



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:12 AM UTC

PDB ID : 2VRQ / pdb_00002vrq
Title : STRUCTURE OF AN INACTIVE MUTANT OF ARABINOFURANOSIDASE FROM THERMOBACILLUS XYLANILYTICUS IN COMPLEX WITH A PENTASACCHARIDE
Authors : Paes, G.; Skov, L.K.; Odonohue, M.J.; Remond, C.; Kastrup, J.S.; Gajhede, M.; Mirza, O.
Deposited on : 2008-04-09
Resolution : 2.00 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

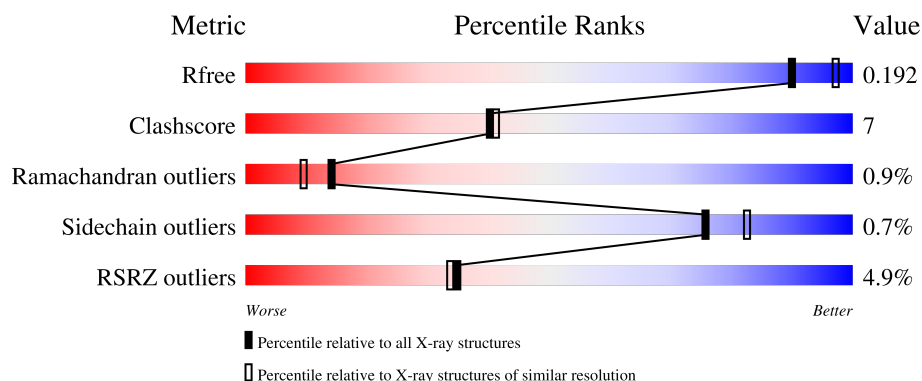
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	496	5% 87% 12% .
1	B	496	6% 88% 11% .
1	C	496	4% 85% 13% ..
2	D	4	25% 75%

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Mol	Chain	Length	Quality of chain
2	E	4	50% 50%
3	F	3	33% 67%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13062 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 ALPHA-L-ARABINOFURANOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	0	0
			3930	2487	694	726	23			
1	B	493	Total	C	N	O	S	0	0	0
			3930	2487	694	726	23			
1	C	493	Total	C	N	O	S	0	0	0
			3930	2487	694	726	23			

There are 3 discrepancies between the modelled and reference sequences:

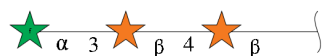
Chain	Residue	Modelled	Actual	Comment	Reference
A	176	GLN	GLU	engineered mutation	UNP O69262
B	176	GLN	GLU	engineered mutation	UNP O69262
C	176	GLN	GLU	engineered mutation	UNP O69262

- Molecule 2 is an oligosaccharide called alpha-L-arabinofuranose-(1-3)-[beta-D-xylopyranose-(1-4)]beta-D-xylopyranose-(1-4)-beta-D-xylopyranose.



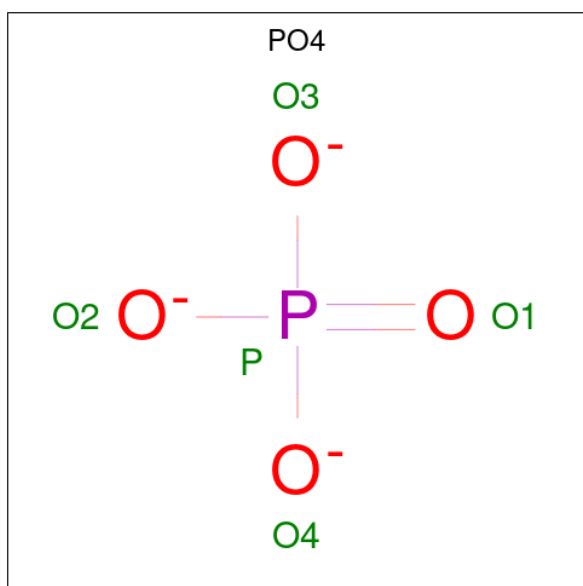
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	D	4	Total	C	O	0	0	0
			37	20	17			
2	E	4	Total	C	O	0	0	0
			37	20	17			

- Molecule 3 is an oligosaccharide called alpha-L-arabinofuranose-(1-3)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	F	3	Total	C	O	0	0	0
			28	15	13			

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		

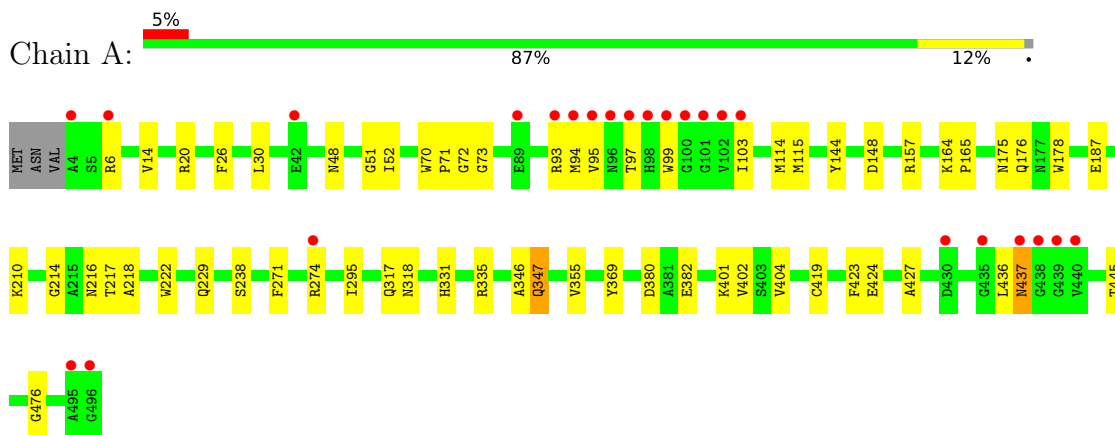
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	413	Total	O	0	0
			413	413		
5	B	376	Total	O	0	0
			376	376		
5	C	371	Total	O	0	0
			371	371		

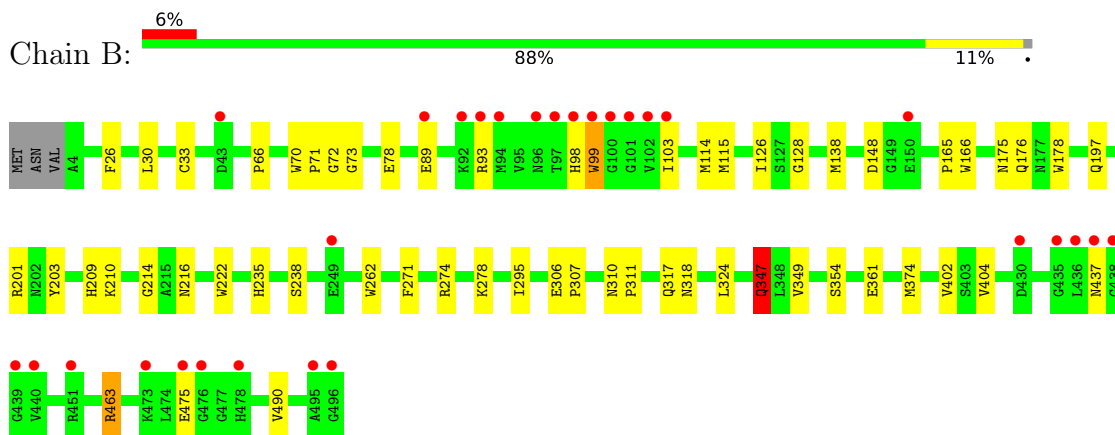
3 Residue-property plots [i](#)

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.

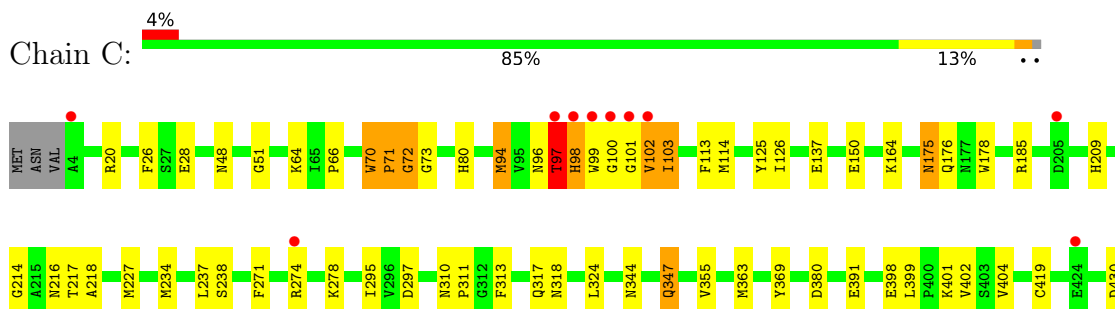
• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE



• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE



• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE





- Molecule 2: alpha-L-arabinofuranose-(1-3)-[beta-D-xylopyranose-(1-4)]beta-D-xylopyranose-(1-4)-beta-D-xylopyranose

Chain D: 25% 75%



- Molecule 2: alpha-L-arabinofuranose-(1-3)-[beta-D-xylopyranose-(1-4)]beta-D-xylopyranose-(1-4)-beta-D-xylopyranose

Chain E: 50% 50%



- Molecule 3: alpha-L-arabinofuranose-(1-3)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose

Chain F: 33% 67%



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	156.80Å 156.80Å 378.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	78.40 – 2.00 78.40 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (78.40-2.00) 100.0 (78.40-2.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.74 (at 2.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.182 , 0.201 0.177 , 0.192	Depositor DCC
R_{free} test set	1847 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.4	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13062	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.65% 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: XYP, AHR, PO4

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	0.55	1/4038 (0.0%)	0.79	0/5482
1	B	0.52	1/4038 (0.0%)	0.76	2/5482 (0.0%)
1	C	0.53	1/4038 (0.0%)	1.23	6/5482 (0.1%)
All	All	0.54	3/12114 (0.0%)	0.95	8/16446 (0.0%)

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
1	B	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	176	GLN	CD-OE1	5.98	1.34	1.23
1	C	176	GLN	CD-OE1	5.84	1.34	1.23
1	B	176	GLN	CD-OE1	5.63	1.34	1.23

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	70	TRP	CA-C-N	48.07	179.93	119.84
1	C	70	TRP	C-N-CA	48.07	179.93	119.84
1	C	70	TRP	C-N-CD	-12.43	74.03	125.00
1	C	72	GLY	N-CA-C	6.83	119.03	110.29
1	C	103	ILE	N-CA-C	6.22	116.88	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	TRP	N-CA-C	5.54	116.34	108.54
1	C	70	TRP	O-C-N	5.47	126.10	121.83
1	B	347	GLN	CB-CA-C	5.07	120.51	110.42

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	346	ALA	Peptide
1	A	72	GLY	Peptide
1	B	72	GLY	Peptide

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	3930	0	3778	48	0
1	B	3930	0	3778	44	0
1	C	3930	0	3778	67	0
2	D	37	0	0	0	0
2	E	37	0	0	0	0
3	F	28	0	0	0	0
4	A	5	0	0	0	0
4	C	5	0	0	0	0
5	A	413	0	0	4	0
5	B	376	0	0	3	0
5	C	371	0	0	7	0
All	All	13062	0	11334	153	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 (153) 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:227:MET:HE3	1:C:237:LEU:HD22	1.35	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:GLN:HE21	1:B:318:ASN:H	1.28	0.79
1:C:71:PRO:HD2	1:C:72:GLY:H	1.49	0.78
1:C:98:HIS:HB2	5:C:752:HOH:O	1.82	0.77
1:A:317:GLN:HE21	1:A:318:ASN:H	1.29	0.76
1:B:114:MET:HE1	1:B:126:ILE:HD11	1.64	0.76
1:C:26:PHE:O	1:C:347:GLN:HA	1.87	0.75
1:B:73:GLY:HA2	1:B:175:ASN:CG	2.12	0.75
1:A:274:ARG:HG2	1:A:274:ARG:HH11	1.50	0.74
1:A:274:ARG:HG2	1:A:274:ARG:NH1	2.00	0.74
1:A:73:GLY:HA2	1:A:175:ASN:CG	2.13	0.74
1:C:271:PHE:HE1	1:C:274:ARG:HD3	1.55	0.71
1:C:317:GLN:HE21	1:C:318:ASN:H	1.37	0.71
1:B:26:PHE:O	1:B:347:GLN:HA	1.91	0.71
1:A:436:LEU:O	1:A:437:ASN:HB2	1.92	0.70
1:B:214:GLY:HA3	1:B:238:SER:O	1.92	0.70
1:C:48:ASN:HD22	1:C:51:GLY:H	1.41	0.68
1:A:331:HIS:O	1:A:335:ARG:HG2	1.94	0.67
1:B:317:GLN:NE2	1:B:318:ASN:H	1.93	0.66
1:C:113:PHE:CD2	1:C:114:MET:HE2	2.29	0.66
1:B:271:PHE:HE1	1:B:274:ARG:HD3	1.61	0.65
1:A:26:PHE:O	1:A:347:GLN:HA	1.97	0.64
1:A:73:GLY:HA2	1:A:175:ASN:ND2	2.13	0.63
1:B:115:MET:HG2	5:B:725:HOH:O	1.98	0.62
1:A:73:GLY:CA	1:A:175:ASN:ND2	2.64	0.61
1:C:113:PHE:HD2	1:C:114:MET:HE2	1.63	0.61
1:C:113:PHE:HD2	1:C:114:MET:CE	2.14	0.60
1:A:214:GLY:HA3	1:A:238:SER:O	2.01	0.60
1:A:274:ARG:HH11	1:A:274:ARG:CG	2.13	0.60
1:C:227:MET:HE2	1:C:227:MET:HA	1.83	0.60
1:C:271:PHE:CE1	1:C:274:ARG:HD3	2.36	0.60
1:B:209:HIS:CE1	1:B:235:HIS:CG	2.91	0.59
1:A:48:ASN:HD22	1:A:51:GLY:H	1.47	0.59
1:C:26:PHE:HE1	1:C:72:GLY:HA2	1.67	0.59
1:C:94:MET:HB2	5:C:812:HOH:O	2.03	0.59
1:A:423:PHE:CE1	1:A:424:GLU:HG3	2.37	0.59
1:C:214:GLY:HA3	1:C:238:SER:O	2.03	0.59
1:C:297:ASP:HA	1:C:344:ASN:HD22	1.68	0.59
1:A:402:VAL:HG12	1:A:404:VAL:HG23	1.86	0.57
1:C:102:VAL:HG22	1:C:313:PHE:HZ	1.70	0.57
1:B:73:GLY:HA2	1:B:175:ASN:OD1	2.02	0.57
1:B:271:PHE:CE1	1:B:274:ARG:HD3	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:ARG:HE	1:C:278:LYS:HZ2	1.52	0.57
1:B:70:TRP:CD1	1:B:71:PRO:HA	2.40	0.57
1:C:227:MET:HE1	1:C:234:MET:SD	2.44	0.56
1:B:30:LEU:HG	1:B:99:TRP:CG	2.40	0.56
1:B:463:ARG:HB2	1:B:463:ARG:HH11	1.71	0.56
1:C:102:VAL:HB	1:C:103:ILE:HD12	1.87	0.56
1:A:97:THR:HG21	1:B:201:ARG:CD	2.36	0.56
1:B:78:GLU:OE2	1:B:98:HIS:ND1	2.40	0.55
1:C:178:TRP:CZ2	1:C:216:ASN:HB2	2.41	0.55
1:C:102:VAL:HG23	1:C:103:ILE:H	1.72	0.55
1:A:30:LEU:HG	1:A:99:TRP:CG	2.42	0.54
1:C:113:PHE:CD2	1:C:114:MET:CE	2.90	0.54
1:A:317:GLN:NE2	1:A:318:ASN:H	2.01	0.54
1:A:93:ARG:HB3	1:A:103:ILE:CG2	2.38	0.53
1:C:96:ASN:O	1:C:98:HIS:N	2.41	0.53
1:A:48:ASN:ND2	1:A:51:GLY:H	2.05	0.53
1:C:274:ARG:NE	1:C:278:LYS:NZ	2.57	0.53
1:C:114:MET:HE3	1:C:126:ILE:HD11	1.90	0.53
1:C:274:ARG:HE	1:C:278:LYS:NZ	2.06	0.52
1:A:73:GLY:C	1:A:175:ASN:ND2	2.68	0.52
1:C:274:ARG:NE	1:C:278:LYS:HZ2	2.08	0.52
1:C:440:VAL:HG23	1:C:474:LEU:HD21	1.93	0.51
1:A:20:ARG:HD3	1:A:380:ASP:OD2	2.11	0.51
1:B:361:GLU:HG2	5:B:702:HOH:O	2.11	0.51
1:C:402:VAL:HA	1:C:419:CYS:O	2.10	0.51
1:A:70:TRP:CD1	1:A:71:PRO:HA	2.45	0.51
1:B:402:VAL:HG12	1:B:404:VAL:HG23	1.93	0.51
1:B:238:SER:HA	1:B:295:ILE:O	2.12	0.50
1:C:96:ASN:OD1	1:C:99:TRP:HB2	2.12	0.50
1:A:144:TYR:O	1:A:157:ARG:HD3	2.12	0.49
1:C:178:TRP:CH2	1:C:216:ASN:HB2	2.47	0.49
1:C:164:LYS:HG2	5:C:807:HOH:O	2.10	0.49
1:C:48:ASN:ND2	1:C:51:GLY:H	2.09	0.49
1:C:98:HIS:C	1:C:100:GLY:H	2.19	0.49
1:C:209:HIS:HD2	5:C:901:HOH:O	1.94	0.49
1:C:430:ASP:HA	1:C:480:ASN:HD22	1.77	0.49
1:A:335:ARG:HH11	1:A:335:ARG:HG3	1.79	0.48
1:C:102:VAL:CG2	1:C:313:PHE:HZ	2.26	0.48
1:A:178:TRP:CZ2	1:A:216:ASN:HB2	2.49	0.48
1:C:217:THR:HG22	1:C:218:ALA:N	2.28	0.48
1:A:164:LYS:HG2	5:A:643:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:THR:HG23	1:C:99:TRP:CD1	2.48	0.47
1:A:210:LYS:HE3	5:A:694:HOH:O	2.14	0.47
1:A:401:LYS:HE3	1:A:427:ALA:HB2	1.96	0.47
1:B:274:ARG:NH2	1:B:278:LYS:HZ2	2.12	0.47
1:C:73:GLY:HA2	1:C:175:ASN:CG	2.40	0.47
1:C:73:GLY:HA2	1:C:175:ASN:OD1	2.15	0.46
1:C:238:SER:HA	1:C:295:ILE:O	2.15	0.46
1:C:317:GLN:NE2	1:C:318:ASN:H	2.07	0.46
1:A:355:VAL:O	1:A:369:TYR:HB2	2.15	0.45
1:A:436:LEU:O	1:A:437:ASN:CB	2.63	0.45
1:B:70:TRP:CG	1:B:71:PRO:HA	2.52	0.45
1:A:187:GLU:HG3	5:A:946:HOH:O	2.16	0.45
1:C:71:PRO:HD2	1:C:72:GLY:N	2.24	0.45
1:A:97:THR:HG21	1:B:201:ARG:HD2	1.98	0.45
1:C:97:THR:O	1:C:99:TRP:N	2.49	0.45
1:B:128:GLY:C	1:B:138:MET:HE2	2.41	0.45
1:A:238:SER:HA	1:A:295:ILE:O	2.17	0.45
1:A:271:PHE:HE1	1:A:274:ARG:HD3	1.82	0.45
1:B:103:ILE:CD1	1:C:150:GLU:O	2.65	0.44
1:A:52:ILE:HD12	1:A:115:MET:HE3	2.00	0.44
1:B:73:GLY:CA	1:B:175:ASN:CG	2.88	0.44
1:C:71:PRO:CD	1:C:72:GLY:N	2.78	0.44
1:C:355:VAL:O	1:C:369:TYR:HB2	2.17	0.44
1:C:80:HIS:HD2	1:C:137:GLU:OE2	2.01	0.44
1:A:94:MET:HE2	1:A:95:VAL:C	2.43	0.43
1:A:229:GLN:OE1	1:C:185:ARG:HD3	2.18	0.43
1:C:398:GLU:HG3	1:C:399:LEU:HG	2.00	0.43
1:B:197:GLN:HB2	1:B:210:LYS:HD3	2.01	0.43
1:A:95:VAL:HG11	1:B:203:TYR:HE1	1.83	0.43
1:A:402:VAL:HA	1:A:419:CYS:O	2.19	0.43
1:B:103:ILE:HD11	5:C:877:HOH:O	2.17	0.43
1:B:317:GLN:NE2	5:B:604:HOH:O	2.48	0.43
1:A:271:PHE:CZ	1:A:274:ARG:NH1	2.87	0.43
1:C:449:SER:O	1:C:486:MET:HE3	2.18	0.43
1:C:20:ARG:HD3	1:C:380:ASP:OD2	2.19	0.43
1:C:98:HIS:C	1:C:100:GLY:N	2.75	0.43
1:C:402:VAL:HG12	1:C:404:VAL:HG23	2.01	0.43
1:A:14:VAL:HG13	1:A:382:GLU:HB3	2.01	0.42
1:A:217:THR:HG22	1:A:218:ALA:N	2.34	0.42
1:C:71:PRO:CD	1:C:72:GLY:H	2.18	0.42
1:B:89:GLU:H	1:B:89:GLU:CD	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:TRP:HB3	1:B:222:TRP:CZ3	2.55	0.42
1:B:262:TRP:CH2	1:B:324:LEU:HD12	2.55	0.42
1:B:178:TRP:CZ2	1:B:216:ASN:HB2	2.54	0.42
1:A:148:ASP:HB3	1:A:165:PRO:HG3	2.02	0.42
1:C:324:LEU:HD23	1:C:419:CYS:HB3	2.01	0.42
1:C:391:GLU:HG3	1:C:401:LYS:CB	2.50	0.42
1:C:455:HIS:O	1:C:463:ARG:HD2	2.19	0.42
1:B:148:ASP:HB3	1:B:165:PRO:HG3	2.02	0.42
1:A:97:THR:HG21	1:B:201:ARG:HG3	2.00	0.42
1:C:28:GLU:HG3	1:C:72:GLY:HA3	2.02	0.42
1:C:70:TRP:CG	1:C:71:PRO:N	2.87	0.41
1:C:310:ASN:HA	1:C:311:PRO:HD2	1.87	0.41
1:A:30:LEU:HD12	1:A:99:TRP:CE2	2.55	0.41
1:A:178:TRP:HB3	1:A:222:TRP:CZ3	2.55	0.41
1:C:114:MET:CE	1:C:126:ILE:HD11	2.50	0.41
1:B:128:GLY:HA3	1:B:138:MET:HE2	2.02	0.41
1:B:347:GLN:O	1:B:354:SER:HA	2.21	0.41
1:A:6:ARG:NH2	5:A:610:HOH:O	2.53	0.41
1:B:310:ASN:HA	1:B:311:PRO:HD2	1.97	0.41
1:B:374:MET:HE2	1:B:490:VAL:HG23	2.01	0.41
1:B:306:GLU:HA	1:B:307:PRO:HD3	1.96	0.41
1:B:33:CYS:HA	1:B:349:VAL:O	2.21	0.41
1:C:102:VAL:HG22	1:C:313:PHE:CZ	2.54	0.41
1:A:271:PHE:CE1	1:A:274:ARG:NH1	2.89	0.40
1:B:93:ARG:HD2	1:B:103:ILE:HG21	2.04	0.40
1:B:178:TRP:CH2	1:B:216:ASN:HB2	2.57	0.40
1:C:80:HIS:HE1	5:C:862:HOH:O	2.02	0.40
1:C:64:LYS:HD3	5:C:753:HOH:O	2.21	0.40
1:B:114:MET:HG2	1:B:166:TRP:CD2	2.56	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	491/496 (99%)	468 (95%)	20 (4%)	3 (1%)	21	17
1	B	491/496 (99%)	469 (96%)	19 (4%)	3 (1%)	21	17
1	C	491/496 (99%)	466 (95%)	18 (4%)	7 (1%)	9	4
All	All	1473/1488 (99%)	1403 (95%)	57 (4%)	13 (1%)	14	9

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	437	ASN
1	B	437	ASN
1	C	71	PRO
1	C	97	THR
1	C	98	HIS
1	A	347	GLN
1	A	476	GLY
1	C	102	VAL
1	C	347	GLN
1	B	347	GLN
1	B	66	PRO
1	C	66	PRO
1	C	101	GLY

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	416/419 (99%)	414 (100%)	2 (0%)	81	87
1	B	416/419 (99%)	414 (100%)	2 (0%)	81	87
1	C	416/419 (99%)	411 (99%)	5 (1%)	63	70
All	All	1248/1257 (99%)	1239 (99%)	9 (1%)	76	82

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	MET
1	A	445	THR
1	B	463	ARG
1	B	475	GLU
1	C	94	MET
1	C	97	THR
1	C	125	TYR
1	C	175	ASN
1	C	363	MET

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	48	ASN
1	A	96	ASN
1	A	155	ASN
1	A	176	GLN
1	A	216	ASN
1	A	221	HIS
1	A	317	GLN
1	A	437	ASN
1	B	176	GLN
1	B	317	GLN
1	B	437	ASN
1	C	21	ASN
1	C	48	ASN
1	C	80	HIS
1	C	155	ASN
1	C	209	HIS
1	C	221	HIS
1	C	235	HIS
1	C	317	GLN
1	C	344	ASN
1	C	437	ASN
1	C	480	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 ⓘ

11 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	D	1	2	10,10,10	0.94	0	14,14,14	0.68	0
2	XYP	D	2	2	9,9,10	1.20	1 (11%)	10,12,14	2.21	4 (40%)
2	AHR	D	3	2	9,9,10	0.77	0	11,12,14	1.30	2 (18%)
2	XYP	D	4	2	9,9,10	1.11	0	10,12,14	1.00	1 (10%)
2	XYP	E	1	2	10,10,10	0.93	0	14,14,14	0.85	0
2	XYP	E	2	2	9,9,10	1.24	0	10,12,14	1.26	2 (20%)
2	AHR	E	3	2	9,9,10	0.99	0	11,12,14	1.42	2 (18%)
2	XYP	E	4	2	9,9,10	0.72	0	10,12,14	0.35	0
3	XYP	F	1	3	10,10,10	0.98	0	14,14,14	0.93	0
3	XYP	F	2	3	9,9,10	1.08	0	10,12,14	1.54	3 (30%)
3	AHR	F	3	3	9,9,10	0.70	0	11,12,14	1.74	2 (18%)

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	D	1	2	-	-	0/1/1/1
2	XYP	D	2	2	-	-	0/1/1/1
2	AHR	D	3	2	-	0/2/15/18	0/1/1/1
2	XYP	D	4	2	-	-	0/1/1/1
2	XYP	E	1	2	-	-	0/1/1/1
2	XYP	E	2	2	-	-	0/1/1/1
2	AHR	E	3	2	-	0/2/15/18	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XYP	E	4	2	-	-	0/1/1/1
3	XYP	F	1	3	-	-	0/1/1/1
3	XYP	F	2	3	-	-	0/1/1/1
3	AHR	F	3	3	-	0/2/15/18	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	XYP	C4-C3	2.03	1.55	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	3	AHR	O2-C2-C3	3.81	118.94	111.43
2	D	2	XYP	C1-C2-C3	3.68	115.00	109.64
3	F	3	AHR	C1-C2-C3	3.29	106.89	101.63
2	D	2	XYP	O2-C2-C3	3.27	116.93	110.15
2	D	2	XYP	C5-C4-C3	2.74	113.63	109.64
2	E	3	AHR	C1-C2-C3	2.71	105.97	101.63
3	F	2	XYP	C5-C4-C3	2.70	113.57	109.64
2	D	3	AHR	C1-C2-C3	2.59	105.77	101.63
3	F	2	XYP	O2-C2-C3	2.42	115.16	110.15
2	E	2	XYP	C5-C4-C3	2.25	112.92	109.64
2	D	4	XYP	C4-C3-C2	-2.19	108.32	110.92
2	D	3	AHR	O2-C2-C3	2.11	115.58	111.43
2	E	3	AHR	O4-C4-C5	2.03	113.06	109.36
3	F	2	XYP	O4-C4-C3	2.03	114.36	110.15
2	E	2	XYP	C4-C3-C2	-2.02	108.52	110.92
2	D	2	XYP	C5-O5-C1	2.01	114.66	111.42

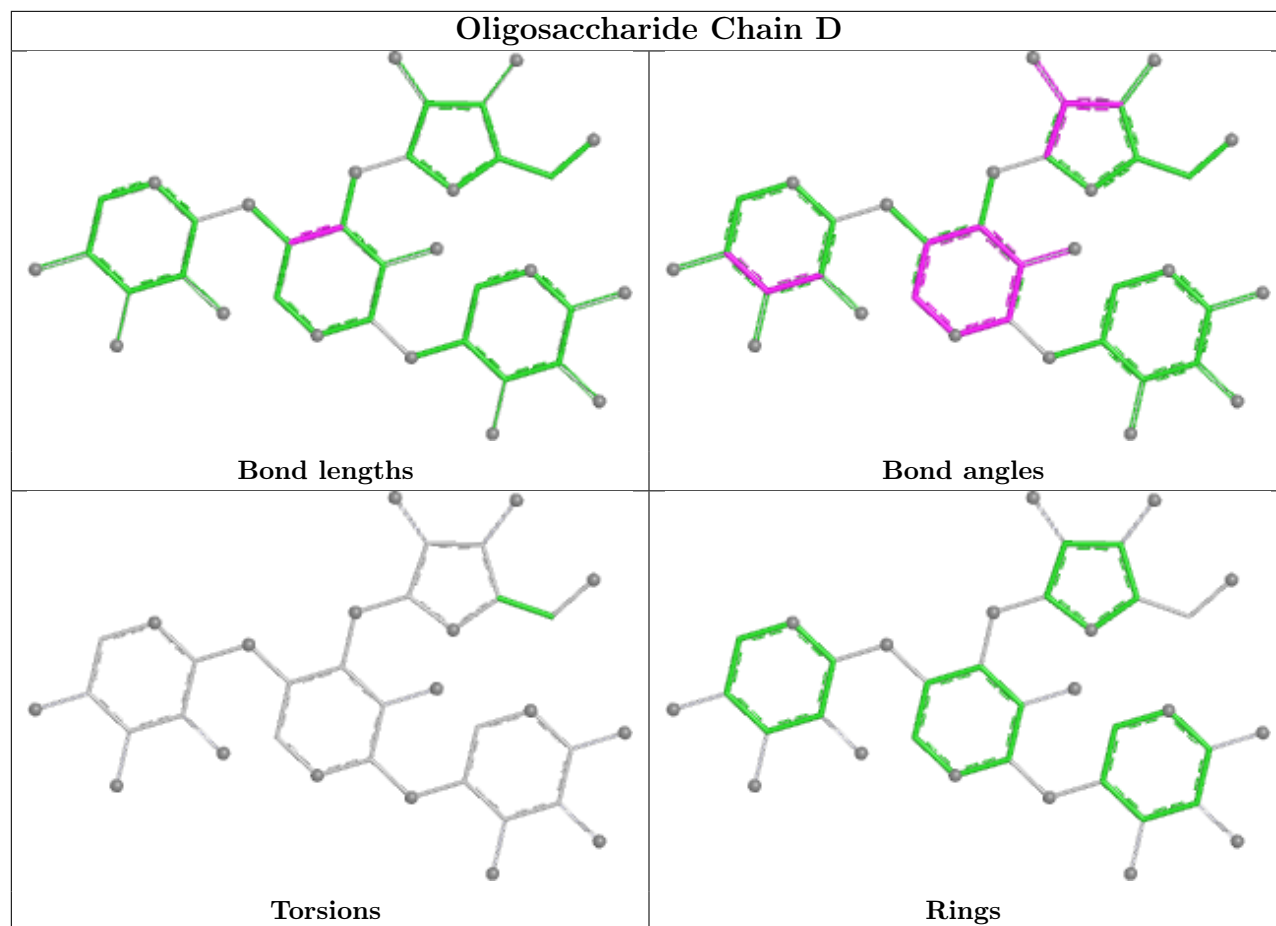
There are no chirality outliers.

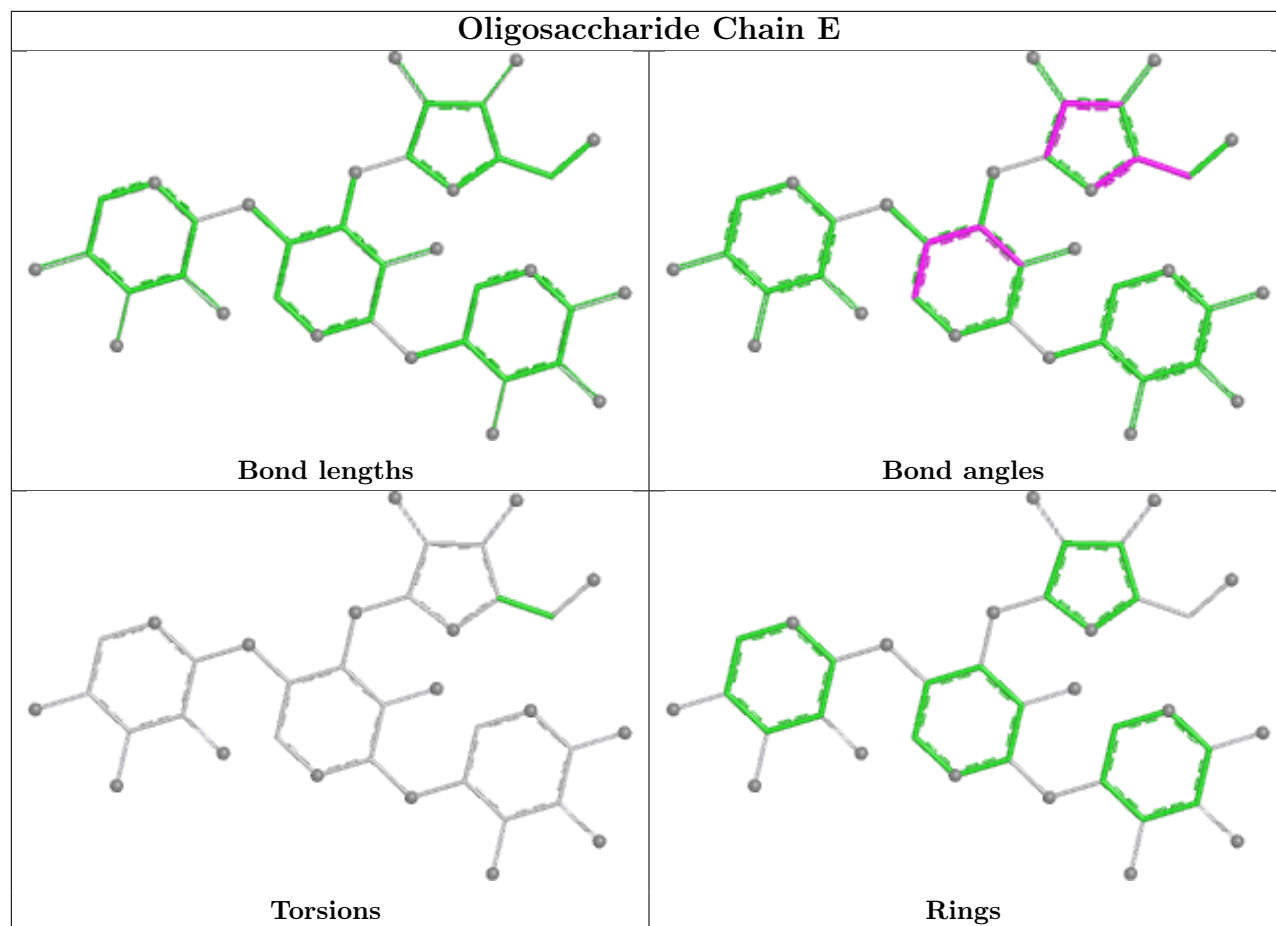
There are no torsion outliers.

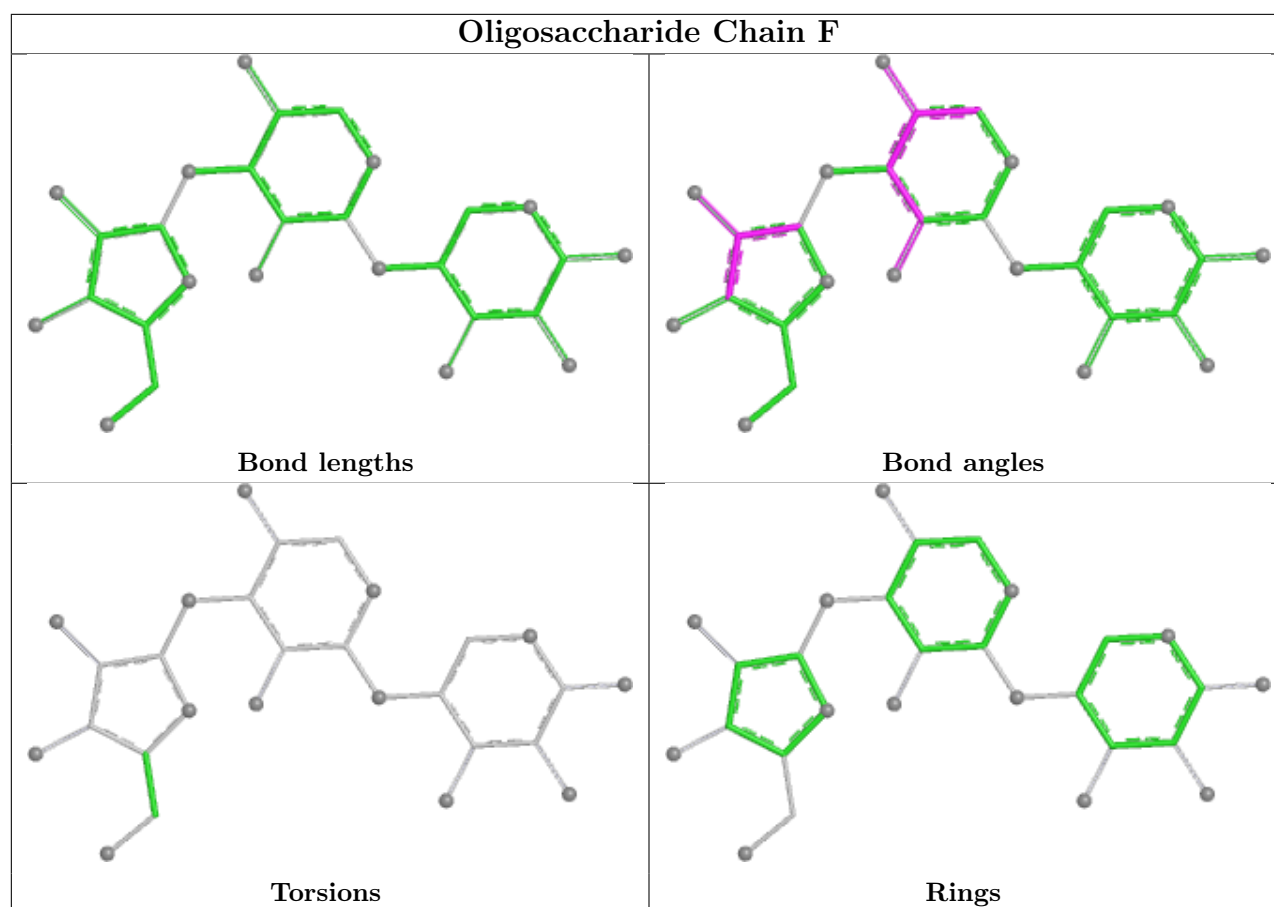
There are no ring outliers.

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](#)

2 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	$\# Z > 2$
4	PO4	C	501	-	4,4,4	0.91	0	6,6,6	0.45	0
4	PO4	A	505	-	4,4,4	0.76	0	6,6,6	0.50	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.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

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	493/496 (99%)	-0.33	24 (4%) 35 34	11, 19, 44, 67	0
1	B	493/496 (99%)	-0.21	29 (5%) 28 27	11, 22, 46, 76	0
1	C	493/496 (99%)	-0.15	20 (4%) 41 40	12, 22, 46, 78	0
All	All	1479/1488 (99%)	-0.23	73 (4%) 35 34	11, 21, 46, 78	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	99	TRP	8.8
1	C	100	GLY	8.2
1	A	103	ILE	8.0
1	C	98	HIS	7.7
1	C	101	GLY	7.3
1	A	94	MET	6.6
1	A	102	VAL	5.8
1	C	102	VAL	5.8
1	A	99	TRP	5.6
1	A	101	GLY	5.5
1	C	496	GLY	5.4
1	A	100	GLY	5.3
1	A	95	VAL	5.1
1	B	496	GLY	4.9
1	A	93	ARG	4.6
1	B	94	MET	4.5
1	C	495	ALA	4.3
1	B	436	LEU	4.2
1	A	98	HIS	4.2
1	A	439	GLY	4.0
1	B	93	ARG	4.0
1	A	97	THR	4.0
1	C	97	THR	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	103	ILE	3.9
1	A	96	ASN	3.7
1	A	496	GLY	3.6
1	C	438	GLY	3.6
1	C	4	ALA	3.6
1	B	102	VAL	3.6
1	A	495	ALA	3.5
1	B	495	ALA	3.5
1	B	89	GLU	3.4
1	B	98	HIS	3.4
1	A	438	GLY	3.4
1	B	100	GLY	3.2
1	B	43	ASP	3.2
1	B	475	GLU	3.1
1	B	96	ASN	3.1
1	B	435	GLY	3.0
1	B	99	TRP	2.9
1	B	101	GLY	2.9
1	A	4	ALA	2.9
1	B	249	GLU	2.8
1	B	439	GLY	2.8
1	B	97	THR	2.8
1	C	436	LEU	2.7
1	C	437	ASN	2.7
1	B	476	GLY	2.6
1	B	437	ASN	2.6
1	A	42	GLU	2.6
1	A	89	GLU	2.5
1	B	92	LYS	2.5
1	A	437	ASN	2.5
1	C	424	GLU	2.5
1	B	478	HIS	2.5
1	C	274	ARG	2.5
1	B	438	GLY	2.4
1	C	476	GLY	2.4
1	B	440	VAL	2.4
1	B	473	LYS	2.3
1	B	430	ASP	2.3
1	C	205	ASP	2.3
1	A	274	ARG	2.2
1	A	440	VAL	2.2
1	B	150	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	433	LEU	2.1
1	A	430	ASP	2.1
1	A	6	ARG	2.1
1	C	439	GLY	2.1
1	C	440	VAL	2.1
1	B	451	ARG	2.0
1	C	474	LEU	2.0
1	A	435	GLY	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](#)

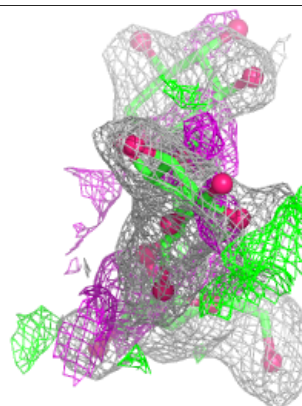
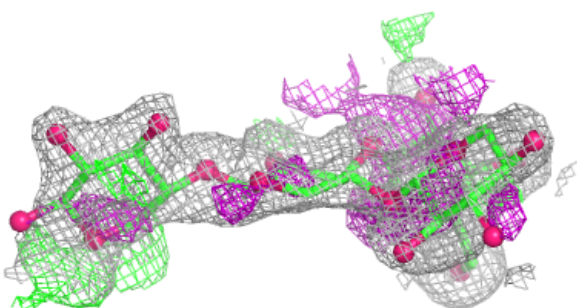
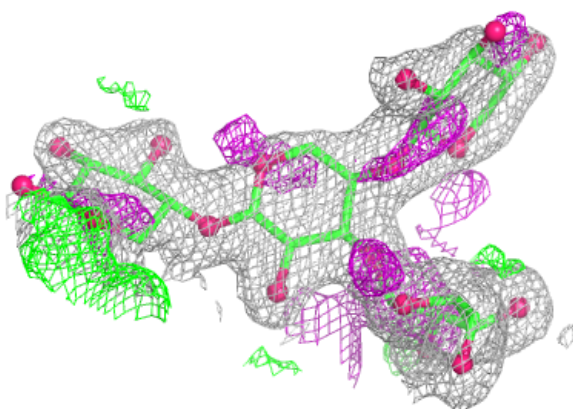
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	XYP	F	1	10/10	0.67	0.18	48,53,63,70	0
2	XYP	E	1	10/10	0.71	0.17	35,47,51,59	0
2	XYP	D	1	10/10	0.73	0.15	39,49,60,65	0
2	XYP	D	4	9/10	0.80	0.16	45,48,55,60	0
2	XYP	E	4	9/10	0.83	0.15	38,41,52,57	0
3	XYP	F	2	9/10	0.86	0.14	36,37,45,48	0
2	AHR	D	3	9/10	0.91	0.12	16,23,29,30	0
2	XYP	D	2	9/10	0.91	0.11	26,32,41,42	0
3	AHR	F	3	9/10	0.91	0.12	19,26,32,35	0
2	AHR	E	3	9/10	0.93	0.10	19,22,27,30	0
2	XYP	E	2	9/10	0.97	0.07	22,28,33,37	0

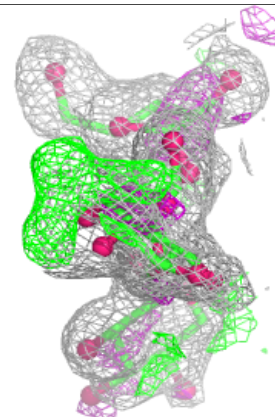
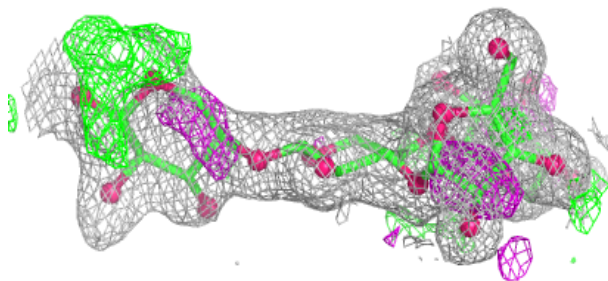
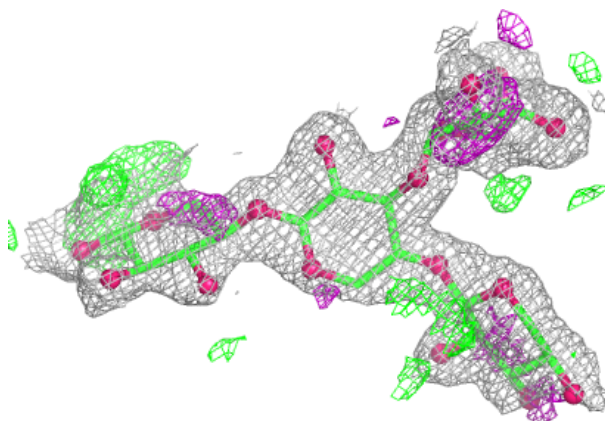
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

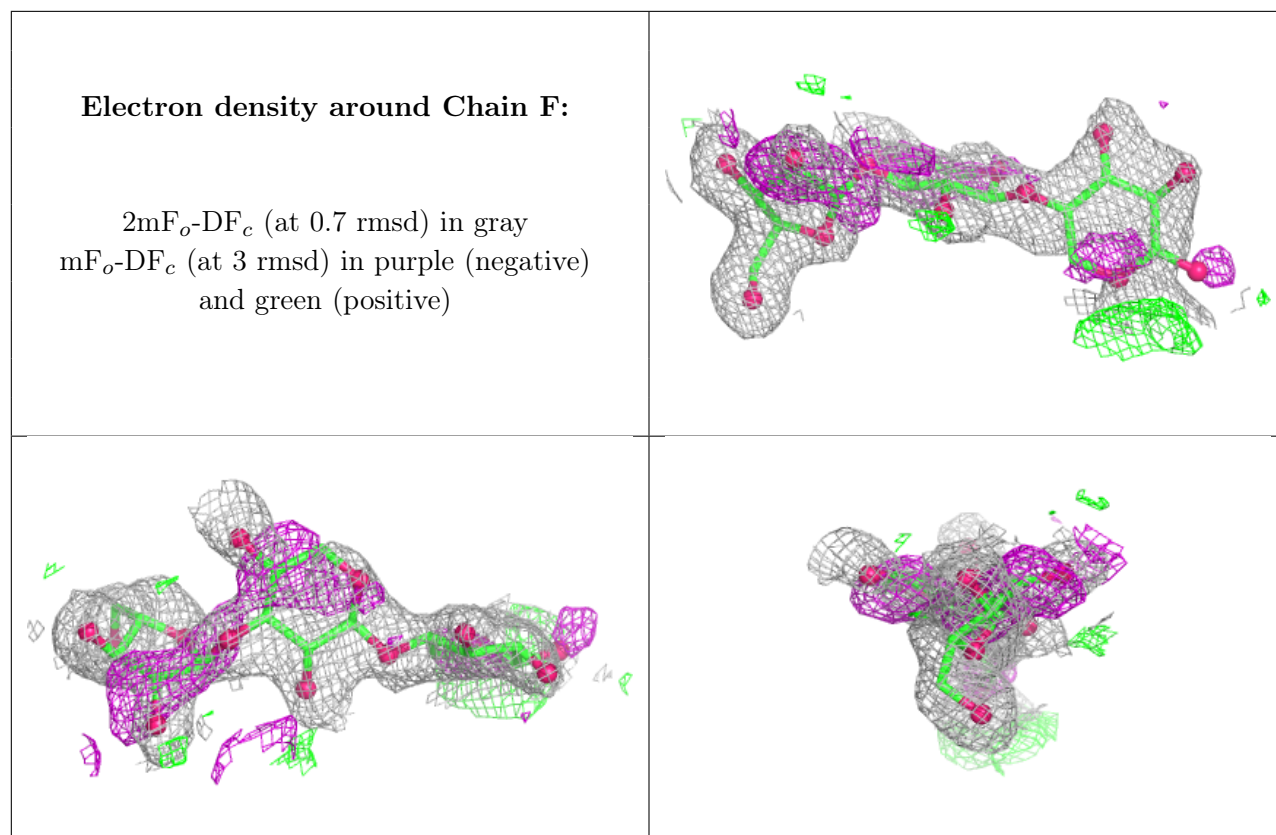
Electron density around Chain D:

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

**Electron density around Chain E:**

$2mF_o-DF_c$ (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	PO4	A	505	5/5	0.93	0.24	26,27,46,52	0
4	PO4	C	501	5/5	0.98	0.06	16,23,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.