



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:19 PM UTC

PDB ID : 6B9O / pdb_00006b9o
Title : Structure of GH 38 Jack Bean alpha-mannosidase
Authors : Howard, E.; Cousido-Siah, A.; Lepage, M.; Bodlenner, A.; Mitschler, A.; Meli, A.; De Riccardis, F.; Izzo, I.; Podjarny, A.; Compain, P.
Deposited on : 2017-10-11
Resolution : 1.84 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

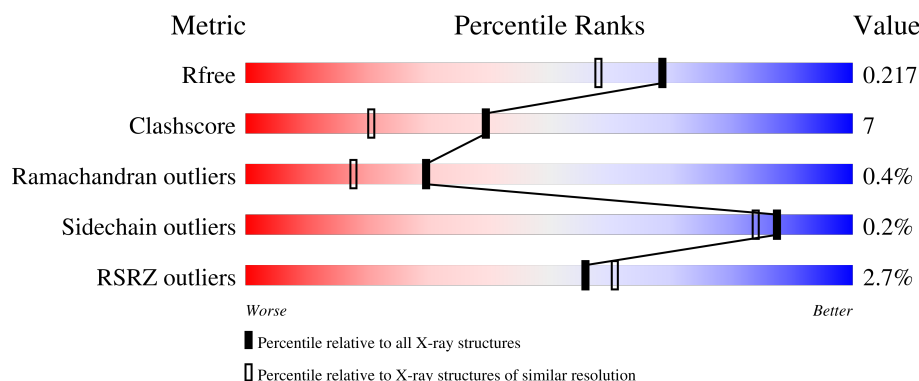
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	981	3% 83% 12% 5%
1	B	981	3% 84% 12% .
2	C	12	17% 50% 33%
2	E	12	8% 75% 17%
3	D	2	50% 50%

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

2 Entry composition [i](#)

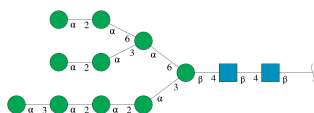
There are 5 unique types of molecules in this entry. The entry contains 17299 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-mannosidase from *Canavalia ensiformis* (jack bean).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	935	Total	C	N	O	S	0	10	0
			7561	4807	1288	1435	31			
1	B	937	Total	C	N	O	S	0	6	0
			7561	4805	1288	1437	31			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	0	1	0
			149	82	2	65			
2	E	12	Total	C	N	O	0	0	0
			138	76	2	60			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	930	Total	O	0	3
			933	933		
5	B	898	Total	O	0	1
			899	899		

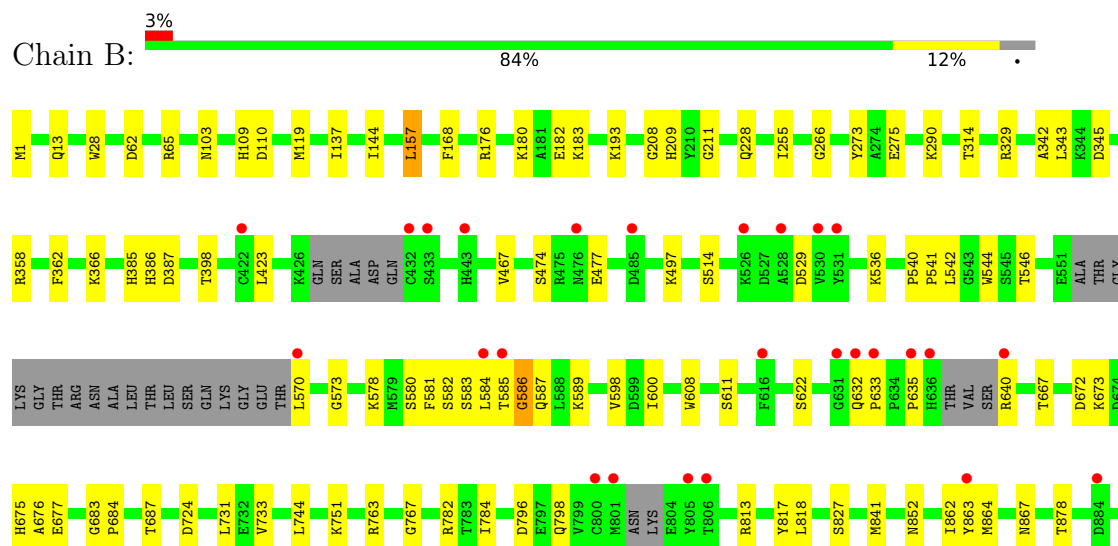
3 Residue-property plots

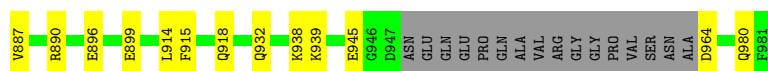
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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-mannosidase from *Canavalia ensiformis* (jack bean)

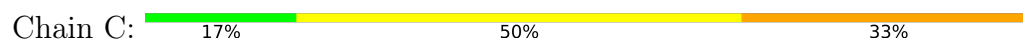


- Molecule 1: Alpha-mannosidase from *Canavalia ensiformis* (jack bean)





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



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



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.31Å 119.67Å 277.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.55 – 1.84 49.55 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.55-1.84) 99.4 (49.55-1.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 1.84Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.175 , 0.216 0.176 , 0.217	Depositor DCC
R_{free} test set	9670 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17299	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 3.26% 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: MAN, BMA, ZN, NAG

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	0.36	0/7777	0.55	1/10540 (0.0%)
1	B	0.35	0/7767	0.54	1/10528 (0.0%)
All	All	0.35	0/15544	0.55	2/21068 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	157	LEU	CA-CB-CG	-5.89	95.69	116.30
1	A	157	LEU	CA-CB-CG	-5.65	96.53	116.30

There are no chirality outliers.

There are no planarity outliers.

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	7561	0	7318	97	0
1	B	7561	0	7298	100	0
2	C	149	0	125	8	0
2	E	138	0	115	5	0
3	D	28	0	25	0	0
3	F	28	0	25	1	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	A	933	0	0	40	2
5	B	899	0	0	52	2
All	All	17299	0	14906	204	2

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 (204) 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:541:PRO:O	1:B:862:ILE:CD1	2.06	1.03
1:B:541:PRO:O	1:B:862:ILE:HD13	1.64	0.97
1:B:813:ARG:H	2:E:1:NAG:H81	1.40	0.84
1:B:632:GLN:NE2	1:B:633:PRO:O	2.13	0.81
1:B:329:ARG:NH2	5:B:1102:HOH:O	2.15	0.80
1:B:585:THR:O	1:B:587:GLN:N	2.18	0.77
1:A:813:ARG:H	2:C:1:NAG:H81	1.52	0.73
1:B:314:THR:HB	5:B:1146:HOH:O	1.90	0.70
1:B:640:ARG:N	5:B:1112:HOH:O	2.26	0.69
1:B:632:GLN:HE22	1:B:635:PRO:HD3	1.58	0.69
1:B:183:LYS:NZ	5:B:1114:HOH:O	2.26	0.69
1:B:541:PRO:O	1:B:862:ILE:HD11	1.91	0.68
1:A:505:MET:SD	5:A:1917:HOH:O	2.52	0.67
1:A:850:MET:SD	5:A:1951:HOH:O	2.53	0.66
1:B:540:PRO:HA	5:B:1103:HOH:O	1.96	0.66
1:B:546:THR:HG22	5:B:1115:HOH:O	1.96	0.65
1:B:675:HIS:NE2	5:B:1101:HOH:O	2.08	0.63
1:A:678:ILE:HG13	5:A:1341:HOH:O	1.98	0.63
1:B:13:GLN:NE2	5:B:1127:HOH:O	2.33	0.61
1:B:183:LYS:N	1:B:183:LYS:HD3	2.16	0.61
1:B:573:GLY:O	1:B:578:LYS:NZ	2.33	0.61
1:A:529:ASP:OD1	1:A:529:ASP:N	2.30	0.60
1:B:290:LYS:NZ	5:B:1125:HOH:O	2.32	0.60
1:B:474:SER:OG	5:B:1103:HOH:O	2.16	0.60
1:A:570:LEU:HD21	5:A:1971:HOH:O	2.00	0.60
1:B:582:SER:O	1:B:586:GLY:HA2	2.02	0.59
1:A:182:GLU:OE2	5:A:1101:HOH:O	2.16	0.59
1:A:880:GLU:HB3	5:A:1166:HOH:O	2.02	0.59
1:B:193:LYS:NZ	1:B:896:GLU:OE1	2.36	0.59
1:B:852:ASN:ND2	5:B:1136:HOH:O	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:MET:HE2	5:B:1260:HOH:O	2.03	0.58
1:A:330:ASN:HB3	1:A:628:ARG:HD3	1.84	0.58
1:B:183:LYS:HE2	5:B:1731:HOH:O	2.04	0.58
1:A:590:ARG:HD2	1:A:599:ASP:OD1	2.04	0.57
1:A:668:ARG:NH2	1:A:679:GLU:OE2	2.37	0.57
1:A:548:PHE:CD1	1:A:850:MET:HG3	2.40	0.56
1:B:183:LYS:HD2	1:B:255:ILE:HA	1.87	0.56
1:B:731:LEU:HD23	5:B:1105:HOH:O	2.05	0.56
1:B:733:VAL:HG22	5:B:1105:HOH:O	2.05	0.56
1:B:570:LEU:HB2	1:B:581:PHE:O	2.06	0.56
1:A:934:LYS:HE3	5:A:1116:HOH:O	2.07	0.55
5:A:1775:HOH:O	2:C:2:NAG:H83	2.06	0.54
1:B:945:GLU:N	5:B:1146:HOH:O	2.40	0.54
5:A:1937:HOH:O	2:C:2:NAG:H81	2.07	0.54
1:A:886:LEU:HB3	5:A:1906:HOH:O	2.08	0.54
1:B:724:ASP:O	5:B:1105:HOH:O	2.18	0.54
1:A:95:LYS:HE2	5:A:1943:HOH:O	2.08	0.54
1:A:888:LEU:HB3	5:A:1166:HOH:O	2.07	0.54
1:A:518:LYS:HE3	1:A:801:MET:HG2	1.89	0.54
1:A:612:SER:HB3	1:A:628:ARG:HG2	1.90	0.54
1:A:44:GLN:NE2	5:A:1113:HOH:O	2.41	0.54
1:A:139:ARG:HD2	5:A:1321:HOH:O	2.08	0.53
2:C:6[B]:MAN:H3	2:C:7:MAN:C1	2.37	0.53
1:A:524:VAL:HG12	5:A:1223:HOH:O	2.07	0.53
1:A:540:PRO:HB2	1:A:862:ILE:HD12	1.91	0.53
5:A:1574:HOH:O	2:C:1:NAG:H83	2.08	0.53
1:A:484:ASN:HA	1:A:530:VAL:HG22	1.90	0.53
1:A:345:ASP:HB2	1:A:763:ARG:CZ	2.40	0.52
5:B:1106:HOH:O	3:F:1:NAG:O7	2.19	0.52
1:A:570:LEU:HD22	1:A:581:PHE:HB2	1.91	0.52
1:A:342:ALA:HB1	2:C:2:NAG:H82	1.91	0.52
1:A:586:GLY:CA	1:A:643:VAL:HG23	2.40	0.52
1:A:888:LEU:HD23	5:A:1166:HOH:O	2.10	0.52
1:B:813:ARG:N	2:E:1:NAG:H81	2.17	0.51
1:A:676:ALA:HB3	1:A:818[A]:LEU:CD2	2.41	0.51
1:B:632:GLN:HG2	1:B:633:PRO:HD2	1.91	0.51
1:A:920:ILE:HD13	1:A:981:PHE:CE1	2.45	0.51
1:A:586:GLY:HA3	1:A:643:VAL:HG23	1.91	0.51
1:A:852:ASN:HB2	5:A:1640:HOH:O	2.11	0.51
1:A:922:GLU:CD	1:A:980:GLN:HE21	2.18	0.51
1:A:160:GLU:OE1	5:A:1102:HOH:O	2.18	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:847:GLN:O	1:A:848:GLU:HG2	2.10	0.50
1:B:497:LYS:HE3	5:B:1943:HOH:O	2.11	0.50
1:B:109:HIS:CE1	5:B:1260:HOH:O	2.63	0.50
1:A:850:MET:HE2	5:A:1119:HOH:O	2.11	0.50
1:A:908:LYS:HE2	1:A:966:VAL:HG11	1.93	0.50
1:B:667:THR:HG22	5:B:1685:HOH:O	2.12	0.50
1:B:157:LEU:HD13	5:B:1260:HOH:O	2.11	0.49
1:B:342:ALA:HB1	2:E:2:NAG:H82	1.94	0.49
1:B:600:ILE:HG12	5:B:1893:HOH:O	2.11	0.49
1:A:851:GLU:CD	1:A:851:GLU:H	2.20	0.49
1:A:587:GLN:NE2	5:A:1147:HOH:O	2.45	0.49
1:A:144:ILE:HA	1:A:168:PHE:HA	1.95	0.49
1:A:726[B]:ARG:HD3	1:A:729:TRP:CH2	2.48	0.49
1:B:751:LYS:NZ	5:B:1159:HOH:O	2.46	0.49
1:B:529:ASP:OD1	1:B:529:ASP:N	2.38	0.49
1:B:673:LYS:HD2	5:B:1101:HOH:O	2.12	0.49
1:A:544:TRP:CH2	1:A:861:GLY:HA3	2.48	0.48
1:B:945:GLU:HG3	5:B:1146:HOH:O	2.13	0.48
1:A:570:LEU:C	1:A:570:LEU:HD23	2.38	0.48
1:B:110:ASP:HB2	1:B:386:HIS:HA	1.96	0.48
1:A:290:LYS:HE2	1:A:290:LYS:HB3	1.68	0.48
1:B:585:THR:O	1:B:585:THR:OG1	2.29	0.48
1:B:964:ASP:N	5:B:1158:HOH:O	2.45	0.48
1:B:273:TYR:CZ	1:B:275:GLU:HB3	2.49	0.48
1:B:542:LEU:O	1:B:862:ILE:HD12	2.14	0.47
1:B:640:ARG:CZ	1:B:640:ARG:HB2	2.44	0.47
1:B:827:SER:OG	1:B:899:GLU:HB3	2.14	0.47
1:A:548:PHE:CG	1:A:850:MET:HG3	2.49	0.47
1:A:813:ARG:N	2:C:1:NAG:H81	2.25	0.47
1:A:829:TRP:CD1	1:A:833:THR:HG1	2.33	0.47
1:A:221:ARG:HA	5:B:1388:HOH:O	2.13	0.47
5:B:1223:HOH:O	2:E:1:NAG:H83	2.14	0.47
1:A:530:VAL:HG23	1:A:531:TYR:CD2	2.50	0.47
1:B:536:LYS:HE2	1:B:672:ASP:OD2	2.15	0.47
1:B:366:LYS:NZ	5:B:1138:HOH:O	2.46	0.47
1:B:193:LYS:HG3	5:B:1542:HOH:O	2.15	0.47
1:B:611:SER:HB3	1:B:622:SER:HB3	1.97	0.47
1:A:936:GLU:HG2	5:A:1934:HOH:O	2.15	0.47
1:B:514:SER:HB2	5:B:1234:HOH:O	2.15	0.47
5:B:1860:HOH:O	2:E:2:NAG:H83	2.15	0.47
1:A:137:ILE:HD13	1:A:932:GLN:HG2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:584:LEU:HB3	5:B:1876:HOH:O	2.15	0.46
1:B:589:LYS:O	5:B:1108:HOH:O	2.21	0.46
1:B:798:GLN:NE2	5:B:1130:HOH:O	2.34	0.46
1:A:528:ALA:HB1	5:A:1917:HOH:O	2.14	0.46
1:A:605:ASN:ND2	1:A:636:HIS:CE1	2.84	0.46
1:A:385:HIS:CD2	1:A:387:ASP:H	2.33	0.46
1:B:137:ILE:HD13	1:B:932:GLN:HG2	1.98	0.46
1:B:497:LYS:HD2	5:B:1873:HOH:O	2.15	0.46
1:B:980:GLN:HB2	5:B:1795:HOH:O	2.14	0.46
1:A:878:THR:HB	1:A:890:ARG:HB2	1.97	0.46
1:B:782:ARG:HD2	1:B:796:ASP:OD1	2.16	0.46
1:A:678:ILE:HG21	1:A:816:TYR:CZ	2.51	0.45
1:B:676:ALA:HB3	1:B:818:LEU:CD2	2.46	0.45
1:B:864:MET:SD	1:B:914:LEU:HD12	2.56	0.45
1:A:110:ASP:HB2	1:A:386:HIS:HA	1.97	0.45
1:A:273:TYR:CZ	1:A:275:GLU:HB3	2.52	0.45
1:A:329:ARG:NH2	1:A:732:GLU:OE2	2.43	0.45
1:A:912:LYS:HE3	5:A:1192:HOH:O	2.17	0.45
1:A:254:LYS:NZ	5:A:1170:HOH:O	2.50	0.44
1:A:782:ARG:HD3	5:A:1442:HOH:O	2.18	0.44
1:B:580:SER:HB2	1:B:589:LYS:HG2	1.98	0.44
1:A:611:SER:HB3	1:A:622:SER:HB3	2.00	0.44
1:B:144:ILE:HA	1:B:168:PHE:HA	1.99	0.44
1:A:570:LEU:HB3	1:A:581:PHE:O	2.17	0.44
1:A:655:VAL:HG22	5:A:1971:HOH:O	2.17	0.44
1:A:912:LYS:HE2	5:A:1914:HOH:O	2.16	0.44
1:B:345:ASP:HB2	1:B:763:ARG:CZ	2.47	0.44
1:B:477:GLU:HG3	5:B:1749:HOH:O	2.16	0.44
1:A:646:VAL:HG13	5:A:1971:HOH:O	2.18	0.44
1:A:670:TYR:HB2	1:A:673:LYS:HG3	2.00	0.44
1:A:922:GLU:OE2	1:A:980:GLN:NE2	2.48	0.44
1:A:591:MET:HB3	5:A:1410:HOH:O	2.18	0.44
1:A:631:GLY:N	5:A:1126:HOH:O	2.51	0.44
1:B:211:GLY:HA2	1:B:266:GLY:O	2.18	0.44
1:A:269:PHE:HE2	5:A:1113:HOH:O	2.01	0.43
1:A:654:GLU:OE2	1:A:668:ARG:HD3	2.18	0.43
1:B:586:GLY:N	5:B:1184:HOH:O	2.51	0.43
1:A:541:PRO:HG2	1:A:862:ILE:HD11	2.00	0.43
1:B:343:LEU:HD22	1:B:398:THR:HG23	1.99	0.43
1:B:541:PRO:HB2	1:B:864:MET:HE2	2.00	0.43
1:A:330:ASN:HB3	1:A:628:ARG:CD	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:922:GLU:HB3	1:A:980:GLN:HG3	2.00	0.43
1:A:211:GLY:HA2	1:A:266:GLY:O	2.19	0.43
1:B:1:MET:SD	5:B:1872:HOH:O	2.62	0.43
1:B:385:HIS:CD2	1:B:387:ASP:H	2.37	0.43
5:A:1202:HOH:O	2:C:6[A]:MAN:H61	2.18	0.43
1:A:486:ALA:HB2	1:A:531:TYR:CE2	2.54	0.43
1:A:751:LYS:HB3	1:A:756:GLU:HG3	2.00	0.43
1:A:707:LYS:NZ	5:A:1154:HOH:O	2.46	0.43
1:A:319[A]:ILE:HG22	1:A:320:ASP:CG	2.43	0.42
1:A:886:LEU:HD12	5:A:1416:HOH:O	2.17	0.42
1:A:37:VAL:HG11	1:B:228:GLN:HG2	2.01	0.42
1:A:666:VAL:HG12	1:A:668:ARG:HG3	2.01	0.42
1:B:62:ASP:HB3	1:B:65:ARG:HD2	2.01	0.42
1:B:209:HIS:HB3	5:B:1821:HOH:O	2.19	0.42
1:B:13:GLN:CD	5:B:1127:HOH:O	2.62	0.42
1:B:862:ILE:HG23	1:B:864:MET:O	2.20	0.42
1:A:922:GLU:HG2	5:A:1187:HOH:O	2.18	0.42
1:B:358:ARG:NH2	5:B:1194:HOH:O	2.52	0.42
1:B:467:VAL:O	1:B:544:TRP:HA	2.20	0.42
1:B:598:VAL:HG12	5:B:1893:HOH:O	2.18	0.42
1:A:491:LYS:HD2	1:A:850:MET:HE1	2.02	0.42
1:A:582:SER:O	1:A:586:GLY:HA2	2.20	0.42
1:B:939:LYS:HE2	5:B:1890:HOH:O	2.20	0.42
1:A:132:GLU:HG3	5:A:1779:HOH:O	2.20	0.42
1:B:176:ARG:HG2	1:B:180:LYS:HE3	2.01	0.42
1:B:744:LEU:HD11	1:B:767:GLY:HA3	2.01	0.42
1:B:687:THR:HG21	1:B:782:ARG:HG2	2.01	0.41
1:B:583:SER:O	5:B:1109:HOH:O	2.21	0.41
1:A:865:ASP:OD2	1:A:913:LYS:HA	2.21	0.41
1:B:477:GLU:HB3	1:B:841[A]:MET:SD	2.60	0.41
1:B:608:TRP:O	1:B:635:PRO:HD2	2.21	0.41
1:B:878:THR:HB	1:B:890:ARG:HB2	2.02	0.41
1:B:362:PHE:CE1	1:B:887:VAL:HG21	2.55	0.41
1:B:635:PRO:HG3	5:B:1822:HOH:O	2.19	0.41
1:B:938:LYS:NZ	5:B:1199:HOH:O	2.53	0.41
1:A:304:TYR:O	1:A:308:LYS:HG2	2.21	0.41
1:B:423:LEU:HD22	1:B:863:TYR:CE1	2.55	0.41
1:B:683:GLY:HA2	1:B:684:PRO:C	2.45	0.41
1:A:587:GLN:HG3	5:A:1135:HOH:O	2.21	0.41
1:A:116:TYR:CD1	1:A:716:ARG:HG3	2.56	0.41
1:A:919:LYS:NZ	5:A:1192:HOH:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:LYS:HD3	1:B:183:LYS:H	1.83	0.41
1:B:915:PHE:HB3	5:B:1363:HOH:O	2.21	0.41
1:B:918:GLN:HB2	5:B:1363:HOH:O	2.21	0.41
1:A:314:THR:HG22	1:A:943:SER:OG	2.21	0.40
1:A:44:GLN:CD	5:A:1113:HOH:O	2.64	0.40
1:A:370:LYS:HD2	1:A:880:GLU:OE2	2.21	0.40
1:A:453:THR:HG22	1:A:463:LEU:HD21	2.04	0.40
1:B:687:THR:HB	1:B:784:ILE:HD11	2.03	0.40
1:B:677:GLU:HG2	1:B:817:TYR:CD2	2.57	0.40
1:B:867:ASN:N	5:B:1129:HOH:O	2.34	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1365:HOH:O	5:B:1804:HOH:O[1_655]	1.93	0.27
5:A:1365:HOH:O	5:B:1133:HOH:O[1_655]	2.19	0.01

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	931/981 (95%)	894 (96%)	34 (4%)	3 (0%)	36	25
1	B	931/981 (95%)	893 (96%)	34 (4%)	4 (0%)	30	18
All	All	1862/1962 (95%)	1787 (96%)	68 (4%)	7 (0%)	30	18

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	586	GLY
1	A	28	TRP

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Mol	Chain	Res	Type
1	B	28	TRP
1	B	208	GLY
1	B	103	ASN
1	A	103	ASN
1	A	208	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	823/849 (97%)	820 (100%)	3 (0%)	84	78
1	B	821/849 (97%)	820 (100%)	1 (0%)	88	85
All	All	1644/1698 (97%)	1640 (100%)	4 (0%)	87	84

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

Mol	Chain	Res	Type
1	A	228	GLN
1	A	818[A]	LEU
1	A	818[B]	LEU
1	B	182	GLU

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	587	GLN
1	A	603	GLN
1	A	605	ASN
1	A	636	HIS
1	A	980	GLN
1	B	131	GLN
1	B	133	GLN
1	B	223	ASN
1	B	313	GLN

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Mol	Chain	Res	Type
1	B	587	GLN
1	B	605	ASN
1	B	632	GLN
1	B	932	GLN

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 ⓘ

29 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	0.84	1 (7%)	17,19,21	0.80	1 (5%)
2	MAN	C	10	2	11,11,12	0.74	0	15,15,17	1.03	1 (6%)
2	MAN	C	11	2	11,11,12	1.05	1 (9%)	15,15,17	1.29	2 (13%)
2	MAN	C	12	2	11,11,12	0.55	0	15,15,17	1.03	1 (6%)
2	NAG	C	2	2	14,14,15	0.63	0	17,19,21	0.81	1 (5%)
2	BMA	C	3	2	11,11,12	0.86	0	15,15,17	0.92	0
2	MAN	C	4	2	11,11,12	0.82	0	15,15,17	1.10	1 (6%)
2	MAN	C	5	2	11,11,12	0.77	0	15,15,17	1.02	0
2	MAN	C	6[A]	2	11,11,12	1.59	3 (27%)	15,15,17	1.66	3 (20%)
2	MAN	C	6[B]	2	11,11,12	1.98	2 (18%)	15,15,17	2.06	4 (26%)
2	MAN	C	7	2	11,11,12	1.90	3 (27%)	15,15,17	1.56	4 (26%)
2	MAN	C	8	2	11,11,12	0.72	0	15,15,17	1.24	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	C	9	2	11,11,12	0.69	0	15,15,17	1.09	2 (13%)
3	NAG	D	1	1,3	14,14,15	0.36	0	17,19,21	0.72	0
3	NAG	D	2	3	14,14,15	0.34	0	17,19,21	1.99	2 (11%)
2	NAG	E	1	1,2	14,14,15	0.66	0	17,19,21	0.77	1 (5%)
2	MAN	E	10	2	11,11,12	0.95	1 (9%)	15,15,17	1.16	1 (6%)
2	MAN	E	11	2	11,11,12	0.85	0	15,15,17	1.10	1 (6%)
2	MAN	E	12	2	11,11,12	0.73	0	15,15,17	0.88	1 (6%)
2	NAG	E	2	2	14,14,15	0.65	0	17,19,21	0.82	1 (5%)
2	BMA	E	3	2	11,11,12	0.63	0	15,15,17	0.98	0
2	MAN	E	4	2	11,11,12	0.78	0	15,15,17	1.06	1 (6%)
2	MAN	E	5	2	11,11,12	0.62	0	15,15,17	1.13	2 (13%)
2	MAN	E	6	2	11,11,12	1.05	0	15,15,17	1.32	2 (13%)
2	MAN	E	7	2	11,11,12	1.59	1 (9%)	15,15,17	1.30	2 (13%)
2	MAN	E	8	2	11,11,12	0.78	0	15,15,17	1.16	2 (13%)
2	MAN	E	9	2	11,11,12	0.86	1 (9%)	15,15,17	1.25	2 (13%)
3	NAG	F	1	1,3	14,14,15	1.52	2 (14%)	17,19,21	1.38	1 (5%)
3	NAG	F	2	3	14,14,15	1.30	2 (14%)	17,19,21	1.42	1 (5%)

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	-	4/6/23/26	0/1/1/1
2	MAN	C	10	2	-	0/2/19/22	0/1/1/1
2	MAN	C	11	2	-	0/2/19/22	0/1/1/1
2	MAN	C	12	2	-	0/2/19/22	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	0/2/19/22	0/1/1/1
2	MAN	C	5	2	-	0/2/19/22	0/1/1/1
2	MAN	C	6[A]	2	-	2/2/19/22	0/1/1/1
2	MAN	C	6[B]	2	-	2/2/19/22	0/1/1/1
2	MAN	C	7	2	-	0/2/19/22	0/1/1/1
2	MAN	C	8	2	-	0/2/19/22	0/1/1/1
2	MAN	C	9	2	-	0/2/19/22	0/1/1/1
3	NAG	D	1	1,3	-	3/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	MAN	E	10	2	-	2/2/19/22	0/1/1/1
2	MAN	E	11	2	-	0/2/19/22	0/1/1/1
2	MAN	E	12	2	-	0/2/19/22	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	MAN	E	4	2	-	1/2/19/22	0/1/1/1
2	MAN	E	5	2	-	0/2/19/22	0/1/1/1
2	MAN	E	6	2	-	2/2/19/22	0/1/1/1
2	MAN	E	7	2	-	0/2/19/22	0/1/1/1
2	MAN	E	8	2	-	0/2/19/22	0/1/1/1
2	MAN	E	9	2	-	0/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	7	MAN	C2-C3	5.28	1.60	1.52
2	E	7	MAN	C2-C3	4.64	1.59	1.52
3	F	1	NAG	O5-C1	4.56	1.51	1.43
3	F	2	NAG	O5-C1	4.36	1.51	1.43
2	C	6[B]	MAN	C2-C3	4.23	1.59	1.52
2	C	6[B]	MAN	O5-C5	4.01	1.51	1.43
3	F	1	NAG	C1-C2	3.23	1.56	1.52
2	C	6[A]	MAN	C2-C3	-3.00	1.47	1.52
2	C	6[A]	MAN	C1-C2	2.63	1.58	1.52
2	C	7	MAN	C4-C3	2.58	1.59	1.52
2	C	6[A]	MAN	C4-C5	2.54	1.58	1.53
2	C	11	MAN	O5-C5	2.54	1.48	1.43
2	C	1	NAG	C1-C2	2.30	1.55	1.52
2	E	9	MAN	O5-C5	2.22	1.47	1.43
2	E	10	MAN	C1-C2	2.09	1.57	1.52
2	C	7	MAN	O5-C1	-2.05	1.40	1.43
3	F	2	NAG	C1-C2	2.03	1.55	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2	NAG	C1-O5-C5	7.43	122.14	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	2	NAG	C1-O5-C5	5.57	119.65	112.19
2	C	6[B]	MAN	C2-C3-C4	5.43	120.41	110.86
3	F	1	NAG	C1-O5-C5	5.32	119.32	112.19
2	C	6[A]	MAN	C1-O5-C5	4.29	117.93	112.19
2	E	6	MAN	C1-O5-C5	3.98	117.52	112.19
2	E	4	MAN	O2-C2-C3	-3.48	102.94	110.15
2	C	6[B]	MAN	C1-C2-C3	3.40	114.59	109.64
2	C	11	MAN	O2-C2-C3	-3.33	103.25	110.15
2	E	9	MAN	C1-O5-C5	3.20	116.47	112.19
2	E	9	MAN	O2-C2-C3	-3.16	103.61	110.15
2	C	6[B]	MAN	O2-C2-C3	-3.10	103.74	110.15
2	C	12	MAN	C1-O5-C5	3.04	116.26	112.19
2	C	7	MAN	C1-C2-C3	3.03	114.05	109.64
2	C	4	MAN	O2-C2-C3	-2.98	103.98	110.15
2	E	11	MAN	O2-C2-C3	-2.95	104.04	110.15
2	E	10	MAN	C1-O5-C5	2.88	116.05	112.19
2	E	7	MAN	C2-C3-C4	2.71	115.62	110.86
2	E	8	MAN	O2-C2-C3	-2.69	104.57	110.15
2	C	9	MAN	C1-O5-C5	2.68	115.78	112.19
2	C	11	MAN	C1-O5-C5	2.65	115.74	112.19
2	E	7	MAN	C1-C2-C3	2.65	113.50	109.64
2	E	5	MAN	C1-O5-C5	2.61	115.68	112.19
2	C	8	MAN	O2-C2-C3	-2.59	104.78	110.15
3	D	2	NAG	C3-C4-C5	2.59	114.92	110.23
2	C	9	MAN	O2-C2-C3	-2.58	104.80	110.15
2	C	7	MAN	O5-C1-C2	2.57	116.92	110.79
2	C	7	MAN	C1-O5-C5	2.56	115.62	112.19
2	C	6[A]	MAN	O5-C1-C2	2.51	116.77	110.79
2	C	7	MAN	C2-C3-C4	2.51	115.27	110.86
2	C	10	MAN	O2-C2-C3	-2.50	104.96	110.15
2	E	2	NAG	C1-O5-C5	2.45	115.47	112.19
2	E	5	MAN	O2-C2-C3	-2.38	105.22	110.15
2	E	12	MAN	C1-O5-C5	2.37	115.37	112.19
2	C	6[A]	MAN	C2-C3-C4	-2.34	106.74	110.86
2	C	6[B]	MAN	O5-C5-C6	2.27	112.09	107.66
2	C	8	MAN	C1-O5-C5	2.23	115.17	112.19
2	C	1	NAG	C1-O5-C5	2.21	115.14	112.19
2	E	1	NAG	C1-O5-C5	2.18	115.11	112.19
2	E	8	MAN	C1-O5-C5	2.16	115.08	112.19
2	C	2	NAG	C1-O5-C5	2.14	115.05	112.19
2	E	6	MAN	O2-C2-C3	-2.04	105.93	110.15

There are no chirality outliers.

All (20) torsion outliers are listed below:

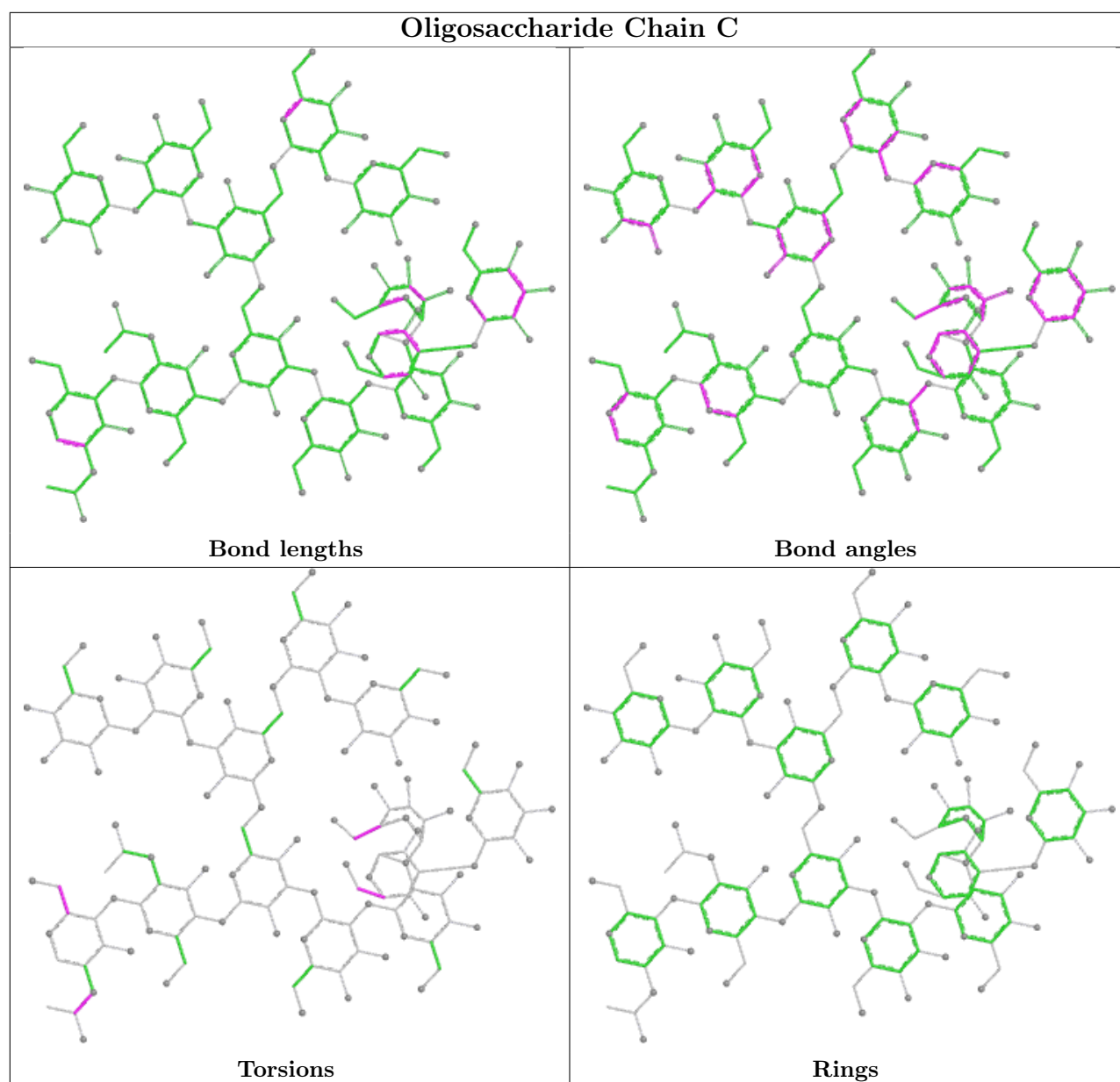
Mol	Chain	Res	Type	Atoms
2	C	6[A]	MAN	O5-C5-C6-O6
2	C	6[B]	MAN	C4-C5-C6-O6
2	C	6[B]	MAN	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
2	C	6[A]	MAN	C4-C5-C6-O6
2	E	6	MAN	O5-C5-C6-O6
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C4-C5-C6-O6
2	E	10	MAN	O5-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
2	E	10	MAN	C4-C5-C6-O6
2	E	6	MAN	C4-C5-C6-O6
3	D	1	NAG	C3-C2-N2-C7
3	D	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	E	4	MAN	C4-C5-C6-O6

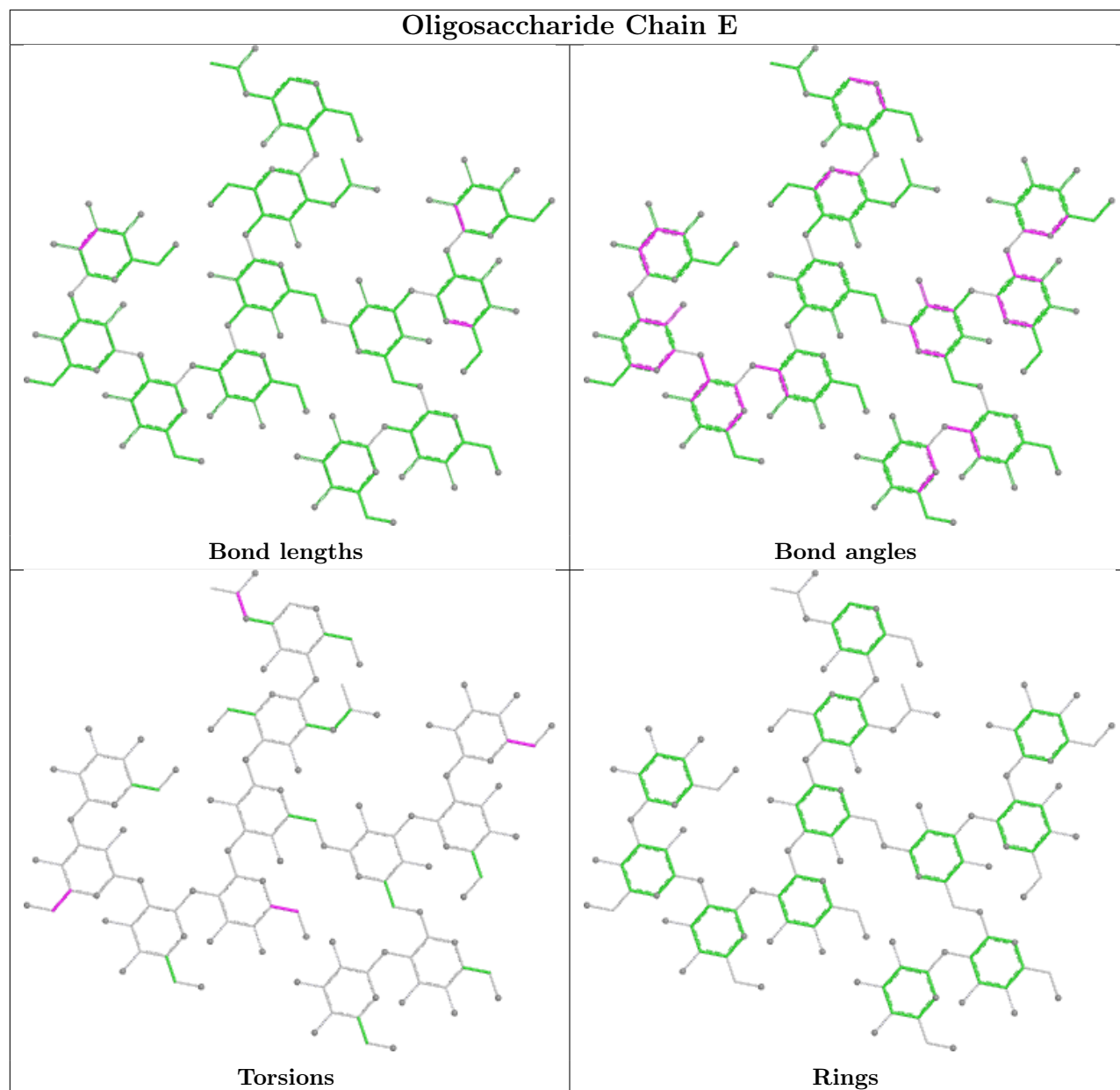
There are no ring outliers.

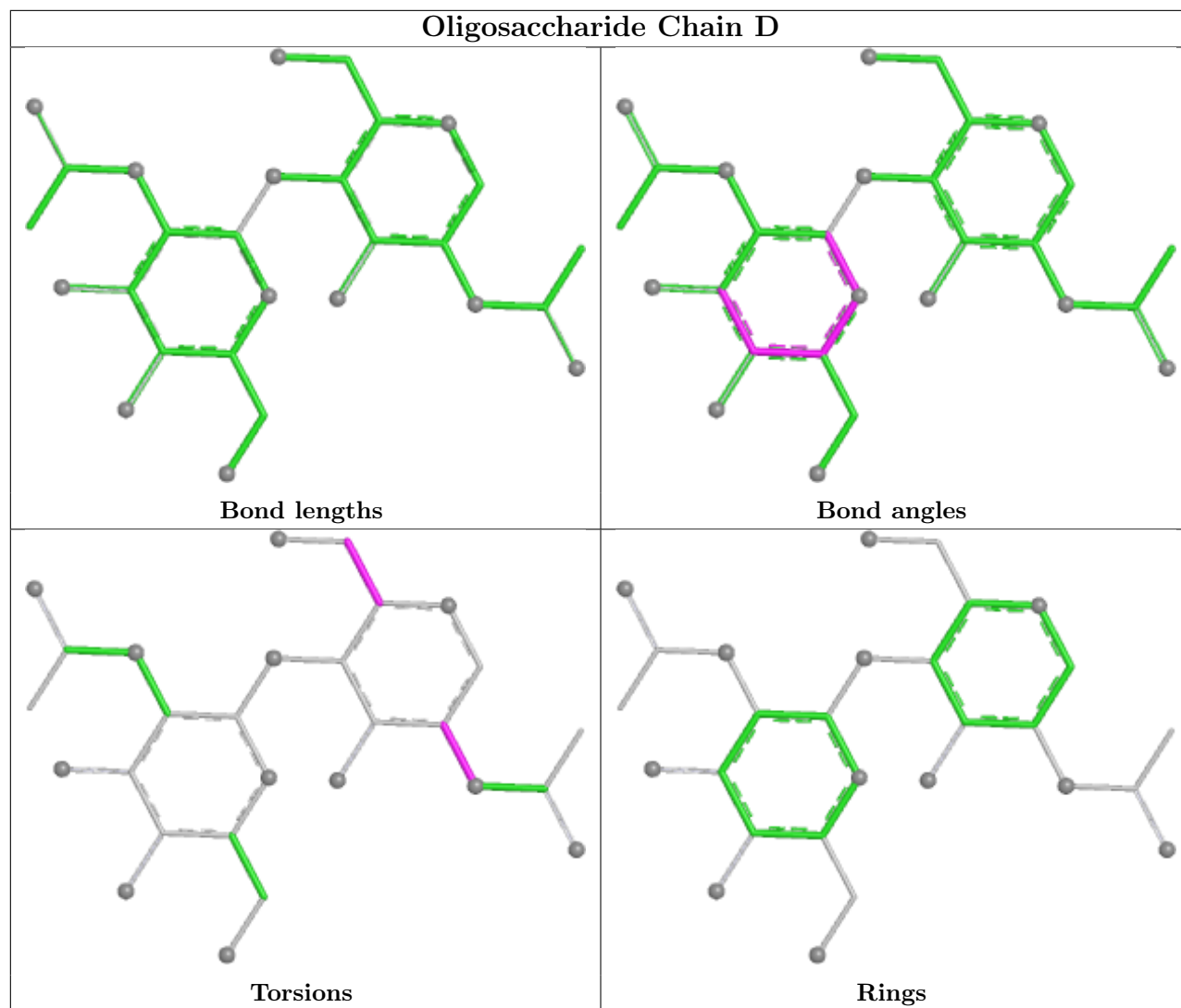
8 monomers are involved in 14 short contacts:

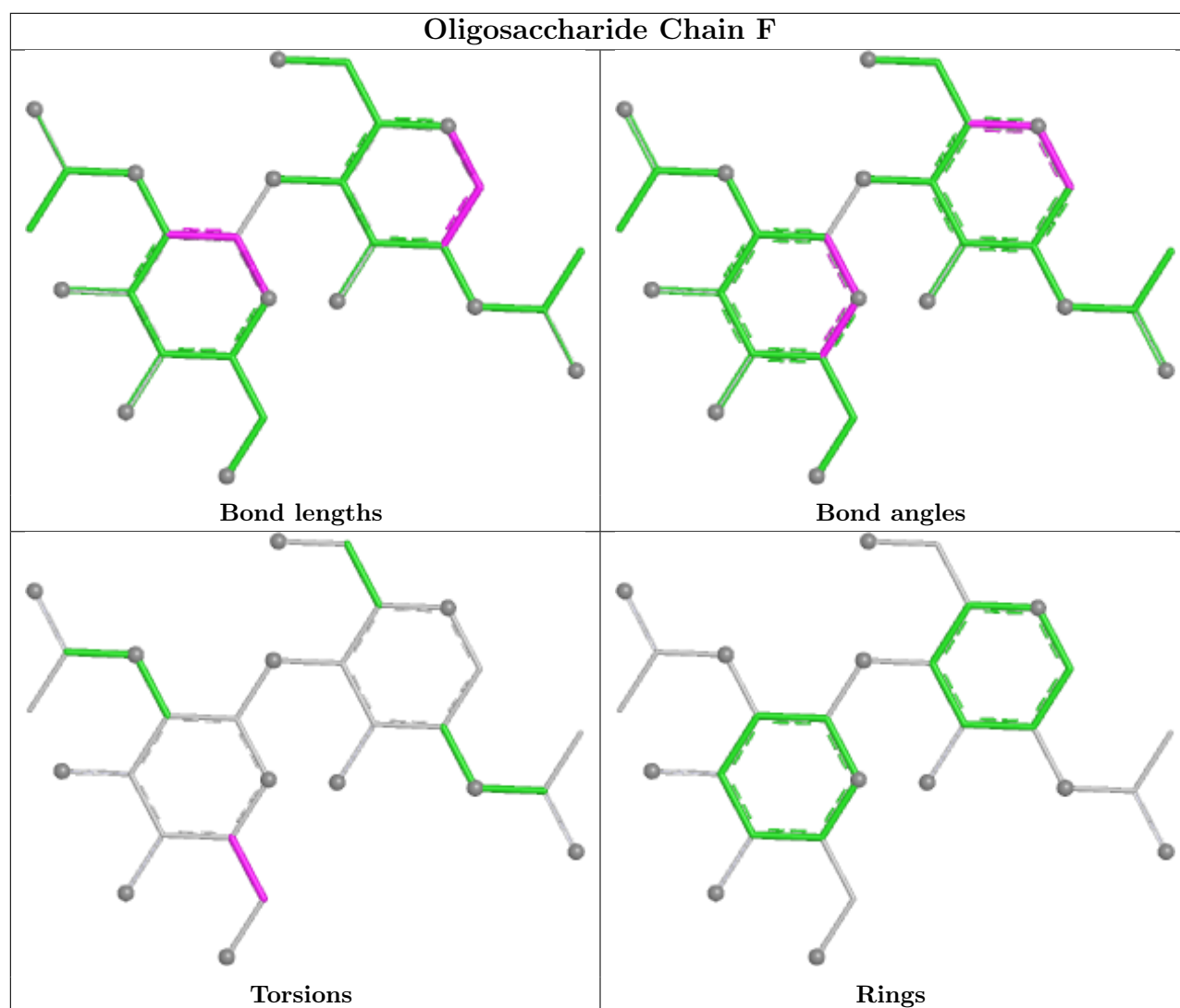
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	3	0
3	F	1	NAG	1	0
2	C	7	MAN	1	0
2	E	1	NAG	3	0
2	C	6[B]	MAN	1	0
2	E	2	NAG	2	0
2	C	6[A]	MAN	1	0
2	C	2	NAG	3	0

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









5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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(Å²)	Q<0.9
1	A	935/981 (95%)	-0.24	25 (2%)	56	61	8, 24, 48, 79	10 (1%)
1	B	937/981 (95%)	-0.13	26 (2%)	55	60	9, 26, 51, 103	6 (0%)
All	All	1872/1962 (95%)	-0.18	51 (2%)	56	61	8, 25, 50, 103	16 (0%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	917	THR	4.9
1	A	570	LEU	4.7
1	B	636	HIS	4.4
1	B	570	LEU	4.3
1	B	805	TYR	4.3
1	B	635	PRO	4.2
1	B	616	PHE	3.8
1	A	801	MET	3.6
1	B	633	PRO	3.5
1	A	636	HIS	3.3
1	B	632	GLN	3.3
1	A	586	GLY	3.1
1	B	422	CYS	2.9
1	A	583	SER	2.9
1	B	585	THR	2.8
1	B	433	SER	2.8
1	A	551	GLU	2.7
1	B	884	ASP	2.7
1	A	633	PRO	2.7
1	A	637	THR	2.7
1	A	635	PRO	2.7
1	B	801	MET	2.6
1	B	800	CYS	2.6
1	A	594	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	432	CYS	2.5
1	A	632	GLN	2.5
1	B	863	TYR	2.5
1	B	476	ASN	2.4
1	B	528	ALA	2.4
1	B	531	TYR	2.4
1	A	806	THR	2.4
1	A	862	ILE	2.4
1	A	643	VAL	2.4
1	B	584	LEU	2.3
1	A	619	TYR	2.3
1	A	964	ASP	2.3
1	A	433	SER	2.3
1	A	630	ASN	2.3
1	B	631	GLY	2.2
1	A	885	GLY	2.2
1	B	640	ARG	2.2
1	A	427	GLN	2.1
1	B	806	THR	2.1
1	A	916	ALA	2.1
1	A	915	PHE	2.1
1	A	432	CYS	2.1
1	B	526	LYS	2.1
1	B	443	HIS	2.1
1	A	531	TYR	2.0
1	B	530	VAL	2.0
1	B	485	ASP	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	NAG	F	2	14/15	0.21	0.24	100,133,145,147	0

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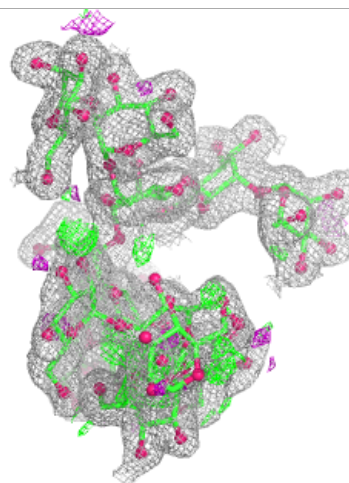
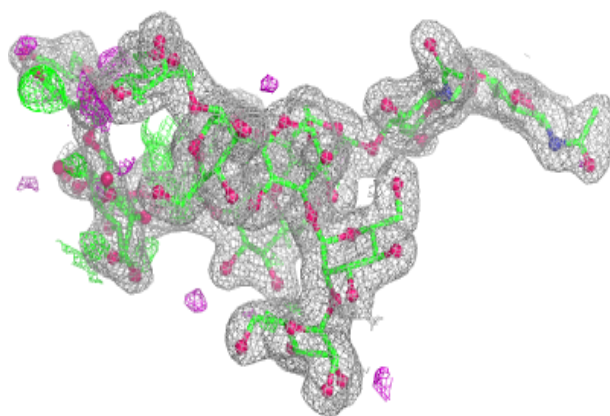
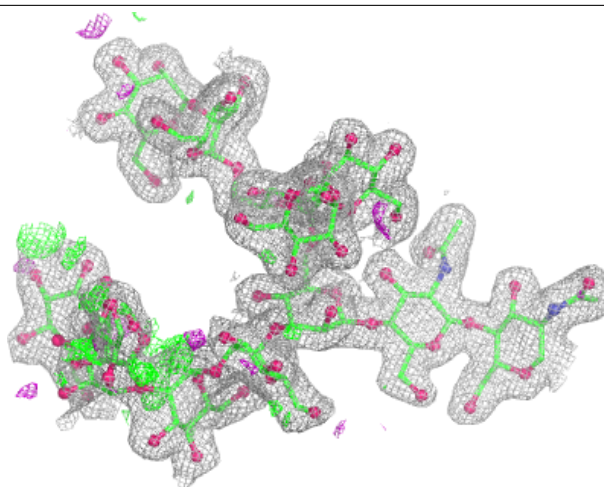
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	F	1	14/15	0.24	0.20	88,106,117,130	0
3	NAG	D	2	14/15	0.33	0.22	76,87,94,95	0
3	NAG	D	1	14/15	0.69	0.14	56,64,74,79	0
2	MAN	C	6[A]	11/12	0.73	0.18	43,46,48,49	11
2	MAN	C	6[B]	11/12	0.73	0.18	45,47,50,50	11
2	MAN	E	6	11/12	0.75	0.15	49,53,58,61	0
2	MAN	C	7	11/12	0.84	0.13	31,43,48,52	0
2	MAN	C	5	11/12	0.86	0.10	27,38,43,44	0
2	MAN	E	7	11/12	0.88	0.13	29,46,52,61	0
2	MAN	E	5	11/12	0.89	0.10	24,39,46,47	0
2	MAN	E	10	11/12	0.89	0.10	34,39,43,45	0
2	MAN	C	4	11/12	0.92	0.09	26,31,40,42	0
2	MAN	E	4	11/12	0.93	0.09	24,30,37,41	0
2	MAN	E	9	11/12	0.94	0.08	28,32,36,37	0
2	MAN	C	10	11/12	0.94	0.07	28,31,38,44	0
2	NAG	E	1	14/15	0.94	0.08	21,24,28,39	0
2	NAG	C	2	14/15	0.95	0.07	21,24,33,38	0
2	NAG	E	2	14/15	0.95	0.07	15,22,31,38	0
2	NAG	C	1	14/15	0.96	0.07	21,24,28,34	0
2	MAN	C	12	11/12	0.97	0.05	17,18,20,21	0
2	MAN	E	8	11/12	0.97	0.05	17,20,24,24	0
2	MAN	C	9	11/12	0.97	0.05	25,25,30,30	0
2	BMA	C	3	11/12	0.97	0.04	18,18,23,24	0
2	MAN	E	12	11/12	0.98	0.05	15,17,20,22	0
2	MAN	C	8	11/12	0.98	0.04	17,19,22,25	0
2	BMA	E	3	11/12	0.98	0.05	18,22,27,29	0
2	MAN	C	11	11/12	0.98	0.04	17,20,24,25	0
2	MAN	E	11	11/12	0.98	0.04	18,21,24,26	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.

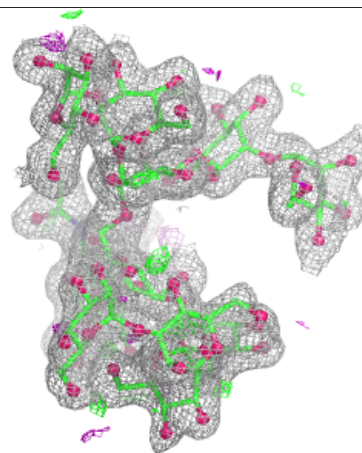
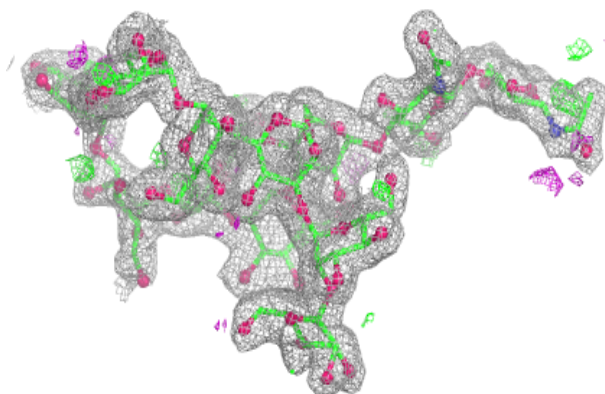
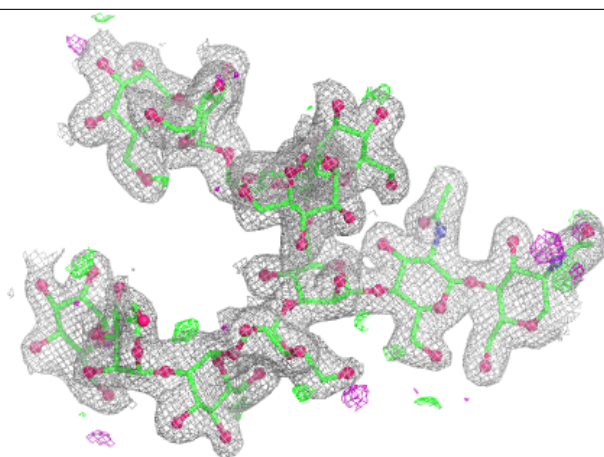
Electron density around Chain C:

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



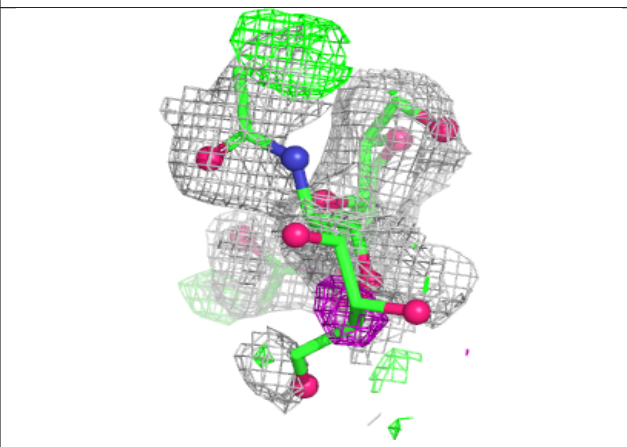
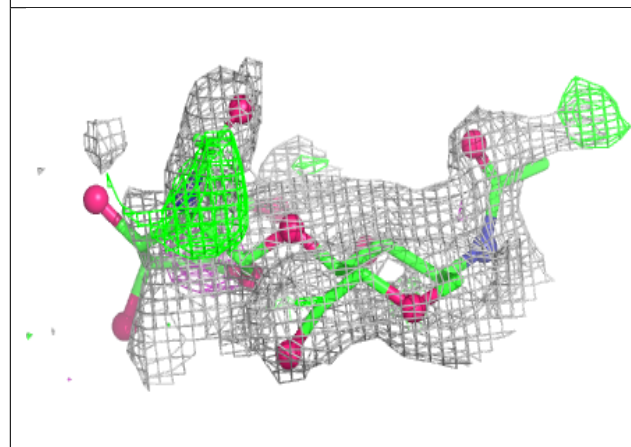
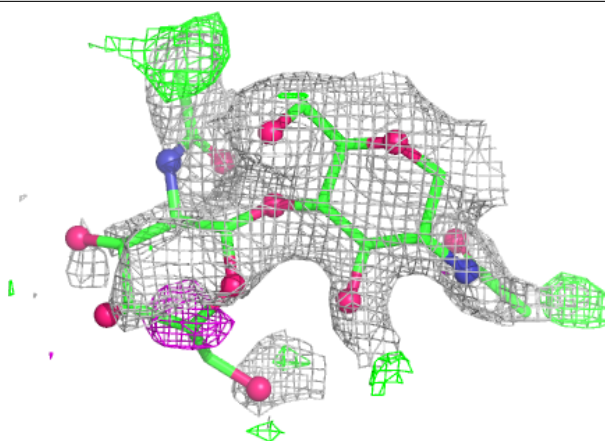
Electron density around Chain E:

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

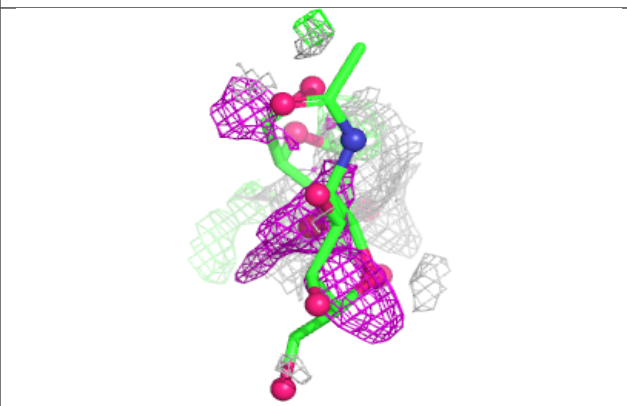
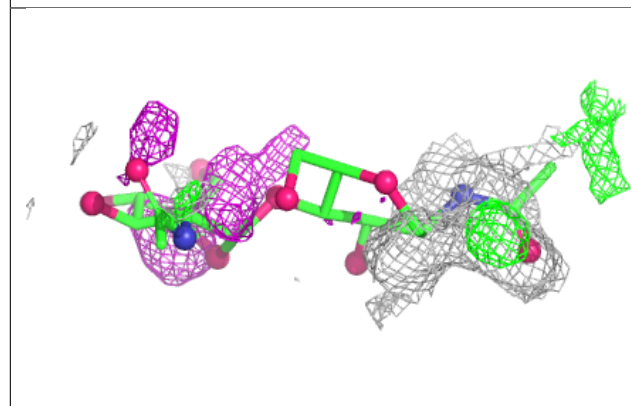
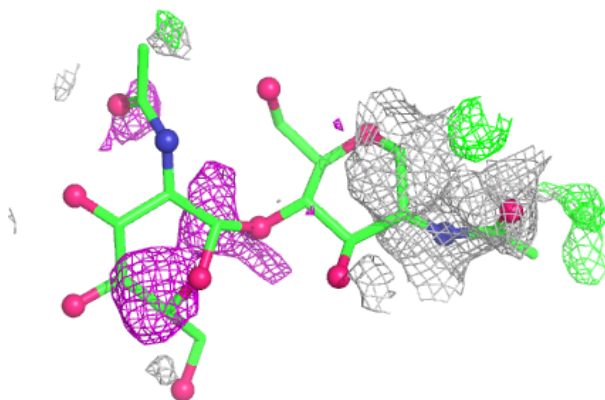


Electron density around Chain D:

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

**Electron density around Chain F:**

$2mF_o-DF_c$ (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($\text{\AA}^2$)	Q<0.9
4	ZN	A	1001	1/1	1.00	0.02	16,16,16,16	0
4	ZN	B	1001	1/1	1.00	0.01	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.