



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 04:21 PM UTC

PDB ID : 1LES / pdb_00001les
Title : LENTIL LECTIN COMPLEXED WITH SUCROSE
Authors : Hamelryck, T.; Loris, R.
Deposited on : 1995-08-23
Resolution : 1.90 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

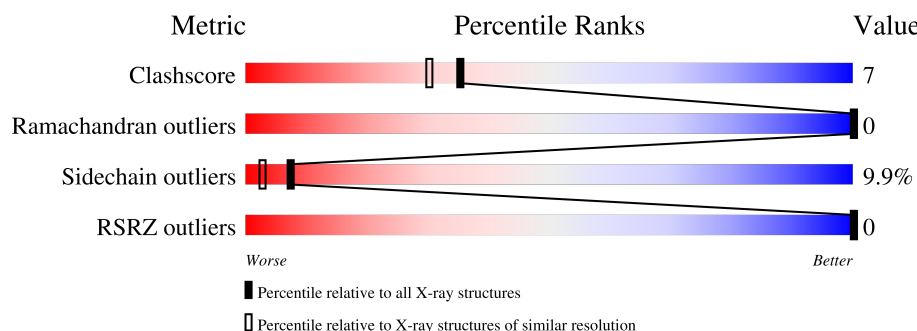
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	181	
1	C	181	
2	B	52	
2	D	52	
3	G	2	
3	H	2	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3838 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 LENTIL LECTIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	0	0	0
			1410	898	232	280			
1	C	181	Total	C	N	O	0	0	0
			1410	898	232	280			

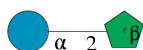
- Molecule 2 is a protein called LENTIL LECTIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	48	Total	C	N	O	0	0	1
			370	239	62	69			
2	D	48	Total	C	N	O	0	0	1
			370	239	62	69			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	35	HIS	GLN	conflict	UNP P02870
B	43	HIS	ASN	conflict	UNP P02870
D	35	HIS	GLN	conflict	UNP P02870
D	43	HIS	ASN	conflict	UNP P02870

- Molecule 3 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	G	2	Total	C	O	0	0	0
			23	12	11			
3	H	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Ca 1	0	0
4	C	1	Total 1	Ca 1	0	0

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Mn 1	0	0
5	C	1	Total 1	Mn 1	0	0

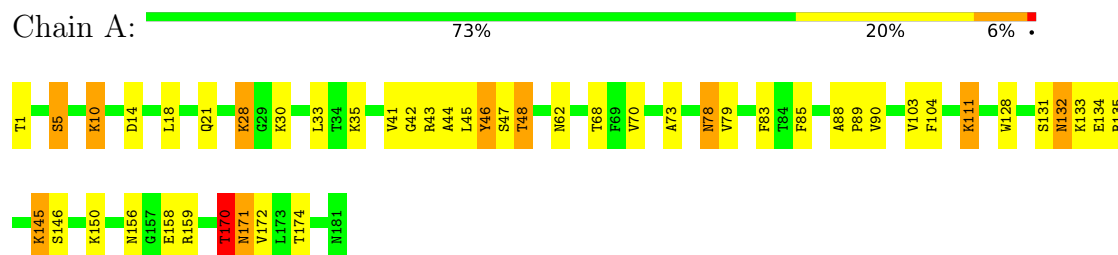
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	107	Total 107	O 107	0	0
6	C	92	Total 92	O 92	0	0
6	B	14	Total 14	O 14	0	0
6	D	15	Total 15	O 15	0	0

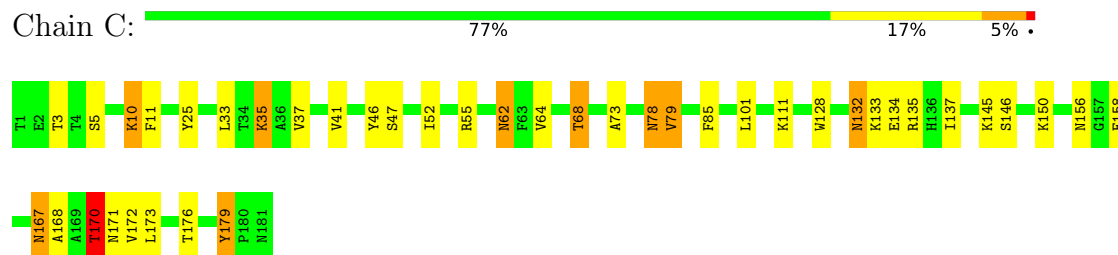
3 Residue-property plots [i](#)

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.

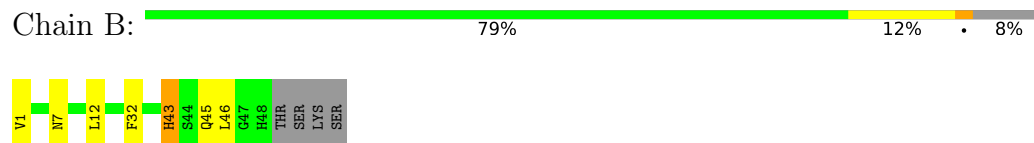
- Molecule 1: LENTIL LECTIN



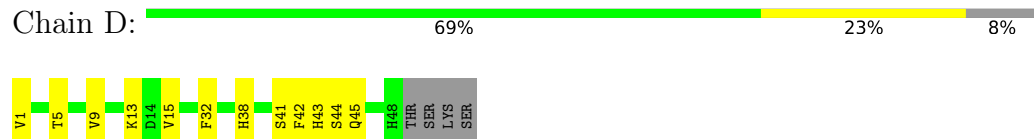
- Molecule 1: LENTIL LECTIN



- Molecule 2: LENTIL LECTIN



- Molecule 2: LENTIL LECTIN



- Molecule 3: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose



GLC1
FRU2

- Molecule 3: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain H:

100%

GLC1
FRU2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.03Å 124.80Å 50.02Å 90.00° 111.80° 90.00°	Depositor
Resolution (Å)	6.00 – 1.90 6.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	75.1 (6.00-1.90) 92.7 (6.00-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 1.90Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.188 , (Not available) 0.171 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.427 for l,-k,h	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	3838	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.63% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN, GLC, FRU, 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	1.04	2/1444 (0.1%)	1.29	12/1970 (0.6%)
1	C	1.07	2/1444 (0.1%)	1.27	13/1970 (0.7%)
2	B	1.26	2/382 (0.5%)	1.33	3/524 (0.6%)
2	D	1.11	0/382	1.20	2/524 (0.4%)
All	All	1.08	6/3652 (0.2%)	1.28	30/4988 (0.6%)

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	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	43	HIS	CB-CG	-7.21	1.40	1.50
1	A	78	ASN	CB-CG	5.72	1.66	1.52
1	A	90	VAL	CA-CB	5.64	1.61	1.54
2	B	43	HIS	CA-C	-5.52	1.45	1.52
1	C	78	ASN	CB-CG	5.36	1.65	1.52
1	C	64	VAL	CA-CB	-5.23	1.47	1.54

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	43	HIS	CB-CA-C	-9.49	93.67	109.80
1	C	170	THR	N-CA-CB	-8.22	99.97	110.90
1	A	170	THR	N-CA-CB	-7.64	100.04	110.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	32	PHE	N-CA-C	7.54	118.56	108.38
2	D	32	PHE	N-CA-C	6.90	117.69	108.38
1	C	85	PHE	N-CA-C	-6.82	98.09	109.07
1	C	179	TYR	CA-C-N	6.70	126.96	119.32
1	C	179	TYR	C-N-CA	6.70	126.96	119.32
2	D	15	VAL	N-CA-C	6.62	117.25	111.56
1	A	48	THR	N-CA-C	6.59	117.95	109.65
1	A	35	LYS	N-CA-C	-6.25	101.53	110.59
1	A	131	SER	N-CA-C	6.22	119.71	111.75
1	A	146	SER	N-CA-C	6.17	118.26	110.24
1	A	85	PHE	N-CA-C	-6.09	99.27	109.07
2	B	43	HIS	CB-CG-ND1	-5.75	114.08	122.70
1	A	47	SER	N-CA-C	5.74	118.48	111.82
1	C	47	SER	N-CA-C	5.67	118.39	111.82
1	C	146	SER	N-CA-C	5.47	117.35	110.24
1	A	30	LYS	CB-CG-CD	-5.36	98.97	111.30
1	C	135	ARG	N-CA-C	-5.32	103.00	110.50
1	C	79	VAL	CB-CA-C	-5.30	102.59	111.29
1	C	137	ILE	N-CA-C	-5.30	100.81	108.71
1	C	62	ASN	N-CA-C	-5.30	101.07	109.76
1	C	64	VAL	CB-CA-C	-5.28	101.58	110.71
1	C	46	TYR	N-CA-C	-5.27	102.48	110.28
1	A	62	ASN	N-CA-C	-5.25	101.15	109.76
1	A	135	ARG	N-CA-C	-5.12	103.28	110.50
1	C	35	LYS	N-CA-C	-5.11	103.18	110.59
1	A	83	PHE	N-CA-CB	-5.04	102.15	111.52
1	A	46	TYR	N-CA-C	-5.02	102.84	110.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	46	TYR	Sidechain

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	1410	0	1363	26	0
1	C	1410	0	1363	24	0
2	B	370	0	350	3	0
2	D	370	0	350	8	0
3	G	23	0	20	0	0
3	H	23	0	21	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	A	107	0	0	3	0
6	B	14	0	0	0	0
6	C	92	0	0	3	0
6	D	15	0	0	0	0
All	All	3838	0	3467	52	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 7.

All (52) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ALA:H	1:A:156:ASN:HD21	1.21	0.86
1:C:170:THR:HG23	1:C:172:VAL:HG23	1.59	0.85
1:C:68:THR:HG22	2:D:38:HIS:HB2	1.58	0.84
1:C:73:ALA:H	1:C:156:ASN:HD21	1.26	0.84
1:A:21:GLN:HE22	1:A:43:ARG:HH21	1.30	0.77
1:A:132:ASN:HD22	1:A:134:GLU:H	1.34	0.76
1:A:73:ALA:H	1:A:156:ASN:ND2	1.89	0.70
1:A:132:ASN:ND2	1:A:134:GLU:H	1.89	0.70
1:A:170:THR:HG23	1:A:172:VAL:HG23	1.74	0.68
1:C:170:THR:HG23	1:C:172:VAL:CG2	2.27	0.65
1:C:73:ALA:H	1:C:156:ASN:ND2	1.94	0.64
1:C:62:ASN:ND2	1:C:168:ALA:H	1.97	0.63
1:A:10:LYS:HE3	1:A:10:LYS:HA	1.82	0.61
1:C:5:SER:HB3	2:D:43:HIS:CD2	2.36	0.60
1:C:132:ASN:C	1:C:132:ASN:HD22	2.08	0.60
1:A:73:ALA:N	1:A:156:ASN:HD21	1.97	0.58
1:A:103:VAL:HG23	1:A:104:PHE:CD2	2.38	0.57
1:C:132:ASN:HD22	1:C:133:LYS:N	2.02	0.57
1:A:174:THR:HG23	6:A:779:HOH:O	2.05	0.56
1:C:128:TRP:HD1	6:C:760:HOH:O	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:THR:CG2	2:D:38:HIS:HB2	2.33	0.56
1:C:73:ALA:N	1:C:156:ASN:HD21	1.99	0.55
1:C:132:ASN:ND2	1:C:134:GLU:H	2.05	0.54
1:A:111:LYS:HD2	6:A:709:HOH:O	2.08	0.54
1:C:179:TYR:O	2:D:1:VAL:HA	2.08	0.53
1:A:5:SER:HB2	2:B:43:HIS:CE1	2.44	0.53
1:C:25:TYR:OH	1:C:35:LYS:HE3	2.10	0.52
1:A:132:ASN:HD22	1:A:132:ASN:C	2.19	0.51
2:D:42:PHE:CE2	2:D:44:SER:HB2	2.46	0.50
1:A:132:ASN:HD22	1:A:133:LYS:N	2.10	0.49
1:A:5:SER:HB2	2:B:43:HIS:ND1	2.27	0.49
1:C:62:ASN:HD22	1:C:168:ALA:H	1.59	0.48
1:A:128:TRP:HB2	1:A:145:LYS:HG2	1.96	0.48
1:C:3:THR:HG22	2:D:45:GLN:HG3	1.96	0.48
1:A:73:ALA:HB2	1:A:79:VAL:HG13	1.95	0.47
1:C:10:LYS:HD2	1:C:11:PHE:O	2.14	0.47
1:C:73:ALA:HB2	1:C:79:VAL:HG13	1.97	0.47
1:C:128:TRP:CD1	6:C:760:HOH:O	2.66	0.47
1:C:167:ASN:HD22	1:C:167:ASN:C	2.23	0.46
1:A:88:ALA:HB1	1:A:89:PRO:HD2	1.99	0.45
1:C:145:LYS:HA	6:C:760:HOH:O	2.16	0.45
1:A:171:ASN:HD22	1:A:171:ASN:HA	1.68	0.44
1:A:14:ASP:HB3	6:A:833:HOH:O	2.17	0.44
1:A:21:GLN:NE2	1:A:43:ARG:HE	2.15	0.43
1:A:1:THR:HA	2:B:46:LEU:O	2.19	0.43
1:C:176:THR:HG23	2:D:5:THR:HG22	2.01	0.42
1:A:44:ALA:C	1:A:45:LEU:HD12	2.45	0.42
1:C:170:THR:O	1:C:171:ASN:HB2	2.21	0.41
2:D:9:VAL:O	2:D:9:VAL:HG23	2.21	0.41
1:A:28:LYS:HE3	1:A:28:LYS:HB2	1.85	0.41
1:A:33:LEU:O	1:A:42:GLY:HA3	2.21	0.40
1:A:88:ALA:HB1	1:A:89:PRO:CD	2.52	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	179/181 (99%)	174 (97%)	5 (3%)	0	100	100
1	C	179/181 (99%)	174 (97%)	5 (3%)	0	100	100
2	B	46/52 (88%)	45 (98%)	1 (2%)	0	100	100
2	D	46/52 (88%)	44 (96%)	2 (4%)	0	100	100
All	All	450/466 (97%)	437 (97%)	13 (3%)	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	156/156 (100%)	139 (89%)	17 (11%)	6	2
1	C	156/156 (100%)	140 (90%)	16 (10%)	7	2
2	B	40/45 (89%)	36 (90%)	4 (10%)	7	3
2	D	40/45 (89%)	38 (95%)	2 (5%)	22	14
All	All	392/402 (98%)	353 (90%)	39 (10%)	7	3

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	10	LYS
1	A	18	LEU
1	A	28	LYS
1	A	41	VAL
1	A	48	THR
1	A	68	THR
1	A	70	VAL

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Mol	Chain	Res	Type
1	A	78	ASN
1	A	111	LYS
1	A	132	ASN
1	A	145	LYS
1	A	150	LYS
1	A	158	GLU
1	A	159	ARG
1	A	170	THR
1	A	171	ASN
1	C	10	LYS
1	C	33	LEU
1	C	37	VAL
1	C	41	VAL
1	C	52	ILE
1	C	55	ARG
1	C	68	THR
1	C	78	ASN
1	C	101	LEU
1	C	111	LYS
1	C	132	ASN
1	C	150	LYS
1	C	158	GLU
1	C	167	ASN
1	C	170	THR
1	C	173	LEU
2	B	1	VAL
2	B	7	ASN
2	B	12	LEU
2	B	45	GLN
2	D	13	LYS
2	D	41	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	59	ASN
1	A	105	ASN
1	A	132	ASN
1	A	142	ASN
1	A	155	GLN
1	A	156	ASN

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Mol	Chain	Res	Type
1	A	161	ASN
1	A	171	ASN
1	A	181	ASN
1	C	62	ASN
1	C	132	ASN
1	C	142	ASN
1	C	156	ASN
1	C	161	ASN
1	C	167	ASN
2	B	7	ASN
2	D	43	HIS

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 ⓘ

4 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	G	1	3	11,11,12	1.97	4 (36%)	15,15,17	2.12	5 (33%)
3	FRU	G	2	3	11,12,12	2.08	2 (18%)	10,18,18	2.07	3 (30%)
3	GLC	H	1	3	11,11,12	1.92	3 (27%)	15,15,17	1.62	3 (20%)
3	FRU	H	2	3	11,12,12	2.94	4 (36%)	10,18,18	1.99	3 (30%)

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	G	1	3	-	0/2/19/22	0/1/1/1
3	FRU	G	2	3	-	0/5/24/24	0/1/1/1
3	GLC	H	1	3	-	2/2/19/22	0/1/1/1
3	FRU	H	2	3	-	1/5/24/24	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	2	FRU	O2-C2	8.33	1.55	1.40
3	G	2	FRU	O2-C2	5.78	1.50	1.40
3	G	1	GLC	C2-C3	-3.82	1.46	1.52
3	H	1	GLC	C1-C2	3.67	1.61	1.52
3	H	2	FRU	C4-C5	-3.28	1.44	1.53
3	H	1	GLC	C4-C5	3.21	1.59	1.53
3	G	2	FRU	C4-C5	-2.85	1.45	1.53
3	H	1	GLC	O5-C5	2.79	1.48	1.43
3	G	1	GLC	O3-C3	2.72	1.49	1.43
3	G	1	GLC	C4-C5	2.53	1.58	1.53
3	H	2	FRU	O5-C5	-2.31	1.38	1.43
3	H	2	FRU	C1-C2	-2.21	1.47	1.52
3	G	1	GLC	O2-C2	-2.04	1.39	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1	GLC	C1-O5-C5	5.05	118.95	112.19
3	G	2	FRU	C6-C5-C4	-4.53	104.41	115.10
3	H	2	FRU	O1-C1-C2	3.68	119.83	111.67
3	H	2	FRU	O6-C6-C5	-3.47	99.51	111.33
3	H	1	GLC	C6-C5-C4	-3.30	104.92	113.02
3	G	1	GLC	O5-C5-C6	3.24	113.97	107.66
3	G	1	GLC	O5-C1-C2	-3.00	103.64	110.79
3	G	2	FRU	O5-C5-C6	2.95	116.94	108.89
3	G	1	GLC	O5-C5-C4	-2.94	103.68	110.83
3	H	1	GLC	O5-C1-C2	-2.51	104.80	110.79
3	G	1	GLC	O4-C4-C3	-2.47	104.56	110.38
3	G	2	FRU	O3-C3-C4	-2.37	104.89	113.25
3	H	2	FRU	O3-C3-C4	-2.31	105.10	113.25
3	H	1	GLC	O2-C2-C3	-2.05	105.91	110.15

There are no chirality outliers.

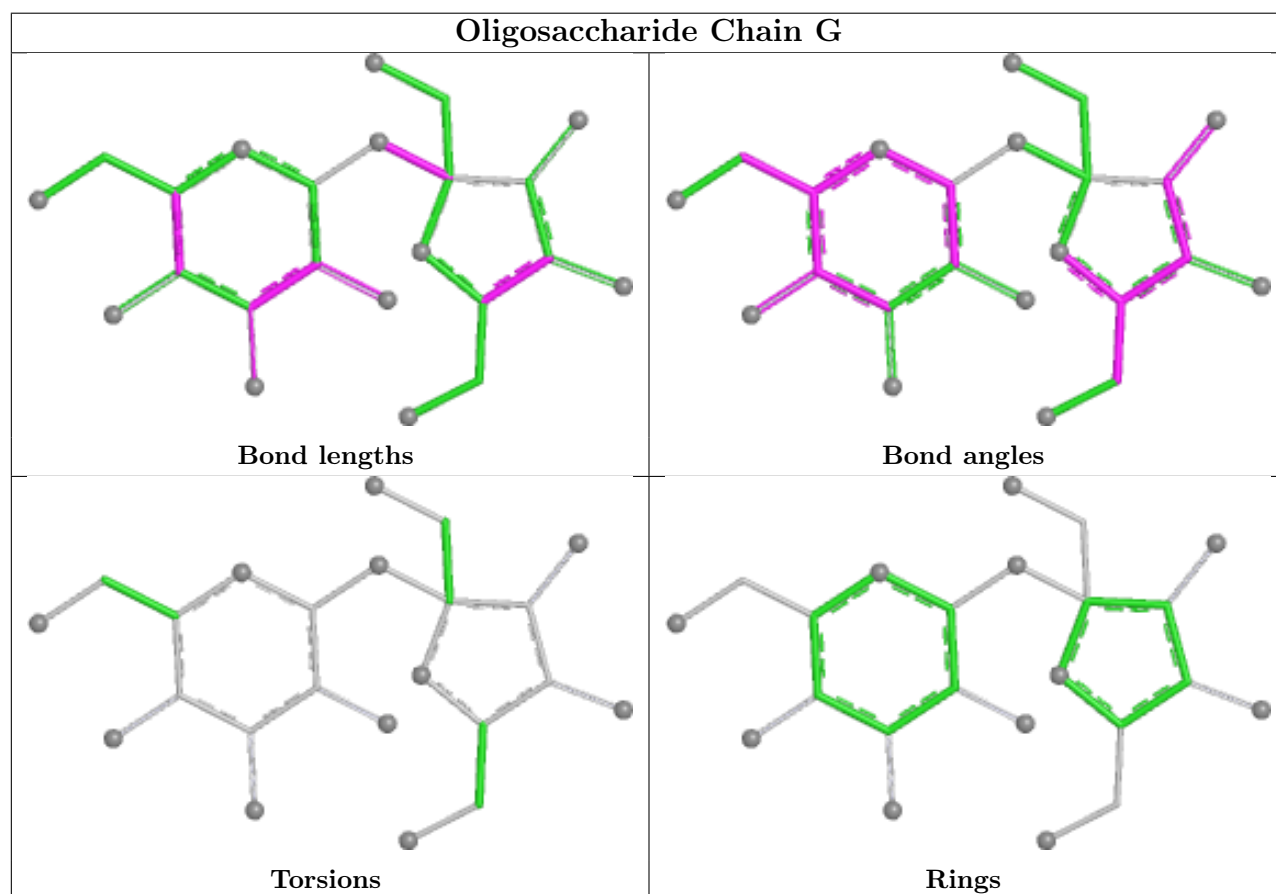
All (3) torsion outliers are listed below:

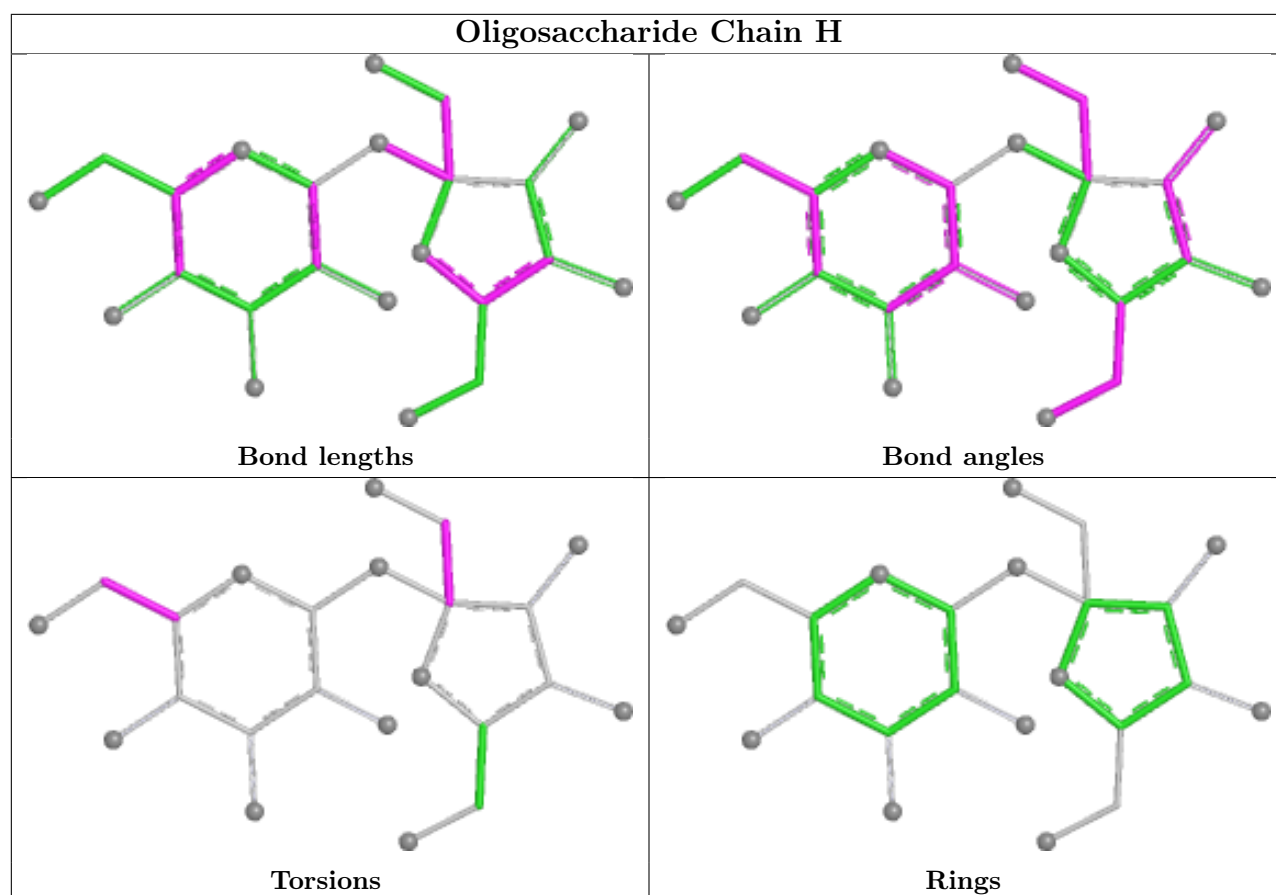
Mol	Chain	Res	Type	Atoms
3	H	1	GLC	O5-C5-C6-O6
3	H	1	GLC	C4-C5-C6-O6
3	H	2	FRU	O1-C1-C2-C3

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 [i](#)

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	181/181 (100%)	-1.42	0 100 100	13, 23, 41, 64	0
1	C	181/181 (100%)	-1.45	0 100 100	13, 22, 41, 68	0
2	B	48/52 (92%)	-1.40	0 100 100	13, 23, 39, 54	0
2	D	48/52 (92%)	-1.41	0 100 100	13, 23, 42, 52	0
All	All	458/466 (98%)	-1.43	0 100 100	13, 23, 42, 68	0

There are no RSRZ outliers to report.

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](#)

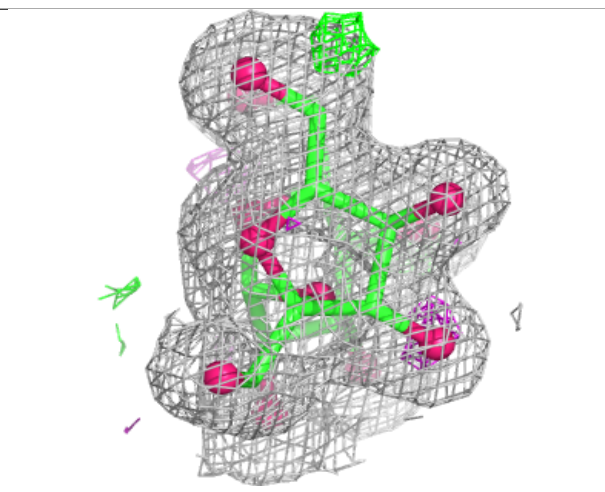
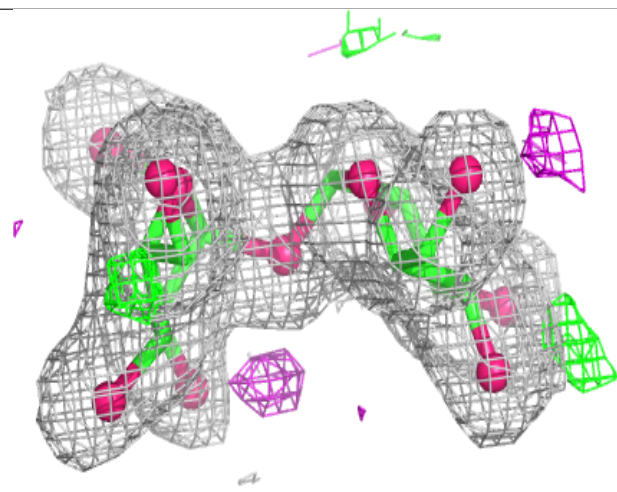
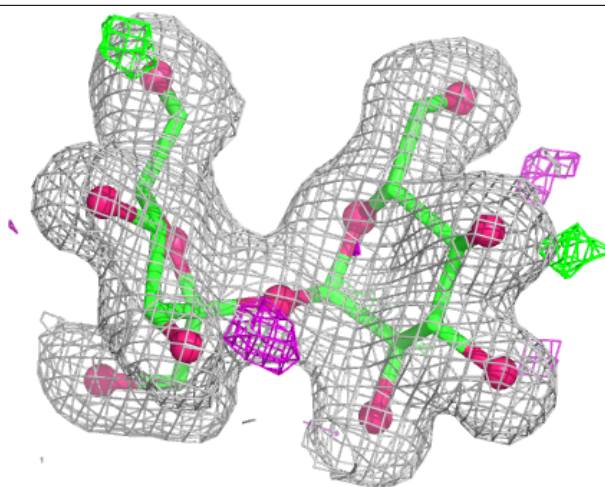
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

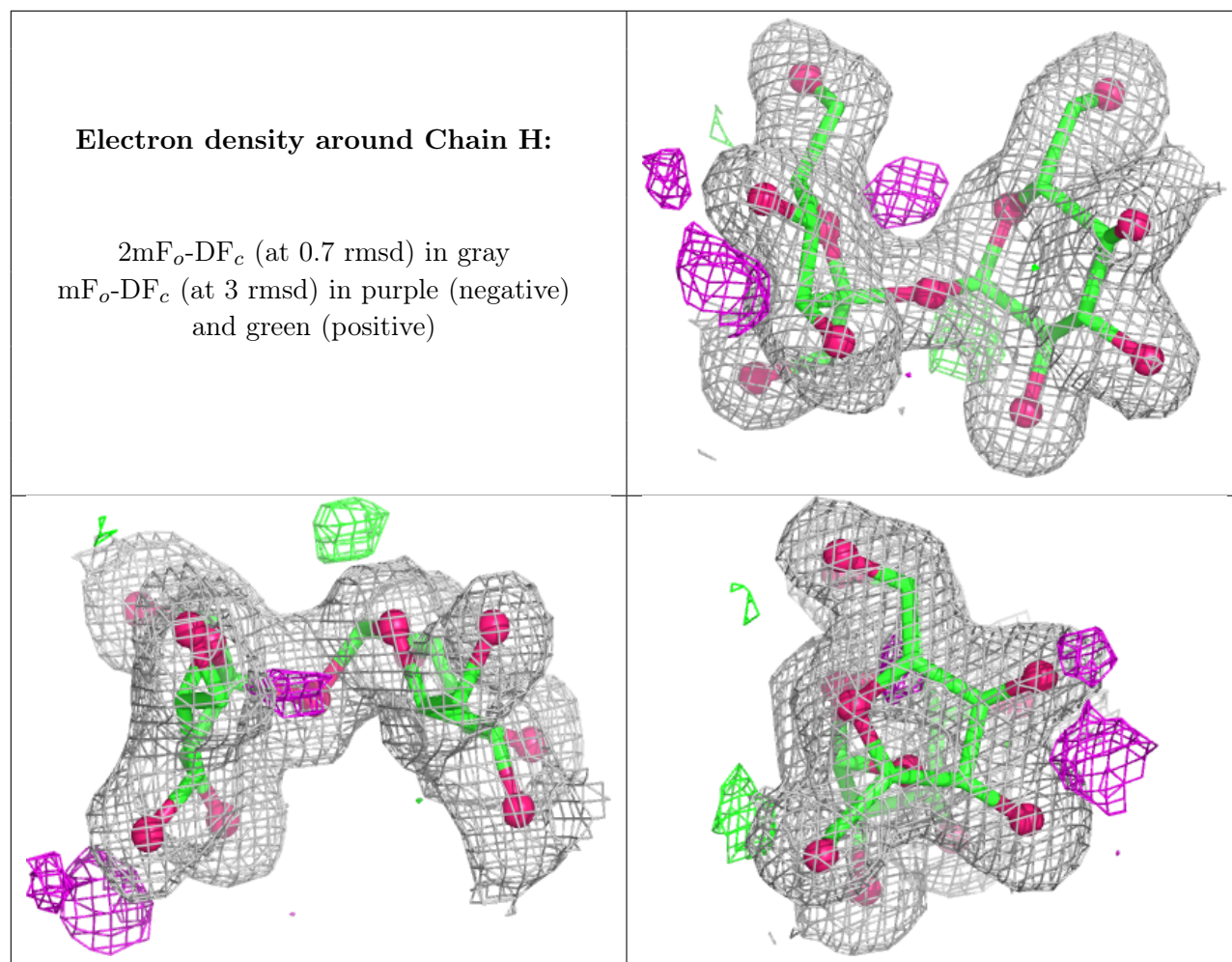
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GLC	G	1	11/12	-	-	12,14,17,17	0
3	FRU	G	2	12/12	-	-	18,22,26,28	0
3	GLC	H	1	11/12	-	-	14,15,16,18	0
3	FRU	H	2	12/12	-	-	19,21,26,28	0

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 G:

$2mF_o-DF_c$ (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($\text{\AA}^2$)	Q<0.9
4	CA	A	203	1/1	1.00	0.01	14,14,14,14	0
4	CA	C	201	1/1	1.00	0.01	15,15,15,15	0
5	MN	A	204	1/1	1.00	0.02	18,18,18,18	0
5	MN	C	202	1/1	1.00	0.01	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.