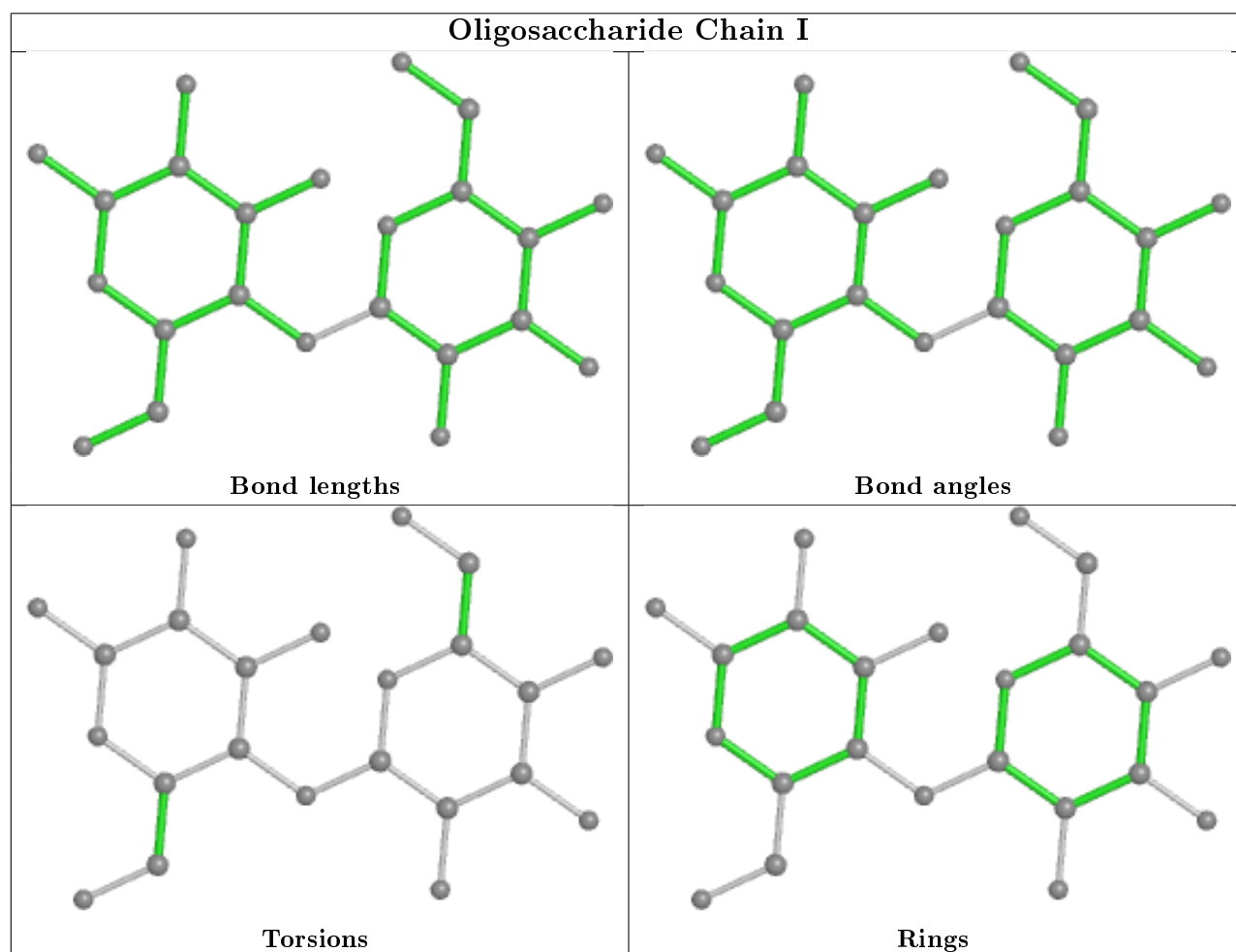
All (5) torsion outliers are listed below:

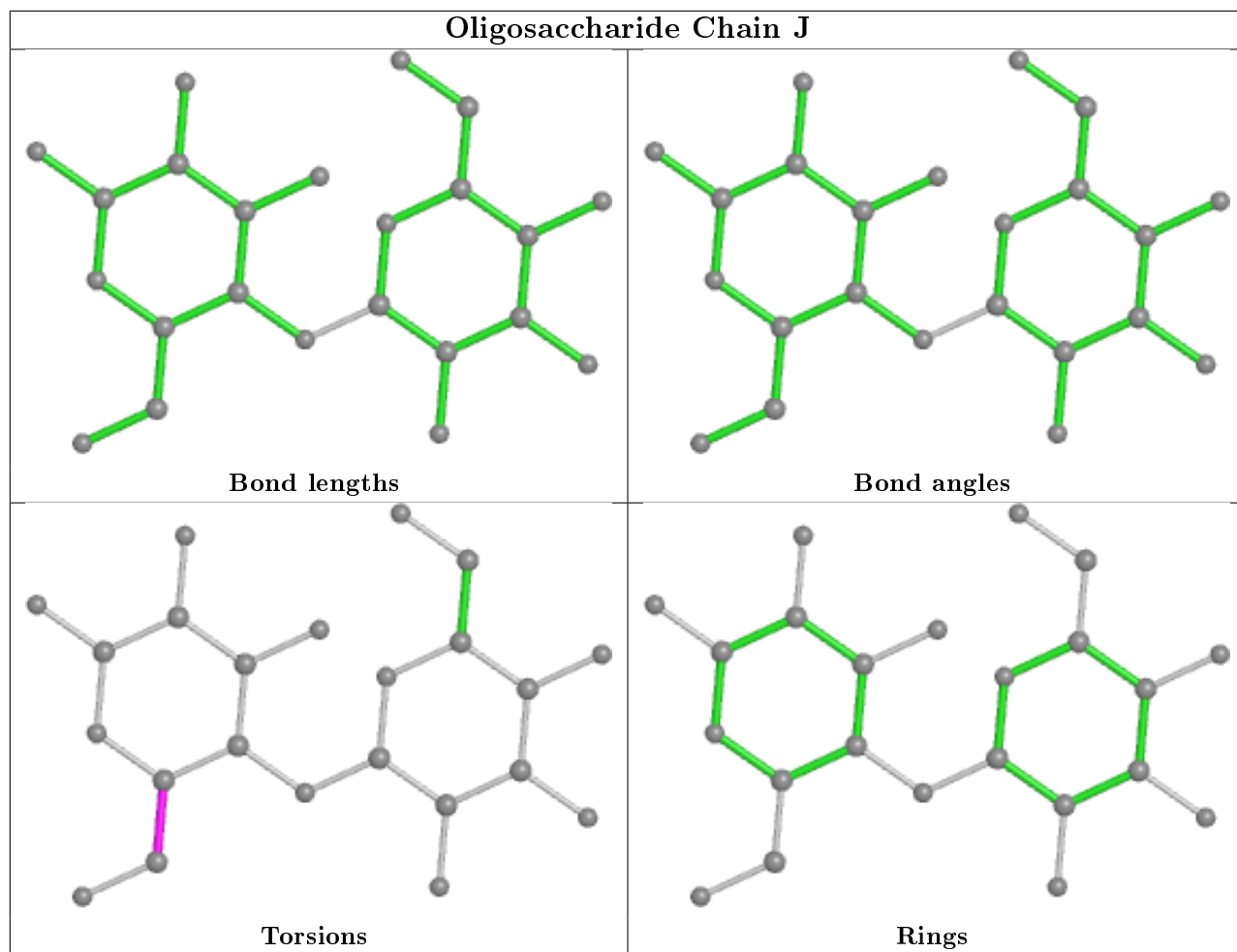
Mol	Chain	Res	Type	Atoms
3	L	1	GLC	C4-C5-C6-O6
3	L	1	GLC	O5-C5-C6-O6
3	J	1	GLC	C4-C5-C6-O6
3	J	1	GLC	O5-C5-C6-O6
3	K	1	GLC	C4-C5-C6-O6

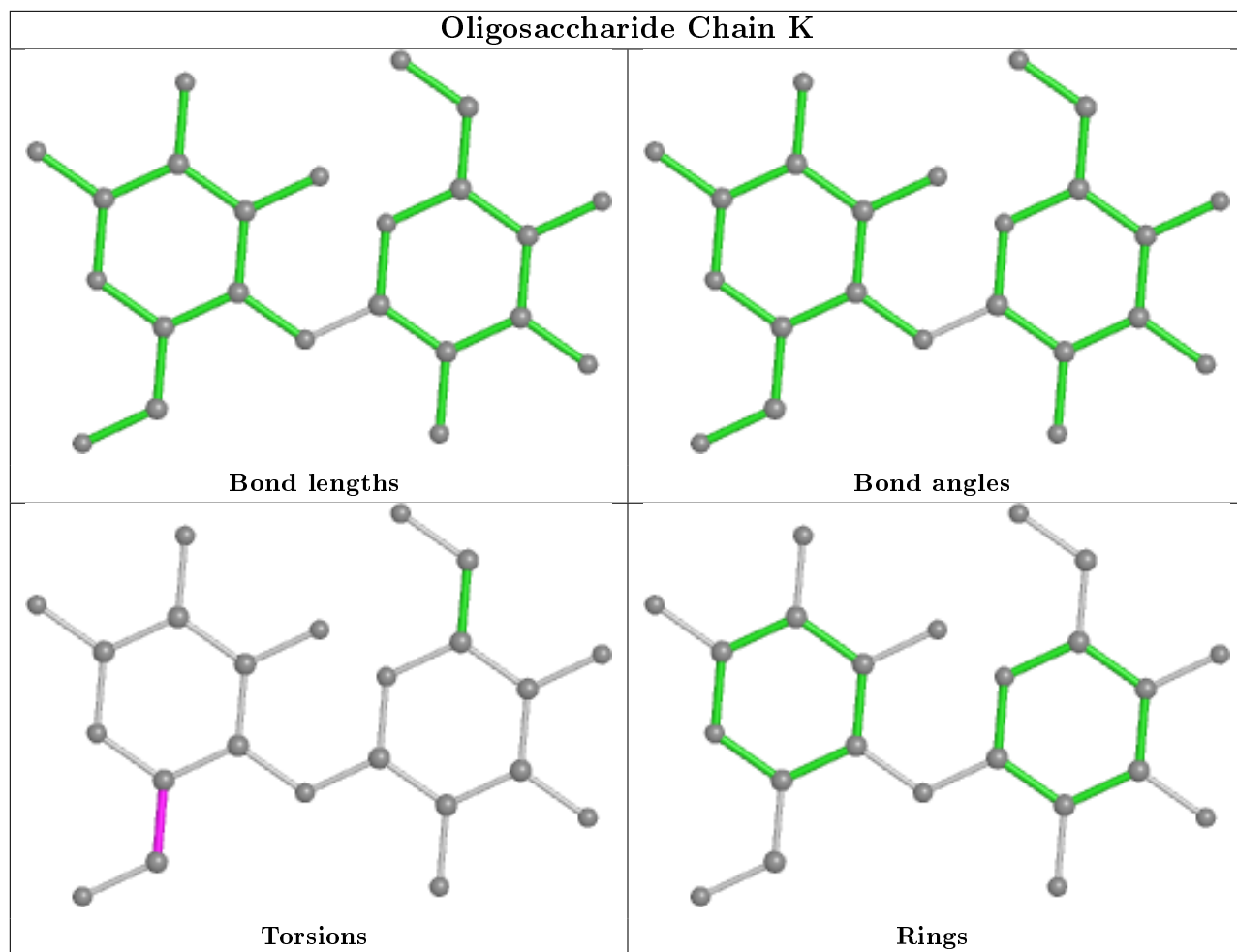
There are no ring outliers.

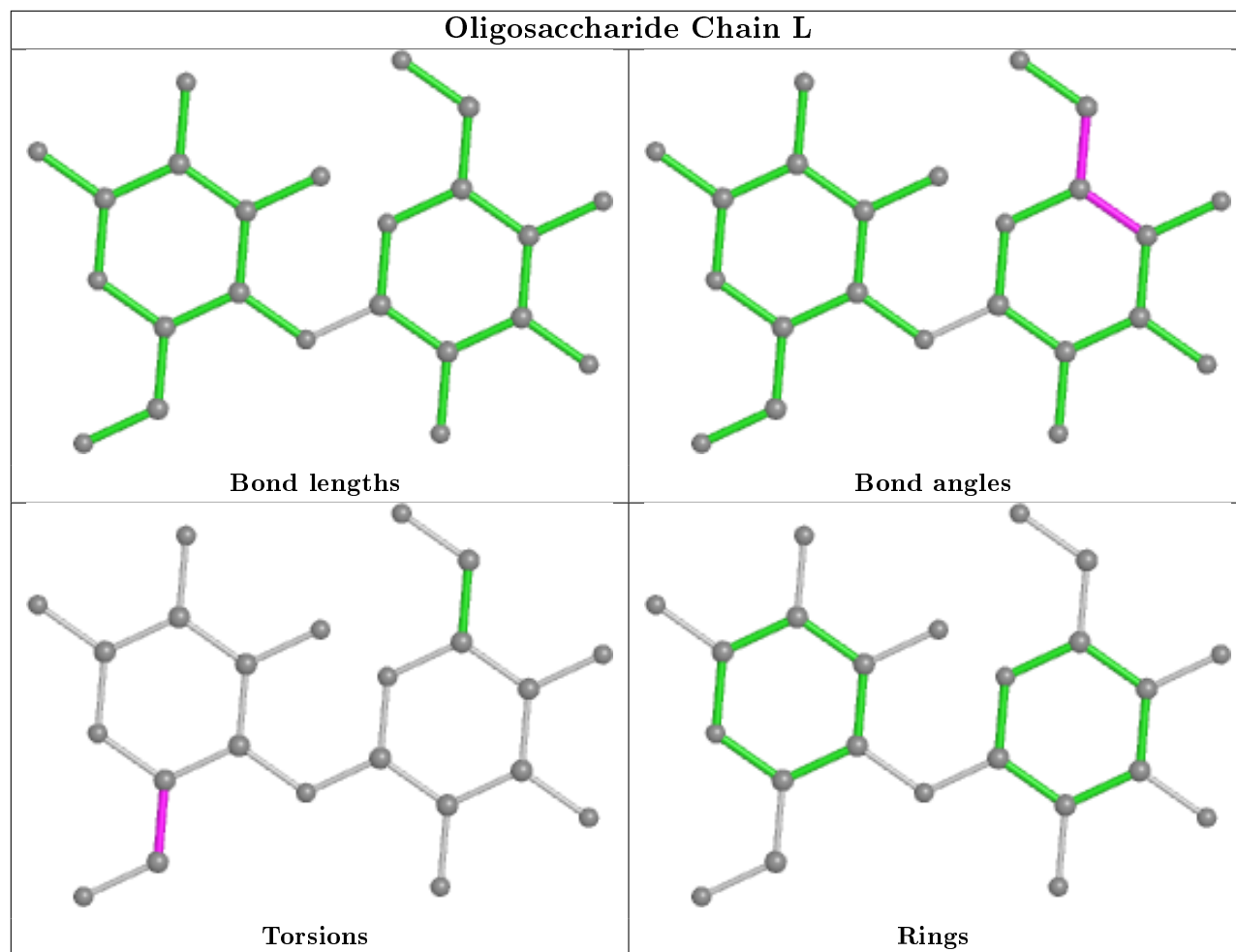
No monomer is involved in short contacts.

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	734/753 (97%)	-0.27	7 (0%) 82 77	50, 81, 127, 176	0
1	B	717/753 (95%)	-0.16	13 (1%) 68 61	51, 87, 167, 209	0
1	C	725/753 (96%)	0.09	52 (7%) 15 8	64, 111, 193, 274	0
1	D	730/753 (96%)	-0.25	11 (1%) 73 68	56, 87, 141, 190	0
2	E	9/17 (52%)	-0.12	0 100 100	61, 71, 105, 107	0
2	F	9/17 (52%)	0.32	1 (11%) 5 3	87, 99, 123, 125	0
2	G	8/17 (47%)	0.35	0 100 100	91, 104, 111, 133	0
2	H	9/17 (52%)	0.21	0 100 100	78, 80, 107, 108	0
All	All	2941/3080 (95%)	-0.14	84 (2%) 51 41	50, 90, 167, 274	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	703	GLU	7.4
1	A	1	MET	7.4
1	C	8	LEU	6.1
1	C	710	GLU	5.9
1	C	548	LEU	5.4
1	C	674	VAL	5.3
1	C	708	MET	5.3
1	C	701	THR	5.2
1	B	720	LEU	5.2
1	B	712	LEU	5.1
1	C	595	VAL	4.9
1	C	552	GLN	4.6
1	C	501	VAL	4.5
1	C	630	GLN	4.2
1	C	700	LEU	3.9
1	C	704	PHE	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	584	ILE	3.8
1	C	103	LYS	3.8
1	C	105	ILE	3.8
1	C	585	PHE	3.8
1	C	721	PRO	3.7
1	C	705	VAL	3.7
1	B	708	MET	3.7
1	C	518	LEU	3.7
1	C	695	TRP	3.6
1	D	627	ARG	3.6
1	A	3	THR	3.5
1	B	645	LEU	3.4
1	C	594	ARG	3.4
1	C	711	ASP	3.3
1	C	104	LEU	3.3
1	C	709	LEU	3.3
1	C	500	HIS	3.2
1	C	682	PHE	3.2
1	C	686	GLU	3.0
1	B	668	ASP	3.0
1	C	6	GLY	3.0
1	D	747	GLU	3.0
1	C	659	HIS	2.9
1	C	723	GLU	2.9
1	C	707	LYS	2.8
1	C	7	LYS	2.7
1	A	4	GLU	2.7
1	C	720	LEU	2.7
1	C	79	GLU	2.6
1	C	676	THR	2.6
1	A	285	LEU	2.6
1	D	21	LEU	2.6
1	B	665	ILE	2.5
1	C	731	LEU	2.5
1	D	569	ALA	2.5
1	C	654	LEU	2.5
1	C	675	SER	2.4
1	C	685	GLU	2.4
1	D	695	TRP	2.4
1	C	563	LEU	2.3
1	A	304	ALA	2.3
1	B	734	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	719	ALA	2.3
1	C	97	ALA	2.3
1	D	748	HIS	2.3
1	B	614	ALA	2.3
1	C	140	LEU	2.3
1	C	683	ALA	2.2
1	B	604	SER	2.2
1	C	304	ALA	2.2
1	C	517	MET	2.2
1	D	629	ARG	2.2
1	B	638	PHE	2.2
1	C	646	ILE	2.2
1	B	735	ILE	2.1
1	D	671	ILE	2.1
1	A	679	GLU	2.1
2	F	121	TYR	2.1
1	C	666	THR	2.1
1	C	502	SER	2.1
1	D	267	ILE	2.1
1	C	636	LEU	2.1
1	C	543	VAL	2.1
1	A	638	PHE	2.1
1	D	302	ALA	2.1
1	B	612	LYS	2.1
1	D	682	PHE	2.0
1	C	638	PHE	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

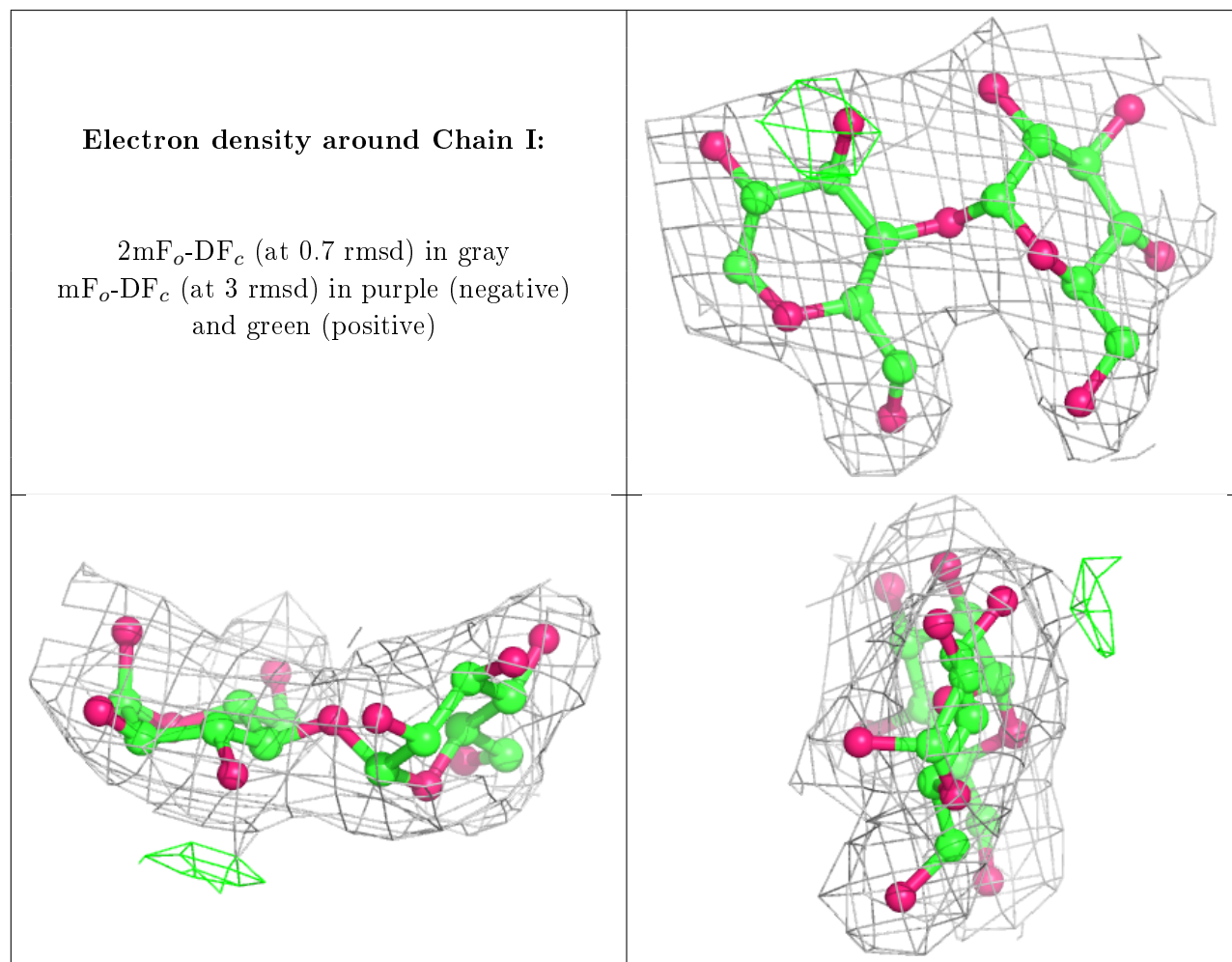
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GLC	J	1	12/12	0.95	0.16	47,55,60,70	0
3	GLC	L	1	12/12	0.96	0.17	57,64,72,72	0

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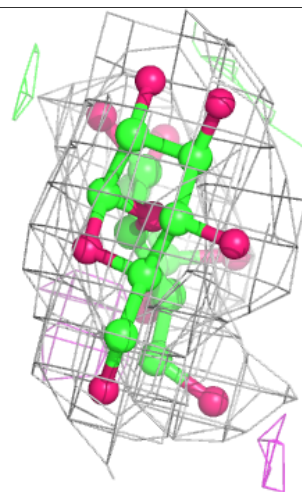
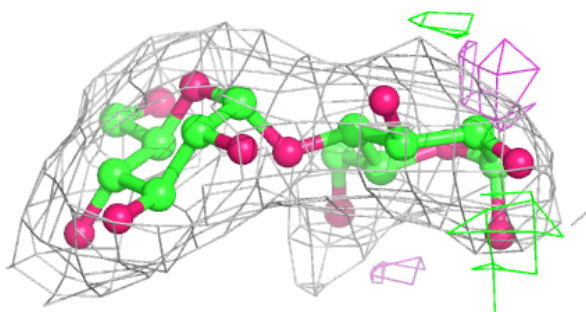
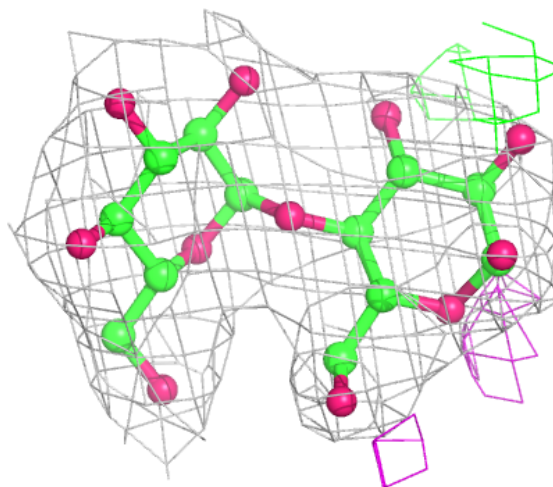
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GLC	K	1	12/12	0.96	0.14	63,84,91,98	0
3	GLC	I	1	12/12	0.96	0.15	55,64,73,82	0
3	GLC	K	2	11/12	0.96	0.16	63,67,86,97	0
3	GLC	I	2	11/12	0.97	0.14	56,60,68,76	0
3	GLC	L	2	11/12	0.97	0.12	55,64,68,69	0
3	GLC	J	2	11/12	0.98	0.16	49,62,67,71	0

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.



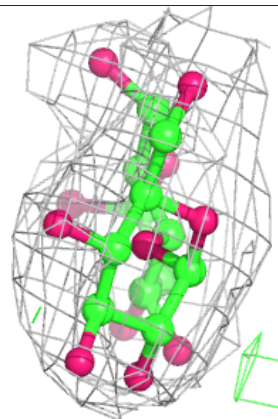
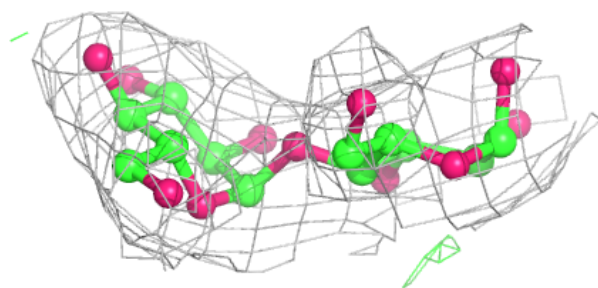
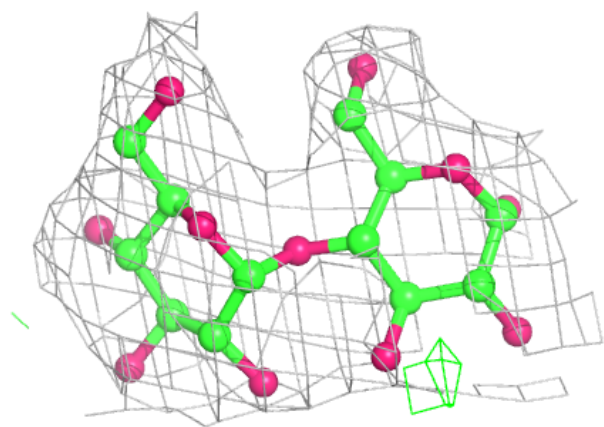
Electron density around Chain J:

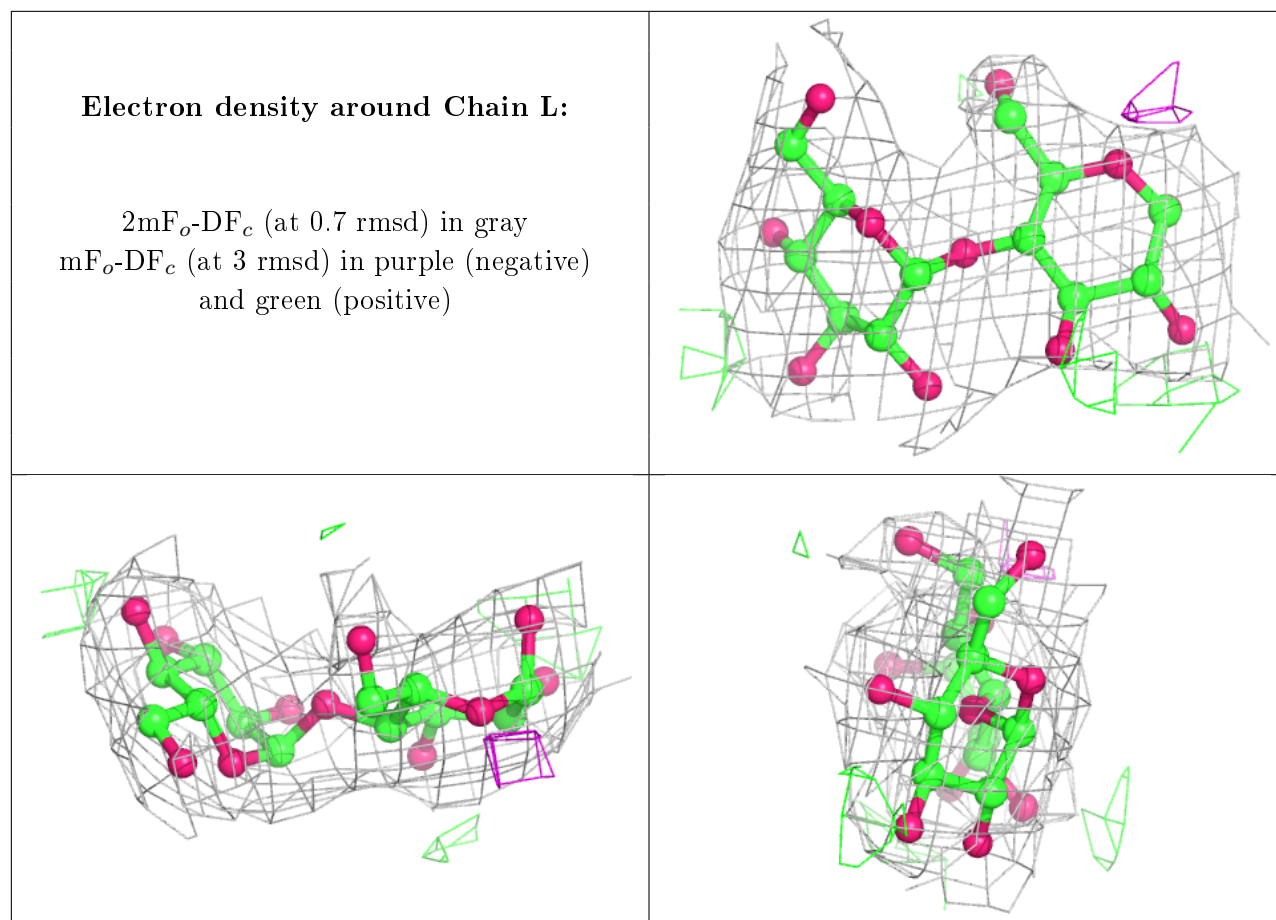
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain K:

$2mF_o-DF_c$ (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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ZN	A	910	1/1	0.96	0.18	75,75,75,75	0
4	ZN	B	910	1/1	0.98	0.23	98,98,98,98	0
4	ZN	D	910	1/1	0.99	0.22	93,93,93,93	0
4	ZN	C	910	1/1	0.99	0.26	84,84,84,84	0

6.5 Other polymers [i](#)

There are no such residues in this entry.