



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:22 PM UTC

PDB ID : 3LIH / pdb_00003lih
Title : Crystal structure of fructosyltransferase (D191A) from *A. japonicus* in complex with raffinose
Authors : Chuankhayan, P.; Chen, C.J.; Chiang, C.M.
Deposited on : 2010-01-24
Resolution : 2.20 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

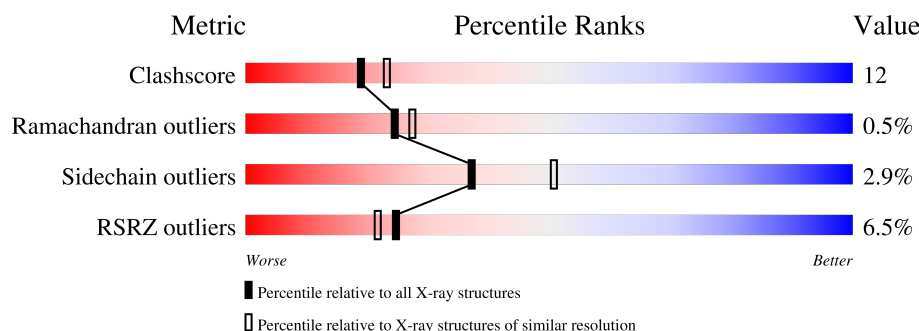
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	634	
2	B	3	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5068 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 Fructosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	634	Total	C	N	O	S	0	0	0
			4883	3091	824	965	3			

- Molecule 2 is an oligosaccharide called alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranoside-(1-2)-beta-D-fructofuranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	3	Total	C	O	0	0	0
			34	18	16			

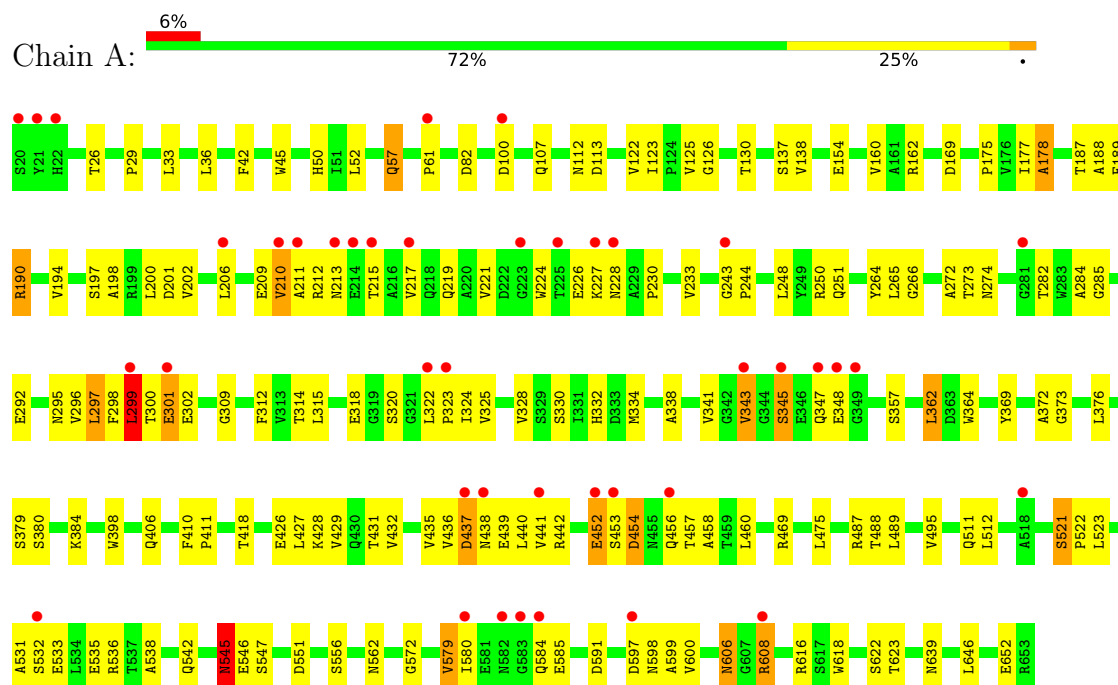
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	151	Total	O	0	0
			151	151		

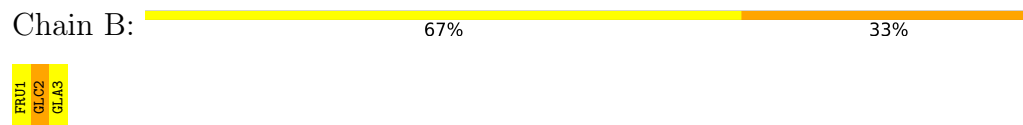
3 Residue-property plots [i](#)

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: Fructosyltransferase



• Molecule 2: alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranose-(1-2)-beta-D-fructofuranos e



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	99.27Å 109.94Å 65.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 30.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	94.5 (30.00-2.20) 94.5 (30.00-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.50 (at 2.19Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.222 , 0.268 0.233 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	38.0	Xtriage
Anisotropy	0.605	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 26.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5068	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: FRU, GLC, GLA

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	0.59	16/5014 (0.3%)	1.07	37/6858 (0.5%)

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	2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	244	PRO	C-N	-10.00	1.20	1.33
1	A	126	GLY	C-N	-9.75	1.22	1.33
1	A	282	THR	C-N	-8.07	1.22	1.33
1	A	125	VAL	C-N	7.66	1.44	1.33
1	A	112	ASN	C-N	7.35	1.43	1.33
1	A	113	ASP	C-N	6.93	1.42	1.34
1	A	639	ASN	C-N	-6.73	1.22	1.33
1	A	545	ASN	C-N	-6.06	1.25	1.33
1	A	300	THR	C-N	6.00	1.43	1.33
1	A	188	ALA	C-N	-5.93	1.23	1.34
1	A	187	THR	C-N	-5.52	1.26	1.33
1	A	189	PHE	C-N	-5.46	1.26	1.33
1	A	274	ASN	C-N	5.34	1.42	1.33
1	A	210	VAL	C-N	-5.25	1.26	1.33
1	A	343	VAL	C-N	-5.21	1.25	1.33
1	A	209	GLU	C-N	5.20	1.40	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	GLY	CA-C-N	8.53	128.47	119.85
1	A	243	GLY	C-N-CA	8.53	128.47	119.85
1	A	343	VAL	O-C-N	8.39	130.03	122.65
1	A	301	GLU	CA-C-N	-7.83	110.09	122.08
1	A	301	GLU	C-N-CA	-7.83	110.09	122.08
1	A	113	ASP	O-C-N	7.55	128.19	121.17
1	A	227	LYS	N-CA-C	7.25	119.26	111.36
1	A	190	ARG	N-CA-C	7.12	118.00	108.38
1	A	292	GLU	CA-C-N	6.84	134.61	121.54
1	A	292	GLU	C-N-CA	6.84	134.61	121.54
1	A	26	THR	N-CA-C	-6.75	104.95	113.72
1	A	357	SER	N-CA-C	-6.74	105.59	113.88
1	A	533	GLU	N-CA-C	-6.73	104.55	112.89
1	A	556	SER	N-CA-C	6.68	119.30	108.41
1	A	122	VAL	N-CA-C	6.61	118.00	108.48
1	A	376	LEU	CA-C-N	6.46	126.48	119.89
1	A	376	LEU	C-N-CA	6.46	126.48	119.89
1	A	299	LEU	N-CA-C	6.41	119.94	110.46
1	A	362	LEU	N-CA-C	-6.16	104.47	112.23
1	A	42	PHE	N-CA-C	6.14	118.77	111.33
1	A	177	ILE	N-CA-C	-6.14	97.15	106.42
1	A	123	ILE	N-CA-C	-5.97	101.09	107.77
1	A	454	ASP	N-CA-C	-5.92	104.72	112.41
1	A	531	ALA	N-CA-C	5.82	118.73	109.24
1	A	469	ARG	N-CA-C	5.71	118.24	111.33
1	A	299	LEU	CA-C-N	5.68	131.14	122.11
1	A	299	LEU	C-N-CA	5.68	131.14	122.11
1	A	112	ASN	O-C-N	5.66	130.27	122.46
1	A	639	ASN	O-C-N	-5.36	116.14	122.58
1	A	639	ASN	CA-C-N	5.36	129.28	122.37
1	A	639	ASN	C-N-CA	5.36	129.28	122.37
1	A	178	ALA	N-CA-C	5.31	119.63	113.20
1	A	138	VAL	N-CA-C	5.12	115.34	108.17
1	A	521	SER	N-CA-C	5.08	116.41	110.31
1	A	495	VAL	N-CA-C	5.07	113.51	107.73
1	A	579	VAL	N-CA-C	5.05	114.69	108.53
1	A	187	THR	N-CA-C	-5.00	107.02	113.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	ASP	Mainchain

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Mol	Chain	Res	Type	Group
1	A	545	ASN	Mainchain

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	4883	0	4610	113	0
2	B	34	0	29	1	0
3	A	151	0	0	1	0
All	All	5068	0	4639	113	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 12.

All (113) 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:299:LEU:O	1:A:309:GLY:HA2	1.56	1.02
1:A:454:ASP:HB2	1:A:456:GLN:NE2	1.94	0.82
1:A:454:ASP:HB2	1:A:456:GLN:HE21	1.42	0.82
1:A:299:LEU:O	1:A:309:GLY:CA	2.28	0.81
1:A:33:LEU:H	1:A:562:ASN:HD21	1.26	0.80
1:A:438:ASN:HD21	1:A:440:LEU:HG	1.46	0.79
1:A:453:SER:HB2	1:A:458:ALA:HB2	1.69	0.75
1:A:435:VAL:HA	1:A:579:VAL:HG12	1.69	0.74
1:A:162:ARG:HG3	1:A:169:ASP:OD2	1.88	0.73
1:A:591:ASP:H	1:A:606:ASN:HD21	1.33	0.72
1:A:438:ASN:HD21	1:A:440:LEU:CG	2.03	0.72
1:A:299:LEU:HG	1:A:428:LYS:HA	1.70	0.72
1:A:606:ASN:HD22	1:A:606:ASN:N	1.89	0.71
1:A:453:SER:CB	1:A:458:ALA:HB2	2.21	0.70
1:A:61:PRO:O	1:A:372:ALA:HB2	1.92	0.70
1:A:295:ASN:ND2	1:A:373:GLY:H	1.92	0.68
1:A:453:SER:HB2	1:A:458:ALA:CB	2.23	0.68
1:A:61:PRO:O	1:A:372:ALA:CB	2.42	0.67
1:A:456:GLN:NE2	1:A:457:THR:H	1.91	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:ASP:H	1:A:606:ASN:ND2	1.93	0.67
1:A:597:ASP:O	1:A:600:VAL:N	2.29	0.66
1:A:50:HIS:ND1	1:A:597:ASP:OD2	2.29	0.65
1:A:206:LEU:HB2	1:A:211:ALA:HB2	1.80	0.63
1:A:523:LEU:C	1:A:523:LEU:HD23	2.24	0.63
1:A:379:SER:O	1:A:384:LYS:HE2	2.01	0.61
1:A:212:ARG:NE	1:A:343:VAL:HG21	2.15	0.61
1:A:33:LEU:H	1:A:562:ASN:ND2	1.99	0.60
1:A:542:GLN:O	1:A:546:GLU:HA	2.04	0.58
1:A:295:ASN:HD21	1:A:373:GLY:H	1.51	0.57
1:A:438:ASN:O	1:A:441:VAL:N	2.26	0.57
1:A:322:LEU:HA	1:A:323:PRO:C	2.30	0.57
1:A:606:ASN:HD22	1:A:606:ASN:H	1.52	0.56
1:A:295:ASN:HD21	1:A:372:ALA:HA	1.70	0.56
1:A:426:GLU:C	1:A:427:LEU:HD12	2.31	0.55
1:A:536:ARG:HH11	1:A:536:ARG:HG3	1.71	0.55
1:A:197:SER:HB2	1:A:298:PHE:CZ	2.42	0.55
1:A:217:VAL:O	1:A:221:VAL:HG23	2.07	0.55
1:A:438:ASN:ND2	1:A:440:LEU:HG	2.18	0.54
1:A:57:GLN:HB2	1:A:418:THR:HB	1.90	0.54
1:A:295:ASN:HB2	1:A:314:THR:OG1	2.07	0.54
1:A:201:ASP:OD2	1:A:230:PRO:HD2	2.09	0.53
1:A:606:ASN:C	1:A:608:ARG:H	2.16	0.53
1:A:248:LEU:HG	1:A:265:LEU:HD12	1.90	0.53
1:A:45:TRP:HB3	1:A:599:ALA:CB	2.39	0.52
1:A:202:VAL:HG11	1:A:217:VAL:HG13	1.91	0.51
1:A:284:ALA:O	1:A:285:GLY:C	2.54	0.51
1:A:315:LEU:N	1:A:315:LEU:HD23	2.26	0.50
1:A:325:VAL:HB	1:A:328:VAL:HB	1.93	0.50
1:A:301:GLU:HB2	1:A:302:GLU:OE2	2.12	0.50
1:A:82:ASP:OD2	1:A:107:GLN:NE2	2.46	0.49
1:A:130:THR:OG1	1:A:160:VAL:HG13	2.12	0.49
1:A:45:TRP:CE2	1:A:616:ARG:HB3	2.47	0.49
1:A:299:LEU:HD22	1:A:312:PHE:CD2	2.48	0.49
1:A:606:ASN:N	1:A:606:ASN:ND2	2.60	0.49
1:A:579:VAL:O	1:A:585:GLU:HA	2.14	0.48
1:A:200:LEU:HD21	1:A:341:VAL:HG11	1.95	0.48
1:A:436:VAL:O	1:A:438:ASN:N	2.46	0.48
1:A:406:GLN:HG3	3:A:729:HOH:O	2.12	0.48
1:A:410:PHE:CG	1:A:411:PRO:HD2	2.49	0.48
1:A:535:GLU:OE1	1:A:622:SER:OG	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:LEU:CD2	1:A:341:VAL:HG11	2.44	0.47
1:A:215:THR:O	1:A:219:GLN:HG3	2.15	0.47
1:A:297:LEU:HD12	1:A:297:LEU:N	2.29	0.47
1:A:438:ASN:O	1:A:439:GLU:C	2.56	0.47
1:A:437:ASP:OD2	1:A:442:ARG:NH2	2.36	0.47
1:A:580:ILE:HA	1:A:584:GLN:O	2.15	0.47
1:A:334:MET:HG3	1:A:362:LEU:HB3	1.96	0.47
1:A:452:GLU:O	1:A:458:ALA:HA	2.15	0.46
1:A:299:LEU:HB3	1:A:429:VAL:HG23	1.96	0.46
1:A:454:ASP:HB2	1:A:456:GLN:HG3	1.97	0.46
1:A:312:PHE:CE1	1:A:338:ALA:HB2	2.49	0.46
1:A:320:SER:HA	1:A:330:SER:OG	2.15	0.46
1:A:542:GLN:O	1:A:546:GLU:CA	2.63	0.46
1:A:29:PRO:HB3	1:A:618:TRP:CE2	2.50	0.46
1:A:511:GLN:C	1:A:512:LEU:HD12	2.41	0.45
1:A:332:HIS:HB3	1:A:369:TYR:CD2	2.52	0.45
1:A:213:ASN:HD22	1:A:215:THR:HB	1.82	0.45
1:A:82:ASP:O	1:A:107:GLN:HG2	2.17	0.45
1:A:318:GLU:OE2	2:B:2:GLC:O3	2.33	0.45
1:A:606:ASN:C	1:A:608:ARG:N	2.71	0.45
1:A:250:ARG:O	1:A:251:GLN:C	2.60	0.45
1:A:154:GLU:OE2	1:A:190:ARG:HD3	2.17	0.44
1:A:213:ASN:ND2	1:A:215:THR:HB	2.32	0.44
1:A:536:ARG:HG3	1:A:536:ARG:NH1	2.32	0.44
1:A:296:VAL:C	1:A:297:LEU:HD12	2.42	0.44
1:A:532:SER:O	1:A:535:GLU:O	2.36	0.44
1:A:538:ALA:HB3	1:A:551:ASP:HB3	2.00	0.44
1:A:438:ASN:OD1	1:A:440:LEU:HB2	2.18	0.43
1:A:224:TRP:CH2	1:A:226:GLU:HG2	2.54	0.43
1:A:488:THR:O	1:A:489:LEU:HD23	2.18	0.43
1:A:297:LEU:N	1:A:297:LEU:CD1	2.82	0.43
1:A:456:GLN:CD	1:A:457:THR:H	2.27	0.43
1:A:437:ASP:OD1	1:A:437:ASP:C	2.62	0.42
1:A:264:TYR:CZ	1:A:266:GLY:HA2	2.55	0.42
1:A:198:ALA:HA	1:A:224:TRP:NE1	2.35	0.42
1:A:322:LEU:HD23	1:A:324:ILE:HG13	2.01	0.42
1:A:431:THR:HA	1:A:460:LEU:O	2.20	0.42
1:A:475:LEU:HD22	1:A:646:LEU:HD22	2.01	0.42
1:A:194:VAL:HG12	1:A:233:VAL:HG22	2.01	0.42
1:A:364:TRP:CE3	1:A:572:GLY:HA2	2.55	0.41
1:A:622:SER:C	1:A:623:THR:HG23	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:VAL:HG12	1:A:210:VAL:O	2.21	0.41
1:A:523:LEU:C	1:A:523:LEU:CD2	2.94	0.41
1:A:175:PRO:HG2	1:A:178:ALA:HB2	2.02	0.41
1:A:453:SER:HB2	1:A:458:ALA:HB1	2.03	0.41
1:A:61:PRO:HD3	1:A:398:TRP:HB2	2.02	0.41
1:A:312:PHE:CD1	1:A:338:ALA:HB2	2.56	0.41
1:A:432:VAL:HB	1:A:460:LEU:HB2	2.02	0.41
1:A:521:SER:HA	1:A:522:PRO:HD3	1.83	0.41
1:A:272:ALA:O	1:A:273:THR:C	2.64	0.40
1:A:380:SER:O	1:A:384:LYS:HG3	2.21	0.40
1:A:454:ASP:C	1:A:456:GLN:H	2.28	0.40
1:A:545:ASN:OD1	1:A:547:SER:OG	2.38	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	632/634 (100%)	590 (93%)	39 (6%)	3 (0%)	24	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	598	ASN
1	A	345	SER
1	A	437	ASP

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	516/516 (100%)	501 (97%)	15 (3%)	37	51

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LEU
1	A	52	LEU
1	A	57	GLN
1	A	137	SER
1	A	228	ASN
1	A	297	LEU
1	A	299	LEU
1	A	345	SER
1	A	347	GLN
1	A	348	GLU
1	A	452	GLU
1	A	487	ARG
1	A	606	ASN
1	A	608	ARG
1	A	652	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	156	GLN
1	A	213	ASN
1	A	295	ASN
1	A	327	GLN
1	A	434	ASN
1	A	438	ASN
1	A	456	GLN
1	A	511	GLN
1	A	524	GLN
1	A	542	GLN
1	A	554	GLN
1	A	562	ASN
1	A	586	GLN

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Mol	Chain	Res	Type
1	A	606	ASN
1	A	639	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 ⓘ

3 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	FRU	B	1	2	11,12,12	1.71	2 (18%)	10,18,18	0.92	0
2	GLC	B	2	2	11,11,12	1.26	1 (9%)	15,15,17	1.14	2 (13%)
2	GLA	B	3	2	11,11,12	1.58	2 (18%)	15,15,17	1.15	2 (13%)

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	FRU	B	1	2	-	3/5/24/24	0/1/1/1
2	GLC	B	2	2	-	2/2/19/22	0/1/1/1
2	GLA	B	3	2	-	0/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	GLA	O5-C1	4.20	1.50	1.43
2	B	1	FRU	O2-C2	3.37	1.46	1.40
2	B	1	FRU	O6-C6	-3.31	1.28	1.42
2	B	2	GLC	O5-C1	3.18	1.49	1.43
2	B	3	GLA	O5-C5	2.69	1.48	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	GLC	C6-C5-C4	-2.55	106.77	113.02
2	B	3	GLA	C1-O5-C5	2.22	115.16	112.19
2	B	3	GLA	C1-C2-C3	2.22	112.87	109.64
2	B	2	GLC	O5-C5-C6	2.04	111.64	107.66

There are no chirality outliers.

All (5) torsion outliers are listed below:

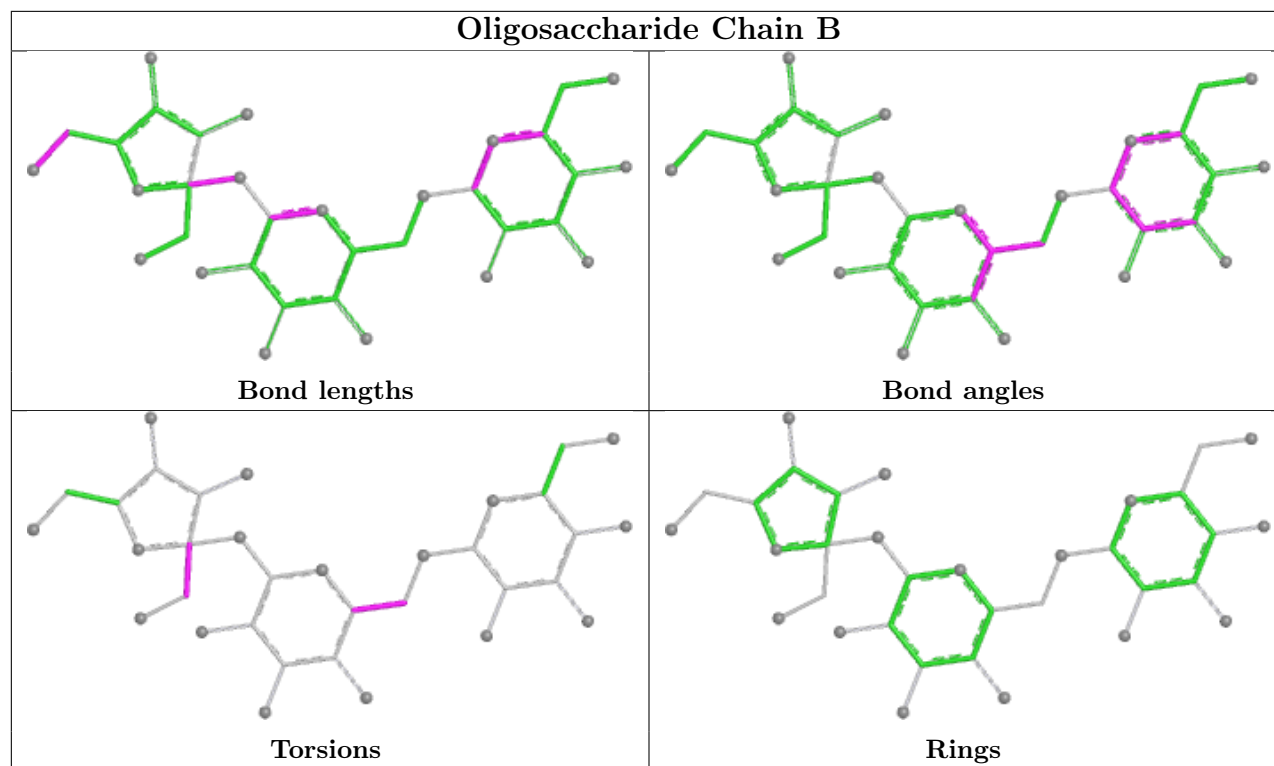
Mol	Chain	Res	Type	Atoms
2	B	1	FRU	O1-C1-C2-C3
2	B	1	FRU	O1-C1-C2-O2
2	B	1	FRU	O1-C1-C2-O5
2	B	2	GLC	O5-C5-C6-O6
2	B	2	GLC	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	GLC	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](#)

There are no ligands in this entry.

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	634/634 (100%)	0.33	41 (6%)	25 22	28, 41, 80, 114	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	210	VAL	4.3
1	A	21	TYR	4.0
1	A	438	ASN	3.5
1	A	345	SER	3.3
1	A	532	SER	3.3
1	A	343	VAL	3.1
1	A	349	GLY	3.0
1	A	225	THR	2.9
1	A	243	GLY	2.8
1	A	582	ASN	2.8
1	A	347	GLN	2.8
1	A	206	LEU	2.7
1	A	211	ALA	2.7
1	A	322	LEU	2.6
1	A	597	ASP	2.6
1	A	453	SER	2.5
1	A	323	PRO	2.5
1	A	608	ARG	2.5
1	A	213	ASN	2.5
1	A	215	THR	2.5
1	A	583	GLY	2.4
1	A	348	GLU	2.4
1	A	20	SER	2.3
1	A	214	GLU	2.3
1	A	452	GLU	2.3
1	A	580	ILE	2.3
1	A	223	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	227	LYS	2.2
1	A	281	GLY	2.2
1	A	441	VAL	2.2
1	A	437	ASP	2.2
1	A	456	GLN	2.2
1	A	22	HIS	2.1
1	A	299	LEU	2.1
1	A	100	ASP	2.1
1	A	228	ASN	2.1
1	A	584	GLN	2.1
1	A	301	GLU	2.0
1	A	61	PRO	2.0
1	A	217	VAL	2.0
1	A	518	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

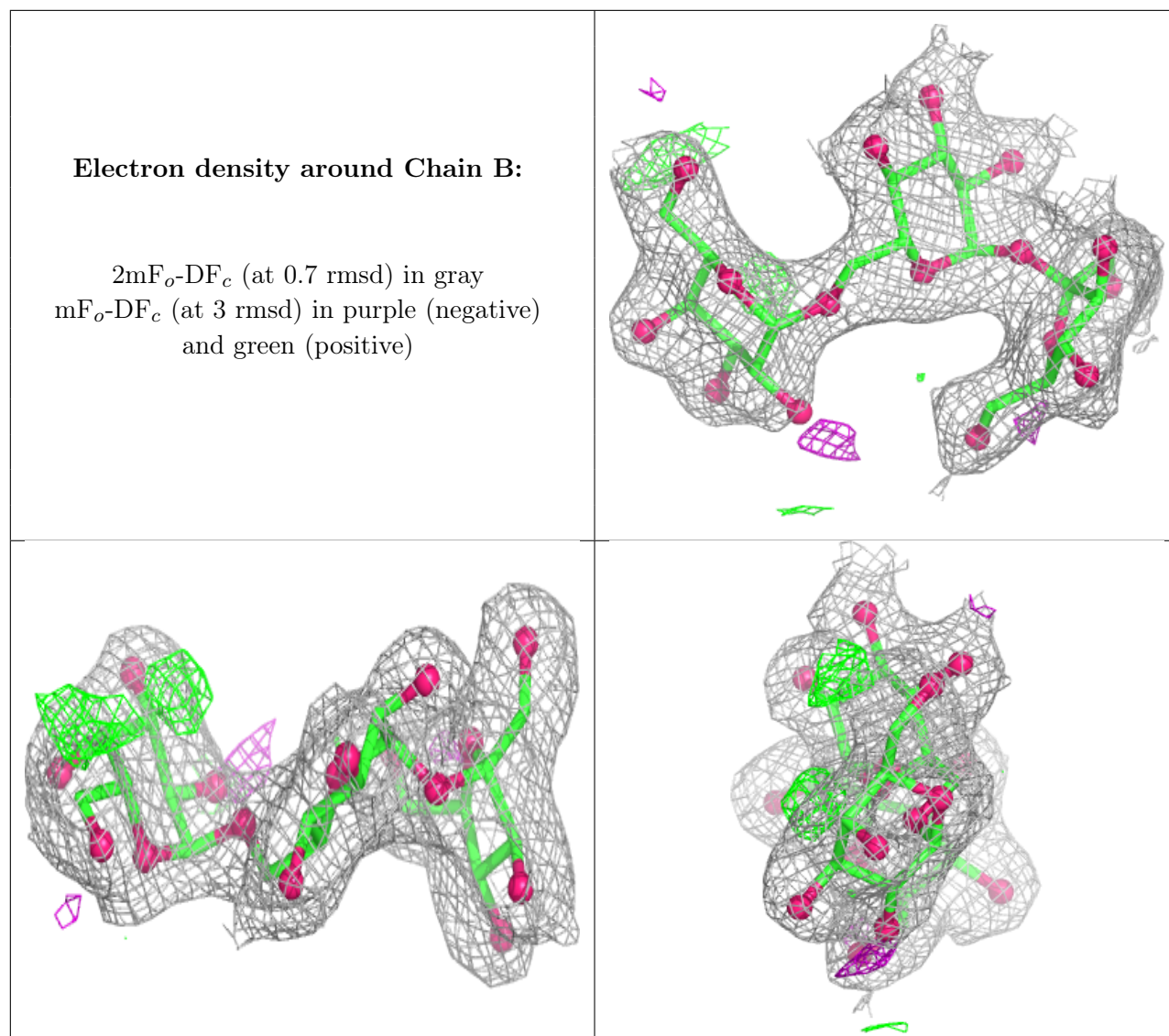
There are no non-standard protein/DNA/RNA residues in this entry.

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($\text{\AA}^2$)	Q<0.9
2	GLA	B	3	11/12	0.81	0.12	41,43,45,47	0
2	GLC	B	2	11/12	0.94	0.07	26,31,36,38	0
2	FRU	B	1	12/12	0.94	0.07	29,32,34,35	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.



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.