



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:15 PM UTC

PDB ID : 4GKX / pdb_00004gkx
Title : Crystal structure of a carbohydrate-binding domain
Authors : Page, R.C.; Zheng, C.; Nix, J.C.; Misra, S.; Zhang, B.
Deposited on : 2012-08-13
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

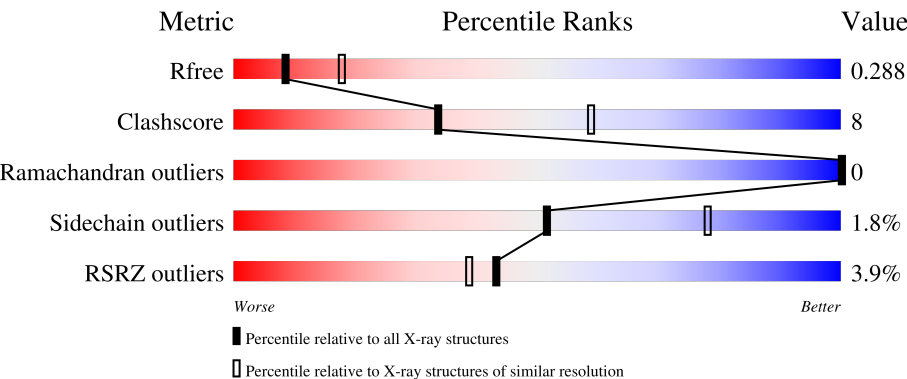
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



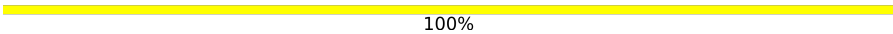
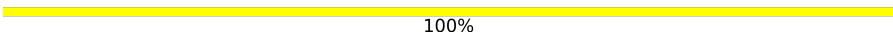
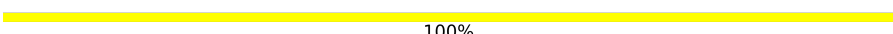
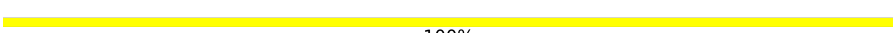
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 $\leq 5\%$. 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	261	2% 56% 31% • 12%
1	B	261	3% 73% 14% • 12%
1	C	261	2% 76% 11% • 12%
1	D	261	2% 72% 16% • 12%
1	E	261	7% 80% 8% • 12%

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Mol	Chain	Length	Quality of chain
1	F	261	
2	G	2	
2	H	2	
2	I	2	
2	J	2	
2	K	2	
2	L	2	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11508 atoms, of which 132 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 Protein ERGIC-53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	0	0
			1807	1147	317	339	4			
1	B	230	Total	C	N	O	S	0	0	0
			1807	1147	317	339	4			
1	C	230	Total	C	N	O	S	0	0	0
			1807	1147	317	339	4			
1	D	230	Total	C	N	O	S	0	1	0
			1816	1153	319	340	4			
1	E	230	Total	C	N	O	S	0	0	0
			1807	1147	317	339	4			
1	F	230	Total	C	N	O	S	0	0	0
			1807	1147	317	339	4			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MET	-	expression tag	UNP P49257
A	11	GLY	-	expression tag	UNP P49257
A	12	SER	-	expression tag	UNP P49257
A	13	SER	-	expression tag	UNP P49257
A	14	HIS	-	expression tag	UNP P49257
A	15	HIS	-	expression tag	UNP P49257
A	16	HIS	-	expression tag	UNP P49257
A	17	HIS	-	expression tag	UNP P49257
A	18	HIS	-	expression tag	UNP P49257
A	19	HIS	-	expression tag	UNP P49257
A	20	SER	-	expression tag	UNP P49257
A	21	SER	-	expression tag	UNP P49257
A	22	GLY	-	expression tag	UNP P49257
A	23	LEU	-	expression tag	UNP P49257
A	24	VAL	-	expression tag	UNP P49257
A	25	PRO	-	expression tag	UNP P49257
A	26	ARG	-	expression tag	UNP P49257

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Chain	Residue	Modelled	Actual	Comment	Reference
A	27	GLY	-	expression tag	UNP P49257
A	28	SER	-	expression tag	UNP P49257
A	29	HIS	-	expression tag	UNP P49257
A	30	MET	-	expression tag	UNP P49257
B	10	MET	-	expression tag	UNP P49257
B	11	GLY	-	expression tag	UNP P49257
B	12	SER	-	expression tag	UNP P49257
B	13	SER	-	expression tag	UNP P49257
B	14	HIS	-	expression tag	UNP P49257
B	15	HIS	-	expression tag	UNP P49257
B	16	HIS	-	expression tag	UNP P49257
B	17	HIS	-	expression tag	UNP P49257
B	18	HIS	-	expression tag	UNP P49257
B	19	HIS	-	expression tag	UNP P49257
B	20	SER	-	expression tag	UNP P49257
B	21	SER	-	expression tag	UNP P49257
B	22	GLY	-	expression tag	UNP P49257
B	23	LEU	-	expression tag	UNP P49257
B	24	VAL	-	expression tag	UNP P49257
B	25	PRO	-	expression tag	UNP P49257
B	26	ARG	-	expression tag	UNP P49257
B	27	GLY	-	expression tag	UNP P49257
B	28	SER	-	expression tag	UNP P49257
B	29	HIS	-	expression tag	UNP P49257
B	30	MET	-	expression tag	UNP P49257
C	10	MET	-	expression tag	UNP P49257
C	11	GLY	-	expression tag	UNP P49257
C	12	SER	-	expression tag	UNP P49257
C	13	SER	-	expression tag	UNP P49257
C	14	HIS	-	expression tag	UNP P49257
C	15	HIS	-	expression tag	UNP P49257
C	16	HIS	-	expression tag	UNP P49257
C	17	HIS	-	expression tag	UNP P49257
C	18	HIS	-	expression tag	UNP P49257
C	19	HIS	-	expression tag	UNP P49257
C	20	SER	-	expression tag	UNP P49257
C	21	SER	-	expression tag	UNP P49257
C	22	GLY	-	expression tag	UNP P49257
C	23	LEU	-	expression tag	UNP P49257
C	24	VAL	-	expression tag	UNP P49257
C	25	PRO	-	expression tag	UNP P49257
C	26	ARG	-	expression tag	UNP P49257

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Chain	Residue	Modelled	Actual	Comment	Reference
C	27	GLY	-	expression tag	UNP P49257
C	28	SER	-	expression tag	UNP P49257
C	29	HIS	-	expression tag	UNP P49257
C	30	MET	-	expression tag	UNP P49257
D	10	MET	-	expression tag	UNP P49257
D	11	GLY	-	expression tag	UNP P49257
D	12	SER	-	expression tag	UNP P49257
D	13	SER	-	expression tag	UNP P49257
D	14	HIS	-	expression tag	UNP P49257
D	15	HIS	-	expression tag	UNP P49257
D	16	HIS	-	expression tag	UNP P49257
D	17	HIS	-	expression tag	UNP P49257
D	18	HIS	-	expression tag	UNP P49257
D	19	HIS	-	expression tag	UNP P49257
D	20	SER	-	expression tag	UNP P49257
D	21	SER	-	expression tag	UNP P49257
D	22	GLY	-	expression tag	UNP P49257
D	23	LEU	-	expression tag	UNP P49257
D	24	VAL	-	expression tag	UNP P49257
D	25	PRO	-	expression tag	UNP P49257
D	26	ARG	-	expression tag	UNP P49257
D	27	GLY	-	expression tag	UNP P49257
D	28	SER	-	expression tag	UNP P49257
D	29	HIS	-	expression tag	UNP P49257
D	30	MET	-	expression tag	UNP P49257
E	10	MET	-	expression tag	UNP P49257
E	11	GLY	-	expression tag	UNP P49257
E	12	SER	-	expression tag	UNP P49257
E	13	SER	-	expression tag	UNP P49257
E	14	HIS	-	expression tag	UNP P49257
E	15	HIS	-	expression tag	UNP P49257
E	16	HIS	-	expression tag	UNP P49257
E	17	HIS	-	expression tag	UNP P49257
E	18	HIS	-	expression tag	UNP P49257
E	19	HIS	-	expression tag	UNP P49257
E	20	SER	-	expression tag	UNP P49257
E	21	SER	-	expression tag	UNP P49257
E	22	GLY	-	expression tag	UNP P49257
E	23	LEU	-	expression tag	UNP P49257
E	24	VAL	-	expression tag	UNP P49257
E	25	PRO	-	expression tag	UNP P49257
E	26	ARG	-	expression tag	UNP P49257

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Chain	Residue	Modelled	Actual	Comment	Reference
E	27	GLY	-	expression tag	UNP P49257
E	28	SER	-	expression tag	UNP P49257
E	29	HIS	-	expression tag	UNP P49257
E	30	MET	-	expression tag	UNP P49257
F	10	MET	-	expression tag	UNP P49257
F	11	GLY	-	expression tag	UNP P49257
F	12	SER	-	expression tag	UNP P49257
F	13	SER	-	expression tag	UNP P49257
F	14	HIS	-	expression tag	UNP P49257
F	15	HIS	-	expression tag	UNP P49257
F	16	HIS	-	expression tag	UNP P49257
F	17	HIS	-	expression tag	UNP P49257
F	18	HIS	-	expression tag	UNP P49257
F	19	HIS	-	expression tag	UNP P49257
F	20	SER	-	expression tag	UNP P49257
F	21	SER	-	expression tag	UNP P49257
F	22	GLY	-	expression tag	UNP P49257
F	23	LEU	-	expression tag	UNP P49257
F	24	VAL	-	expression tag	UNP P49257
F	25	PRO	-	expression tag	UNP P49257
F	26	ARG	-	expression tag	UNP P49257
F	27	GLY	-	expression tag	UNP P49257
F	28	SER	-	expression tag	UNP P49257
F	29	HIS	-	expression tag	UNP P49257
F	30	MET	-	expression tag	UNP P49257

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	2	Total	C	H	O	0	0	0
			45	12	22	11			
2	H	2	Total	C	H	O	0	0	0
			45	12	22	11			
2	I	2	Total	C	H	O	0	0	0
			45	12	22	11			
2	J	2	Total	C	H	O	0	0	0
			45	12	22	11			

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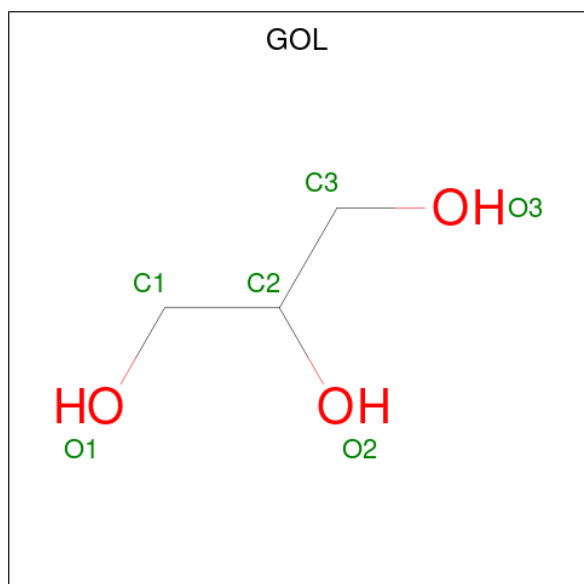
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	K	2	Total	C	H	O	0	0	0
			45	12	22	11			
2	L	2	Total	C	H	O	0	0	0
			45	12	22	11			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Ca	0	0
			2	2		
3	B	2	Total	Ca	0	0
			2	2		
3	C	2	Total	Ca	0	0
			2	2		
3	D	2	Total	Ca	0	0
			2	2		
3	E	2	Total	Ca	0	0
			2	2		
3	F	2	Total	Ca	0	0
			2	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

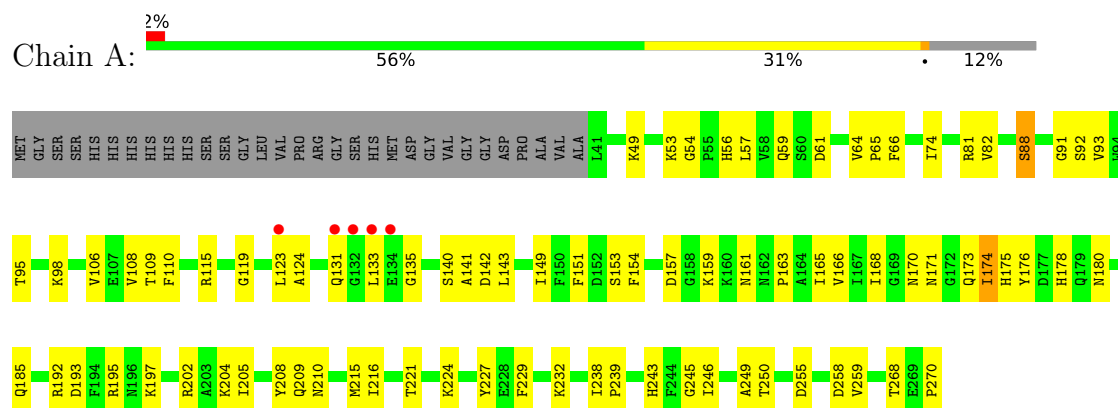
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	61	Total	O	0	0
			61	61		
5	B	62	Total	O	0	0
			62	62		
5	C	64	Total	O	0	0
			64	64		
5	D	57	Total	O	0	0
			57	57		
5	E	42	Total	O	0	0
			42	42		
5	F	47	Total	O	0	0
			47	47		

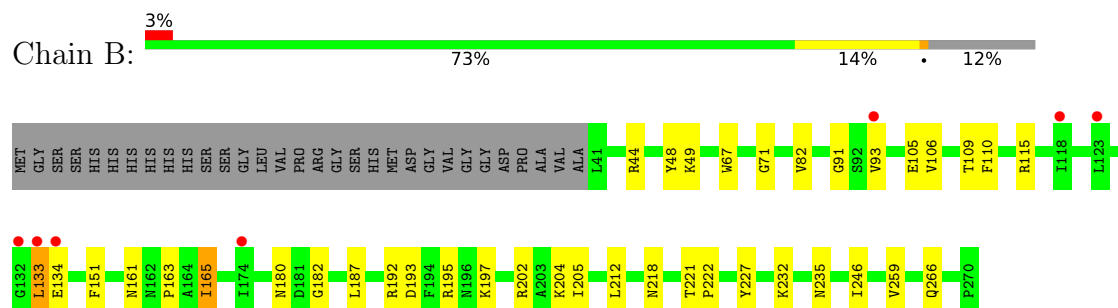
3 Residue-property plots [i](#)

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.

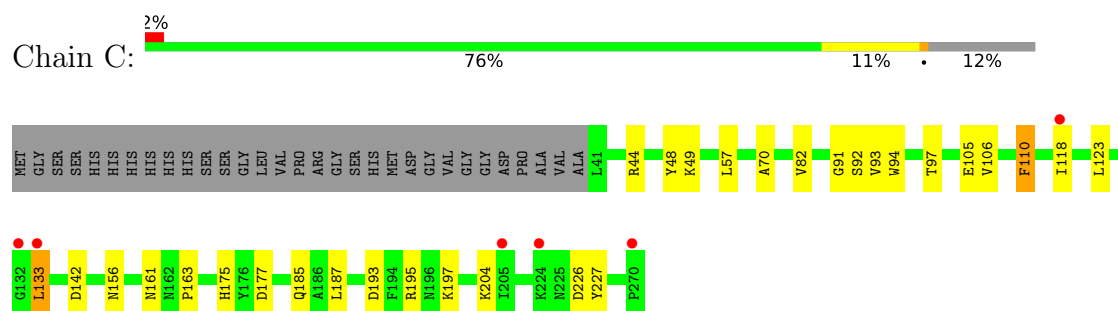
• Molecule 1: Protein ERGIC-53



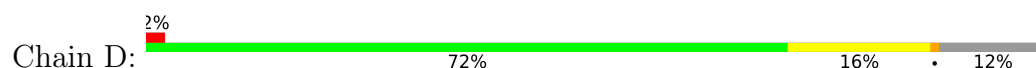
• Molecule 1: Protein ERGIC-53

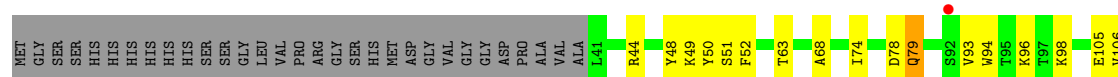


• Molecule 1: Protein ERGIC-53

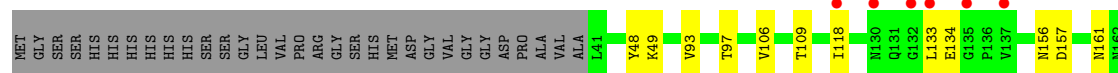
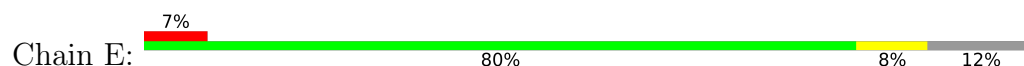


• Molecule 1: Protein ERGIC-53

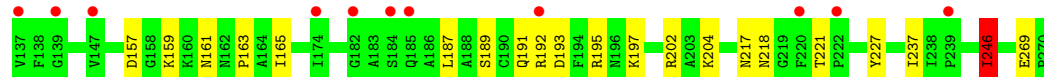
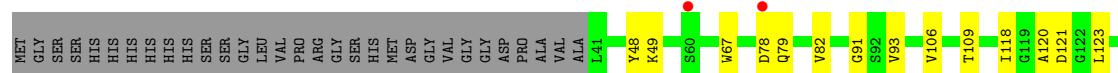
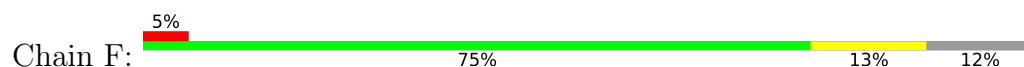




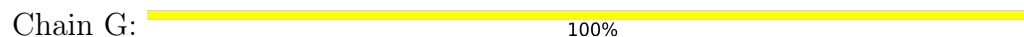
• Molecule 1: Protein ERGIC-53



• Molecule 1: Protein ERGIC-53



• Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose



• Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose



• Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose



• Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain J:  100%

MAN1
MAN2

- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain K:  50%  50%

MAN1
MAN2

- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain L:  100%

MAN1
MAN2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	39.28Å 111.50Å 111.39Å 60.09° 89.21° 86.25°	Depositor
Resolution (Å)	37.19 – 2.70 37.19 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.8 (37.19-2.70) 93.4 (37.19-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.246 , 0.286 0.247 , 0.288	Depositor DCC
R_{free} test set	2172 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	61.8	Xtriage
Anisotropy	0.114	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 36.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.000 for -h,-k+l,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11508	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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: GOL, MAN, CA

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	1.34	4/1858 (0.2%)	1.14	10/2523 (0.4%)
1	B	0.79	1/1858 (0.1%)	0.89	3/2523 (0.1%)
1	C	0.68	0/1858	0.88	2/2523 (0.1%)
1	D	0.73	2/1867 (0.1%)	0.90	1/2534 (0.0%)
1	E	0.60	0/1858	0.84	0/2523
1	F	0.62	1/1858 (0.1%)	0.86	1/2523 (0.0%)
All	All	0.83	8/11157 (0.1%)	0.92	17/15149 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	154	PHE	C-O	-6.08	1.16	1.24
1	A	64	VAL	CA-C	-5.99	1.48	1.52
1	A	106	VAL	C-O	-5.65	1.18	1.24
1	B	165	ILE	C-O	-5.62	1.18	1.24
1	D	125	ILE	C-O	-5.33	1.18	1.24
1	F	246	ILE	C-O	-5.11	1.18	1.24
1	D	79	GLN	C-O	-5.08	1.18	1.23
1	A	221	THR	C-O	-5.01	1.19	1.24

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	SER	N-CA-C	8.27	121.41	111.82
1	A	108	VAL	N-CA-C	6.56	117.33	107.75
1	F	187	LEU	N-CA-C	-5.81	106.03	113.23
1	B	187	LEU	N-CA-C	-5.76	106.09	113.23
1	A	57	LEU	N-CA-C	5.67	117.26	111.14
1	A	54	GLY	CA-C-O	-5.66	116.55	122.38
1	B	133	LEU	N-CA-C	5.50	118.03	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	ALA	N-CA-C	5.49	117.42	109.07
1	A	110	PHE	N-CA-C	5.41	117.01	108.96
1	B	110	PHE	N-CA-C	5.39	116.98	108.96
1	C	187	LEU	N-CA-C	-5.38	106.22	112.89
1	A	174	ILE	CB-CA-C	-5.35	103.89	111.28
1	A	88	SER	N-CA-C	5.28	118.52	111.24
1	A	133	LEU	N-CA-C	5.24	117.79	111.40
1	D	187	LEU	N-CA-C	-5.16	106.84	113.23
1	A	245	GLY	N-CA-C	5.09	120.10	111.19
1	C	110	PHE	N-CA-C	5.08	116.02	108.60

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	1807	0	1713	59	0
1	B	1807	0	1713	30	0
1	C	1807	0	1713	17	0
1	D	1816	0	1725	30	0
1	E	1807	0	1713	12	0
1	F	1807	0	1713	24	0
2	G	23	22	21	0	0
2	H	23	22	21	0	0
2	I	23	22	21	1	0
2	J	23	22	21	0	0
2	K	23	22	21	1	0
2	L	23	22	21	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
4	A	30	0	40	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	6	0	8	1	0
4	C	6	0	8	0	0
5	A	61	0	0	1	0
5	B	62	0	0	2	0
5	C	64	0	0	1	0
5	D	57	0	0	1	0
5	E	42	0	0	0	0
5	F	47	0	0	0	0
All	All	11376	132	10472	165	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 8.

All (165) 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:177:ASP:H	1:C:185:GLN:HE22	1.22	0.86
1:A:49:LYS:HD3	4:A:307:GOL:H11	1.60	0.81
1:A:74:ILE:HD11	1:D:74:ILE:HD11	1.63	0.80
1:A:81:ARG:HG3	1:A:258:ASP:OD1	1.84	0.78
1:A:170:ASN:OD1	1:A:174:ILE:HG13	1.86	0.75
1:A:95:THR:O	1:A:131:GLN:NE2	2.21	0.73
1:A:149:ILE:HD13	1:A:205:ILE:HD13	1.71	0.72
1:A:268:THR:O	1:A:270:PRO:HD3	1.90	0.70
1:D:163:PRO:HG2	1:D:193:ASP:HA	1.73	0.70
1:D:48:TYR:HB3	1:F:237:ILE:HG13	1.75	0.68
1:F:67:TRP:CE3	1:F:93:VAL:HG12	2.29	0.68
1:F:67:TRP:HE3	1:F:93:VAL:HG12	1.59	0.68
1:B:163:PRO:HG2	1:B:193:ASP:HA	1.76	0.68
1:C:94:TRP:CD1	1:C:133:LEU:HD23	2.28	0.67
1:A:143:LEU:HD11	1:A:175:HIS:CE1	2.29	0.67
1:B:266:GLN:CG	5:B:405:HOH:O	2.42	0.67
1:B:266:GLN:HG3	5:B:405:HOH:O	1.94	0.67
1:C:163:PRO:HG2	1:C:193:ASP:HA	1.77	0.66
1:D:202:ARG:HB2	1:D:217:ASN:HB3	1.76	0.66
1:F:163:PRO:HG2	1:F:193:ASP:HA	1.78	0.65
1:F:165:ILE:O	1:F:189:SER:HB2	1.96	0.65
1:A:159:LYS:HB2	1:A:161:ASN:OD1	1.97	0.64
1:E:163:PRO:HG2	1:E:193:ASP:HA	1.78	0.64
1:A:53:LYS:NZ	1:B:235:ASN:O	2.31	0.63
1:A:229:PHE:CE1	4:A:305:GOL:H12	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:GLY:HA3	1:A:140:SER:O	1.99	0.62
1:C:226:ASP:HB3	5:C:447:HOH:O	1.98	0.62
1:A:123:LEU:HG	1:A:124:ALA:N	2.16	0.60
1:F:161:ASN:C	1:F:191:GLN:HE22	2.10	0.60
1:B:133:LEU:HD12	1:B:134:GLU:N	2.17	0.60
1:A:143:LEU:HD11	1:A:175:HIS:ND1	2.17	0.59
1:D:167:ILE:HD11	1:D:214:VAL:HG21	1.84	0.59
1:B:93:VAL:HB	1:B:246:ILE:HG22	1.83	0.59
1:A:171:ASN:HD21	1:A:173:GLN:HG3	1.67	0.59
1:A:135:GLY:HA3	1:A:140:SER:C	2.27	0.59
1:F:93:VAL:HB	1:F:246:ILE:CG2	2.33	0.59
1:A:74:ILE:CD1	1:D:74:ILE:HD11	2.33	0.59
1:D:78:ASP:O	1:D:79:GLN:HB3	2.03	0.58
1:B:93:VAL:HB	1:B:246:ILE:CG2	2.34	0.58
1:A:135:GLY:HA3	1:A:141:ALA:HA	1.85	0.57
1:C:110:PHE:CD2	1:C:123:LEU:HD21	2.39	0.57
1:F:109:THR:HG23	1:F:202:ARG:HG3	1.86	0.57
1:B:133:LEU:HD12	1:B:133:LEU:C	2.30	0.56
1:D:135:GLY:HA3	1:D:140:SER:O	2.06	0.56
1:B:221:THR:OG1	1:B:222:PRO:HD2	2.06	0.56
1:D:167:ILE:HD11	1:D:214:VAL:CG2	2.37	0.55
1:A:229:PHE:CZ	4:A:305:GOL:H12	2.40	0.55
1:F:193:ASP:O	1:F:197:LYS:HE3	2.07	0.55
1:A:232:LYS:HG2	4:A:305:GOL:H2	1.88	0.55
1:D:48:TYR:CB	1:F:237:ILE:HG13	2.37	0.55
1:B:67:TRP:CE3	1:B:93:VAL:HG12	2.41	0.54
1:F:157:ASP:OD2	1:F:159:LYS:HD2	2.08	0.54
1:A:98:LYS:HE2	1:A:243:HIS:NE2	2.23	0.54
1:C:94:TRP:HD1	1:C:133:LEU:HD23	1.73	0.54
1:E:133:LEU:HD12	1:E:134:GLU:N	2.22	0.54
1:F:192:ARG:HD2	1:F:218:ASN:HD22	1.72	0.53
1:A:208:TYR:CD2	1:A:209:GLN:HG3	2.44	0.53
1:D:135:GLY:C	1:D:140:SER:O	2.52	0.53
1:B:109:THR:HG23	1:B:202:ARG:CG	2.40	0.52
1:D:135:GLY:CA	1:D:140:SER:O	2.58	0.52
1:B:67:TRP:HE3	1:B:93:VAL:HG12	1.74	0.52
1:B:48:TYR:OH	1:B:49:LYS:HE2	2.10	0.52
1:E:133:LEU:HD12	1:E:134:GLU:CB	2.40	0.52
1:A:151:PHE:CE2	1:A:165:ILE:HD13	2.45	0.52
1:E:193:ASP:O	1:E:197:LYS:HE3	2.09	0.52
1:C:193:ASP:O	1:C:197:LYS:HE3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:ALA:HB2	1:D:96[B]:LYS:HD3	1.92	0.51
1:F:93:VAL:HB	1:F:246:ILE:HG22	1.92	0.51
1:A:166:VAL:HG12	1:A:168:ILE:HG13	1.92	0.51
1:A:215:MET:O	1:A:216:ILE:HD13	2.09	0.51
1:E:48:TYR:OH	1:E:49:LYS:HE2	2.11	0.51
1:A:174:ILE:HG22	1:A:175:HIS:N	2.24	0.51
1:D:48:TYR:OH	1:D:49:LYS:HE2	2.11	0.51
1:A:178:HIS:O	1:A:180:ASN:N	2.44	0.50
1:B:193:ASP:O	1:B:197:LYS:HE3	2.12	0.50
1:B:71:GLY:HA2	4:B:304:GOL:H12	1.94	0.50
1:B:151:PHE:CE2	1:B:165:ILE:HD13	2.47	0.49
1:F:78:ASP:OD1	1:F:79:GLN:HG2	2.12	0.49
1:A:98:LYS:HG3	1:A:243:HIS:CD2	2.47	0.49
1:A:59:GLN:HB2	1:A:61:ASP:OD1	2.13	0.49
1:A:142:ASP:CG	1:A:143:LEU:HD12	2.38	0.49
1:E:109:THR:HG23	1:E:202:ARG:HG2	1.94	0.49
1:B:195:ARG:O	1:B:197:LYS:HE2	2.13	0.48
1:F:48:TYR:OH	1:F:49:LYS:HE2	2.14	0.48
1:A:123:LEU:HG	1:A:124:ALA:H	1.79	0.47
1:F:269:GLU:O	1:F:269:GLU:HG3	2.14	0.47
1:A:193:ASP:O	1:A:197:LYS:HE3	2.15	0.47
1:A:171:ASN:ND2	1:A:173:GLN:HG3	2.29	0.47
1:A:176:TYR:HA	1:A:185:GLN:OE1	2.15	0.47
1:C:161:ASN:OD1	1:C:161:ASN:N	2.48	0.47
1:F:195:ARG:O	1:F:197:LYS:HE2	2.15	0.47
1:C:48:TYR:OH	1:C:49:LYS:HE2	2.14	0.46
1:A:238:ILE:HB	1:A:239:PRO:CD	2.46	0.46
1:D:108:VAL:HG21	1:D:125:ILE:HD13	1.98	0.46
1:F:192:ARG:HD2	1:F:218:ASN:ND2	2.29	0.46
1:A:88:SER:HA	1:A:250:THR:O	2.15	0.46
1:A:109:THR:HG23	1:A:202:ARG:HG2	1.97	0.46
1:A:209:GLN:O	1:A:210:ASN:HB2	2.16	0.46
1:A:224:LYS:HB2	1:A:224:LYS:HE2	1.80	0.45
1:B:212:LEU:O	1:B:232:LYS:HA	2.16	0.45
1:A:142:ASP:O	1:A:143:LEU:HB2	2.15	0.45
1:B:204:LYS:HB2	1:B:227:TYR:CE2	2.51	0.45
1:E:161:ASN:OD1	1:E:161:ASN:N	2.49	0.45
1:A:178:HIS:CE1	5:A:447:HOH:O	2.69	0.45
1:A:157:ASP:OD2	1:A:159:LYS:HG3	2.17	0.45
1:A:195:ARG:O	1:A:197:LYS:HE2	2.16	0.45
1:E:212:LEU:O	1:E:232:LYS:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:ARG:HD2	1:B:218:ASN:HD22	1.82	0.44
1:D:68:ALA:HB2	1:D:96[B]:LYS:CD	2.47	0.44
1:A:208:TYR:CE2	1:A:209:GLN:HG3	2.51	0.44
1:D:195:ARG:O	1:D:197:LYS:HE2	2.17	0.44
1:A:56:HIS:O	1:A:56:HIS:CG	2.70	0.44
1:B:246:ILE:CD1	1:B:259:VAL:HG21	2.47	0.44
1:C:123:LEU:HD12	1:C:123:LEU:C	2.42	0.44
1:C:195:ARG:O	1:C:197:LYS:HE2	2.18	0.44
1:D:221:THR:OG1	1:D:222:PRO:HD2	2.17	0.44
1:A:115:ARG:HH22	1:D:113:THR:HG21	1.82	0.44
1:A:161:ASN:OD1	1:A:161:ASN:N	2.51	0.44
1:D:161:ASN:OD1	1:D:161:ASN:N	2.50	0.44
1:A:65:PRO:O	1:A:66:PHE:HB2	2.18	0.44
1:F:161:ASN:OD1	1:F:161:ASN:N	2.50	0.44
1:A:204:LYS:HB2	1:A:227:TYR:CE2	2.53	0.44
1:D:51:SER:HB3	1:D:263:LEU:HA	2.00	0.43
1:E:133:LEU:HD12	1:E:134:GLU:HB2	2.00	0.43
1:F:202:ARG:HB2	1:F:217:ASN:HB3	2.00	0.43
1:C:44:ARG:HD3	1:C:105:GLU:OE2	2.18	0.43
1:C:204:LYS:HB2	1:C:227:TYR:CE2	2.53	0.43
1:A:82:VAL:O	1:A:91:GLY:HA3	2.19	0.43
1:A:49:LYS:HD3	4:A:307:GOL:C1	2.40	0.42
1:A:119:GLY:HA3	1:A:255:ASP:OD2	2.18	0.42
1:A:229:PHE:CD1	1:A:229:PHE:C	2.97	0.42
1:D:44:ARG:HD3	1:D:105:GLU:OE2	2.18	0.42
1:D:50:TYR:OH	1:D:98:LYS:O	2.27	0.42
1:E:157:ASP:HB3	1:E:180:ASN:HA	2.01	0.42
1:A:151:PHE:CD1	1:A:151:PHE:N	2.88	0.42
1:A:163:PRO:HB2	1:A:192:ARG:O	2.19	0.42
1:B:106:VAL:CG2	1:B:205:ILE:HB	2.49	0.42
1:E:195:ARG:O	1:E:197:LYS:HE2	2.19	0.42
1:D:204:LYS:HB2	1:D:227:TYR:CE2	2.55	0.42
1:F:120:ALA:HA	1:F:121:ASP:HA	1.81	0.42
1:A:216:ILE:HD12	1:A:216:ILE:HG23	1.86	0.41
1:B:109:THR:HG23	1:B:202:ARG:HG3	2.01	0.41
1:A:91:GLY:O	1:A:92:SER:HB3	2.19	0.41
1:F:82:VAL:O	1:F:91:GLY:HA3	2.20	0.41
1:B:133:LEU:HD12	1:B:134:GLU:HG2	2.01	0.41
1:B:161:ASN:OD1	1:B:161:ASN:N	2.51	0.41
1:D:197:LYS:HD2	1:D:220:PHE:CE1	2.55	0.41
1:E:156:ASN:ND2	2:K:2:MAN:O4	2.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:VAL:O	1:C:91:GLY:HA3	2.21	0.41
1:D:63:THR:HB	5:D:408:HOH:O	2.19	0.41
1:C:156:ASN:ND2	2:I:2:MAN:O4	2.48	0.41
1:B:44:ARG:HD3	1:B:105:GLU:OE2	2.21	0.41
1:B:192:ARG:HG3	1:B:218:ASN:ND2	2.36	0.41
1:D:246:ILE:CD1	1:D:259:VAL:HG21	2.50	0.41
1:B:109:THR:HG23	1:B:202:ARG:HG2	2.02	0.41
1:B:180:ASN:C	1:B:182:GLY:N	2.78	0.41
1:C:142:ASP:OD2	1:C:175:HIS:HB2	2.21	0.41
1:F:123:LEU:HD12	1:F:123:LEU:C	2.46	0.41
1:A:74:ILE:HD11	1:D:74:ILE:CD1	2.40	0.40
1:B:82:VAL:O	1:B:91:GLY:HA3	2.21	0.40
1:D:52:PHE:CD1	1:D:52:PHE:C	2.99	0.40
1:A:246:ILE:CD1	1:A:259:VAL:HG21	2.52	0.40
1:C:70:ALA:HB3	1:C:92:SER:HB2	2.02	0.40
1:D:94:TRP:HB3	1:D:131:GLN:HG2	2.04	0.40
1:F:204:LYS:HB2	1:F:227:TYR:CE2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	228/261 (87%)	217 (95%)	11 (5%)	0	100	100
1	B	228/261 (87%)	216 (95%)	12 (5%)	0	100	100
1	C	228/261 (87%)	220 (96%)	8 (4%)	0	100	100
1	D	229/261 (88%)	221 (96%)	8 (4%)	0	100	100
1	E	228/261 (87%)	215 (94%)	13 (6%)	0	100	100
1	F	228/261 (87%)	217 (95%)	11 (5%)	0	100	100
All	All	1369/1566 (87%)	1306 (95%)	63 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	188/211 (89%)	187 (100%)	1 (0%)	81	92
1	B	188/211 (89%)	187 (100%)	1 (0%)	81	92
1	C	188/211 (89%)	182 (97%)	6 (3%)	34	64
1	D	189/211 (90%)	185 (98%)	4 (2%)	47	75
1	E	188/211 (89%)	184 (98%)	4 (2%)	47	75
1	F	188/211 (89%)	184 (98%)	4 (2%)	47	75
All	All	1129/1266 (89%)	1109 (98%)	20 (2%)	51	78

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	VAL
1	B	115	ARG
1	C	57	LEU
1	C	93	VAL
1	C	97	THR
1	C	106	VAL
1	C	118	ILE
1	C	133	LEU
1	D	93	VAL
1	D	106	VAL
1	D	117	ARG
1	D	118	ILE
1	E	93	VAL
1	E	97	THR
1	E	106	VAL
1	E	118	ILE
1	F	106	VAL
1	F	118	ILE
1	F	221	THR

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Mol	Chain	Res	Type
1	F	246	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	HIS
1	A	178	HIS
1	A	218	ASN
1	A	225	ASN
1	B	69	HIS
1	B	179	GLN
1	B	180	ASN
1	B	191	GLN
1	B	209	GLN
1	B	210	ASN
1	B	218	ASN
1	C	89	GLN
1	C	185	GLN
1	C	218	ASN
1	D	89	GLN
1	D	218	ASN
1	E	56	HIS
1	E	89	GLN
1	F	43	HIS
1	F	178	HIS
1	F	191	GLN
1	F	218	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

12 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	MAN	G	1	2	12,12,12	0.58	0	17,17,17	1.58	4 (23%)
2	MAN	G	2	2	11,11,12	0.83	0	15,15,17	1.94	5 (33%)
2	MAN	H	1	2	12,12,12	0.57	0	17,17,17	1.59	4 (23%)
2	MAN	H	2	2	11,11,12	0.83	0	15,15,17	1.94	5 (33%)
2	MAN	I	1	2	12,12,12	0.58	0	17,17,17	1.59	4 (23%)
2	MAN	I	2	2	11,11,12	0.83	0	15,15,17	1.94	5 (33%)
2	MAN	J	1	2	12,12,12	0.58	0	17,17,17	1.59	4 (23%)
2	MAN	J	2	2	11,11,12	0.83	0	15,15,17	1.94	5 (33%)
2	MAN	K	1	2	12,12,12	0.57	0	17,17,17	1.59	4 (23%)
2	MAN	K	2	2	11,11,12	0.84	0	15,15,17	1.94	5 (33%)
2	MAN	L	1	2	12,12,12	0.58	0	17,17,17	1.58	4 (23%)
2	MAN	L	2	2	11,11,12	0.83	0	15,15,17	1.95	5 (33%)

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	MAN	G	1	2	-	2/2/22/22	0/1/1/1
2	MAN	G	2	2	-	0/2/19/22	0/1/1/1
2	MAN	H	1	2	-	2/2/22/22	0/1/1/1
2	MAN	H	2	2	-	0/2/19/22	0/1/1/1
2	MAN	I	1	2	-	2/2/22/22	0/1/1/1
2	MAN	I	2	2	-	0/2/19/22	0/1/1/1
2	MAN	J	1	2	-	2/2/22/22	0/1/1/1
2	MAN	J	2	2	-	0/2/19/22	0/1/1/1
2	MAN	K	1	2	-	2/2/22/22	0/1/1/1
2	MAN	K	2	2	-	0/2/19/22	0/1/1/1
2	MAN	L	1	2	-	2/2/22/22	0/1/1/1
2	MAN	L	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	2	MAN	C2-C3-C4	3.60	117.19	110.86
2	L	2	MAN	C2-C3-C4	3.59	117.18	110.86
2	I	2	MAN	C2-C3-C4	3.58	117.17	110.86
2	H	2	MAN	C2-C3-C4	3.58	117.16	110.86
2	G	2	MAN	C2-C3-C4	3.58	117.16	110.86
2	K	2	MAN	C2-C3-C4	3.57	117.15	110.86
2	J	2	MAN	O5-C5-C6	-3.49	100.86	107.66
2	G	2	MAN	O5-C5-C6	-3.49	100.87	107.66
2	I	2	MAN	O5-C5-C6	-3.49	100.88	107.66
2	L	2	MAN	O5-C5-C6	-3.48	100.89	107.66
2	K	2	MAN	O5-C5-C6	-3.48	100.90	107.66
2	H	2	MAN	O5-C5-C6	-3.47	100.91	107.66
2	L	2	MAN	O5-C1-C2	-2.92	103.81	110.79
2	H	2	MAN	O5-C1-C2	-2.91	103.84	110.79
2	G	2	MAN	O5-C1-C2	-2.91	103.85	110.79
2	I	2	MAN	O5-C1-C2	-2.91	103.85	110.79
2	J	2	MAN	O5-C1-C2	-2.89	103.89	110.79
2	K	2	MAN	O5-C1-C2	-2.89	103.89	110.79
2	H	2	MAN	C1-C2-C3	2.88	113.84	109.64
2	L	2	MAN	C1-C2-C3	2.88	113.84	109.64
2	J	2	MAN	C1-C2-C3	2.87	113.83	109.64
2	G	2	MAN	C1-C2-C3	2.87	113.82	109.64
2	K	2	MAN	C1-C2-C3	2.85	113.80	109.64
2	I	2	MAN	C1-C2-C3	2.84	113.78	109.64
2	J	1	MAN	C1-C2-C3	2.62	115.70	110.36
2	L	1	MAN	C1-C2-C3	2.62	115.70	110.36
2	H	1	MAN	C1-C2-C3	2.61	115.69	110.36
2	G	1	MAN	C1-C2-C3	2.61	115.69	110.36
2	I	1	MAN	C1-C2-C3	2.60	115.66	110.36
2	K	1	MAN	C1-C2-C3	2.60	115.66	110.36
2	H	1	MAN	O5-C5-C6	2.59	112.85	106.44
2	J	1	MAN	O5-C5-C6	2.58	112.84	106.44
2	I	1	MAN	O5-C5-C6	2.58	112.83	106.44
2	K	1	MAN	O5-C5-C6	2.58	112.83	106.44
2	L	1	MAN	O5-C5-C6	2.58	112.83	106.44
2	G	1	MAN	O5-C5-C6	2.57	112.81	106.44
2	I	2	MAN	O3-C3-C2	-2.42	105.12	110.05
2	G	2	MAN	O3-C3-C2	-2.41	105.14	110.05
2	K	2	MAN	O3-C3-C2	-2.41	105.14	110.05
2	L	2	MAN	O3-C3-C2	-2.40	105.15	110.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	2	MAN	O3-C3-C2	-2.40	105.16	110.05
2	H	2	MAN	O3-C3-C2	-2.39	105.17	110.05
2	K	1	MAN	O5-C1-C2	2.39	114.50	110.30
2	L	1	MAN	O5-C1-C2	2.38	114.48	110.30
2	I	1	MAN	O5-C1-C2	2.38	114.48	110.30
2	G	1	MAN	O5-C1-C2	2.38	114.48	110.30
2	J	1	MAN	O5-C1-C2	2.37	114.47	110.30
2	H	1	MAN	O5-C1-C2	2.37	114.46	110.30
2	K	1	MAN	O2-C2-C3	-2.28	104.99	110.38
2	L	1	MAN	O2-C2-C3	-2.28	105.01	110.38
2	I	1	MAN	O2-C2-C3	-2.27	105.03	110.38
2	J	1	MAN	O2-C2-C3	-2.27	105.03	110.38
2	G	1	MAN	O2-C2-C3	-2.26	105.04	110.38
2	H	1	MAN	O2-C2-C3	-2.26	105.05	110.38

There are no chirality outliers.

All (12) torsion outliers are listed below:

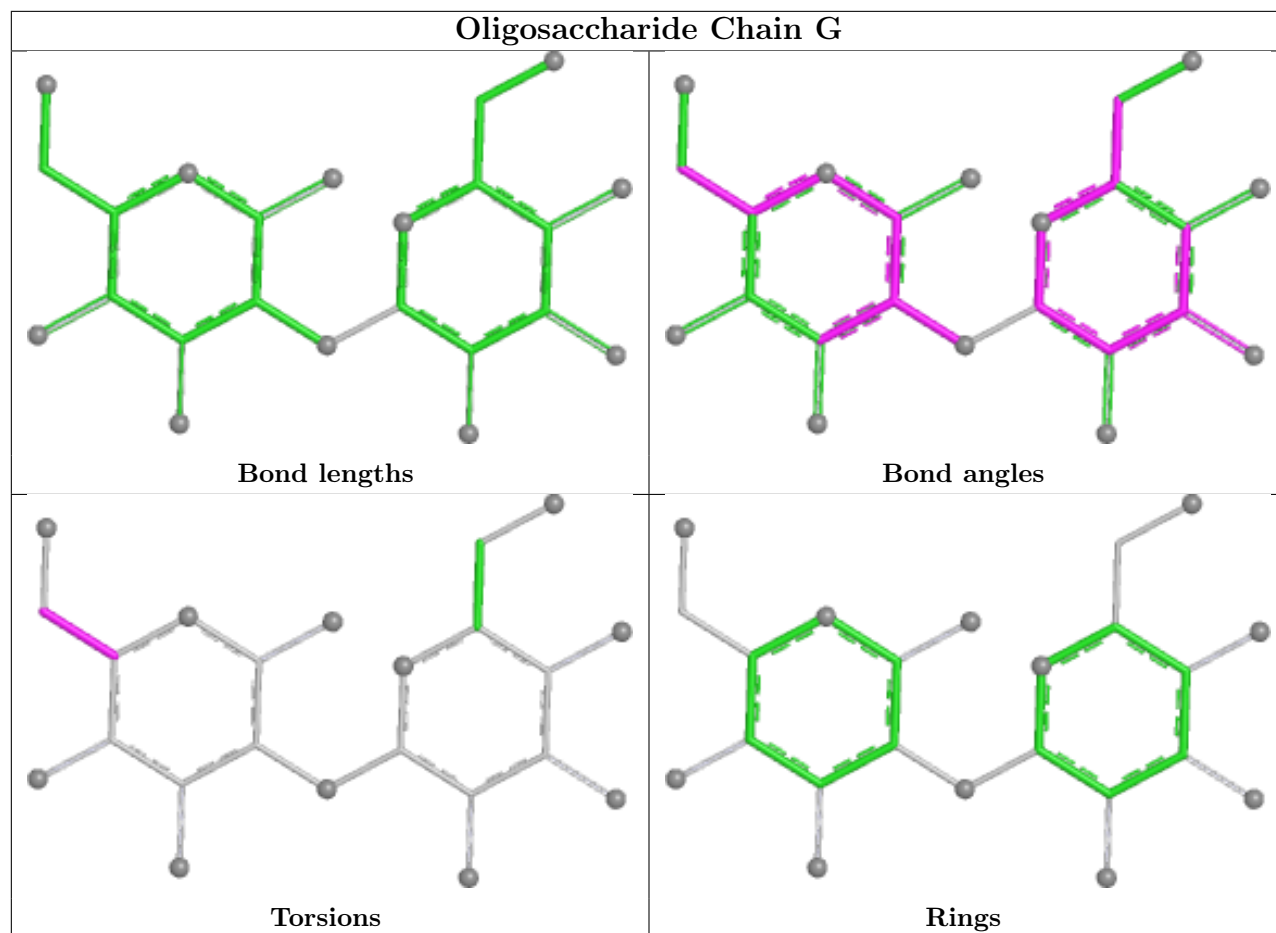
Mol	Chain	Res	Type	Atoms
2	G	1	MAN	C4-C5-C6-O6
2	H	1	MAN	C4-C5-C6-O6
2	I	1	MAN	C4-C5-C6-O6
2	J	1	MAN	C4-C5-C6-O6
2	K	1	MAN	C4-C5-C6-O6
2	L	1	MAN	C4-C5-C6-O6
2	G	1	MAN	O5-C5-C6-O6
2	H	1	MAN	O5-C5-C6-O6
2	I	1	MAN	O5-C5-C6-O6
2	J	1	MAN	O5-C5-C6-O6
2	K	1	MAN	O5-C5-C6-O6
2	L	1	MAN	O5-C5-C6-O6

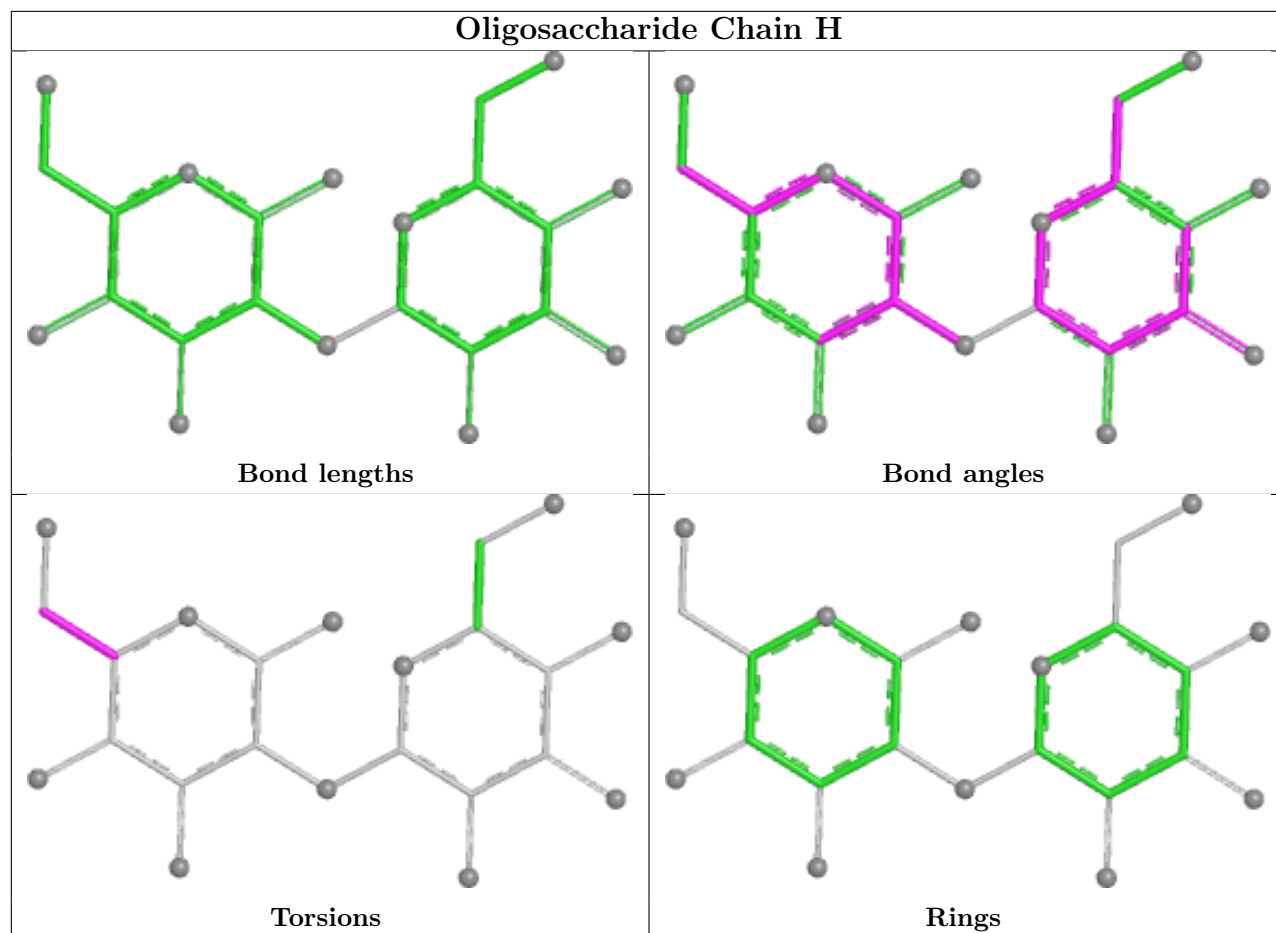
There are no ring outliers.

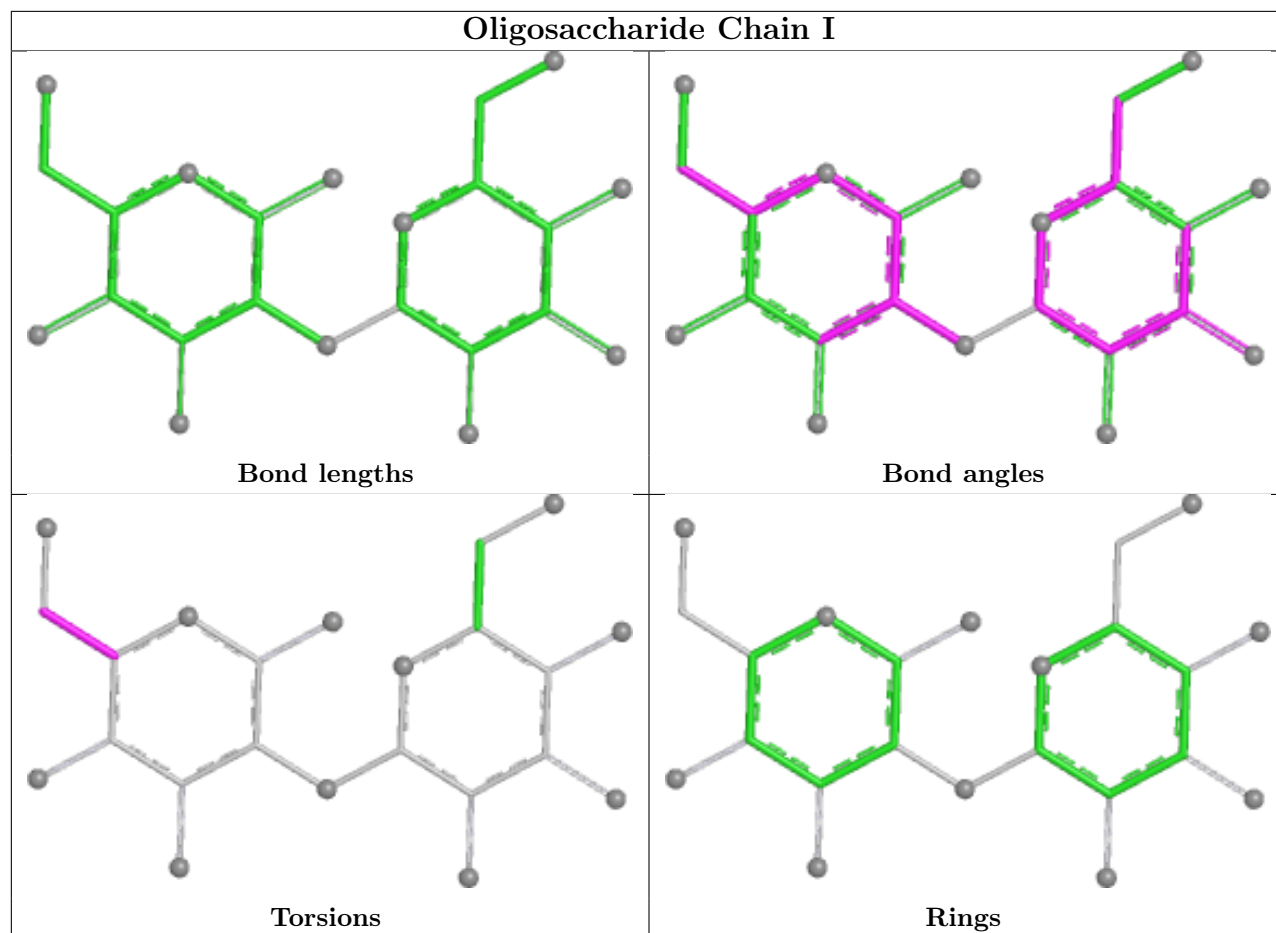
2 monomers are involved in 2 short contacts:

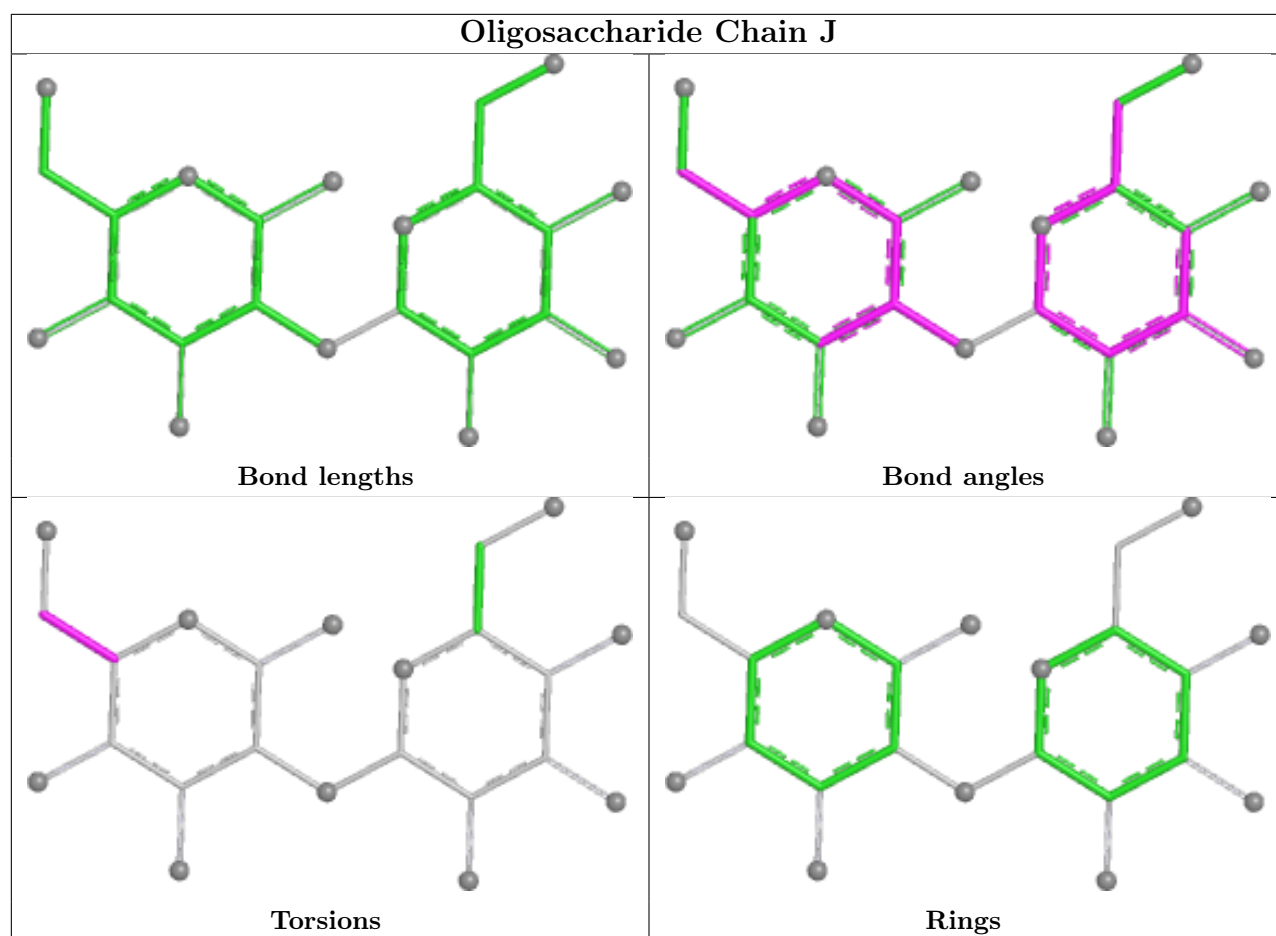
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	2	MAN	1	0
2	I	2	MAN	1	0

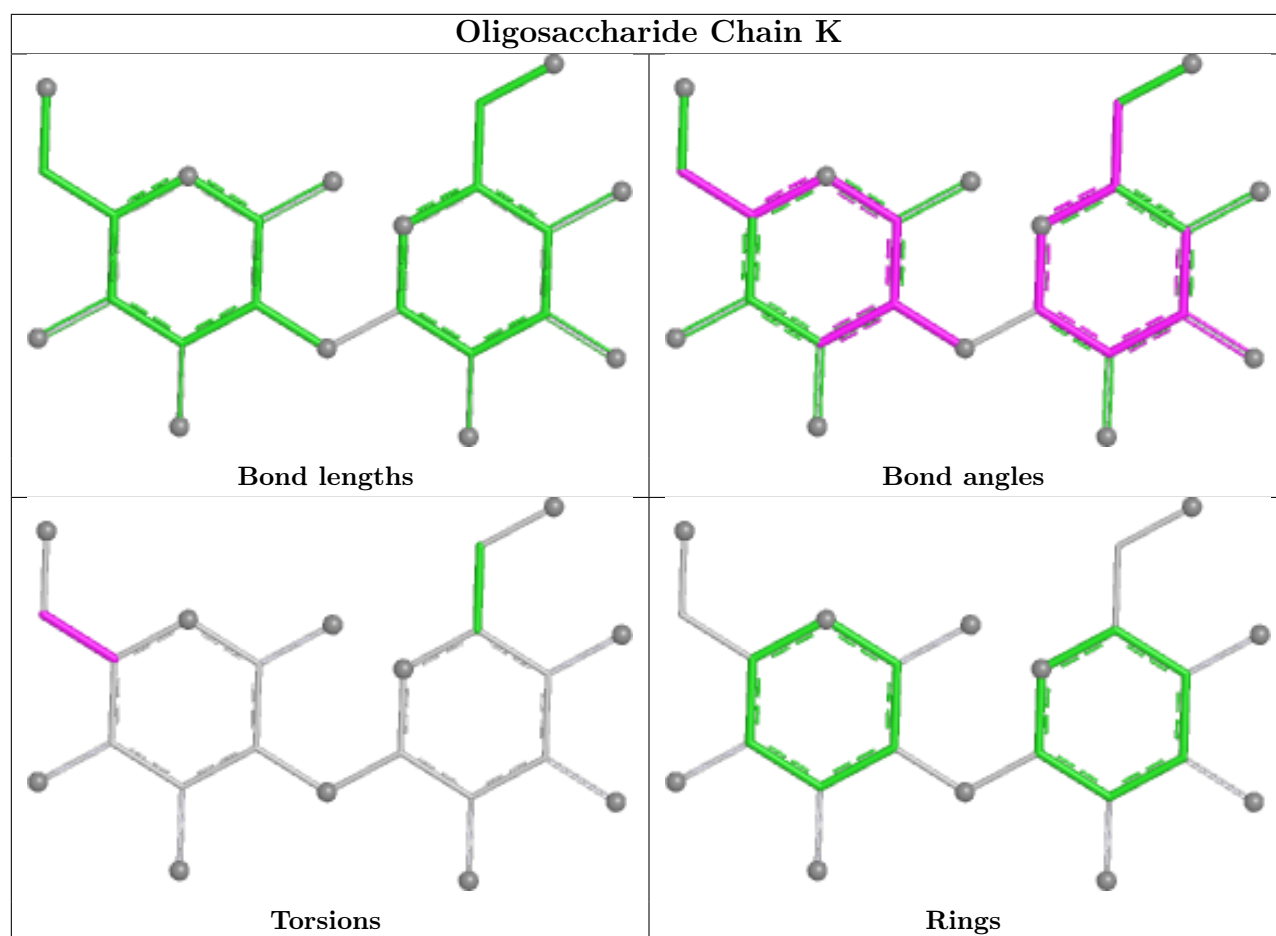
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

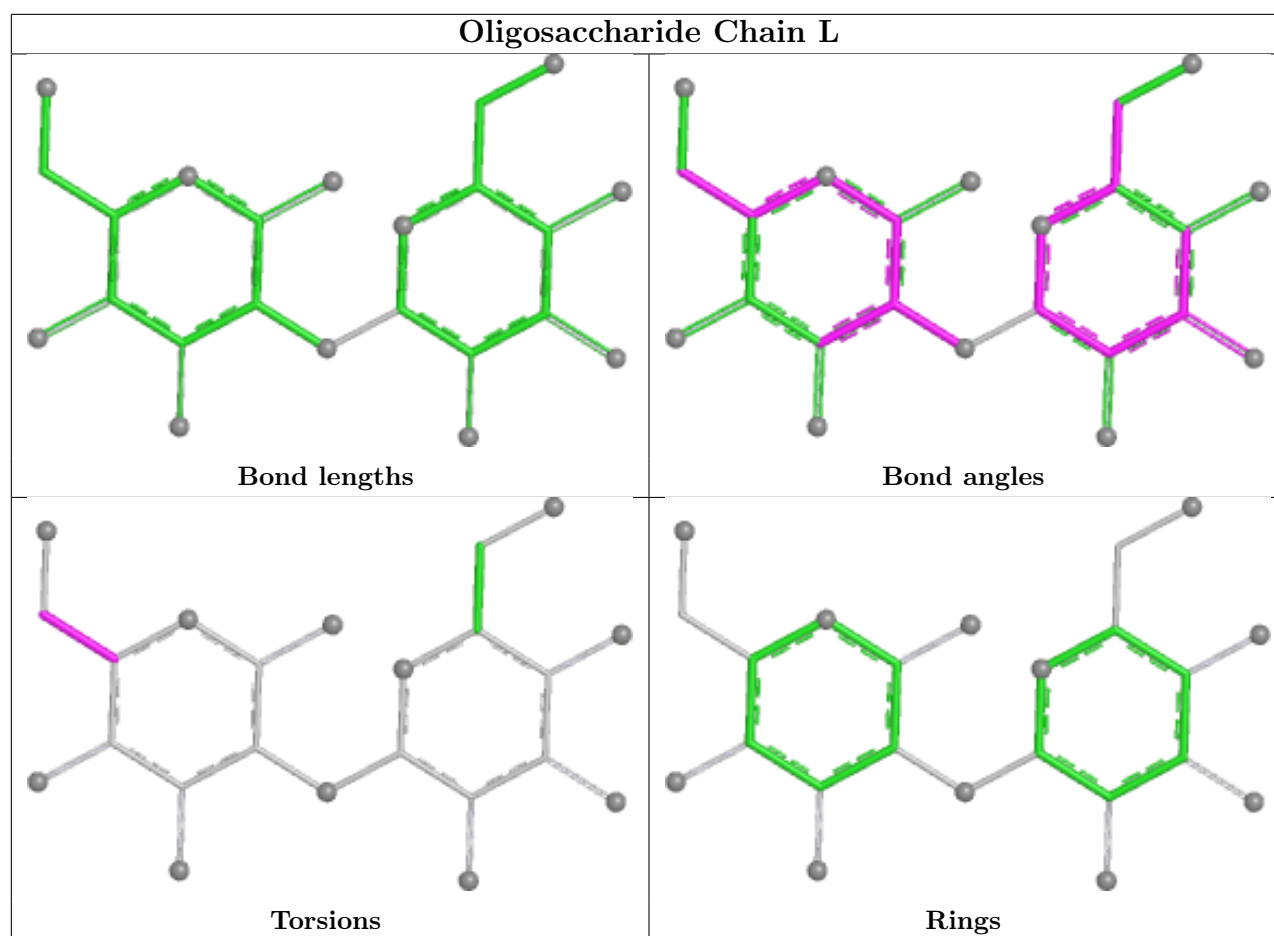












5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 12 are monoatomic - leaving 7 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	GOL	B	304	-	5,5,5	0.30	0	5,5,5	0.47	0
4	GOL	A	308	-	5,5,5	0.23	0	5,5,5	0.38	0
4	GOL	A	305	-	5,5,5	0.26	0	5,5,5	0.31	0
4	GOL	A	304	-	5,5,5	0.21	0	5,5,5	0.35	0
4	GOL	C	304	-	5,5,5	0.17	0	5,5,5	0.33	0
4	GOL	A	306	-	5,5,5	0.24	0	5,5,5	0.21	0
4	GOL	A	307	-	5,5,5	0.19	0	5,5,5	0.33	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	304	-	-	2/4/4/4	-
4	GOL	A	308	-	-	0/4/4/4	-
4	GOL	A	305	-	-	2/4/4/4	-
4	GOL	A	304	-	-	0/4/4/4	-
4	GOL	C	304	-	-	2/4/4/4	-
4	GOL	A	306	-	-	4/4/4/4	-
4	GOL	A	307	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	305	GOL	C1-C2-C3-O3
4	A	305	GOL	O2-C2-C3-O3
4	A	306	GOL	O1-C1-C2-C3
4	A	307	GOL	O1-C1-C2-O2
4	A	307	GOL	O1-C1-C2-C3
4	B	304	GOL	C1-C2-C3-O3
4	C	304	GOL	C1-C2-C3-O3
4	A	306	GOL	O1-C1-C2-O2
4	A	306	GOL	O2-C2-C3-O3
4	A	306	GOL	C1-C2-C3-O3
4	B	304	GOL	O2-C2-C3-O3
4	C	304	GOL	O2-C2-C3-O3
4	A	307	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	304	GOL	1	0
4	A	305	GOL	3	0
4	A	307	GOL	2	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, 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	230/261 (88%)	0.30	5 (2%) 62 59	23, 36, 57, 83	0
1	B	230/261 (88%)	0.30	7 (3%) 52 49	20, 37, 61, 93	0
1	C	230/261 (88%)	0.44	6 (2%) 57 54	24, 44, 69, 88	0
1	D	230/261 (88%)	0.57	5 (2%) 62 59	20, 46, 82, 96	1 (0%)
1	E	230/261 (88%)	0.63	18 (7%) 19 16	30, 49, 82, 103	0
1	F	230/261 (88%)	0.66	13 (5%) 29 26	26, 50, 82, 114	0
All	All	1380/1566 (88%)	0.48	54 (3%) 43 39	20, 44, 77, 114	1 (0%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	LEU	4.8
1	B	133	LEU	4.7
1	C	132	GLY	4.2
1	E	174	ILE	3.7
1	E	216	ILE	3.6
1	D	221	THR	3.6
1	B	134	GLU	3.5
1	B	132	GLY	3.2
1	B	123	LEU	3.1
1	E	241	GLN	3.0
1	F	220	PHE	3.0
1	C	133	LEU	3.0
1	D	222	PRO	2.9
1	E	185	GLN	2.9
1	B	93	VAL	2.9
1	F	60	SER	2.8
1	E	133	LEU	2.8
1	F	185	GLN	2.7
1	F	182	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	132	GLY	2.7
1	C	224	LYS	2.7
1	A	131	GLN	2.7
1	E	182	GLY	2.6
1	F	192	ARG	2.6
1	F	174	ILE	2.6
1	C	270	PRO	2.6
1	E	130	ASN	2.5
1	E	135	GLY	2.5
1	F	184	SER	2.4
1	F	78	ASP	2.4
1	F	239	PRO	2.4
1	E	230	CYS	2.4
1	E	179	GLN	2.4
1	F	147	VAL	2.4
1	C	118	ILE	2.4
1	C	205	ILE	2.4
1	E	218	ASN	2.3
1	E	245	GLY	2.3
1	F	137	VAL	2.3
1	D	92	SER	2.2
1	A	134	GLU	2.2
1	F	222	PRO	2.2
1	B	118	ILE	2.2
1	A	123	LEU	2.1
1	E	202	ARG	2.1
1	D	137	VAL	2.1
1	D	148	GLY	2.1
1	E	132	GLY	2.1
1	E	137	VAL	2.1
1	F	139	GLY	2.1
1	B	174	ILE	2.1
1	E	220	PHE	2.1
1	E	223	ASP	2.0
1	E	118	ILE	2.0

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

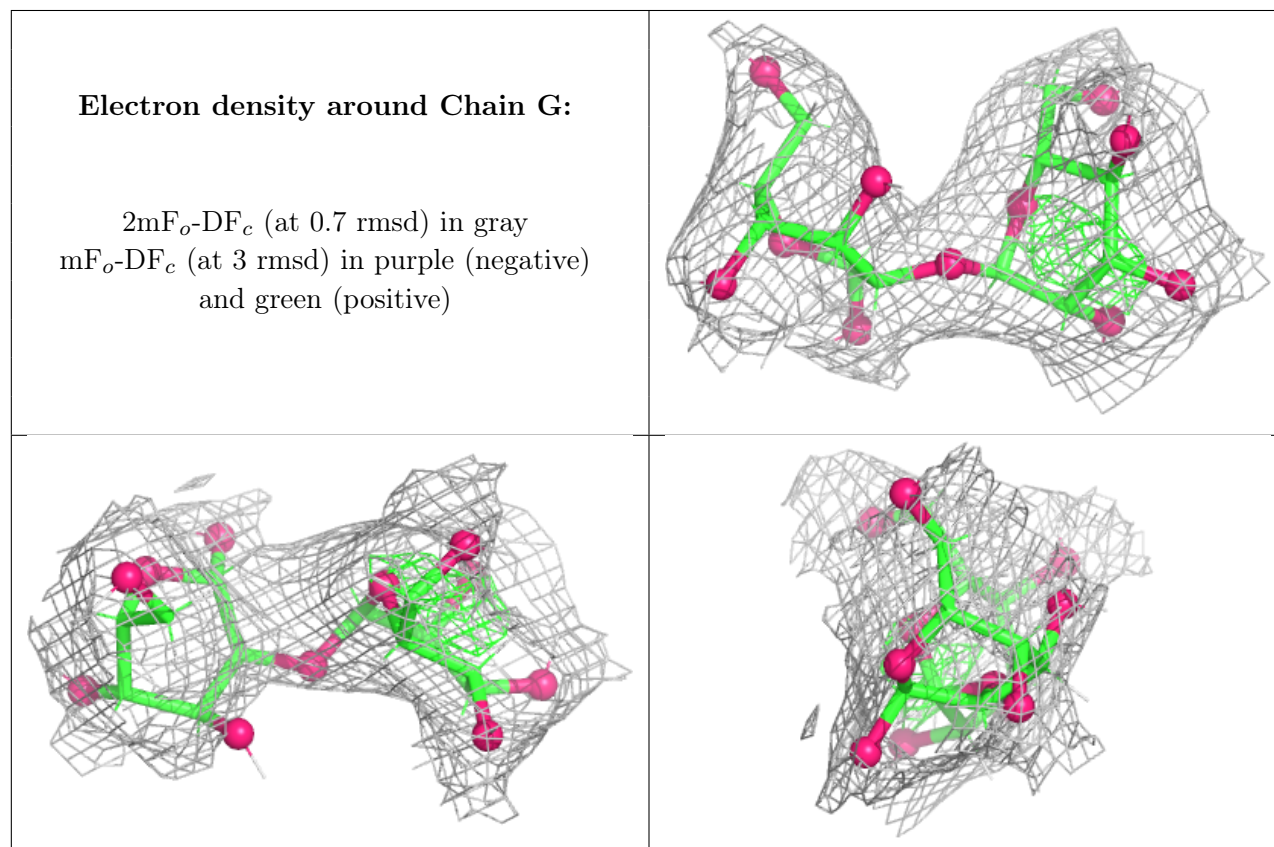
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [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.

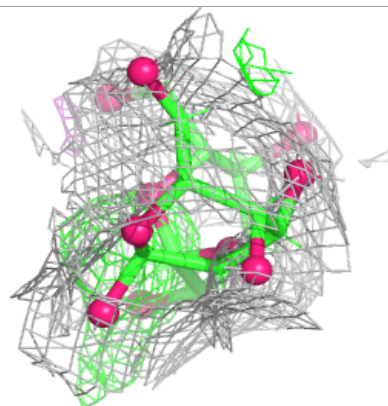
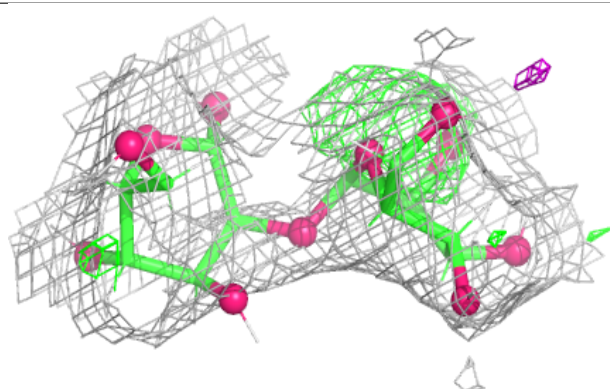
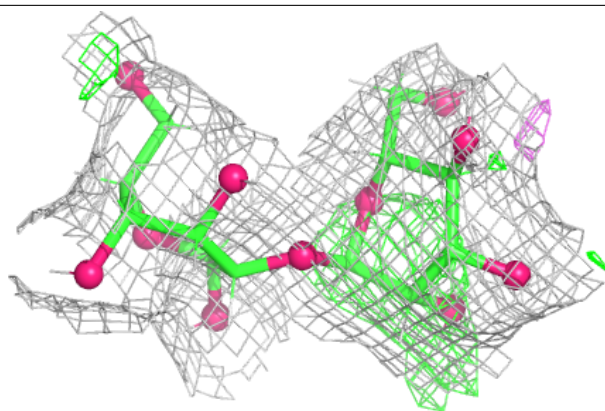
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MAN	K	1	12/12	0.58	0.22	67,82,98,101	23
2	MAN	L	1	12/12	0.65	0.22	61,67,80,91	23
2	MAN	H	1	12/12	0.66	0.16	58,76,92,93	23
2	MAN	I	1	12/12	0.67	0.18	53,63,75,83	23
2	MAN	J	1	12/12	0.72	0.18	56,70,79,88	23
2	MAN	L	2	11/12	0.72	0.23	47,57,69,71	22
2	MAN	G	1	12/12	0.73	0.14	50,60,72,75	23
2	MAN	J	2	11/12	0.74	0.26	46,59,68,74	22
2	MAN	K	2	11/12	0.79	0.20	53,62,69,74	22
2	MAN	I	2	11/12	0.80	0.22	38,51,60,62	22
2	MAN	G	2	11/12	0.83	0.16	36,44,52,56	22
2	MAN	H	2	11/12	0.84	0.16	39,51,59,63	22

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.

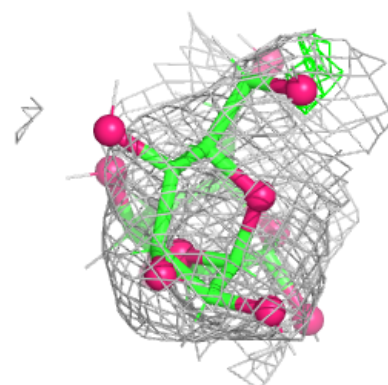
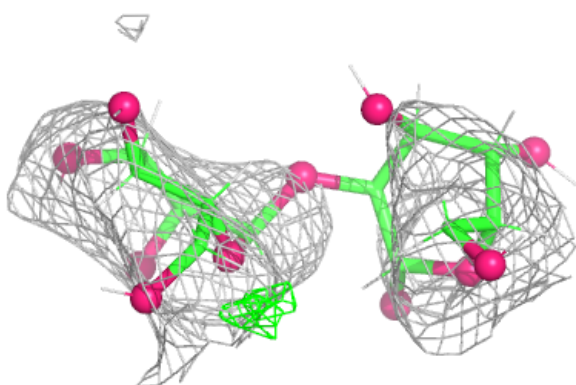
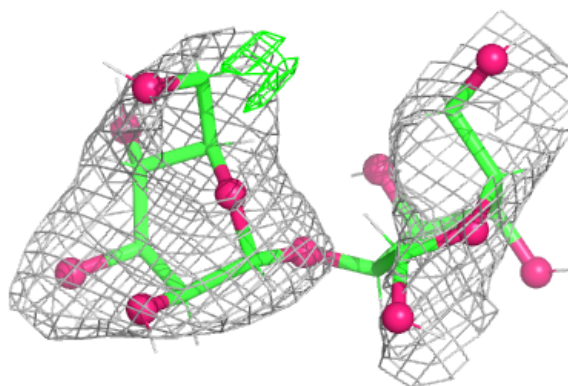


Electron density around Chain H:

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

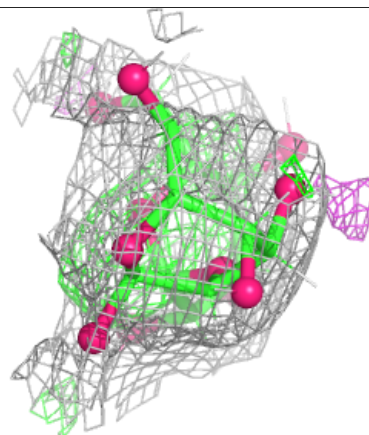
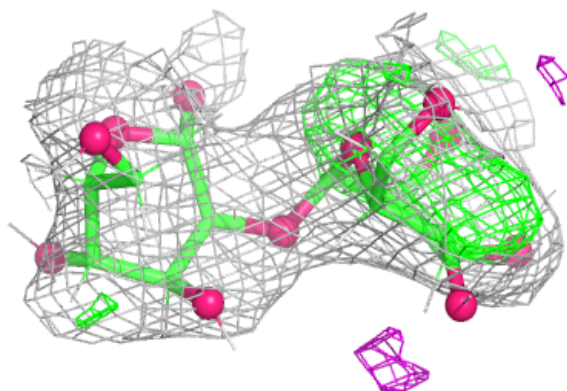
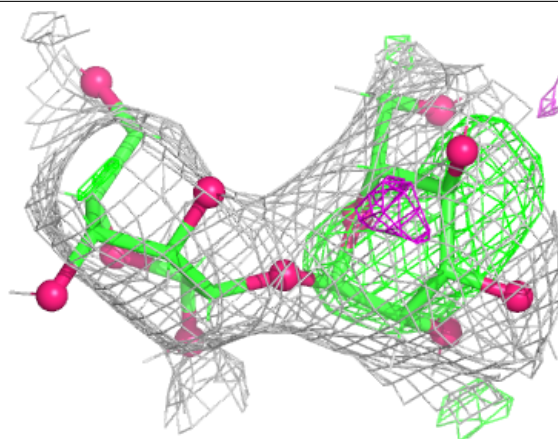
**Electron density around Chain I:**

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



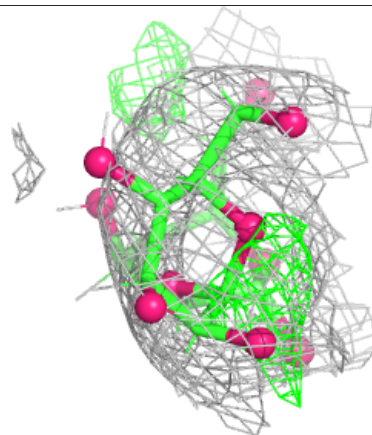
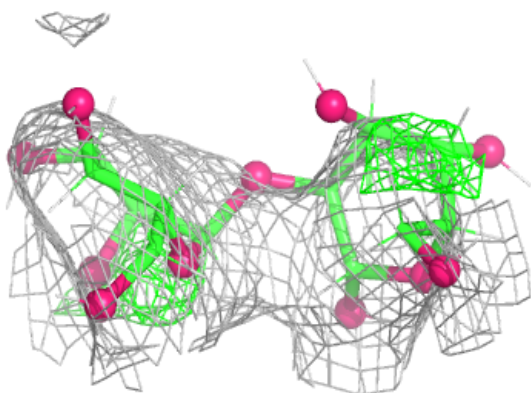
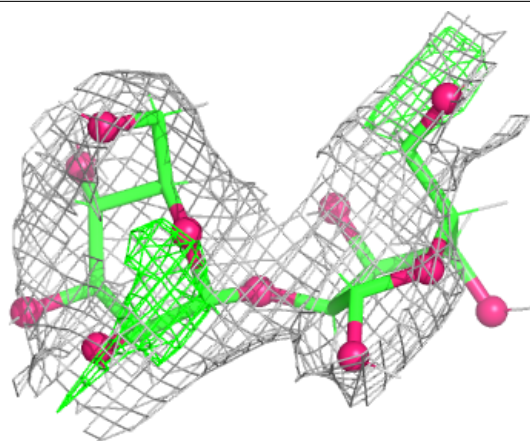
Electron density around Chain J:

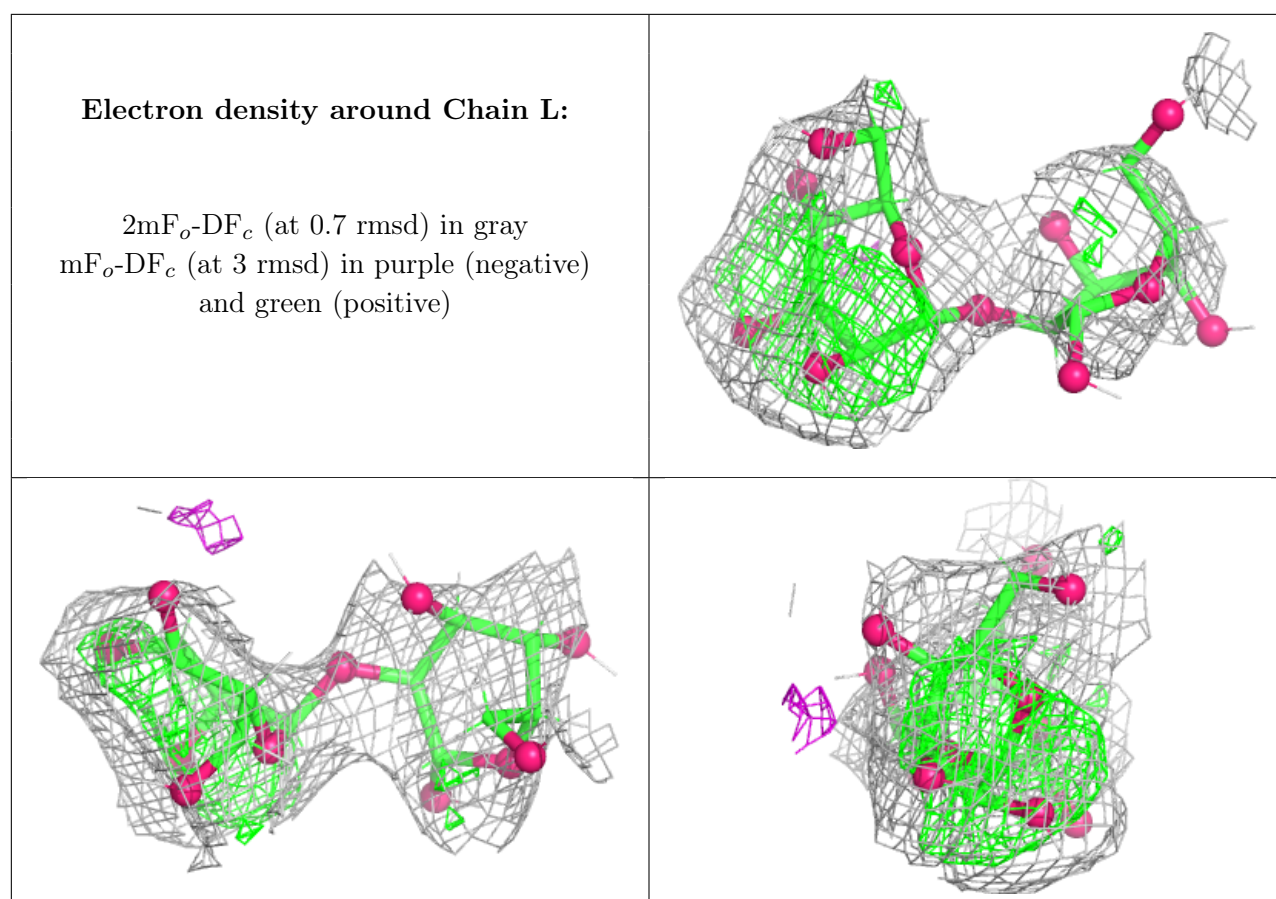
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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	GOL	C	304	6/6	0.56	0.16	67,78,82,82	0
4	GOL	B	304	6/6	0.62	0.15	61,63,70,74	0
4	GOL	A	306	6/6	0.68	0.13	51,61,71,77	0
4	GOL	A	308	6/6	0.70	0.15	48,57,60,67	0
4	GOL	A	304	6/6	0.72	0.11	62,77,79,86	0
4	GOL	A	307	6/6	0.79	0.12	44,54,56,56	0
3	CA	D	301	1/1	0.83	0.12	72,72,72,72	0
4	GOL	A	305	6/6	0.87	0.12	37,38,45,51	0
3	CA	E	301	1/1	0.89	0.09	57,57,57,57	0
3	CA	F	301	1/1	0.90	0.08	67,67,67,67	0
3	CA	B	301	1/1	0.91	0.06	45,45,45,45	0
3	CA	A	301	1/1	0.93	0.08	43,43,43,43	0
3	CA	D	302	1/1	0.94	0.06	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	F	302	1/1	0.94	0.05	78,78,78,78	0
3	CA	A	302	1/1	0.94	0.04	46,46,46,46	0
3	CA	C	302	1/1	0.96	0.06	63,63,63,63	0
3	CA	C	301	1/1	0.96	0.05	56,56,56,56	0
3	CA	E	302	1/1	0.96	0.11	63,63,63,63	0
3	CA	B	302	1/1	0.99	0.07	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.