Full wwPDB X-ray Structure Validation Report ⓘ

Oct 18, 2023 – 02:47 PM EDT

PDB ID	:	1ISW
Title	:	Crystal structure of xylanase from Streptomyces olivaceoviridis E-86 complexed with xylobiose
Authors	:	Fujimoto, Z.; Kuno, A.; Kaneko, S.; Kobayashi, H.; Kusakabe, I.; Mizuno, H.
Deposited on	:	2001-12-27
Resolution	:	2.10 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:

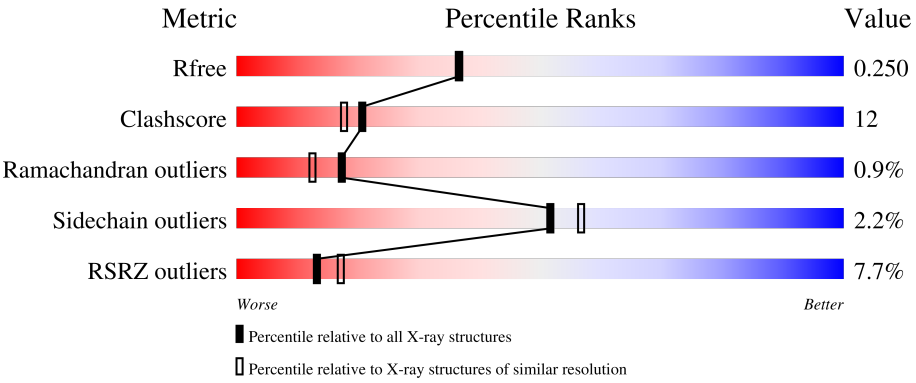
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	436	5% 79% 20% .
1	B	436	10% 79% 20% .
2	C	2	100%
2	D	2	50% 50%
2	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
2	F	2	50% 50%
2	G	2	50% 50%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7466 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.

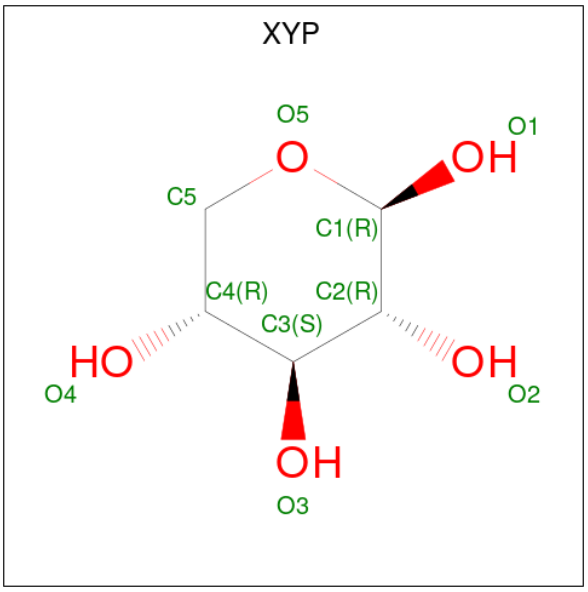
- Molecule 1 is a protein called endo-1,4-beta-D-xylanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	436	Total	C	N	O	S	0	4	0
			3297	2024	600	656	17			
1	B	436	Total	C	N	O	S	0	5	0
			3294	2024	597	656	17			

- Molecule 2 is an oligosaccharide called beta-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				
2	E	2	Total	C	O		0	0	0
			19	10	9				
2	F	2	Total	C	O		0	0	0
			19	10	9				
2	G	2	Total	C	O		0	0	0
			19	10	9				

- Molecule 3 is beta-D-xylopyranose (three-letter code: XYP) (formula: $C_5H_{10}O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 10 5 5	0	0
3	A	1	Total C O 10 5 5	0	0
3	B	1	Total C O 10 5 5	0	0
3	B	1	Total C O 10 5 5	0	0

- Molecule 4 is water.

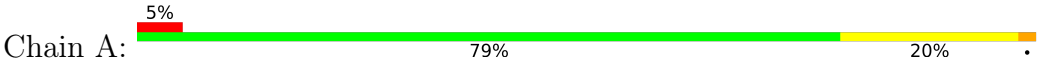
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	357	Total O 357 357	0	0
4	B	383	Total O 383 383	0	0

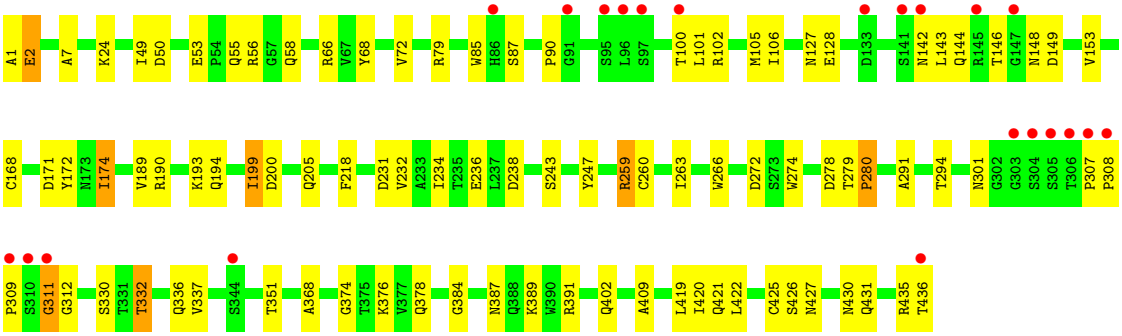
3 Residue-property plots (i)

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: endo-1,4-beta-D-xylanase

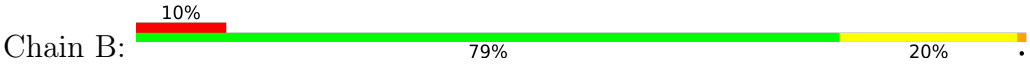
Chain A:

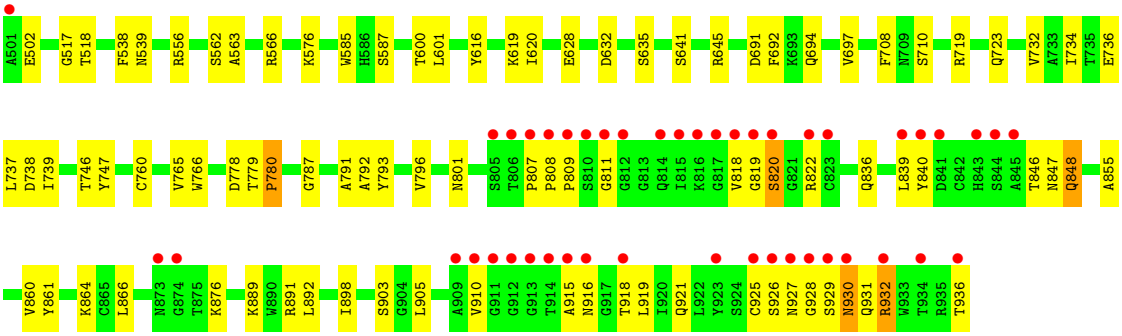




• Molecule 1: endo-1,4-beta-D-xylanase

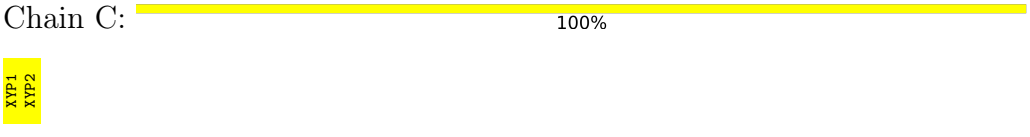
Chain B:





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

Chain C:



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

Chain D:

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

Chain E:

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

Chain F:

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

Chain G:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	74.96Å 94.26Å 137.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.74 – 2.10 44.70 – 2.00	Depositor EDS
% Data completeness (in resolution range)	87.8 (29.74-2.10) 86.5 (44.70-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.69 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.194 , 0.242 0.205 , 0.250	Depositor DCC
R_{free} test set	5881 reflections (10.18%)	wwPDB-VP
Wilson B-factor (Å^2)	24.1	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 62.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7466	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	0.32	0/3390	0.60	0/4600
1	B	0.32	0/3395	0.60	0/4607
All	All	0.32	0/6785	0.60	0/9207

There are no bond length outliers.

There are no bond angle outliers.

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	3297	0	3098	76	0
1	B	3294	0	3084	83	0
2	C	19	0	0	1	0
2	D	19	0	0	1	0
2	E	19	0	0	0	0
2	F	19	0	0	1	0
2	G	19	0	0	0	0
3	A	20	0	0	0	0
3	B	20	0	0	0	0
4	A	357	0	0	8	0
4	B	383	0	0	7	0
All	All	7466	0	6182	159	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 (159) 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:50:ASP:HA	1:A:90:PRO:HG3	1.43	1.00
1:A:24[B]:LYS:NZ	1:A:24[B]:LYS:HE2	1.33	0.98
1:A:24[B]:LYS:NZ	1:A:24[B]:LYS:HE3	1.33	0.97
1:B:925:CYS:HA	1:B:931:GLN:HE22	1.36	0.88
1:A:144:GLN:HE22	1:A:148:ASN:HA	1.37	0.88
1:B:616:TYR:HB3	1:B:620[A]:ILE:HD13	1.56	0.85
1:B:616:TYR:HB3	1:B:620[A]:ILE:CD1	2.09	0.81
1:A:144:GLN:NE2	1:A:148:ASN:HA	2.04	0.72
1:A:2:GLU:HB2	1:A:301:ASN:OD1	1.90	0.72
1:A:234:ILE:HD12	1:A:263:ILE:HG12	1.72	0.72
1:B:708:PHE:HB2	1:B:739[A]:ILE:HD12	1.71	0.72
1:A:49:ILE:HG22	1:A:90:PRO:HD3	1.71	0.72
1:A:174:ILE:H	1:A:174:ILE:HD13	1.55	0.70
1:A:55:GLN:HB3	1:A:58:GLN:HB2	1.74	0.70
1:A:66:ARG:NH1	4:A:1818:HOH:O	2.25	0.69
1:B:929:SER:HA	1:B:932:ARG:HE	1.57	0.68
1:B:839:LEU:HD12	1:B:918:THR:OG1	1.94	0.67
1:A:236:GLU:HG2	1:A:266:TRP:CE3	2.30	0.67
1:B:818:VAL:HG23	1:B:819:GLY:N	2.09	0.67
1:B:819:GLY:HA3	1:B:929:SER:HB2	1.77	0.67
1:B:903:SER:HB2	1:B:905:LEU:HD13	1.76	0.66
1:A:1:ALA:HB1	1:A:7:ALA:HB1	1.78	0.65
1:B:818:VAL:HG23	1:B:819:GLY:H	1.62	0.64
1:B:836:GLN:HB2	1:B:876:LYS:HE3	1.80	0.64
1:A:102:ARG:O	1:A:106:ILE:HG12	1.99	0.63
1:B:719:ARG:O	1:B:723:GLN:HG3	2.00	0.62
1:B:915:ALA:O	1:B:918:THR:HG23	2.01	0.59
1:A:238:ASP:HB2	1:A:280:PRO:HB2	1.85	0.59
1:B:502:GLU:HB2	1:B:801:ASN:OD1	2.03	0.58
1:A:435:ARG:HD3	4:A:1726:HOH:O	2.03	0.57
1:B:556:ARG:HH12	1:B:600:THR:HG22	1.69	0.57
1:A:426:SER:O	1:A:427:ASN:HB2	2.05	0.57
1:B:635:SER:HB3	4:B:1335:HOH:O	2.03	0.57
1:A:231:ASP:OD1	1:A:259:ARG:HG3	2.04	0.56
1:B:778:ASP:O	1:B:779:THR:C	2.43	0.56
1:B:925:CYS:HA	1:B:931:GLN:NE2	2.12	0.56
1:B:910:VAL:HG23	1:B:919:LEU:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ILE:HG22	1:A:90:PRO:CD	2.36	0.55
1:B:891:ARG:NH1	4:B:1330:HOH:O	2.28	0.55
1:A:190:ARG:O	1:A:194:GLN:HG3	2.06	0.55
1:A:1:ALA:CB	1:A:7:ALA:HB1	2.37	0.55
1:B:738:ASP:HB2	1:B:780:PRO:HB2	1.88	0.54
1:B:846:THR:C	1:B:848:GLN:H	2.10	0.54
1:B:822:ARG:HG2	1:B:822:ARG:HH11	1.72	0.54
1:A:311:GLY:HA2	1:A:351:THR:HG23	1.89	0.54
1:B:820:SER:OG	1:B:839:LEU:HD13	2.08	0.54
1:A:172:TYR:HB3	1:A:205:GLN:OE1	2.07	0.54
1:A:105:MET:HE3	1:A:143:LEU:HD22	1.90	0.53
1:B:930:ASN:HD22	1:B:930:ASN:N	2.06	0.53
1:B:860:VAL:O	1:B:861:TYR:HB2	2.09	0.53
1:B:563:ALA:HA	1:B:566:ARG:NH1	2.24	0.53
1:A:24[A]:LYS:HE2	1:A:272:ASP:OD1	2.09	0.52
1:A:278:ASP:O	1:A:279:THR:C	2.48	0.52
1:B:739[B]:ILE:HD13	1:B:746:THR:HG22	1.90	0.52
1:A:199:ILE:HD13	1:A:200:ASP:N	2.25	0.51
1:A:279:THR:N	1:A:280:PRO:HD3	2.25	0.51
1:A:128:GLU:HA	1:A:171:ASP:CG	2.31	0.51
1:B:619:LYS:C	1:B:620[A]:ILE:HD12	2.31	0.51
1:B:840:TYR:CD1	1:B:916:ASN:HB3	2.46	0.51
1:A:24[B]:LYS:HE3	1:A:24[B]:LYS:HZ1	1.24	0.50
1:A:409:ALA:HB3	1:A:430:ASN:HB2	1.92	0.50
1:A:425:CYS:HA	1:A:431:GLN:OE1	2.12	0.50
1:A:101:LEU:HD21	1:A:142:ASN:OD1	2.11	0.50
1:A:294:THR:OG1	1:A:391:ARG:HD2	2.12	0.50
1:B:820:SER:C	1:B:822:ARG:H	2.15	0.49
1:A:100:THR:HG23	4:A:1661:HOH:O	2.12	0.49
1:A:105:MET:CE	1:A:143:LEU:HD22	2.42	0.49
1:B:926:SER:C	1:B:928:GLY:H	2.15	0.49
1:A:53:GLU:OE2	1:A:56:ARG:HA	2.13	0.49
1:A:1:ALA:HB3	4:A:1637:HOH:O	2.11	0.49
1:B:619:LYS:HE2	4:B:1281:HOH:O	2.11	0.49
1:B:860:VAL:O	1:B:864:LYS:HB2	2.12	0.49
1:B:925:CYS:CA	1:B:931:GLN:HE22	2.18	0.49
1:A:336:GLN:HB2	1:A:376:LYS:HE2	1.95	0.49
1:B:616:TYR:HB3	1:B:620[A]:ILE:HD11	1.90	0.49
1:B:708:PHE:HB2	1:B:739[A]:ILE:CD1	2.42	0.49
1:A:24[B]:LYS:HE2	1:A:24[B]:LYS:HZ3	1.23	0.49
1:A:142:ASN:O	1:A:146:THR:HG23	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:779:THR:N	1:B:780:PRO:HD3	2.28	0.48
1:A:330:SER:OG	1:A:332:THR:HG23	2.13	0.48
1:B:585:TRP:CE2	1:B:587:SER:HB3	2.48	0.48
1:A:199:ILE:HD13	1:A:199:ILE:C	2.34	0.48
1:B:628:GLU:OE2	2:F:1:XYP:O1	2.31	0.48
1:A:85:TRP:CE2	1:A:87:SER:HB3	2.49	0.48
1:B:818:VAL:CG2	1:B:819:GLY:N	2.76	0.48
1:A:127:ASN:ND2	1:A:128:GLU:HG3	2.29	0.48
1:A:149:ASP:O	1:A:153:VAL:HG23	2.13	0.48
1:A:368:ALA:HB1	1:A:422:LEU:HD11	1.96	0.48
1:B:556:ARG:HH12	1:B:600:THR:CG2	2.26	0.48
1:A:374:GLY:HA2	1:A:421:GLN:OE1	2.13	0.47
1:A:2:GLU:CB	1:A:301:ASN:OD1	2.62	0.47
1:B:855:ALA:O	1:B:889:LYS:HD2	2.13	0.47
1:A:236:GLU:HG2	1:A:266:TRP:CZ3	2.49	0.47
1:A:387:ASN:HB3	1:A:402:GLN:OE1	2.15	0.47
1:A:189:VAL:O	1:A:193:LYS:HG2	2.15	0.47
1:A:391:ARG:NH2	4:A:1822:HOH:O	2.48	0.47
1:B:738:ASP:O	1:B:739[A]:ILE:HD13	2.13	0.47
1:B:926:SER:C	1:B:928:GLY:N	2.69	0.46
1:B:930:ASN:HD22	1:B:930:ASN:H	1.63	0.46
1:A:308:PRO:HA	1:A:309:PRO:HD3	1.83	0.46
1:B:708:PHE:HD2	1:B:739[A]:ILE:CD1	2.28	0.46
1:B:819:GLY:CA	1:B:929:SER:HB2	2.45	0.46
1:B:926:SER:O	1:B:928:GLY:N	2.49	0.46
1:A:291:ALA:HA	1:A:391:ARG:HG2	1.98	0.45
1:B:539:ASN:HA	1:B:576:LYS:HG3	1.98	0.45
1:B:562:SER:O	1:B:566:ARG:HG3	2.17	0.45
1:B:791:ALA:HA	1:B:891:ARG:HG2	1.99	0.44
1:B:818:VAL:CG2	1:B:819:GLY:H	2.28	0.44
1:B:517:GLY:HA2	1:B:538:PHE:HB3	1.99	0.44
1:B:921:GLN:CD	4:B:1324:HOH:O	2.55	0.44
1:B:840:TYR:HA	1:B:916:ASN:OD1	2.17	0.44
1:A:66:ARG:NH1	4:A:1623:HOH:O	2.51	0.44
1:A:174:ILE:HD13	1:A:174:ILE:N	2.28	0.44
1:A:171:ASP:O	1:A:205:GLN:HG3	2.18	0.44
1:A:307:PRO:HA	1:A:308:PRO:HD3	1.92	0.44
1:B:787:GLY:HA3	4:B:1152:HOH:O	2.18	0.44
1:B:932:ARG:HG2	1:B:932:ARG:HH11	1.82	0.43
1:A:384:GLY:O	1:A:389:LYS:HE2	2.18	0.43
1:B:866:LEU:HD12	1:B:866:LEU:HA	1.94	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:691:ASP:HA	1:B:694:GLN:HG2	2.01	0.43
1:B:732:VAL:HG23	1:B:760:CYS:HA	2.00	0.43
1:A:79:ARG:HB3	1:A:79:ARG:NH1	2.33	0.43
1:B:518:THR:HA	1:B:765:VAL:O	2.19	0.43
1:B:692:PHE:HB3	1:B:697:VAL:HB	2.00	0.43
1:B:739[B]:ILE:HD13	1:B:746:THR:CG2	2.49	0.43
1:B:566:ARG:NH1	4:B:1118:HOH:O	2.52	0.42
2:C:1:XYP:O1	2:D:2:XYP:O4	2.37	0.42
1:B:807:PRO:HA	1:B:808:PRO:HD3	1.84	0.42
1:B:641:SER:O	1:B:645:ARG:HG3	2.19	0.42
1:B:819:GLY:HA3	1:B:929:SER:CB	2.48	0.42
1:B:892:LEU:CD2	1:B:898:ILE:HG12	2.49	0.42
1:B:738:ASP:C	1:B:739[A]:ILE:HD13	2.40	0.42
1:B:929:SER:HA	1:B:932:ARG:NE	2.30	0.42
1:A:309:PRO:HB3	1:A:436:THR:O	2.19	0.42
1:B:793:TYR:CD1	1:B:793:TYR:C	2.93	0.42
1:B:892:LEU:HD23	1:B:898:ILE:HG12	2.01	0.42
1:A:49:ILE:CG2	1:A:90:PRO:HD3	2.47	0.42
1:A:378:GLN:NE2	1:B:710:SER:HB2	2.35	0.42
1:A:259:ARG:HB2	4:A:1556:HOH:O	2.19	0.42
1:B:734:ILE:CG2	1:B:737:LEU:HB2	2.50	0.41
1:B:736:GLU:HG2	1:B:766:TRP:CE3	2.55	0.41
1:B:910:VAL:HA	4:B:1324:HOH:O	2.20	0.41
1:A:419:LEU:C	1:A:420[A]:ILE:HD12	2.41	0.41
1:A:232:VAL:HG23	1:A:260:CYS:HA	2.03	0.41
1:A:68:TYR:CZ	1:A:72:VAL:HG21	2.55	0.41
1:A:218:PHE:HB2	4:A:1691:HOH:O	2.21	0.41
1:B:792:ALA:O	1:B:796:VAL:HG23	2.21	0.41
1:B:822:ARG:HB2	1:B:839:LEU:HB3	2.03	0.41
1:B:936:THR:HG22	1:B:936:THR:OXT	2.20	0.41
1:B:929:SER:CA	1:B:932:ARG:HE	2.28	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	438/436 (100%)	417 (95%)	18 (4%)	3 (1%)	22 18
1	B	439/436 (101%)	415 (94%)	19 (4%)	5 (1%)	14 9
All	All	877/872 (101%)	832 (95%)	37 (4%)	8 (1%)	17 12

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	311	GLY
1	B	820	SER
1	B	927	ASN
1	A	2	GLU
1	B	811	GLY
1	B	809	PRO
1	B	847	ASN
1	A	312	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	345/341 (101%)	337 (98%)	8 (2%)	50 55
1	B	346/341 (102%)	339 (98%)	7 (2%)	55 60
All	All	691/682 (101%)	676 (98%)	15 (2%)	52 57

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

Mol	Chain	Res	Type
1	A	174	ILE
1	A	199	ILE
1	A	243	SER
1	A	247	TYR
1	A	259	ARG

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Mol	Chain	Res	Type
1	A	274	TRP
1	A	280	PRO
1	A	332	THR
1	B	601	LEU
1	B	632	ASP
1	B	747	TYR
1	B	780	PRO
1	B	848	GLN
1	B	930	ASN
1	B	932	ARG

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	55	GLN
1	A	69	ASN
1	A	173	ASN
1	A	284	ASN
1	A	314	GLN
1	B	511	GLN
1	B	603	GLN
1	B	673	ASN
1	B	784	ASN
1	B	814	GLN
1	B	927	ASN
1	B	930	ASN
1	B	931	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

10 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	0.59	0	14,14,14	0.57	0
2	XYP	C	2	2	9,9,10	0.69	0	10,12,14	0.73	1 (10%)
2	XYP	D	1	2	10,10,10	0.62	0	14,14,14	0.55	0
2	XYP	D	2	2	9,9,10	0.80	0	10,12,14	0.72	1 (10%)
2	XYP	E	1	2	10,10,10	0.72	0	14,14,14	0.55	0
2	XYP	E	2	2	9,9,10	0.69	0	10,12,14	0.75	1 (10%)
2	XYP	F	1	2	10,10,10	0.64	0	14,14,14	0.63	0
2	XYP	F	2	2	9,9,10	0.73	0	10,12,14	0.70	0
2	XYP	G	1	2	10,10,10	0.60	0	14,14,14	0.56	0
2	XYP	G	2	2	9,9,10	0.82	0	10,12,14	0.72	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	XYP	C	2	2	-	-	0/1/1/1
2	XYP	D	1	2	-	-	0/1/1/1
2	XYP	D	2	2	-	-	0/1/1/1
2	XYP	E	1	2	-	-	0/1/1/1
2	XYP	E	2	2	-	-	0/1/1/1
2	XYP	F	1	2	-	-	0/1/1/1
2	XYP	F	2	2	-	-	0/1/1/1
2	XYP	G	1	2	-	-	0/1/1/1
2	XYP	G	2	2	-	-	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	E	2	XYP	C4-C3-C2	-2.16	108.36	110.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	2	XYP	C4-C3-C2	-2.09	108.44	110.92
2	D	2	XYP	C4-C3-C2	-2.04	108.49	110.92
2	G	2	XYP	C4-C3-C2	-2.03	108.51	110.92

There are no chirality outliers.

There are no torsion outliers.

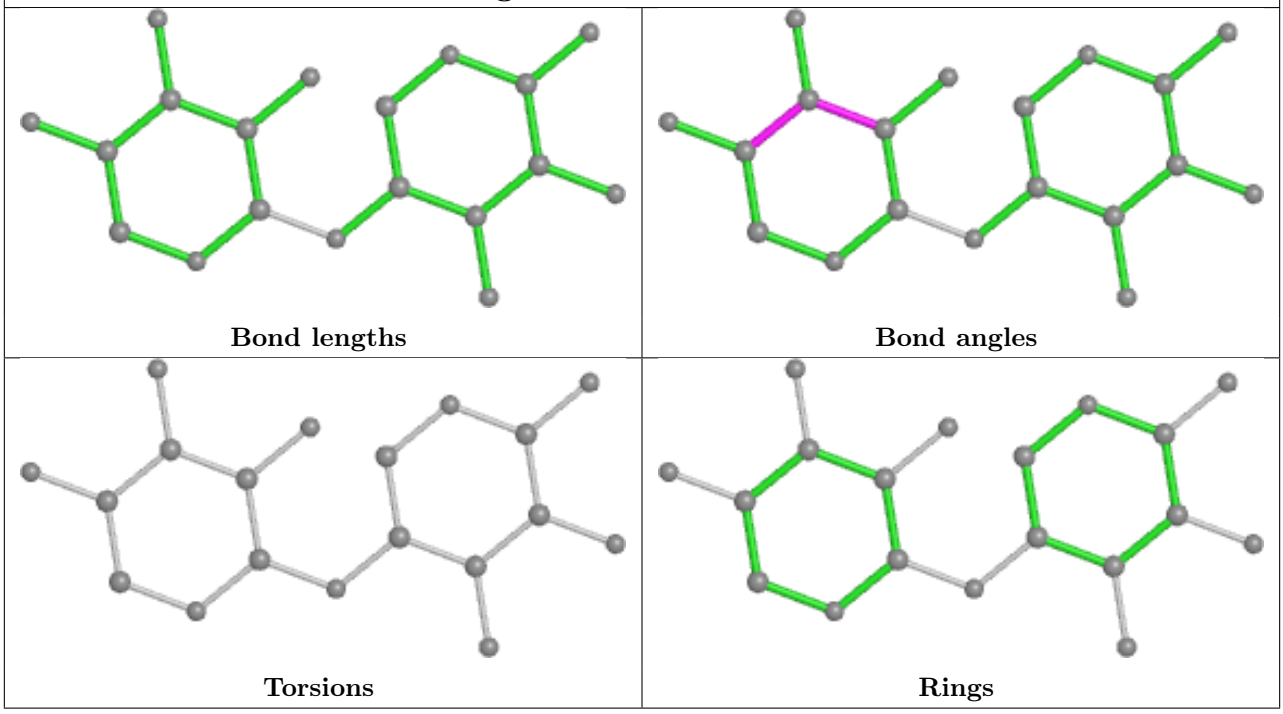
There are no ring outliers.

3 monomers are involved in 2 short contacts:

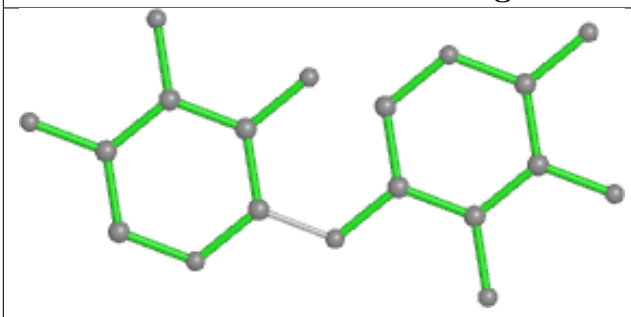
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1	XYP	1	0
2	C	1	XYP	1	0
2	D	2	XYP	1	0

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

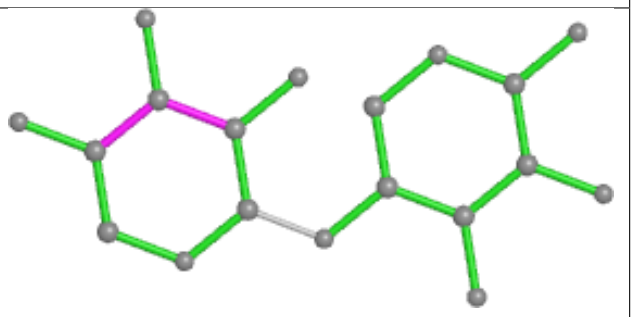
Oligosaccharide Chain C



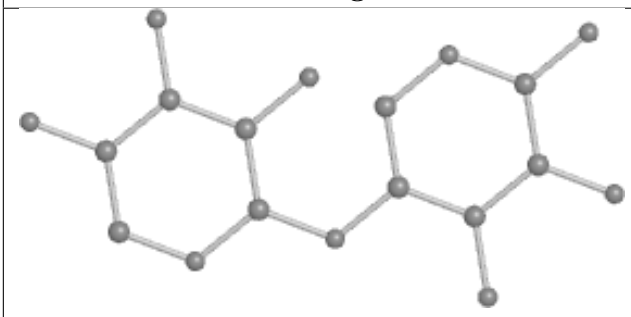
Oligosaccharide Chain D



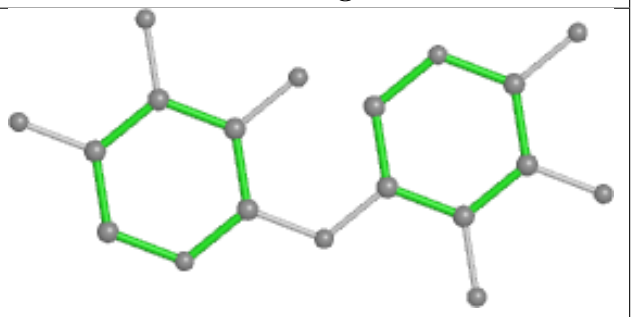
Bond lengths



Bond angles

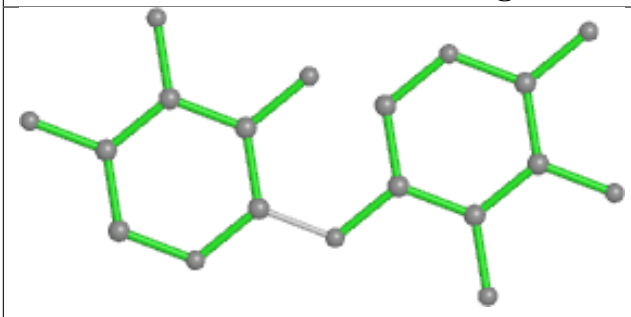


Torsions

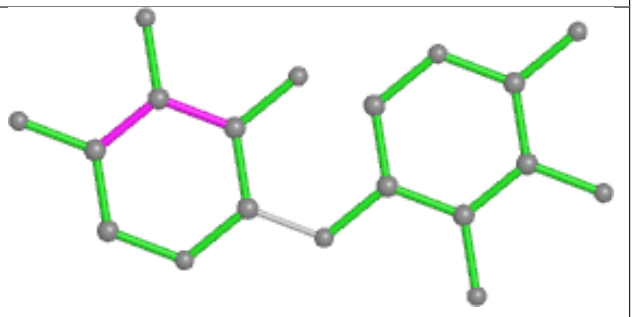


Rings

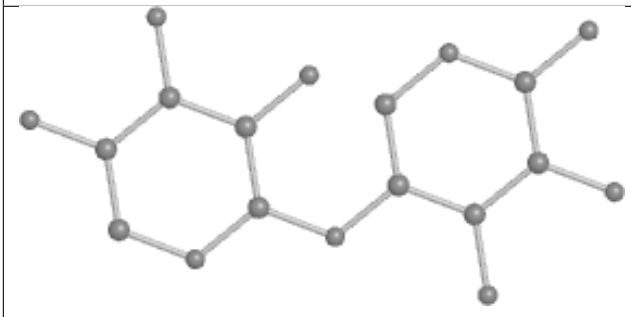
Oligosaccharide Chain E



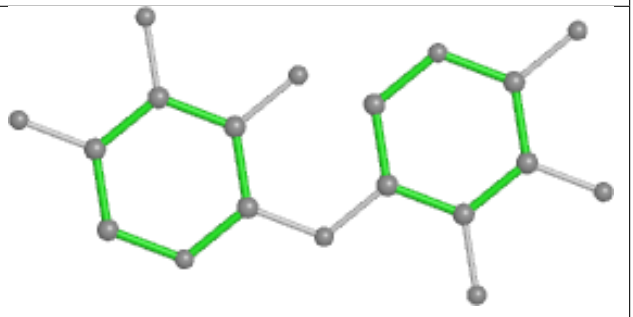
Bond lengths



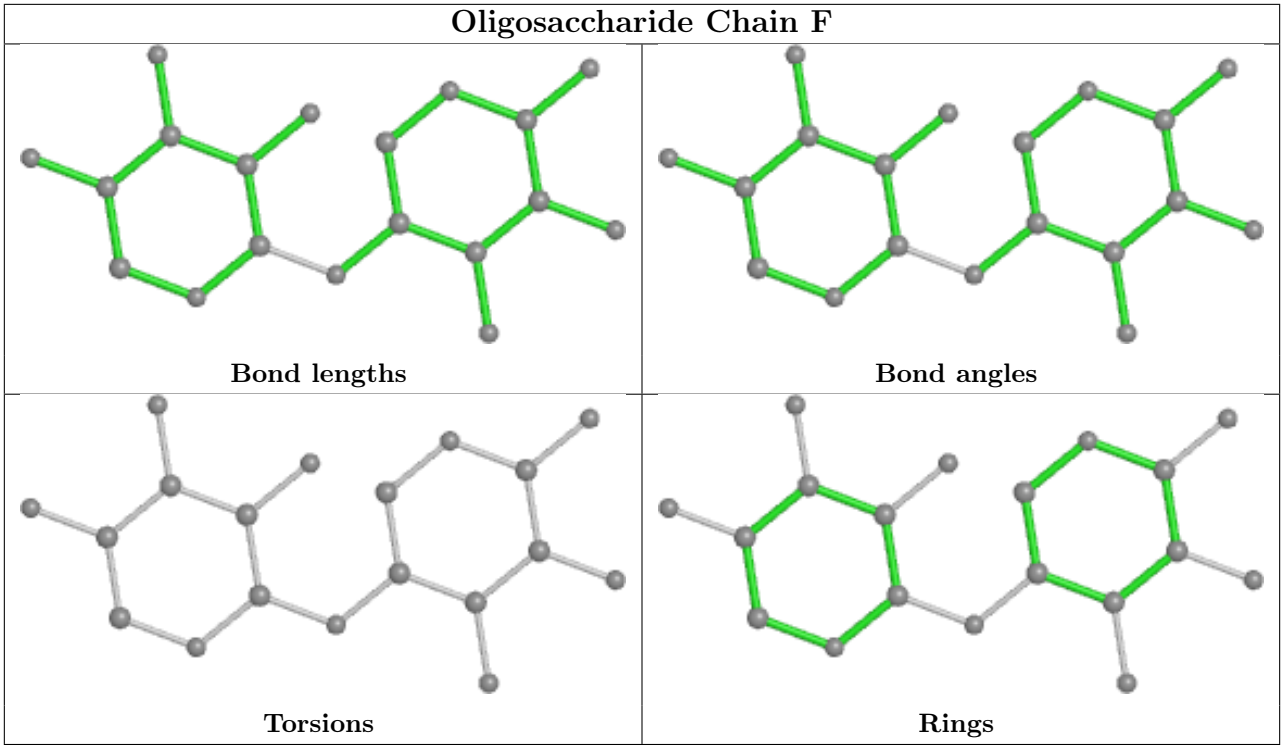
Bond angles

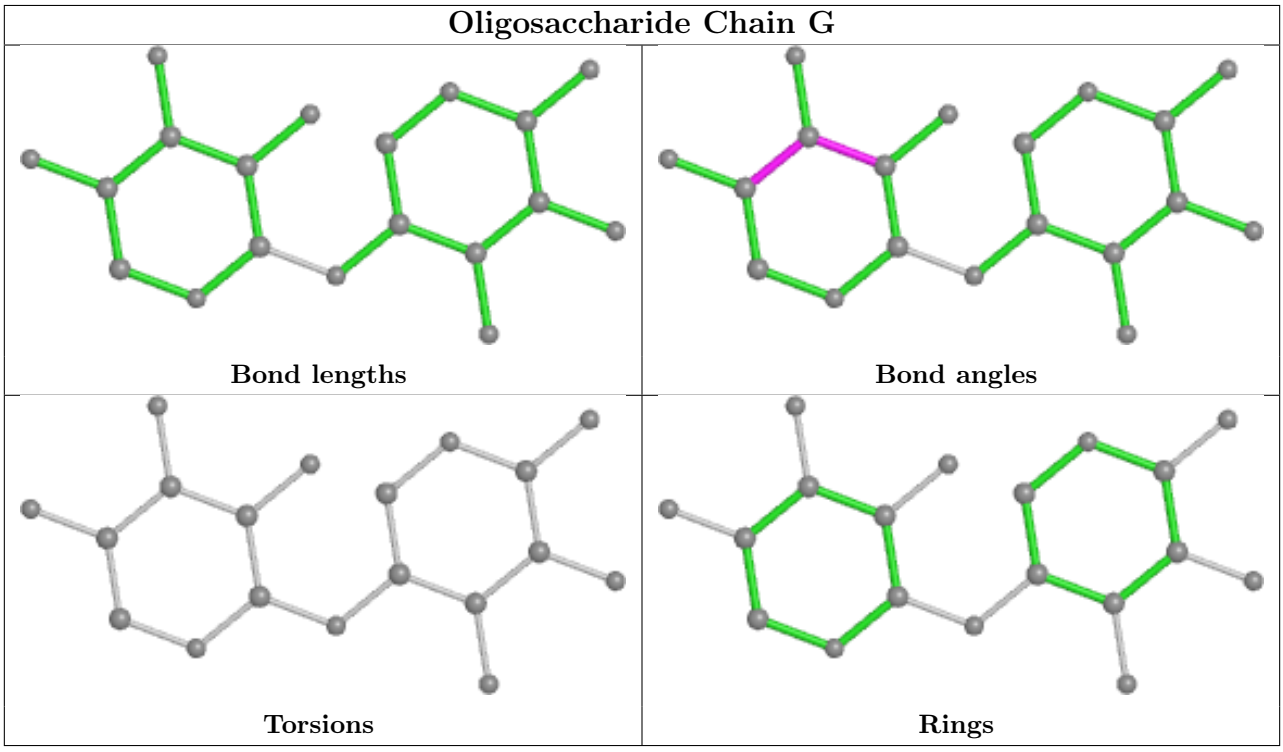


Torsions



Rings





5.6 Ligand geometry (i)

4 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	XYP	B	971	-	10,10,10	0.71	0	14,14,14	0.56	0
3	XYP	A	1461	-	10,10,10	0.58	0	14,14,14	0.56	0
3	XYP	A	1491	-	10,10,10	0.68	0	14,14,14	0.56	0
3	XYP	B	961	-	10,10,10	0.69	0	14,14,14	0.56	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	XYP	B	971	-	-	-	0/1/1/1
3	XYP	A	1461	-	-	-	0/1/1/1
3	XYP	A	1491	-	-	-	0/1/1/1
3	XYP	B	961	-	-	-	0/1/1/1

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

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, 95^{th} 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	436/436 (100%)	0.32	22 (5%) 28 34	14, 23, 43, 83	0
1	B	436/436 (100%)	0.48	45 (10%) 6 8	14, 22, 65, 77	0
All	All	872/872 (100%)	0.40	67 (7%) 13 17	14, 23, 57, 83	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	306	THR	10.8
1	A	305	SER	10.1
1	B	820	SER	7.2
1	A	304	SER	6.8
1	B	819	GLY	6.7
1	B	822	ARG	6.2
1	B	936	THR	6.1
1	B	818	VAL	6.0
1	A	303	GLY	5.8
1	A	436	THR	5.5
1	B	923	TYR	5.5
1	A	307	PRO	5.5
1	B	915	ALA	5.2
1	B	809	PRO	5.2
1	B	805	SER	4.8
1	A	309	PRO	4.8
1	A	308	PRO	4.7
1	B	928	GLY	4.6
1	B	810	SER	4.5
1	B	844	SER	4.5
1	B	806	THR	4.4
1	B	932	ARG	4.3
1	B	911	GLY	4.3
1	B	823	CYS	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	934	THR	4.1
1	B	845	ALA	4.1
1	A	100	THR	3.9
1	B	927	ASN	3.9
1	A	95	SER	3.9
1	B	843	HIS	3.8
1	A	97	SER	3.8
1	B	816	LYS	3.8
1	B	839	LEU	3.7
1	A	311	GLY	3.7
1	B	912	GLY	3.7
1	B	914	THR	3.6
1	B	929	SER	3.5
1	B	807	PRO	3.4
1	B	925	CYS	3.3
1	B	815	ILE	3.2
1	B	811	GLY	3.2
1	B	501	ALA	3.1
1	B	926	SER	3.1
1	B	918	THR	3.1
1	B	840	TYR	3.0
1	A	133	ASP	2.9
1	A	147	GLY	2.8
1	B	841	ASP	2.8
1	B	808	PRO	2.8
1	B	817	GLY	2.8
1	B	909	ALA	2.8
1	A	344	SER	2.8
1	B	874	GLY	2.7
1	B	814	GLN	2.4
1	A	91	GLY	2.4
1	B	910	VAL	2.3
1	A	145	ARG	2.3
1	A	141	SER	2.3
1	A	310	SER	2.3
1	B	916	ASN	2.3
1	A	86	HIS	2.2
1	B	930	ASN	2.2
1	B	873	ASN	2.1
1	B	913	GLY	2.1
1	A	142	ASN	2.1
1	A	96	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	812	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)

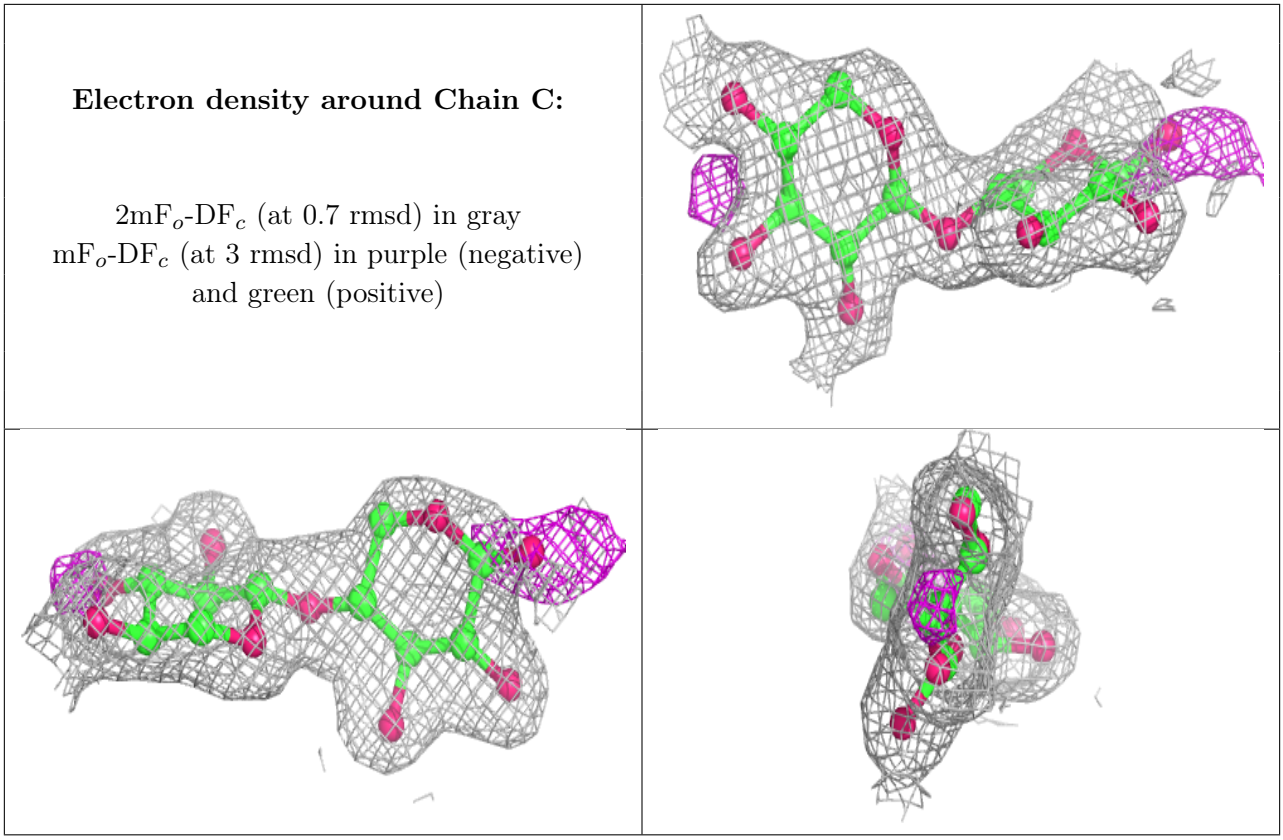
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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	XYP	D	2	9/10	0.74	0.30	60,60,62,62	0
2	XYP	G	2	9/10	0.80	0.23	48,50,51,52	0
2	XYP	E	1	10/10	0.81	0.33	37,51,52,55	0
2	XYP	G	1	10/10	0.81	0.23	42,45,46,46	0
2	XYP	D	1	10/10	0.81	0.24	57,59,60,61	0
2	XYP	C	1	10/10	0.88	0.21	26,33,36,41	0
2	XYP	E	2	9/10	0.91	0.11	23,27,29,30	0
2	XYP	F	1	10/10	0.92	0.14	16,22,27,32	0
2	XYP	C	2	9/10	0.96	0.10	22,24,26,28	0
2	XYP	F	2	9/10	0.96	0.11	16,17,19,19	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.

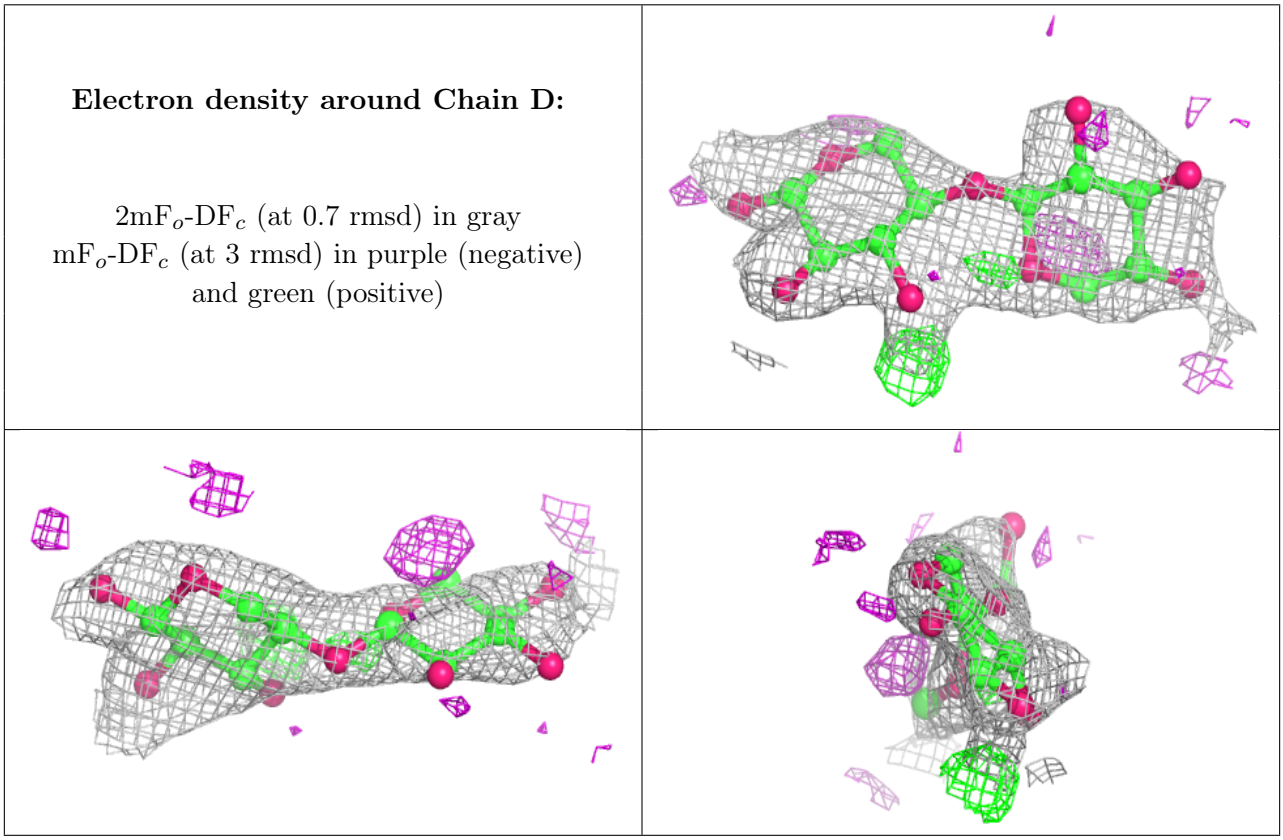
Electron density around Chain C:

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



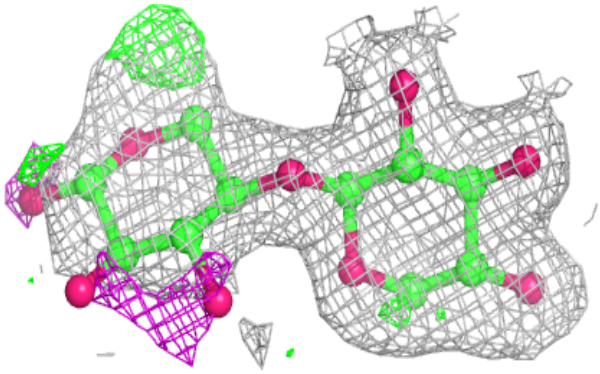
Electron density around Chain D:

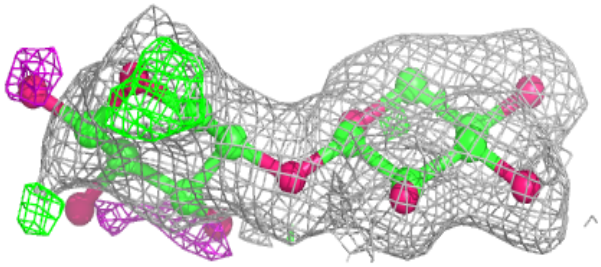
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

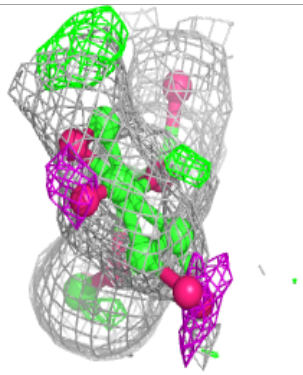


Electron density around Chain E:

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

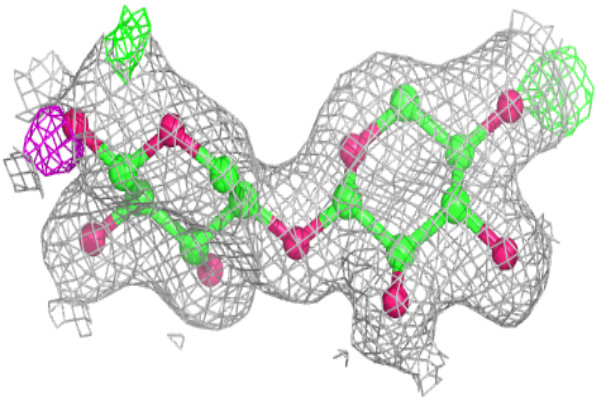


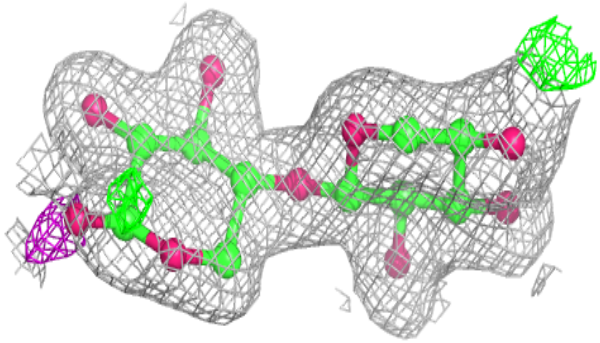


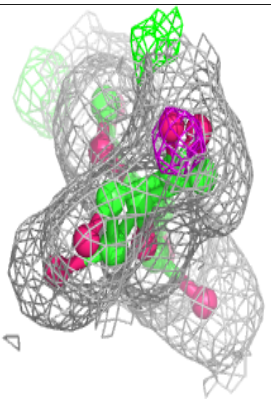


Electron density around Chain F:

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

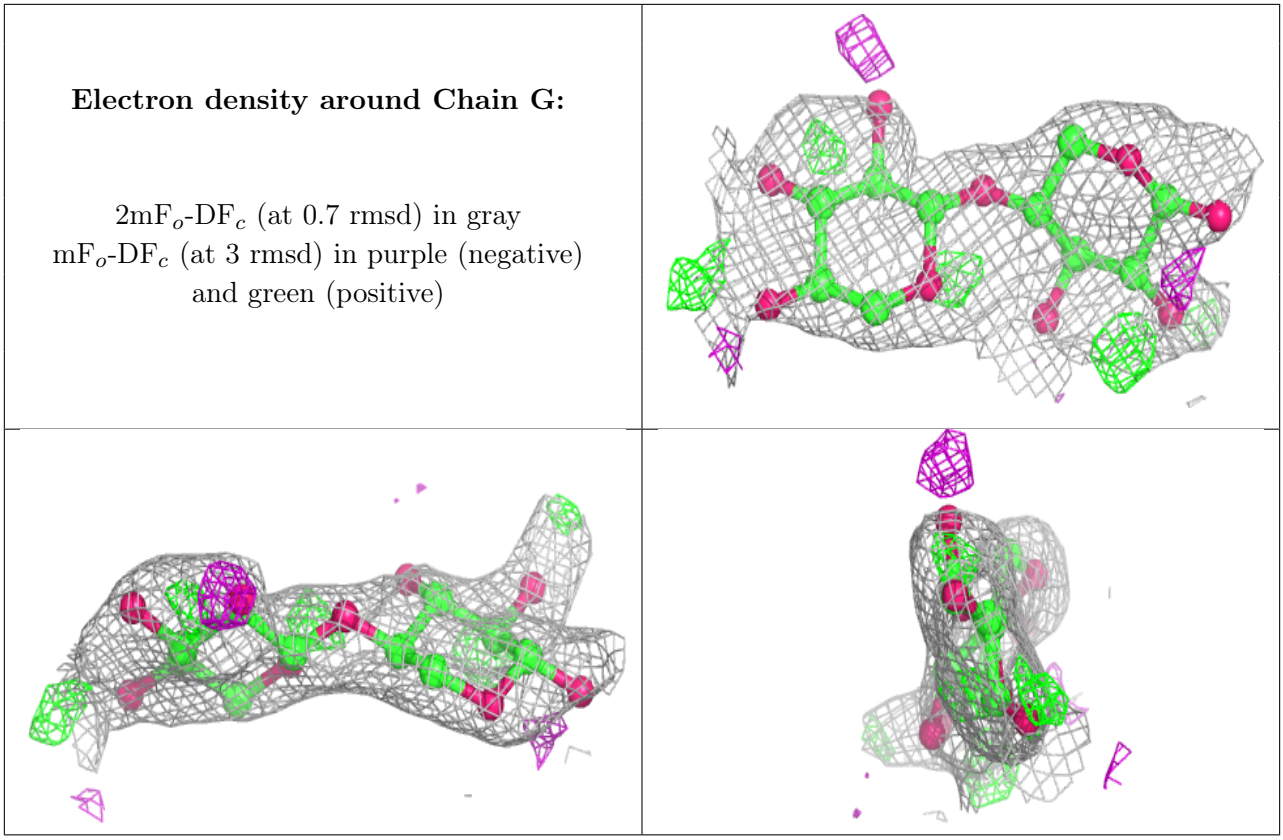






Electron density around Chain G:

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



6.4 Ligands

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, 95^{th} 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	XYP	A	1491	10/10	0.70	0.20	57,58,59,60	0
3	XYP	B	961	10/10	0.72	0.20	64,65,66,66	0
3	XYP	B	971	10/10	0.72	0.31	54,55,55,55	0
3	XYP	A	1461	10/10	0.93	0.12	28,30,31,33	0

6.5 Other polymers

There are no such residues in this entry.