



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:25 PM UTC

PDB ID : 1VBR / pdb_00001vbr
Title : Crystal structure of complex xylanase 10B from *Thermotoga maritima* with xylobiose
Authors : Ihsanawati; Kumasaka, T.; Kaneko, T.; Nakamura, S.; Tanaka, N.
Deposited on : 2004-03-02
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

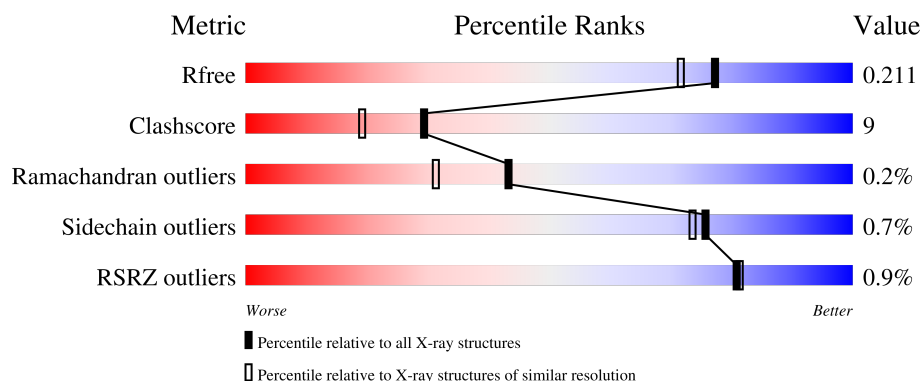
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	328	80% 17% ..
1	B	328	79% 18% ..
2	C	2	100%
2	D	2	50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6114 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 endo-1,4-beta-xylanase B.

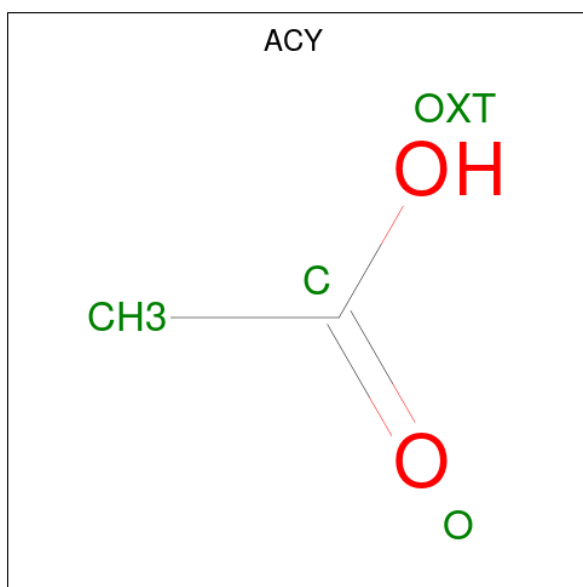
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2703	1751	447	496	9			
1	B	324	Total	C	N	O	S	0	0	0
			2703	1751	447	496	9			

- Molecule 2 is an oligosaccharide called alpha-D-xylopyranose-(1-4)-beta-D-xylopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			19	10	9			
2	D	2	Total	C	O	0	0	0
			19	10	9			

- Molecule 3 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



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

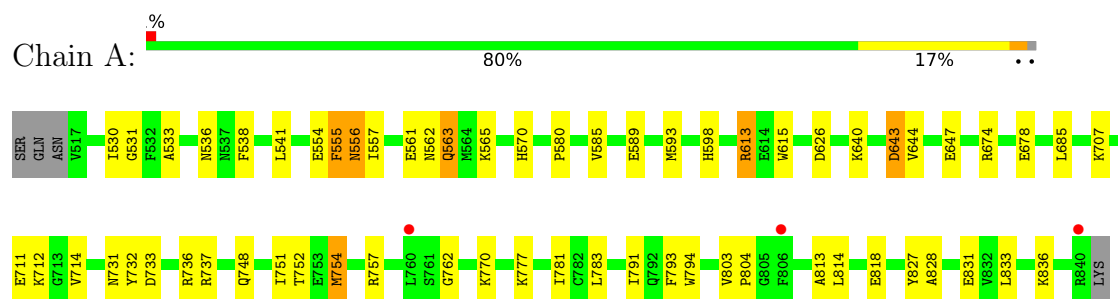
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	333	Total	O	0	0
			333	333		
4	B	313	Total	O	0	0
			313	313		

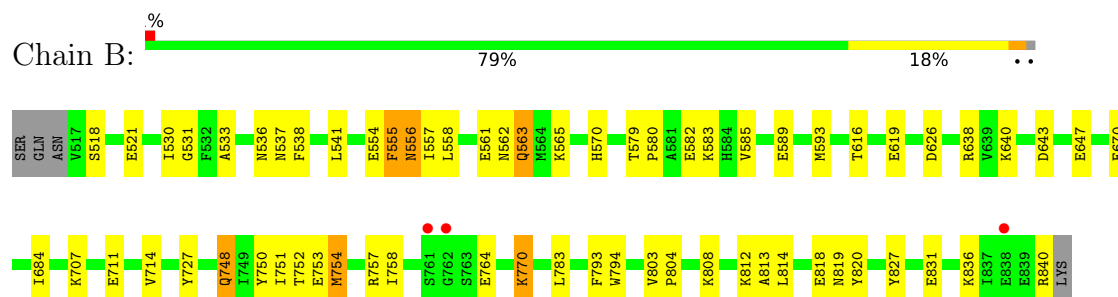
3 Residue-property plots [i](#)

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.

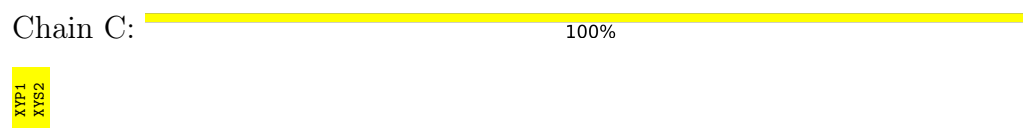
- Molecule 1: endo-1,4-beta-xylanase B



- Molecule 1: endo-1,4-beta-xylanase B



- Molecule 2: alpha-D-xylopyranose-(1-4)-beta-D-xylopyranose



- Molecule 2: alpha-D-xylopyranose-(1-4)-beta-D-xylopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.46Å 58.43Å 61.29Å 70.43° 84.33° 68.77°	Depositor
Resolution (Å)	29.58 – 1.80 29.58 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.58-1.80) 97.4 (29.58-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 1.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.181 , 0.215 0.175 , 0.211	Depositor DCC
R_{free} test set	6490 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6114	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.13% 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: ACY, XYS, XYP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.35	0/2778	0.87	13/3760 (0.3%)
1	B	0.36	0/2778	0.89	11/3760 (0.3%)
All	All	0.35	0/5556	0.88	24/7520 (0.3%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	752	THR	N-CA-C	10.12	125.75	113.41
1	A	752	THR	N-CA-C	9.86	125.44	113.41
1	B	751	ILE	N-CA-C	-7.45	96.41	107.51
1	B	570	HIS	CA-C-N	6.58	126.52	119.28
1	B	570	HIS	C-N-CA	6.58	126.52	119.28
1	A	754	MET	N-CA-C	6.56	119.88	110.10
1	A	751	ILE	N-CA-C	-6.56	97.74	107.51
1	B	554	GLU	N-CA-C	6.52	121.37	113.41
1	B	563	GLN	N-CA-C	6.17	121.89	113.37
1	A	554	GLU	N-CA-C	6.12	121.02	113.50
1	B	754	MET	N-CA-C	6.08	119.09	109.96
1	B	555	PHE	N-CA-C	6.07	119.43	109.72
1	A	555	PHE	N-CA-C	6.06	119.42	109.85
1	A	563	GLN	N-CA-C	6.05	121.72	113.37
1	A	748	GLN	N-CA-C	-5.80	101.10	110.32
1	A	570	HIS	CA-C-N	5.70	125.55	119.28
1	A	570	HIS	C-N-CA	5.70	125.55	119.28
1	B	814	LEU	N-CA-C	5.47	118.60	109.96
1	B	748	GLN	N-CA-C	-5.44	101.67	110.32
1	A	814	LEU	N-CA-C	5.26	118.27	109.96
1	A	643	ASP	N-CA-C	-5.14	98.94	107.99
1	A	732	TYR	N-CA-C	5.05	116.78	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	731	ASN	N-CA-C	-5.04	99.56	108.23
1	B	643	ASP	N-CA-C	-5.03	98.52	107.98

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	2703	0	2626	42	0
1	B	2703	0	2626	49	0
2	C	19	0	8	3	0
2	D	19	0	8	4	0
3	A	16	0	12	0	0
3	B	8	0	6	1	0
4	A	333	0	0	4	0
4	B	313	0	0	6	0
All	All	6114	0	5286	92	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 9.

All (92) 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:783:LEU:HD13	1:A:836:LYS:HG3	1.54	0.88
1:B:561:GLU:OE2	2:D:2:XYS:H51	1.75	0.87
1:A:562:ASN:HD21	2:C:2:XYS:H3	1.43	0.83
1:B:536:ASN:HD21	1:B:562:ASN:HB2	1.45	0.82
1:B:783:LEU:HD13	1:B:836:LYS:HG3	1.62	0.81
1:A:536:ASN:HD21	1:A:562:ASN:HB2	1.49	0.78
1:B:818:GLU:H	1:B:818:GLU:CD	1.94	0.76
1:A:530:ILE:H	1:A:556:ASN:HD21	1.35	0.74
1:A:561:GLU:OE2	2:C:2:XYS:H51	1.87	0.74
1:B:707:LYS:O	1:B:711:GLU:HG3	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:562:ASN:HD21	2:D:2:XYS:H3	1.53	0.71
1:A:818:GLU:CD	1:A:818:GLU:H	2.00	0.70
1:A:674:ARG:O	1:A:678:GLU:HG3	1.94	0.68
1:B:537:ASN:ND2	1:B:563:GLN:HE22	1.92	0.67
1:A:707:LYS:O	1:A:711:GLU:HG3	1.95	0.66
1:B:530:ILE:H	1:B:556:ASN:HD21	1.43	0.66
1:B:579:THR:HB	1:B:580:PRO:HD3	1.77	0.66
1:A:770:LYS:HA	1:A:770:LYS:HE2	1.77	0.65
1:B:626:ASP:HB2	4:B:1080:HOH:O	1.98	0.64
1:B:670:GLU:HG3	1:B:714:VAL:CG2	2.30	0.62
1:B:754:MET:HE3	1:B:793:PHE:CZ	2.35	0.62
1:B:827:TYR:O	1:B:831:GLU:HG3	2.01	0.60
1:A:754:MET:HE3	1:A:793:PHE:CZ	2.37	0.59
1:B:561:GLU:CD	2:D:2:XYS:H51	2.27	0.59
1:A:803:VAL:HB	1:A:804:PRO:HD3	1.85	0.58
1:A:827:TYR:O	1:A:831:GLU:HG3	2.04	0.57
1:A:556:ASN:H	1:A:556:ASN:HD22	1.50	0.57
1:B:754:MET:HE3	1:B:793:PHE:CE1	2.39	0.57
1:B:770:LYS:HE3	1:B:770:LYS:HA	1.86	0.56
1:B:531:GLY:HA3	1:B:557:ILE:HG23	1.86	0.56
1:B:670:GLU:HG3	1:B:714:VAL:HG22	1.87	0.56
1:A:531:GLY:HA2	1:A:555:PHE:HB3	1.90	0.54
1:A:556:ASN:HA	1:A:593:MET:HG2	1.89	0.54
1:A:626:ASP:HB2	4:A:998:HOH:O	2.07	0.54
1:B:582:GLU:OE2	1:B:638:ARG:NH1	2.40	0.54
1:A:754:MET:HE3	1:A:793:PHE:CE1	2.43	0.53
1:B:583:LYS:HE3	4:B:894:HOH:O	2.08	0.53
1:A:613:ARG:HD2	1:A:615:TRP:CE2	2.44	0.52
1:B:638:ARG:HD3	4:B:860:HOH:O	2.09	0.52
1:B:640:LYS:HG2	4:B:1000:HOH:O	2.08	0.52
1:A:531:GLY:HA3	1:A:557:ILE:HG23	1.89	0.52
1:B:556:ASN:HA	1:B:593:MET:HG2	1.92	0.52
1:A:640:LYS:HG2	4:A:1001:HOH:O	2.10	0.52
1:B:757:ARG:HA	1:B:813:ALA:HB2	1.91	0.52
1:A:757:ARG:HA	1:A:813:ALA:HB2	1.92	0.51
1:B:803:VAL:HB	1:B:804:PRO:HD3	1.92	0.51
1:B:562:ASN:ND2	1:B:565:LYS:NZ	2.59	0.51
1:B:647:GLU:OE2	2:D:1:XYP:O1	2.28	0.51
1:B:531:GLY:HA2	1:B:555:PHE:HB3	1.94	0.50
1:B:808:LYS:HG3	4:B:929:HOH:O	2.12	0.50
1:A:783:LEU:HD11	1:A:833:LEU:HD23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:558:LEU:C	1:B:558:LEU:HD23	2.38	0.49
1:A:530:ILE:H	1:A:556:ASN:ND2	2.05	0.49
1:A:585:VAL:O	1:A:589:GLU:HG3	2.12	0.49
1:A:533:ALA:HB2	1:A:794:TRP:CE3	2.48	0.49
1:B:670:GLU:HG3	1:B:714:VAL:HG21	1.95	0.48
1:A:556:ASN:HD22	1:A:556:ASN:N	2.11	0.47
1:B:537:ASN:HD21	1:B:563:GLN:HE22	1.59	0.47
1:A:647:GLU:OE2	2:C:1:XYP:O1	2.32	0.47
1:A:736:ARG:HD2	4:A:1052:HOH:O	2.13	0.47
1:B:727:TYR:HB3	1:B:758:ILE:HG22	1.97	0.47
1:B:684:ILE:HD13	1:B:748:GLN:HE22	1.78	0.47
1:B:518:SER:OG	1:B:521:GLU:HG3	2.14	0.47
1:A:538:PHE:HA	1:A:541:LEU:HG	1.97	0.46
1:B:764:GLU:OE1	1:B:812:LYS:HE3	2.15	0.46
1:A:644:VAL:HG21	1:A:685:LEU:HB3	1.97	0.46
1:B:783:LEU:CD1	1:B:836:LYS:HG3	2.41	0.46
1:A:562:ASN:HD22	1:A:565:LYS:NZ	2.14	0.46
1:A:712:LYS:HG3	1:A:714:VAL:HG23	1.98	0.46
1:A:562:ASN:ND2	1:A:565:LYS:HZ1	2.13	0.45
1:B:818:GLU:CD	1:B:818:GLU:N	2.69	0.45
1:B:748:GLN:NE2	1:B:750:TYR:OH	2.49	0.45
1:B:836:LYS:O	1:B:840:ARG:HG3	2.17	0.45
1:A:783:LEU:CD1	1:A:836:LYS:HG3	2.38	0.45
1:A:536:ASN:ND2	1:A:563:GLN:HG3	2.32	0.44
1:B:530:ILE:H	1:B:556:ASN:ND2	2.10	0.44
1:B:533:ALA:HB2	1:B:794:TRP:CE3	2.52	0.44
1:A:580:PRO:HG2	4:A:845:HOH:O	2.16	0.44
1:A:598:HIS:HD2	1:A:643:ASP:OD2	2.00	0.44
1:B:536:ASN:ND2	1:B:563:GLN:HG3	2.31	0.44
3:B:842:ACY:H2	4:B:876:HOH:O	2.17	0.43
1:B:538:PHE:HA	1:B:541:LEU:HG	2.02	0.42
1:B:585:VAL:O	1:B:589:GLU:HG3	2.19	0.42
1:B:562:ASN:HD22	1:B:565:LYS:NZ	2.17	0.42
1:A:777:LYS:HZ1	1:A:781:ILE:HD11	1.85	0.41
1:A:733:ASP:O	1:A:737:ARG:HG3	2.19	0.41
1:B:819:ASN:O	1:B:820:TYR:HB2	2.19	0.41
1:B:616:THR:OG1	1:B:619:GLU:HG3	2.21	0.41
1:B:753:GLU:HG2	1:B:794:TRP:CE3	2.56	0.41
1:A:562:ASN:ND2	1:A:565:LYS:NZ	2.68	0.41
1:A:754:MET:HE1	1:A:791:ILE:CG2	2.51	0.40
1:A:828:ALA:HA	1:A:831:GLU:OE1	2.22	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	322/328 (98%)	311 (97%)	10 (3%)	1 (0%)	36	25
1	B	322/328 (98%)	310 (96%)	12 (4%)	0	100	100
All	All	644/656 (98%)	621 (96%)	22 (3%)	1 (0%)	43	31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	762	GLY

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	287/291 (99%)	285 (99%)	2 (1%)	76	73
1	B	287/291 (99%)	285 (99%)	2 (1%)	76	73
All	All	574/582 (99%)	570 (99%)	4 (1%)	76	73

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	556	ASN
1	A	613	ARG

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Mol	Chain	Res	Type
1	B	556	ASN
1	B	770	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	536	ASN
1	A	556	ASN
1	A	562	ASN
1	A	577	ASN
1	A	598	HIS
1	A	792	GLN
1	B	536	ASN
1	B	537	ASN
1	B	556	ASN
1	B	562	ASN
1	B	577	ASN
1	B	748	GLN
1	B	819	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 ⓘ

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
2	XYP	C	1	2	10,10,10	0.63	0	14,14,14	0.41	0
2	XYS	C	2	2	9,9,10	0.65	0	10,12,14	0.87	0
2	XYP	D	1	2	10,10,10	0.63	0	14,14,14	0.42	0
2	XYS	D	2	2	9,9,10	0.73	0	10,12,14	1.13	1 (10%)

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	XYP	C	1	2	-	-	0/1/1/1
2	XYS	C	2	2	-	-	0/1/1/1
2	XYP	D	1	2	-	-	0/1/1/1
2	XYS	D	2	2	-	-	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	XYS	C1-C2-C3	2.54	113.34	109.64

There are no chirality outliers.

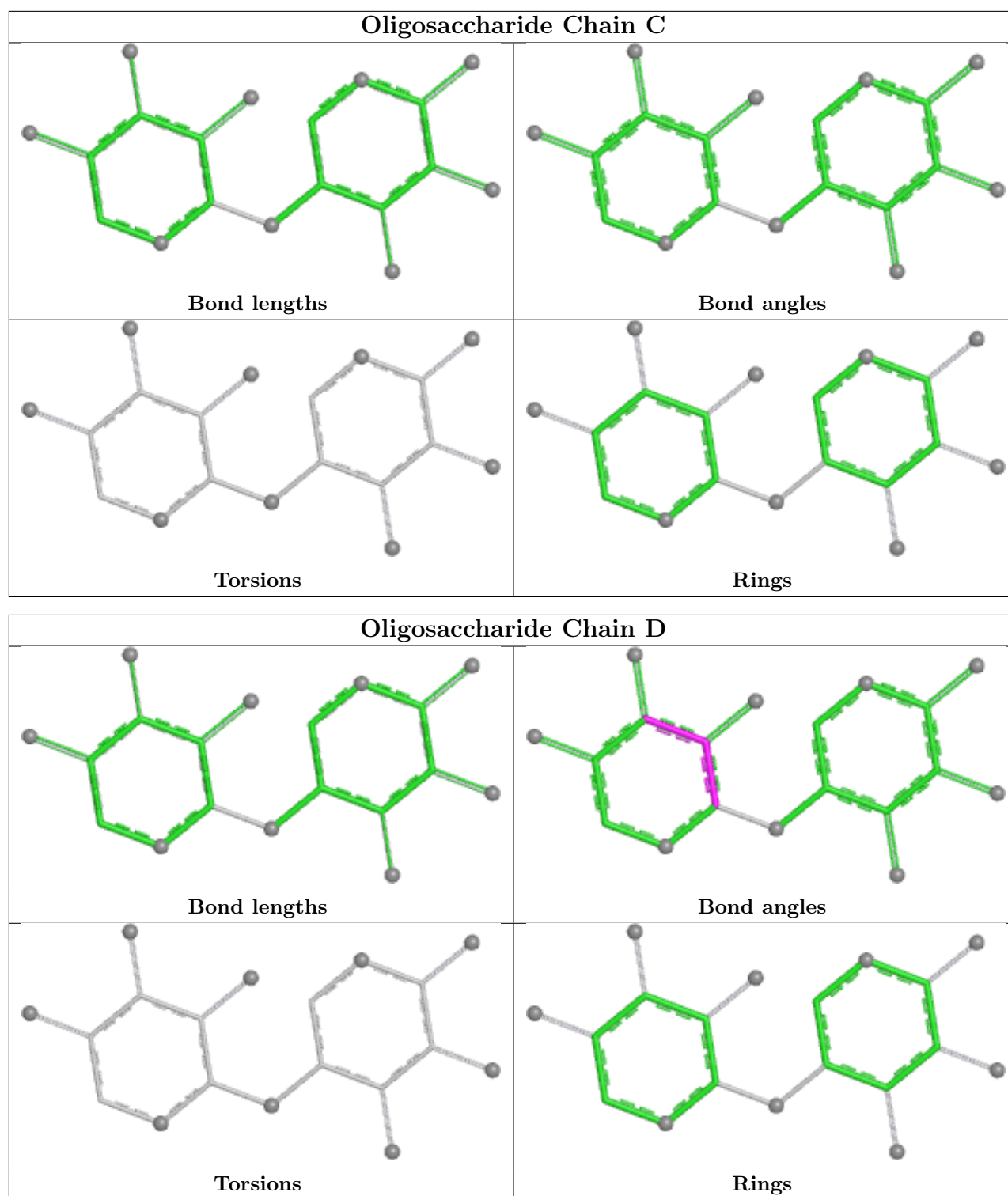
There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	XYP	1	0
2	D	1	XYP	1	0
2	C	2	XYS	2	0
2	D	2	XYS	3	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 ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACY	A	6	-	3,3,3	1.55	0	3,3,3	1.37	0
3	ACY	A	3	-	3,3,3	1.52	0	3,3,3	1.39	0
3	ACY	A	5	-	3,3,3	1.52	0	3,3,3	1.36	0
3	ACY	B	842	-	3,3,3	1.50	0	3,3,3	1.36	0
3	ACY	A	842	-	3,3,3	1.51	0	3,3,3	1.37	0
3	ACY	B	2	-	3,3,3	1.57	0	3,3,3	1.35	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	842	ACY	1	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 [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	324/328 (98%)	-0.17	3 (0%) 81 81	12, 19, 33, 53	0
1	B	324/328 (98%)	-0.16	3 (0%) 81 81	11, 19, 31, 48	0
All	All	648/656 (98%)	-0.17	6 (0%) 81 81	11, 19, 32, 53	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	762	GLY	4.3
1	A	840	ARG	2.7
1	A	760	LEU	2.6
1	A	806	PHE	2.5
1	B	838	GLU	2.3
1	B	761	SER	2.2

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](#)

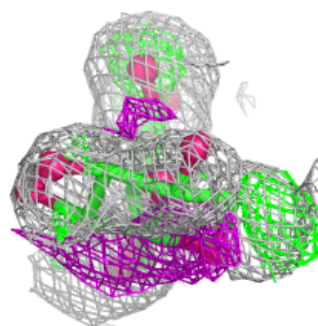
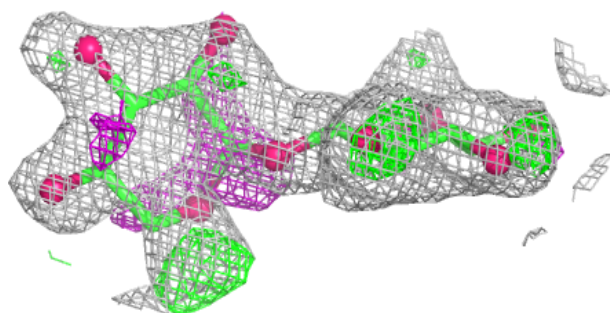
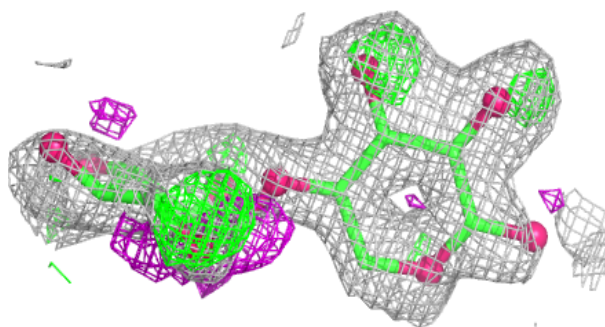
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	XYS	C	2	9/10	0.61	0.23	52,53,54,54	0
2	XYS	D	2	9/10	0.70	0.19	39,40,41,42	0
2	XYP	C	1	10/10	0.77	0.16	42,47,49,51	0
2	XYP	D	1	10/10	0.88	0.11	23,32,36,37	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.

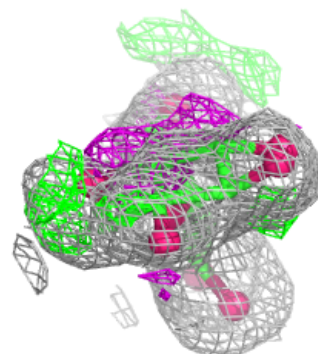
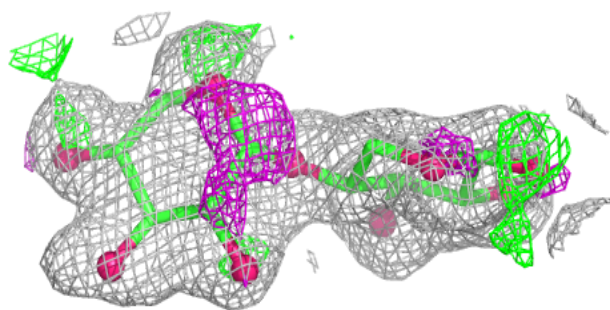
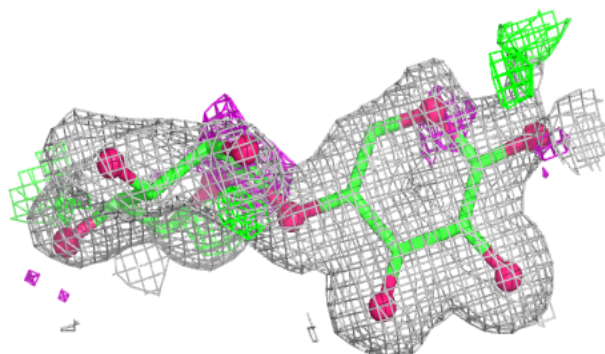
Electron density around Chain C:

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



Electron density around Chain D:

$2mF_o-DF_c$ (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
3	ACY	A	5	4/4	0.73	0.20	44,44,45,46	0
3	ACY	B	2	4/4	0.73	0.21	45,46,46,46	0
3	ACY	B	842	4/4	0.80	0.19	44,44,44,46	0
3	ACY	A	3	4/4	0.84	0.15	38,38,39,41	0
3	ACY	A	6	4/4	0.84	0.19	45,46,47,47	0
3	ACY	A	842	4/4	0.88	0.15	29,31,31,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.