Full wwPDB X-ray Structure Validation Report (i)

Dec 17, 2023 – 10:08 AM EST

PDB ID	:	1CF5
Title	:	BETA-MOMORCHARIN STRUCTURE AT 2.55 A
Authors	:	Yuan, Y.-R.; He, Y.-N.; Xiong, J.-P.; Xia, Z.-X.
Deposited on	:	1999-03-24
Resolution	:	2.55 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 (i) 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 (i)) were used in the production of this report:

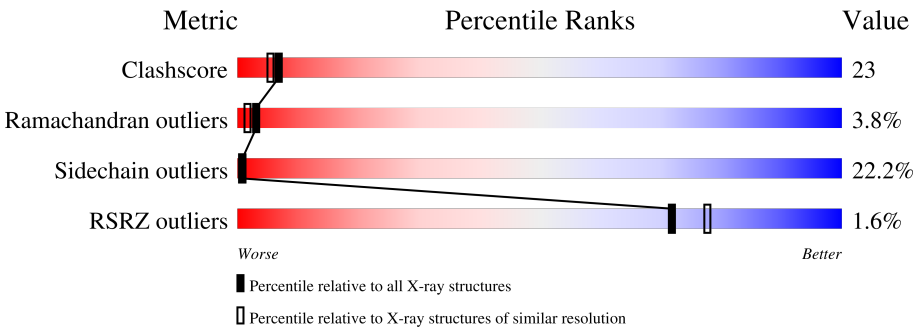
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.55 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	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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	249	2%, 46%, 40%, 12%, •
1	B	249	2%, 43%, 42%, 14%, •
2	C	6	67%, 33%
2	D	6	67%, 33%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	XYP	C	4	-	-	-	X
2	MAN	C	5	-	-	-	X
2	FUC	C	6	X	-	-	-
2	XYP	D	4	-	-	-	X
2	FUC	D	6	X	-	-	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4188 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.

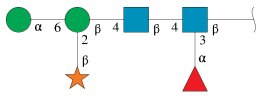
- Molecule 1 is a protein called PROTEIN (BETA-MOMORCHARIN).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	249	1991	1282	333	376	0	0	0
1	B	249	1991	1282	333	376	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	ASN	ASP	conflict	UNP P29339
A	219	GLN	GLU	conflict	UNP P29339

- Molecule 2 is an oligosaccharide called beta-D-xylopyranose-(1-2)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	6	69	39	2	28	0	0	0
2	D	6	69	39	2	28	0	0	0

- Molecule 3 is water.

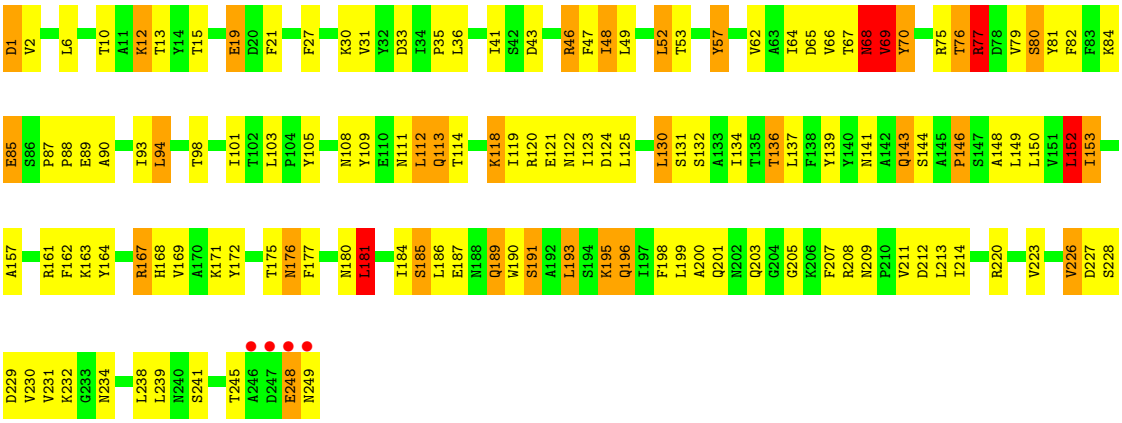
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	31	31	31	0	0
3	B	37	37	37	0	0

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.

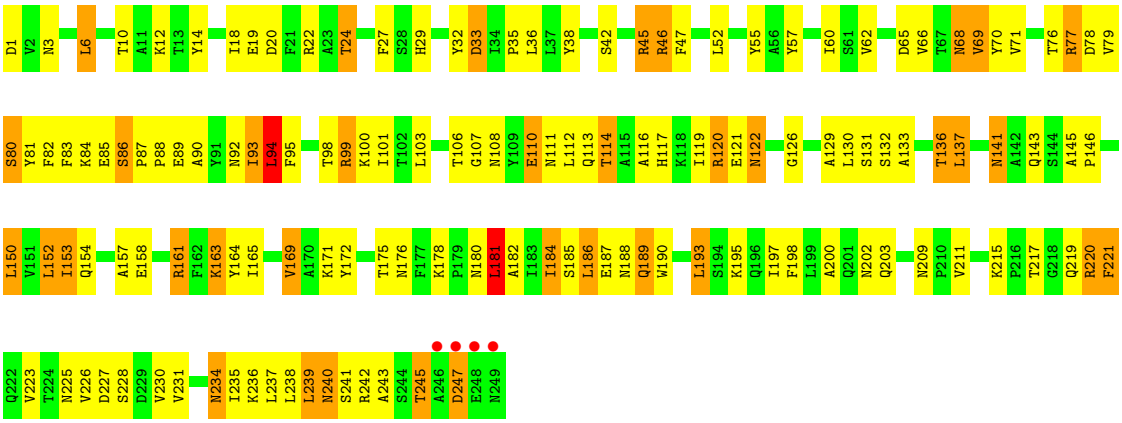
• Molecule 1: PROTEIN (BETA-MOMORCHARIN)

Chain A:

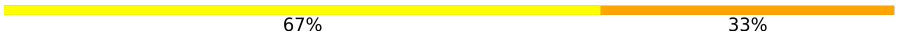


• Molecule 1: PROTEIN (BETA-MOMORCHARIN)

Chain B:



• Molecule 2: beta-D-xylopyranose-(1-2)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C: 

• Molecule 2: beta-D-xylopyranose-(1-2)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	49.09Å 50.58Å 61.12Å 72.98° 78.39° 76.97°	Depositor
Resolution (Å)	10.00 – 2.55 32.81 – 2.43	Depositor EDS
% Data completeness (in resolution range)	79.3 (10.00-2.55) 79.4 (32.81-2.43)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.82 (at 2.42Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.172 , 0.278 0.191 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	27.2	Xtriage
Anisotropy	0.669	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 94.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4188	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.98% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, XYP, MAN, BMA, FUC

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	0.61	0/2034	0.77	2/2767 (0.1%)
1	B	0.61	0/2034	0.78	2/2767 (0.1%)
All	All	0.61	0/4068	0.77	4/5534 (0.1%)

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	7
1	B	0	10
All	All	0	17

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	152	LEU	CA-CB-CG	6.64	130.58	115.30
1	A	152	LEU	CA-CB-CG	6.40	130.01	115.30
1	A	181	LEU	CA-CB-CG	6.03	129.18	115.30
1	B	181	LEU	CA-CB-CG	5.93	128.93	115.30

There are no chirality outliers.

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	TYR	Sidechain
1	A	139	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	A	168	HIS	Sidechain
1	A	176	ASN	Mainchain
1	A	198	PHE	Sidechain
1	A	70	TYR	Sidechain
1	A	77	ARG	Sidechain
1	B	120	ARG	Sidechain
1	B	14	TYR	Sidechain
1	B	161	ARG	Sidechain
1	B	220	ARG	Sidechain
1	B	38	TYR	Sidechain
1	B	55	TYR	Sidechain
1	B	70	TYR	Sidechain
1	B	77	ARG	Sidechain
1	B	80	SER	Mainchain
1	B	94	LEU	Mainchain

5.2 Too-close contacts (i)

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	1991	0	2003	96	0
1	B	1991	0	2003	95	0
2	C	69	0	51	1	0
2	D	69	0	51	1	0
3	A	31	0	0	0	0
3	B	37	0	0	1	0
All	All	4188	0	4108	191	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 23.

All (191) 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:B:226:VAL:HG23	1:B:231:VAL:HG11	1.25	1.17
1:B:189:GLN:HE21	1:B:189:GLN:HA	1.32	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:PHE:HD2	1:A:93:ILE:HD11	1.33	0.92
1:A:241:SER:O	1:A:245:THR:HG22	1.73	0.88
1:B:180:ASN:O	1:B:181:LEU:HB2	1.74	0.88
1:B:190:TRP:CZ3	1:B:239:LEU:HG	2.11	0.85
1:A:230:VAL:O	1:A:234:ASN:HB3	1.76	0.85
1:A:196:GLN:HE21	1:A:209:ASN:H	1.28	0.81
1:B:172:TYR:HD2	1:B:175:THR:HG23	1.46	0.79
1:A:48:ILE:HG23	1:A:66:VAL:HG12	1.66	0.75
1:A:47:PHE:CD2	1:A:93:ILE:HD11	2.21	0.75
1:A:172:TYR:HB3	1:A:175:THR:O	1.85	0.75
1:B:132:SER:O	1:B:136:THR:HG23	1.85	0.74
1:A:6:LEU:HG	1:A:131:SER:HB2	1.71	0.73
1:B:231:VAL:HA	1:B:235:ILE:HD12	1.69	0.73
1:B:6:LEU:HD12	1:B:131:SER:HB2	1.70	0.72
1:A:109:TYR:CE1	1:A:153:ILE:HD11	2.26	0.71
1:B:113:GLN:HE22	1:B:120:ARG:HG2	1.55	0.71
1:A:118:LYS:HG2	1:A:123:ILE:HD11	1.73	0.71
1:B:163:LYS:HB2	1:B:236:LYS:HE2	1.72	0.70
1:A:205:GLY:O	1:A:226:VAL:HG12	1.92	0.69
1:A:196:GLN:NE2	1:A:209:ASN:H	1.89	0.69
1:B:80:SER:HB3	1:B:98:THR:HG21	1.74	0.68
1:B:215:LYS:HD2	1:B:219:GLN:HB2	1.74	0.68
1:A:2:VAL:HG11	1:A:21:PHE:HB2	1.76	0.66
1:A:193:LEU:HD13	1:A:211:VAL:HG11	1.77	0.66
1:B:80:SER:HG	1:B:82:PHE:HE1	1.43	0.66
1:A:149:LEU:O	1:A:153:ILE:HG23	1.96	0.66
1:A:48:ILE:HG12	1:A:64:ILE:HB	1.77	0.66
1:A:190:TRP:O	1:A:191:SER:HB2	1.96	0.66
1:B:126:GLY:HA3	1:B:176:ASN:HD22	1.61	0.66
1:A:109:TYR:O	1:A:113:GLN:HB2	1.95	0.65
1:B:242:ARG:O	1:B:245:THR:HG22	1.96	0.65
1:B:20:ASP:O	1:B:24:THR:HG23	1.96	0.65
1:B:165:ILE:O	1:B:169:VAL:HG12	1.97	0.64
1:A:48:ILE:HG23	1:A:66:VAL:CG1	2.26	0.64
1:A:75:ARG:HD2	1:A:98:THR:HG21	1.79	0.64
1:B:3:ASN:HD21	2:D:1:NAG:H2	1.63	0.63
1:A:113:GLN:HG3	1:A:119:ILE:HA	1.79	0.63
1:B:164:TYR:H	1:B:234:ASN:ND2	1.96	0.63
1:A:113:GLN:HG2	1:A:119:ILE:HG22	1.81	0.63
1:B:113:GLN:NE2	1:B:120:ARG:HG2	2.13	0.63
1:A:113:GLN:NE2	1:A:120:ARG:H	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:SER:OG	1:B:82:PHE:HE1	1.81	0.63
1:B:89:GLU:O	1:B:93:ILE:HG23	1.97	0.62
1:B:193:LEU:HD13	1:B:211:VAL:HG11	1.81	0.62
1:A:132:SER:O	1:A:136:THR:HG23	2.00	0.62
1:B:230:VAL:HG23	1:B:231:VAL:H	1.63	0.62
1:A:164:TYR:H	1:A:234:ASN:ND2	1.97	0.61
1:A:185:SER:O	1:A:189:GLN:HB2	2.00	0.61
1:A:87:PRO:HG2	1:A:90:ALA:HB2	1.82	0.61
1:A:214:ILE:HG13	1:A:220:ARG:HG3	1.83	0.61
1:B:189:GLN:HA	1:B:189:GLN:NE2	2.11	0.61
1:A:65:ASP:HB3	1:A:68:ASN:HD21	1.66	0.60
1:A:162:PHE:CE2	1:A:190:TRP:HB2	2.36	0.60
1:A:6:LEU:HD21	1:A:130:LEU:HD13	1.84	0.60
1:B:87:PRO:HG2	1:B:90:ALA:HB2	1.84	0.60
1:A:181:LEU:HD13	1:A:214:ILE:HG21	1.82	0.60
1:A:90:ALA:O	1:A:94:LEU:HB2	2.03	0.59
1:B:225:ASN:HB3	1:B:227:ASP:H	1.67	0.59
1:A:161:ARG:HG2	1:A:190:TRP:CE2	2.39	0.58
1:B:141:ASN:ND2	1:B:143:GLN:H	2.02	0.57
1:A:119:ILE:HD12	1:A:121:GLU:HB2	1.86	0.57
1:B:164:TYR:H	1:B:234:ASN:HD21	1.50	0.57
1:B:217:THR:OG1	1:B:219:GLN:HG3	2.05	0.57
1:A:46:ARG:HH11	1:A:46:ARG:HG3	1.70	0.57
1:A:113:GLN:HE22	1:A:120:ARG:HD3	1.69	0.57
1:B:95:PHE:O	1:B:98:THR:HG22	2.05	0.57
1:A:27:PHE:HB2	1:A:36:LEU:O	2.04	0.56
1:A:141:ASN:HB3	1:A:144:SER:OG	2.05	0.56
1:A:113:GLN:CG	1:A:119:ILE:HA	2.35	0.56
1:B:113:GLN:HE22	1:B:120:ARG:HH11	1.54	0.56
1:B:157:ALA:O	1:B:161:ARG:HD3	2.05	0.55
1:B:226:VAL:HA	1:B:231:VAL:HG21	1.89	0.55
1:B:10:THR:OG1	1:B:12:LYS:HB3	2.07	0.54
1:B:120:ARG:NH2	1:B:188:ASN:OD1	2.40	0.54
1:A:109:TYR:CZ	1:A:153:ILE:HD11	2.42	0.54
1:B:6:LEU:CD1	1:B:131:SER:HB2	2.39	0.53
1:B:80:SER:CB	1:B:98:THR:HG21	2.38	0.53
1:A:69:VAL:HB	1:A:161:ARG:HH11	1.74	0.53
1:B:18:ILE:O	1:B:22:ARG:HG3	2.09	0.53
1:B:161:ARG:HG2	1:B:190:TRP:NE1	2.23	0.53
1:A:124:ASP:HA	1:A:177:PHE:O	2.10	0.52
1:B:180:ASN:O	1:B:181:LEU:CB	2.53	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:LEU:HD12	1:B:238:LEU:HB3	1.92	0.52
1:A:84:LYS:O	1:A:85:GLU:HB2	2.10	0.51
1:B:153:ILE:O	1:B:153:ILE:HG13	2.07	0.51
1:B:90:ALA:HA	1:B:94:LEU:HD22	1.93	0.51
1:B:113:GLN:NE2	1:B:120:ARG:H	2.08	0.51
1:A:137:LEU:HD13	1:A:148:ALA:HB1	1.92	0.51
1:A:196:GLN:HE21	1:A:209:ASN:N	2.04	0.51
1:B:65:ASP:HB3	1:B:68:ASN:ND2	2.26	0.51
1:B:146:PRO:HD2	3:B:939:HOH:O	2.11	0.51
1:A:76:THR:CG2	1:A:81:TYR:HE1	2.24	0.51
1:B:240:ASN:ND2	1:B:243:ALA:H	2.08	0.51
1:A:27:PHE:CE2	1:A:30:LYS:HG2	2.46	0.51
1:A:57:TYR:HB3	2:C:2:NAG:H81	1.93	0.51
1:B:230:VAL:HG23	1:B:231:VAL:N	2.27	0.50
1:B:32:TYR:O	1:B:33:ASP:HB2	2.11	0.50
1:B:197:ILE:HD13	1:B:231:VAL:HG13	1.94	0.50
1:A:195:LYS:NZ	1:A:248:GLU:HA	2.26	0.50
1:A:68:ASN:ND2	1:A:70:TYR:H	2.10	0.50
1:A:195:LYS:HE3	1:A:208:ARG:NH2	2.27	0.50
1:A:6:LEU:HD21	1:A:130:LEU:CD1	2.42	0.49
1:B:195:LYS:O	1:B:198:PHE:HB3	2.12	0.49
1:A:10:THR:OG1	1:A:12:LYS:HD3	2.13	0.49
1:A:12:LYS:HG2	1:A:13:THR:N	2.28	0.49
1:B:197:ILE:O	1:B:200:ALA:HB3	2.12	0.48
1:B:221:PHE:CD1	1:B:221:PHE:N	2.79	0.48
1:A:65:ASP:OD1	1:A:67:THR:HB	2.12	0.48
1:A:125:LEU:O	1:A:176:ASN:HB3	2.13	0.48
1:B:171:LYS:HB3	1:B:171:LYS:HE2	1.59	0.48
1:A:52:LEU:HD23	1:A:152:LEU:HD21	1.95	0.48
1:B:193:LEU:CD1	1:B:211:VAL:HG11	2.44	0.48
1:A:130:LEU:O	1:A:134:ILE:HD12	2.13	0.48
1:A:157:ALA:O	1:A:161:ARG:HD3	2.13	0.48
1:B:110:GLU:O	1:B:114:THR:HG23	2.14	0.48
1:B:35:PRO:HG2	1:B:237:LEU:HD23	1.95	0.47
1:B:184:ILE:O	1:B:187:GLU:N	2.48	0.47
1:A:1:ASP:HB3	1:A:49:LEU:O	2.15	0.47
1:B:46:ARG:O	1:B:66:VAL:HG22	2.14	0.47
1:B:141:ASN:HD22	1:B:143:GLN:H	1.63	0.47
1:B:172:TYR:HD2	1:B:175:THR:CG2	2.22	0.46
1:B:83:PHE:O	1:B:86:SER:HB2	2.16	0.46
1:B:129:ALA:H	1:B:176:ASN:HD21	1.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:ALA:O	1:B:137:LEU:HB2	2.15	0.46
1:B:137:LEU:HA	1:B:137:LEU:HD12	1.61	0.46
1:A:41:ILE:HG23	1:A:46:ARG:HG2	1.98	0.46
1:B:83:PHE:CE1	1:B:107:GLY:HA2	2.51	0.46
1:B:182:ALA:O	1:B:186:LEU:HB2	2.16	0.46
1:A:15:THR:O	1:A:19:GLU:HB2	2.16	0.45
1:A:163:LYS:O	1:A:167:ARG:HG2	2.16	0.45
1:B:46:ARG:HB3	1:B:47:PHE:CD2	2.51	0.45
1:B:172:TYR:CD2	1:B:175:THR:HG23	2.38	0.45
1:B:65:ASP:HB3	1:B:68:ASN:HD21	1.81	0.45
1:B:172:TYR:HB3	1:B:175:THR:O	2.16	0.45
1:A:31:VAL:HB	1:A:238:LEU:CD2	2.47	0.45
1:A:196:GLN:OE1	1:A:249:ASN:ND2	2.50	0.45
1:A:46:ARG:HG3	1:A:46:ARG:NH1	2.32	0.45
1:B:236:LYS:O	1:B:237:LEU:HD23	2.16	0.45
1:A:27:PHE:HE2	1:A:30:LYS:HG2	1.80	0.45
1:A:79:VAL:HG11	1:A:101:ILE:HD12	1.99	0.45
1:B:120:ARG:HG2	1:B:120:ARG:HH11	1.82	0.45
1:A:6:LEU:HG	1:A:131:SER:CB	2.45	0.44
1:A:195:LYS:HZ1	1:A:248:GLU:HA	1.81	0.44
1:A:6:LEU:CG	1:A:131:SER:HB2	2.45	0.44
1:B:161:ARG:HG2	1:B:190:TRP:CE2	2.53	0.44
1:A:195:LYS:O	1:A:199:LEU:HD13	2.18	0.44
1:A:161:ARG:HG2	1:A:190:TRP:NE1	2.33	0.44
1:B:113:GLN:NE2	1:B:120:ARG:CG	2.81	0.44
1:B:113:GLN:HG3	1:B:150:LEU:HD21	2.00	0.44
1:A:199:LEU:O	1:A:201:GLN:N	2.51	0.44
1:B:154:GLN:HA	1:B:158:GLU:HB2	1.99	0.44
1:A:46:ARG:O	1:A:66:VAL:HG13	2.17	0.43
1:A:43:ASP:OD2	1:A:89:GLU:HG2	2.18	0.43
1:A:226:VAL:HA	1:A:231:VAL:HG11	2.00	0.43
1:B:84:LYS:O	1:B:106:THR:HG22	2.19	0.43
1:A:76:THR:O	1:A:77:ARG:HB2	2.17	0.43
1:B:126:GLY:HA3	1:B:176:ASN:ND2	2.30	0.43
1:A:30:LYS:HD3	1:A:35:PRO:HG3	2.00	0.43
1:A:137:LEU:HD13	1:A:148:ALA:CB	2.48	0.43
1:B:197:ILE:CD1	1:B:231:VAL:HG13	2.49	0.43
1:B:90:ALA:O	1:B:94:LEU:HB2	2.19	0.43
1:B:108:ASN:OD1	1:B:108:ASN:C	2.57	0.43
1:A:53:THR:CG2	1:A:57:TYR:HA	2.49	0.42
1:A:87:PRO:HA	1:A:88:PRO:HD2	1.69	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:VAL:HB	1:A:238:LEU:HD22	2.00	0.42
1:B:80:SER:HB3	1:B:98:THR:CG2	2.46	0.42
1:B:87:PRO:HA	1:B:88:PRO:HD3	1.93	0.42
1:B:230:VAL:HG23	1:B:231:VAL:HG23	2.01	0.42
1:A:171:LYS:HE2	1:A:171:LYS:HB3	1.83	0.42
1:B:116:ALA:O	1:B:117:HIS:HB2	2.20	0.42
1:B:122:ASN:HD22	1:B:178:LYS:NZ	2.18	0.42
1:A:79:VAL:HG12	1:A:80:SER:N	2.34	0.41
1:A:163:LYS:NZ	1:A:167:ARG:NH1	2.68	0.41
1:A:180:ASN:O	1:A:181:LEU:HG	2.20	0.41
1:B:27:PHE:HD2	1:B:29:HIS:O	2.03	0.41
1:B:42:SER:HB3	1:B:45:ARG:NH1	2.35	0.41
1:B:76:THR:O	1:B:78:ASP:N	2.54	0.41
1:B:79:VAL:HG12	1:B:80:SER:N	2.36	0.41
1:A:213:LEU:C	1:A:214:ILE:HD12	2.41	0.41
1:A:80:SER:HG	1:A:82:PHE:HE1	1.67	0.41
1:A:164:TYR:H	1:A:234:ASN:HD21	1.67	0.41
1:B:81:TYR:HA	1:B:101:ILE:O	2.21	0.41
1:A:189:GLN:HG2	1:A:211:VAL:CG1	2.51	0.41
1:B:60:ILE:O	1:B:60:ILE:HG13	2.21	0.41
1:A:48:ILE:CG2	1:A:66:VAL:HG12	2.44	0.40
1:A:184:ILE:O	1:A:187:GLU:HB2	2.21	0.40
1:A:196:GLN:HG2	1:A:207:PHE:HB3	2.03	0.40
1:B:145:ALA:O	1:B:146:PRO:C	2.59	0.40
1:A:143:GLN:O	1:A:146:PRO:HD2	2.21	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	247/249 (99%)	213 (86%)	24 (10%)	10 (4%)	3 1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	247/249 (99%)	219 (89%)	19 (8%)	9 (4%)	3 2
All	All	494/498 (99%)	432 (87%)	43 (9%)	19 (4%)	3 2

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	85	GLU
1	A	191	SER
1	B	77	ARG
1	B	228	SER
1	A	68	ASN
1	A	112	LEU
1	A	181	LEU
1	B	181	LEU
1	B	185	SER
1	B	99	ARG
1	B	184	ILE
1	A	57	TYR
1	A	228	SER
1	B	247	ASP
1	A	200	ALA
1	B	57	TYR
1	A	77	ARG
1	A	69	VAL
1	B	69	VAL

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	218/218 (100%)	172 (79%)	46 (21%)	1 1
1	B	218/218 (100%)	167 (77%)	51 (23%)	1 0
All	All	436/436 (100%)	339 (78%)	97 (22%)	1 1

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ASP
1	A	12	LYS
1	A	19	GLU
1	A	33	ASP
1	A	46	ARG
1	A	48	ILE
1	A	52	LEU
1	A	62	VAL
1	A	68	ASN
1	A	69	VAL
1	A	76	THR
1	A	77	ARG
1	A	80	SER
1	A	94	LEU
1	A	103	LEU
1	A	108	ASN
1	A	111	ASN
1	A	112	LEU
1	A	113	GLN
1	A	114	THR
1	A	118	LYS
1	A	122	ASN
1	A	130	LEU
1	A	136	THR
1	A	143	GLN
1	A	146	PRO
1	A	150	LEU
1	A	152	LEU
1	A	153	ILE
1	A	167	ARG
1	A	169	VAL
1	A	185	SER
1	A	186	LEU
1	A	189	GLN
1	A	193	LEU
1	A	195	LYS
1	A	196	GLN
1	A	203	GLN
1	A	212	ASP
1	A	223	VAL
1	A	226	VAL
1	A	227	ASP
1	A	229	ASP

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Mol	Chain	Res	Type
1	A	232	LYS
1	A	239	LEU
1	A	248	GLU
1	B	1	ASP
1	B	6	LEU
1	B	19	GLU
1	B	24	THR
1	B	33	ASP
1	B	45	ARG
1	B	46	ARG
1	B	52	LEU
1	B	62	VAL
1	B	68	ASN
1	B	69	VAL
1	B	71	VAL
1	B	85	GLU
1	B	86	SER
1	B	92	ASN
1	B	93	ILE
1	B	94	LEU
1	B	99	ARG
1	B	100	LYS
1	B	103	LEU
1	B	110	GLU
1	B	111	ASN
1	B	112	LEU
1	B	114	THR
1	B	119	ILE
1	B	121	GLU
1	B	122	ASN
1	B	130	LEU
1	B	136	THR
1	B	137	LEU
1	B	141	ASN
1	B	150	LEU
1	B	152	LEU
1	B	153	ILE
1	B	163	LYS
1	B	169	VAL
1	B	186	LEU
1	B	189	GLN
1	B	193	LEU

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Mol	Chain	Res	Type
1	B	202	ASN
1	B	203	GLN
1	B	209	ASN
1	B	220	ARG
1	B	221	PHE
1	B	223	VAL
1	B	234	ASN
1	B	239	LEU
1	B	240	ASN
1	B	241	SER
1	B	245	THR
1	B	247	ASP

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	68	ASN
1	A	113	GLN
1	A	122	ASN
1	A	143	GLN
1	A	154	GLN
1	A	196	GLN
1	A	234	ASN
1	A	249	ASN
1	B	3	ASN
1	B	29	HIS
1	B	68	ASN
1	B	113	GLN
1	B	117	HIS
1	B	122	ASN
1	B	141	ASN
1	B	176	ASN
1	B	189	GLN
1	B	202	ASN
1	B	209	ASN
1	B	222	GLN
1	B	234	ASN
1	B	240	ASN

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)

12 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	NAG	C	1	1,2	14,14,15	1.28	2 (14%)	17,19,21	2.00	7 (41%)
2	NAG	C	2	2	14,14,15	1.80	3 (21%)	17,19,21	3.15	10 (58%)
2	BMA	C	3	2	11,11,12	2.94	5 (45%)	15,15,17	1.74	4 (26%)
2	XYP	C	4	2	9,9,10	3.12	6 (66%)	10,12,14	1.90	3 (30%)
2	MAN	C	5	2	11,11,12	2.95	6 (54%)	15,15,17	1.50	2 (13%)
2	FUC	C	6	2	10,10,11	1.59	1 (10%)	14,14,16	1.26	2 (14%)
2	NAG	D	1	1,2	14,14,15	1.96	6 (42%)	17,19,21	3.21	9 (52%)
2	NAG	D	2	2	14,14,15	2.78	4 (28%)	17,19,21	1.64	4 (23%)
2	BMA	D	3	2	11,11,12	2.90	6 (54%)	15,15,17	3.03	7 (46%)
2	XYP	D	4	2	9,9,10	3.10	2 (22%)	10,12,14	1.59	2 (20%)
2	MAN	D	5	2	11,11,12	3.30	3 (27%)	15,15,17	2.33	6 (40%)
2	FUC	D	6	2	10,10,11	1.52	2 (20%)	14,14,16	1.18	1 (7%)

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	NAG	C	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	2/2/19/22	0/1/1/1
2	XYP	C	4	2	-	-	0/1/1/1
2	MAN	C	5	2	-	1/2/19/22	1/1/1/1
2	FUC	C	6	2	4/4/4/5	-	0/1/1/1
2	NAG	D	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	D	2	2	-	3/6/23/26	0/1/1/1
2	BMA	D	3	2	-	2/2/19/22	0/1/1/1
2	XYP	D	4	2	-	-	0/1/1/1
2	MAN	D	5	2	-	2/2/19/22	0/1/1/1
2	FUC	D	6	2	4/4/4/5	-	0/1/1/1

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	5	MAN	C2-C3	7.41	1.63	1.52
2	D	5	MAN	C2-C3	7.31	1.63	1.52
2	D	4	XYP	C4-C3	7.09	1.63	1.52
2	C	4	XYP	C2-C3	6.57	1.62	1.52
2	D	2	NAG	C1-C2	6.40	1.61	1.52
2	D	3	BMA	C2-C3	6.32	1.61	1.52
2	C	3	BMA	C4-C5	6.30	1.66	1.53
2	D	5	MAN	C1-C2	5.51	1.64	1.52
2	D	5	MAN	C4-C3	5.09	1.65	1.52
2	D	2	NAG	C3-C2	4.89	1.62	1.52
2	C	3	BMA	C4-C3	4.86	1.64	1.52
2	D	4	XYP	C2-C3	4.55	1.59	1.52
2	C	2	NAG	C4-C3	4.40	1.63	1.52
2	D	2	NAG	O5-C1	-4.29	1.36	1.43
2	C	4	XYP	C4-C3	4.02	1.58	1.52
2	C	6	FUC	C4-C5	4.00	1.61	1.52
2	D	3	BMA	C1-C2	3.97	1.61	1.52
2	D	1	NAG	C4-C5	3.94	1.61	1.53
2	D	3	BMA	O2-C2	3.89	1.51	1.43
2	D	2	NAG	C4-C3	3.75	1.61	1.52
2	C	2	NAG	C2-N2	3.64	1.52	1.46
2	C	3	BMA	O2-C2	3.41	1.50	1.43
2	C	5	MAN	O5-C5	3.14	1.49	1.43
2	C	5	MAN	O5-C1	3.07	1.48	1.43
2	D	1	NAG	C1-C2	3.04	1.56	1.52
2	D	1	NAG	O3-C3	2.89	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	XYP	C1-C2	2.88	1.58	1.52
2	C	1	NAG	C3-C2	2.78	1.58	1.52
2	C	4	XYP	C5-C4	2.74	1.58	1.52
2	C	5	MAN	C1-C2	2.74	1.58	1.52
2	C	2	NAG	C3-C2	2.70	1.58	1.52
2	D	3	BMA	C6-C5	2.69	1.60	1.51
2	D	6	FUC	C1-C2	2.64	1.58	1.52
2	C	3	BMA	C6-C5	2.57	1.60	1.51
2	D	3	BMA	C4-C5	2.57	1.58	1.53
2	C	3	BMA	C2-C3	2.53	1.56	1.52
2	D	1	NAG	O5-C5	2.43	1.48	1.43
2	C	1	NAG	C4-C3	2.33	1.58	1.52
2	D	3	BMA	C4-C3	2.32	1.58	1.52
2	C	5	MAN	C4-C3	2.31	1.58	1.52
2	D	1	NAG	C4-C3	2.18	1.57	1.52
2	C	5	MAN	C4-C5	2.14	1.57	1.53
2	C	4	XYP	O5-C5	2.13	1.47	1.42
2	C	4	XYP	O5-C1	2.10	1.47	1.42
2	D	6	FUC	O5-C1	2.09	1.47	1.43
2	D	1	NAG	C6-C5	2.04	1.58	1.51

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	1	NAG	C1-O5-C5	8.38	123.54	112.19
2	D	3	BMA	C1-C2-C3	-7.67	100.24	109.67
2	D	1	NAG	C2-N2-C7	6.74	132.49	122.90
2	C	2	NAG	O5-C1-C2	5.88	120.58	111.29
2	D	3	BMA	C1-O5-C5	5.87	120.14	112.19
2	D	5	MAN	C6-C5-C4	5.76	126.50	113.00
2	C	2	NAG	C2-N2-C7	4.79	129.73	122.90
2	C	2	NAG	C3-C4-C5	4.62	118.49	110.24
2	C	4	XYP	C1-C2-C3	4.46	115.15	109.67
2	C	2	NAG	O5-C5-C6	4.35	114.02	107.20
2	C	1	NAG	C1-C2-N2	4.29	117.81	110.49
2	C	5	MAN	C1-O5-C5	3.98	117.59	112.19
2	D	5	MAN	C1-C2-C3	3.94	114.51	109.67
2	C	2	NAG	O4-C4-C5	-3.78	99.91	109.30
2	C	2	NAG	C4-C3-C2	3.76	116.53	111.02
2	D	1	NAG	C1-C2-N2	3.56	116.56	110.49
2	C	2	NAG	O7-C7-C8	-3.49	115.58	122.06
2	D	2	NAG	C3-C4-C5	3.43	116.36	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	NAG	C2-N2-C7	3.38	127.71	122.90
2	C	3	BMA	C1-C2-C3	-3.37	105.53	109.67
2	C	2	NAG	O5-C5-C4	-3.33	102.73	110.83
2	D	2	NAG	O4-C4-C5	-3.22	101.30	109.30
2	D	3	BMA	O2-C2-C1	3.07	115.43	109.15
2	D	3	BMA	O2-C2-C3	3.02	116.19	110.14
2	D	3	BMA	C3-C4-C5	2.97	115.53	110.24
2	D	4	XYP	O3-C3-C2	-2.83	104.57	109.99
2	D	1	NAG	C4-C3-C2	-2.82	106.88	111.02
2	D	1	NAG	C8-C7-N2	-2.82	111.32	116.10
2	C	3	BMA	O2-C2-C1	2.82	114.92	109.15
2	D	5	MAN	C3-C4-C5	-2.78	105.27	110.24
2	D	5	MAN	O4-C4-C3	2.67	116.52	110.35
2	C	2	NAG	C8-C7-N2	2.64	120.57	116.10
2	D	6	FUC	O5-C1-C2	2.63	114.83	110.77
2	C	1	NAG	O7-C7-C8	-2.62	117.20	122.06
2	C	1	NAG	C1-O5-C5	2.54	115.64	112.19
2	C	5	MAN	O3-C3-C2	2.49	114.75	109.99
2	D	1	NAG	C3-C4-C5	2.42	114.55	110.24
2	D	2	NAG	C2-N2-C7	2.40	126.33	122.90
2	C	1	NAG	C3-C4-C5	2.38	114.48	110.24
2	C	3	BMA	C6-C5-C4	2.37	118.56	113.00
2	C	3	BMA	C3-C4-C5	2.36	114.45	110.24
2	D	2	NAG	O4-C4-C3	2.32	115.72	110.35
2	D	5	MAN	O2-C2-C3	-2.29	105.55	110.14
2	D	1	NAG	O4-C4-C3	-2.27	105.10	110.35
2	D	4	XYP	C5-C4-C3	2.24	112.42	109.67
2	C	4	XYP	C5-O5-C1	2.21	114.92	111.52
2	D	5	MAN	O5-C5-C4	-2.19	105.49	110.83
2	C	2	NAG	C6-C5-C4	2.16	118.06	113.00
2	C	6	FUC	O5-C5-C4	2.15	113.37	109.52
2	D	1	NAG	O6-C6-C5	2.14	118.63	111.29
2	D	3	BMA	O5-C5-C6	-2.13	103.87	107.20
2	D	1	NAG	O5-C1-C2	-2.09	107.98	111.29
2	C	6	FUC	O4-C4-C5	2.05	114.21	109.67
2	D	3	BMA	O3-C3-C2	2.05	113.92	109.99
2	C	1	NAG	O4-C4-C5	-2.02	104.28	109.30
2	C	1	NAG	C8-C7-N2	2.02	119.51	116.10
2	C	4	XYP	O2-C2-C3	-2.01	106.12	110.14

All (8) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	6	FUC	C2
2	C	6	FUC	C4
2	C	6	FUC	C3
2	C	6	FUC	C1
2	D	6	FUC	C2
2	D	6	FUC	C4
2	D	6	FUC	C3
2	D	6	FUC	C1

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	NAG	C3-C2-N2-C7
2	C	3	BMA	O5-C5-C6-O6
2	D	3	BMA	C4-C5-C6-O6
2	D	5	MAN	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	D	3	BMA	O5-C5-C6-O6
2	C	1	NAG	C1-C2-N2-C7
2	C	3	BMA	C4-C5-C6-O6
2	D	5	MAN	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C1-C2-N2-C7
2	D	2	NAG	C4-C5-C6-O6
2	C	5	MAN	O5-C5-C6-O6

All (1) ring outliers are listed below:

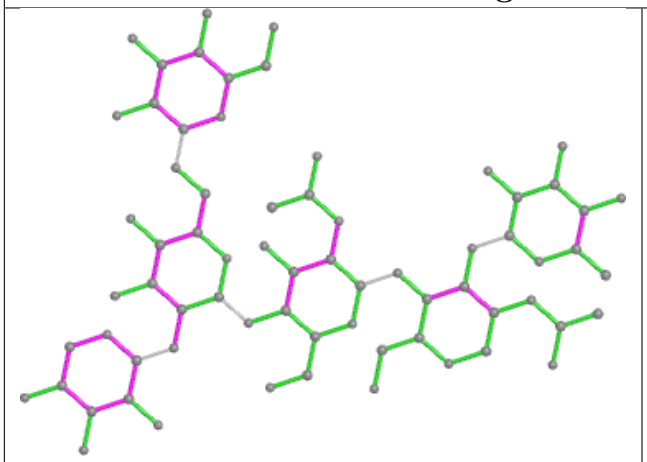
Mol	Chain	Res	Type	Atoms
2	C	5	MAN	C1-C2-C3-C4-C5-O5

2 monomers are involved in 2 short contacts:

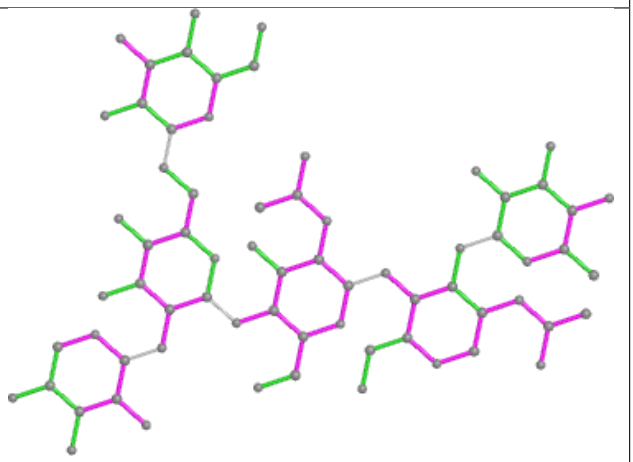
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	NAG	1	0
2	C	2	NAG	1	0

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

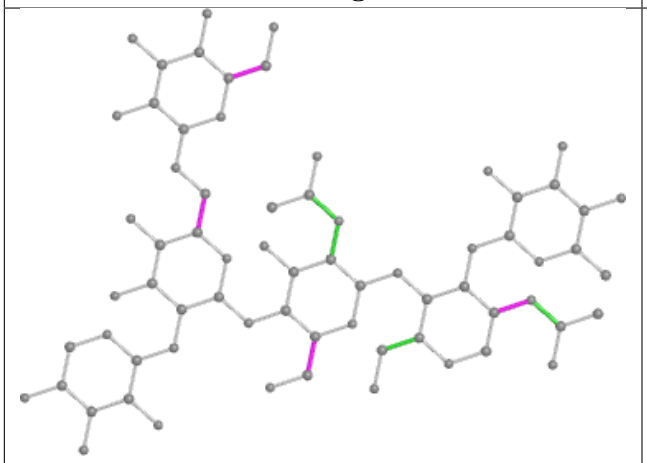
Oligosaccharide Chain C



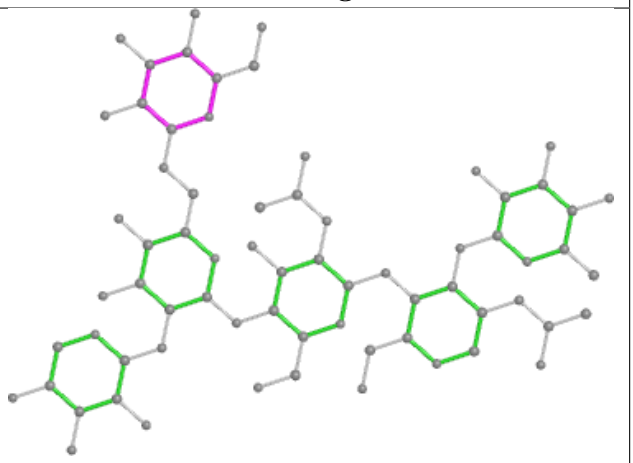
Bond lengths



Bond angles

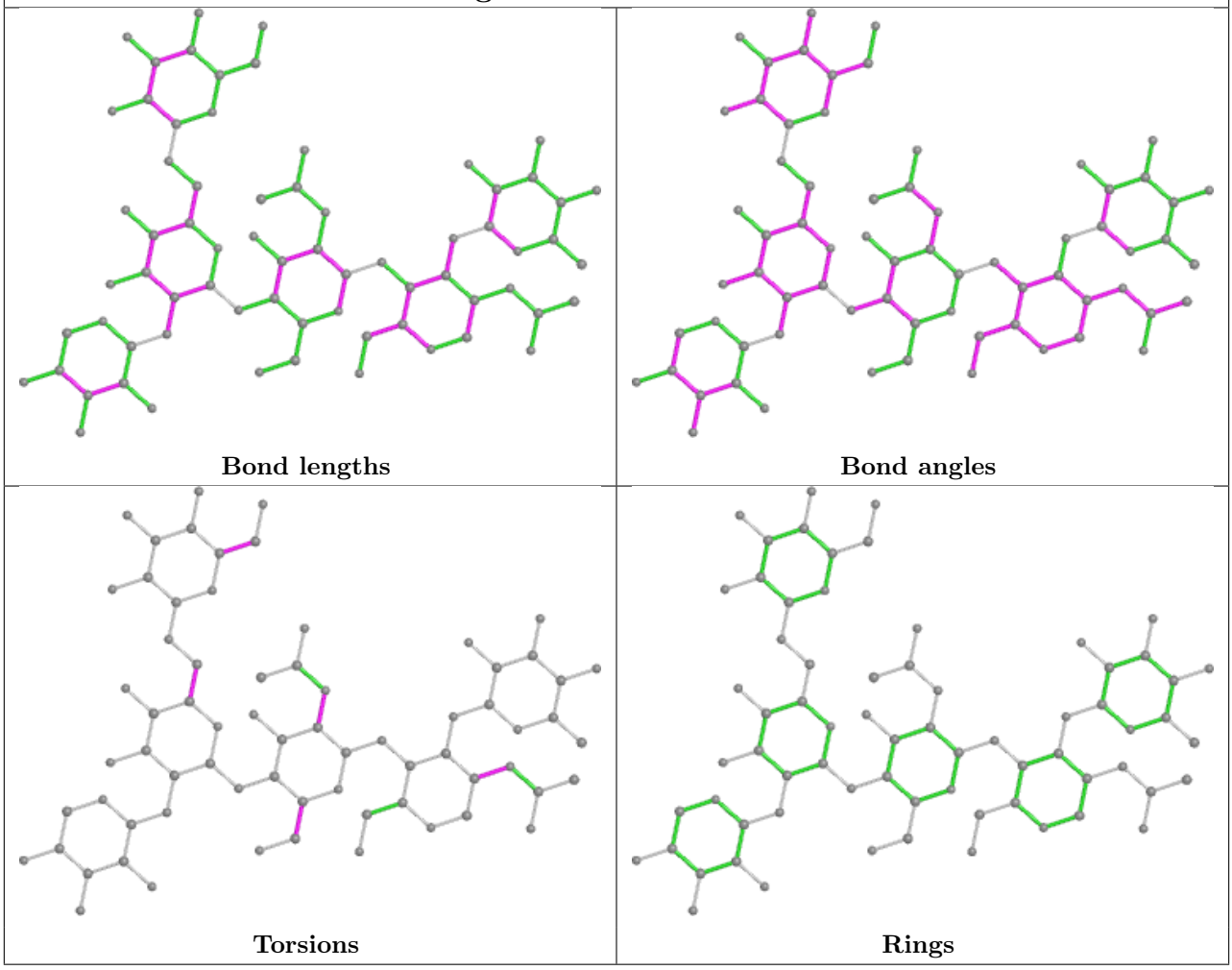


Torsions



Rings

Oligosaccharide Chain D



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 (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	249/249 (100%)	-0.21	4 (1%)	72	78	3, 17, 49, 96	0
1	B	249/249 (100%)	-0.36	4 (1%)	72	78	2, 16, 51, 100	0
All	All	498/498 (100%)	-0.29	8 (1%)	72	78	2, 16, 50, 100	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	249	ASN	9.3
1	A	247	ASP	7.6
1	A	248	GLU	7.2
1	B	248	GLU	6.3
1	B	247	ASP	4.5
1	B	249	ASN	4.4
1	B	246	ALA	3.4
1	A	246	ALA	2.5

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	MAN	C	5	11/12	0.51	0.41	97,101,103,103	0

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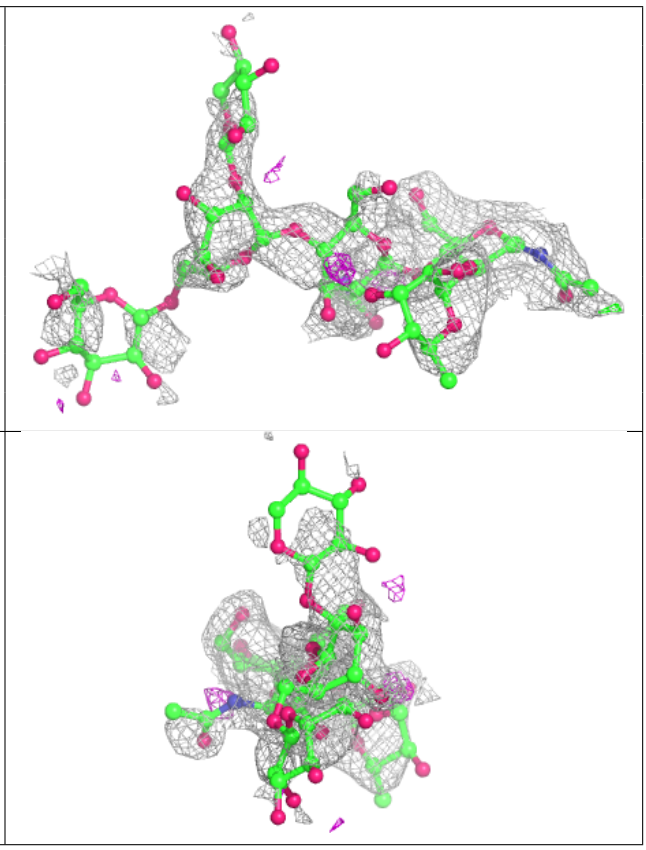
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	XYP	D	4	9/10	0.57	0.45	89,94,96,98	0
2	XYP	C	4	9/10	0.59	0.50	103,104,106,107	0
2	BMA	D	3	11/12	0.64	0.26	95,98,101,102	0
2	FUC	C	6	10/11	0.69	0.38	85,92,96,98	0
2	BMA	C	3	11/12	0.71	0.28	94,99,102,103	0
2	NAG	C	2	14/15	0.71	0.28	86,93,95,96	0
2	NAG	D	2	14/15	0.76	0.24	75,79,87,90	0
2	MAN	D	5	11/12	0.76	0.32	93,98,100,101	0
2	FUC	D	6	10/11	0.76	0.21	86,89,93,94	0
2	NAG	D	1	14/15	0.77	0.24	54,62,72,79	0
2	NAG	C	1	14/15	0.81	0.17	57,70,78,79	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.

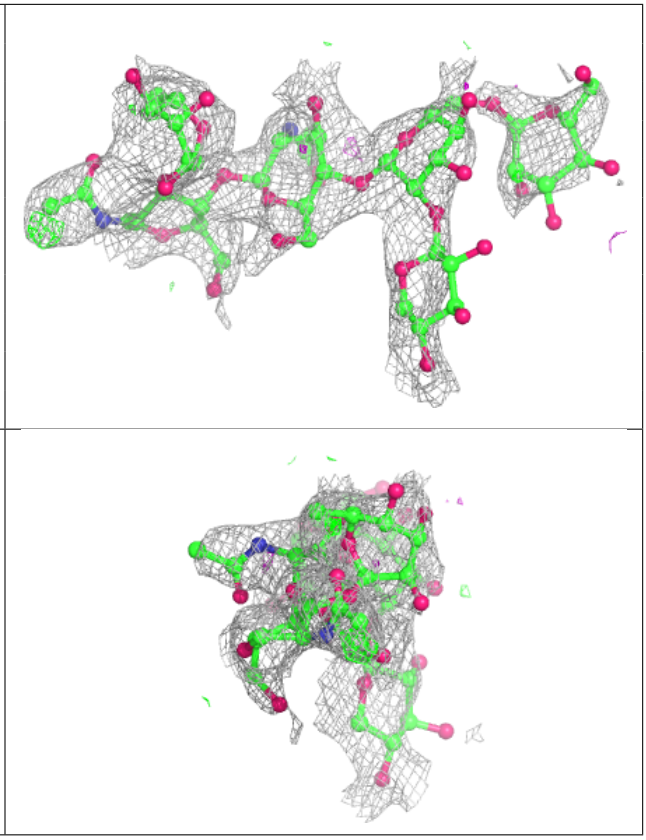
Electron density around Chain C:

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.