



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 05:37 PM EDT

PDB ID : 3GXP / pdb_00003gxp
Title : Crystal structure of acid-alpha-galactosidase A complexed with galactose at pH 4.5
Authors : Lieberman, R.L.
Deposited on : 2009-04-02
Resolution : 2.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

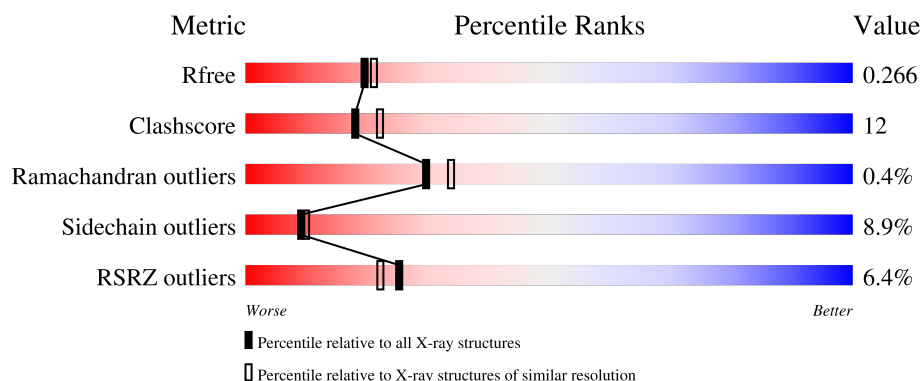
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	398	5% 70% 24% 5%
1	B	398	8% 66% 27% 5%
2	C	3	67% 33%
3	D	2	50% 50%
3	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
3	F	2	 100%
3	G	2	 50% 50%

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	MAN	C	3	X	-	-	-
6	TAM	A	6744	-	X	X	-

2 Entry composition [i](#)

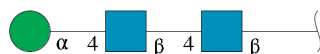
There are 9 unique types of molecules in this entry. The entry contains 6666 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-galactosidase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	0	0
			3122	1988	534	574	26			
1	B	391	Total	C	N	O	S	0	0	0
			3131	1993	536	576	26			

- Molecule 2 is an oligosaccharide called alpha-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	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



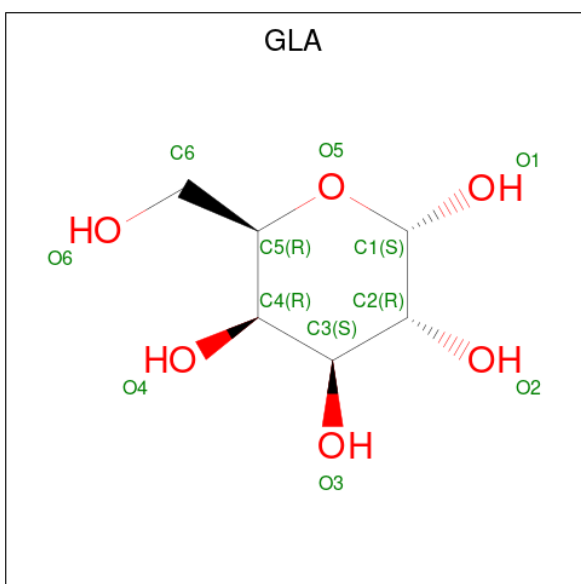
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



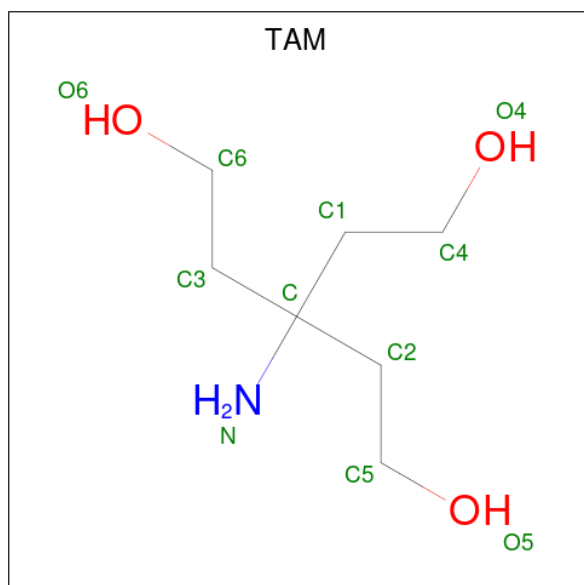
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is alpha-D-galactopyranose (CCD ID: GLA) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			12	6	6		
5	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 6 is TRIS(HYDROXYETHYL)AMINOMETHANE (CCD ID: TAM) (formula: $C_7H_{17}NO_3$).



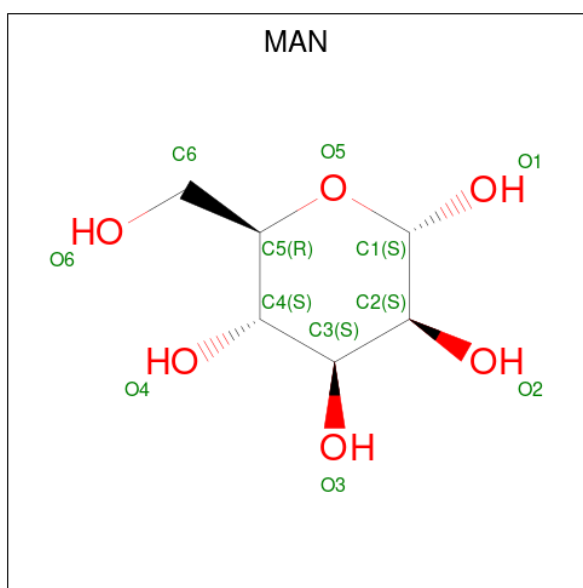
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			11	7	1	3		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			11	6	5		
8	B	1	Total	C	O	0	0
			11	6	5		

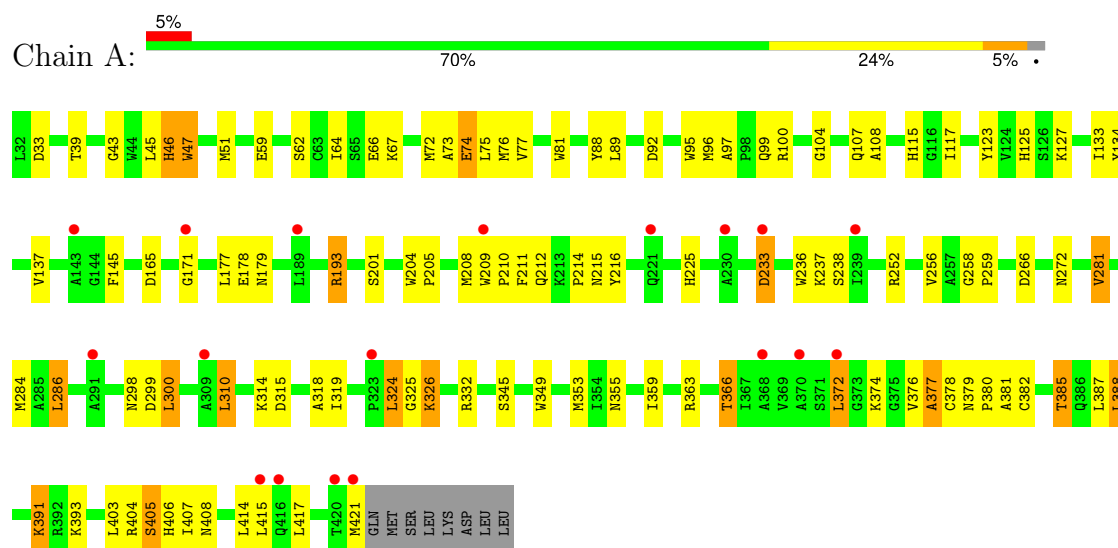
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	86	Total 86	O 86	0	0
9	B	71	Total 71	O 71	0	0

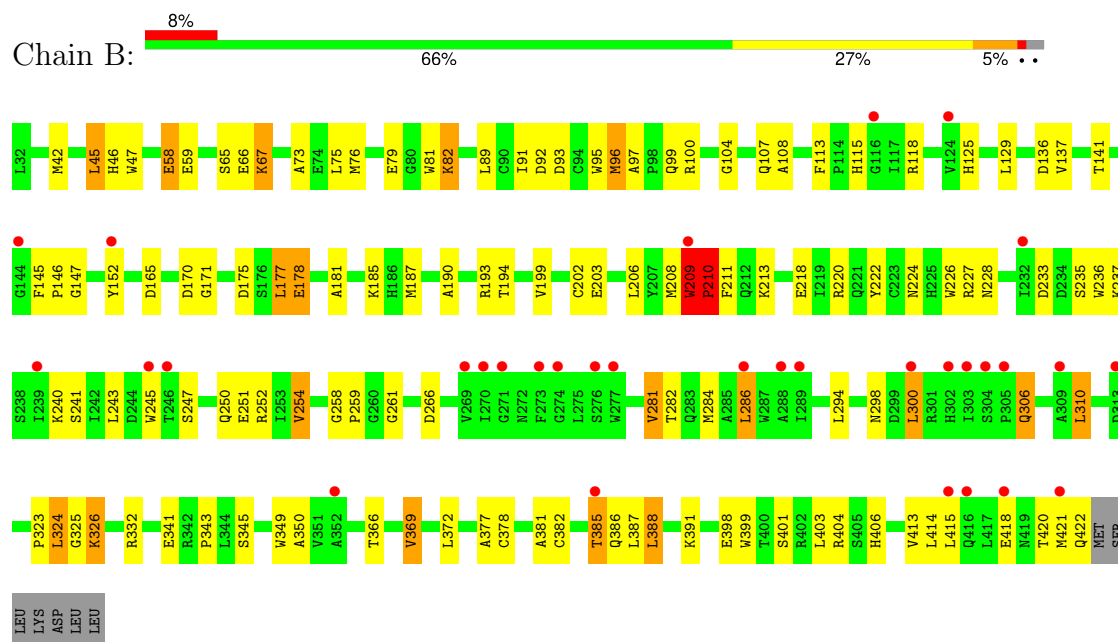
3 Residue-property plots [i](#)

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: Alpha-galactosidase A



• Molecule 1: Alpha-galactosidase A



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

Chain C:  67% 33%

MAG1
MAG2
MAG3

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

Chain D:  50% 50%

MAG1
MAG2

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

Chain E:  50% 50%

MAG1
MAG2

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

Chain F:  100%

MAG1
MAG2

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

Chain G:  50% 50%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	89.51Å 89.51Å 215.93Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.15 – 2.20 38.15 – 2.20	Depositor EDS
% Data completeness (in resolution range)	81.9 (38.15-2.20) 81.9 (38.15-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.201 , 0.265 0.205 , 0.266	Depositor DCC
R_{free} test set	2144 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6666	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, NAG, TAM, SO4, GLA

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	1.14	4/3209 (0.1%)	1.23	13/4358 (0.3%)
1	B	1.14	5/3218 (0.2%)	1.18	9/4370 (0.2%)
All	All	1.14	9/6427 (0.1%)	1.21	22/8728 (0.3%)

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	B	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	209	TRP	CA-C	7.38	1.61	1.53
1	B	281	VAL	CA-CB	6.98	1.63	1.54
1	B	210	PRO	N-CA	-6.33	1.39	1.47
1	B	385	THR	CA-CB	6.05	1.62	1.53
1	A	345	SER	C-O	5.50	1.30	1.23
1	B	282	THR	C-O	-5.41	1.17	1.24
1	B	199	VAL	CA-CB	5.21	1.60	1.54
1	A	318	ALA	C-O	-5.18	1.18	1.24
1	A	407	ILE	CA-CB	5.08	1.60	1.54

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	377	ALA	N-CA-C	-7.77	98.30	109.96
1	A	281	VAL	N-CA-CB	7.38	119.18	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	47	TRP	N-CA-C	7.08	118.64	111.07
1	A	39	THR	CA-C-N	-7.00	112.76	119.90
1	A	39	THR	C-N-CA	-7.00	112.76	119.90
1	A	47	TRP	N-CA-C	6.99	119.49	111.11
1	A	286	LEU	CA-CB-CG	-6.89	92.19	116.30
1	B	281	VAL	N-CA-CB	6.86	119.87	110.54
1	A	377	ALA	CA-C-N	-6.45	113.72	125.66
1	A	377	ALA	C-N-CA	-6.45	113.72	125.66
1	B	99	GLN	N-CA-C	6.26	116.54	108.34
1	B	413	VAL	N-CA-C	5.56	116.11	107.99
1	B	79	GLU	N-CA-C	5.55	119.75	112.92
1	B	211	PHE	N-CA-C	-5.52	105.34	111.36
1	B	209	TRP	CA-C-O	5.43	124.85	119.75
1	A	319	ILE	N-CA-C	5.39	115.53	110.30
1	A	46	HIS	N-CA-C	5.36	118.04	111.82
1	B	323	PRO	N-CA-C	5.33	120.52	113.86
1	A	233	ASP	CB-CA-C	-5.31	100.15	109.71
1	A	145	PHE	N-CA-C	-5.20	104.07	110.31
1	B	286	LEU	CA-CB-CG	-5.11	98.41	116.30
1	A	393	LYS	N-CA-C	-5.04	102.40	109.96

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	209	TRP	Peptide
1	B	210	PRO	Peptide

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	3122	0	2982	70	0
1	B	3131	0	2987	82	0
2	C	39	0	34	1	0
3	D	28	0	25	3	0
3	E	28	0	24	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	28	0	25	1	0
3	G	28	0	22	1	0
4	A	5	0	0	0	0
4	B	15	0	0	0	0
5	A	12	0	12	0	0
5	B	12	0	12	4	0
6	A	11	0	17	11	0
7	A	28	0	25	2	0
8	B	22	0	20	0	0
9	A	86	0	0	6	0
9	B	71	0	0	2	0
All	All	6666	0	6185	155	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 (155) 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:215:ASN:ND2	3:D:1:NAG:C1	2.09	1.15
1:A:215:ASN:HD21	3:D:1:NAG:C1	1.72	0.99
1:A:236:TRP:H	6:A:6744:TAM:H52	1.31	0.94
6:A:6744:TAM:H22	1:B:236:TRP:HB3	1.47	0.94
1:A:66:GLU:OE1	1:A:115:HIS:HD2	1.51	0.93
1:B:170:ASP:OD2	5:B:3681:GLA:H1	1.73	0.86
1:B:65:SER:HB2	1:B:67:LYS:NZ	1.90	0.86
1:A:66:GLU:OE1	1:A:115:HIS:CD2	2.32	0.81
1:B:324:LEU:HD13	1:B:326:LYS:HG3	1.63	0.81
1:B:65:SER:HB2	1:B:67:LYS:HZ3	1.46	0.78
1:B:125:HIS:HE1	1:B:165:ASP:OD2	1.65	0.77
1:A:377:ALA:O	1:A:378:CYS:HB2	1.84	0.76
1:A:379:ASN:HD22	1:A:380:PRO:HA	1.49	0.76
1:A:237:LYS:H	6:A:6744:TAM:HN1	1.32	0.73
1:A:208:MET:CE	1:A:212:GLN:HG3	2.19	0.73
1:B:209:TRP:HE3	1:B:210:PRO:HG3	1.51	0.72
1:B:66:GLU:OE1	1:B:115:HIS:HD2	1.73	0.72
1:A:237:LYS:N	6:A:6744:TAM:N	2.36	0.71
1:A:51:MET:O	1:A:64:ILE:HD11	1.91	0.71
1:A:236:TRP:N	6:A:6744:TAM:H52	2.03	0.70
1:B:306:GLN:H	1:B:306:GLN:HE21	1.37	0.70
1:A:46:HIS:CD2	1:A:92:ASP:H	2.11	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:LEU:H	1:B:177:LEU:HD12	1.59	0.68
1:B:46:HIS:CD2	1:B:92:ASP:H	2.12	0.67
1:B:208:MET:O	1:B:210:PRO:HA	1.96	0.66
1:B:236:TRP:CE2	1:B:240:LYS:HD2	2.31	0.66
1:A:97:ALA:HB3	1:A:107:GLN:HG3	1.77	0.66
1:B:306:GLN:HE21	1:B:306:GLN:N	1.94	0.65
1:A:125:HIS:HE1	1:A:165:ASP:OD2	1.80	0.65
1:A:193:ARG:HG3	1:A:193:ARG:O	1.96	0.65
1:A:377:ALA:O	1:A:378:CYS:CB	2.42	0.65
1:B:45:LEU:HD22	1:B:92:ASP:HB2	1.80	0.64
1:A:372:LEU:HD23	1:A:417:LEU:HD21	1.81	0.62
1:B:213:LYS:HD3	1:B:252:ARG:NH1	2.14	0.62
1:B:349:TRP:CD1	1:B:377:ALA:HB2	2.36	0.61
1:B:181:ALA:O	1:B:185:LYS:HG3	2.01	0.61
1:A:349:TRP:CD1	1:A:377:ALA:HB2	2.37	0.60
1:A:237:LYS:H	6:A:6744:TAM:HN2	1.46	0.60
1:A:406:HIS:HE1	1:B:59:GLU:OE2	1.85	0.60
1:A:377:ALA:O	9:A:3476:HOH:O	2.17	0.60
9:B:6763:HOH:O	3:G:2:NAG:H2	2.02	0.60
1:B:326:LYS:HB3	1:B:326:LYS:NZ	2.18	0.59
7:A:498:NAG:H4	7:A:430:NAG:C1	2.32	0.59
1:A:403:LEU:HD21	1:A:415:LEU:CD1	2.33	0.59
1:B:388:LEU:HB2	1:B:414:LEU:HB3	1.85	0.58
6:A:6744:TAM:H32	1:B:237:LYS:HB2	1.85	0.58
1:A:72:MET:SD	1:A:300:LEU:HB2	2.43	0.58
1:B:306:GLN:H	1:B:306:GLN:NE2	2.02	0.58
1:A:100:ARG:HD3	1:A:104:GLY:O	2.04	0.57
1:A:379:ASN:HD22	1:A:380:PRO:CA	2.17	0.57
1:A:236:TRP:H	6:A:6744:TAM:C5	2.10	0.57
1:A:385:THR:HG23	9:A:3471:HOH:O	2.05	0.56
1:B:137:VAL:HG12	1:B:171:GLY:HA2	1.86	0.56
1:B:236:TRP:NE1	1:B:240:LYS:HD2	2.20	0.56
1:A:233:ASP:HB2	1:A:238:SER:HB2	1.86	0.56
1:A:212:GLN:HB3	9:A:3474:HOH:O	2.06	0.56
1:B:209:TRP:CE3	1:B:210:PRO:HG3	2.37	0.56
1:A:272:ASN:HD21	1:A:298:ASN:HA	1.71	0.56
1:B:82:LYS:HB2	1:B:129:LEU:HD21	1.87	0.55
1:B:284:MET:HG2	1:B:310:LEU:HD22	1.89	0.55
1:A:403:LEU:HD21	1:A:415:LEU:HD13	1.89	0.55
1:A:324:LEU:HD13	1:A:326:LYS:HG3	1.89	0.54
3:F:1:NAG:H61	3:F:2:NAG:H82	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:TRP:HB3	1:A:205:PRO:HD3	1.89	0.54
1:A:421:MET:HA	9:A:3415:HOH:O	2.07	0.54
1:B:378:CYS:HA	1:B:381:ALA:C	2.34	0.53
1:A:388:LEU:HB2	1:A:414:LEU:HB3	1.90	0.53
1:B:95:TRP:CZ2	1:B:96:MET:HE2	2.43	0.53
1:B:136:ASP:CB	1:B:141:THR:HA	2.39	0.53
1:B:228:ASN:HB3	1:B:245:TRP:CH2	2.42	0.53
1:B:381:ALA:HB3	1:B:422:GLN:HB3	1.90	0.52
1:A:72:MET:O	1:A:76:MET:HG3	2.08	0.52
1:B:243:LEU:O	1:B:247:SER:HB2	2.09	0.51
1:A:214:PRO:HG2	1:A:216:TYR:CE2	2.46	0.51
1:A:74:GLU:OE2	1:A:123:TYR:OH	2.29	0.51
1:B:66:GLU:OE1	1:B:115:HIS:CD2	2.60	0.50
1:B:152:TYR:HD1	1:B:187:MET:HE2	1.76	0.50
1:B:341:GLU:HA	1:B:350:ALA:O	2.11	0.50
1:B:42:MET:HE3	1:B:294:LEU:HD12	1.94	0.50
1:A:210:PRO:HD2	1:A:211:PHE:CD2	2.47	0.49
1:B:250:GLN:O	1:B:254:VAL:HG13	2.12	0.49
1:A:117:ILE:HG21	1:A:133:ILE:HD12	1.95	0.49
1:B:224:ASN:O	1:B:261:GLY:HA2	2.13	0.49
1:A:237:LYS:N	6:A:6744:TAM:HN2	2.05	0.49
1:B:404:ARG:HH21	1:B:406:HIS:HE1	1.59	0.49
1:B:76:MET:HA	1:B:81:TRP:HB2	1.94	0.49
1:B:108:ALA:HB1	1:B:113:PHE:HB2	1.95	0.48
7:A:498:NAG:C4	7:A:430:NAG:C1	2.92	0.48
1:B:203:GLU:OE2	5:B:3681:GLA:H2	2.14	0.48
9:A:3405:HOH:O	1:B:58:GLU:HG3	2.14	0.48
1:B:136:ASP:HB2	1:B:141:THR:HA	1.96	0.48
1:B:203:GLU:HA	1:B:227:ARG:HB2	1.96	0.48
1:B:404:ARG:HH21	1:B:406:HIS:CE1	2.32	0.48
1:A:108:ALA:HB2	1:A:117:ILE:HG12	1.95	0.48
1:A:378:CYS:HA	1:A:382:CYS:HB3	1.96	0.47
1:A:315:ASP:OD2	1:A:391:LYS:HE2	2.13	0.47
1:A:359:ILE:HD12	1:B:233:ASP:HB3	1.96	0.47
1:B:210:PRO:N	9:B:6761:HOH:O	2.46	0.47
1:B:403:LEU:HD21	1:B:415:LEU:HD11	1.97	0.47
1:A:77:VAL:HG21	1:A:127:LYS:HB3	1.95	0.47
1:B:65:SER:HB2	1:B:67:LYS:HZ2	1.75	0.47
1:B:218:GLU:O	1:B:222:TYR:HD2	1.98	0.47
1:A:73:ALA:HB2	1:A:89:LEU:HD22	1.96	0.46
1:A:137:VAL:HG12	1:A:171:GLY:HA2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:MET:HG2	1:A:310:LEU:HD22	1.98	0.46
1:A:134:TYR:CD2	1:A:134:TYR:C	2.94	0.46
6:A:6744:TAM:H62	6:A:6744:TAM:H21	1.70	0.46
1:B:46:HIS:HD2	1:B:92:ASP:H	1.61	0.46
1:A:379:ASN:ND2	1:A:380:PRO:HA	2.26	0.45
1:B:220:ARG:NH2	1:B:259:PRO:O	2.49	0.45
1:A:67:LYS:HA	1:A:67:LYS:HE3	1.98	0.45
1:A:208:MET:HE2	1:A:208:MET:HB3	1.79	0.45
1:B:145:PHE:HB3	1:B:146:PRO:CD	2.46	0.45
1:A:378:CYS:HA	1:A:381:ALA:C	2.41	0.45
1:B:178:GLU:H	1:B:178:GLU:HG2	1.49	0.45
1:A:95:TRP:CE2	1:A:96:MET:HG3	2.52	0.45
1:B:45:LEU:CD2	1:B:92:ASP:HB2	2.45	0.45
1:B:298:ASN:OD1	1:B:300:LEU:HD13	2.17	0.45
1:B:175:ASP:N	3:E:1:NAG:O7	2.50	0.44
1:B:213:LYS:HD3	1:B:252:ARG:CZ	2.47	0.44
1:B:202:CYS:O	1:B:226:TRP:HA	2.17	0.44
1:B:170:ASP:OD2	5:B:3681:GLA:C1	2.57	0.44
1:B:190:ALA:O	1:B:194:THR:HG23	2.18	0.44
1:B:258:GLY:HA2	1:B:325:GLY:O	2.17	0.44
1:A:215:ASN:HD22	3:D:1:NAG:C1	2.17	0.44
1:B:250:GLN:HB2	1:B:254:VAL:CG1	2.48	0.44
1:B:100:ARG:NH2	1:B:147:GLY:O	2.50	0.43
1:A:43:GLY:HA3	1:A:88:TYR:O	2.17	0.43
1:B:97:ALA:HB3	1:B:107:GLN:HG3	2.01	0.43
1:A:201:SER:HB2	1:A:225:HIS:CE1	2.54	0.43
1:B:100:ARG:HD2	1:B:104:GLY:O	2.18	0.43
1:B:369:VAL:HB	1:B:401:SER:O	2.17	0.43
1:A:76:MET:HA	1:A:81:TRP:HB2	2.00	0.43
1:B:386:GLN:HA	1:B:415:LEU:HD23	2.01	0.43
1:B:326:LYS:HB2	1:B:343:PRO:HG2	2.01	0.42
1:B:382:CYS:HA	1:B:418:GLU:O	2.19	0.42
1:B:177:LEU:H	1:B:177:LEU:CD1	2.30	0.42
1:B:91:ILE:HG22	1:B:92:ASP:O	2.19	0.42
1:B:398:GLU:O	1:B:399:TRP:C	2.62	0.42
1:A:355:ASN:HB3	1:A:408:ASN:O	2.20	0.42
1:A:33:ASP:O	2:C:1:NAG:H4	2.19	0.42
1:A:252:ARG:NH2	9:A:6748:HOH:O	2.51	0.41
1:A:353:MET:HE1	1:A:405:SER:HB2	2.01	0.41
1:B:73:ALA:HB2	1:B:89:LEU:HD22	2.02	0.41
1:B:125:HIS:CE1	1:B:165:ASP:OD2	2.58	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:6744:TAM:H61	1:B:235:SER:HB2	2.01	0.41
1:A:366:THR:HG22	1:A:404:ARG:HG3	2.01	0.41
1:A:258:GLY:HA2	1:A:325:GLY:O	2.21	0.41
1:A:258:GLY:O	1:A:259:PRO:C	2.62	0.41
1:A:177:LEU:HD11	1:A:211:PHE:HB3	2.03	0.41
1:B:93:ASP:HB2	5:B:3681:GLA:C6	2.51	0.41
1:B:145:PHE:HB3	1:B:146:PRO:HD2	2.04	0.40
1:B:203:GLU:O	1:B:206:LEU:HB3	2.21	0.40
1:A:72:MET:HE1	1:A:299:ASP:OD1	2.21	0.40
1:A:47:TRP:HB2	1:A:92:ASP:CG	2.46	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	388/398 (98%)	370 (95%)	17 (4%)	1 (0%)	36	42
1	B	389/398 (98%)	369 (95%)	18 (5%)	2 (0%)	24	27
All	All	777/796 (98%)	739 (95%)	35 (4%)	3 (0%)	30	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	ASP
1	B	210	PRO
1	B	266	ASP

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	331/339 (98%)	303 (92%)	28 (8%)	10	11
1	B	332/339 (98%)	301 (91%)	31 (9%)	8	9
All	All	663/678 (98%)	604 (91%)	59 (9%)	9	10

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LEU
1	A	59	GLU
1	A	62	SER
1	A	74	GLU
1	A	75	LEU
1	A	99	GLN
1	A	178	GLU
1	A	179	ASN
1	A	193	ARG
1	A	256	VAL
1	A	281	VAL
1	A	286	LEU
1	A	300	LEU
1	A	310	LEU
1	A	314	LYS
1	A	324	LEU
1	A	326	LYS
1	A	332	ARG
1	A	363	ARG
1	A	366	THR
1	A	372	LEU
1	A	374	LYS
1	A	376	VAL
1	A	385	THR
1	A	387	LEU
1	A	388	LEU
1	A	391	LYS

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Mol	Chain	Res	Type
1	A	405	SER
1	B	45	LEU
1	B	58	GLU
1	B	67	LYS
1	B	75	LEU
1	B	82	LYS
1	B	96	MET
1	B	118	ARG
1	B	177	LEU
1	B	178	GLU
1	B	193	ARG
1	B	241	SER
1	B	251	GLU
1	B	254	VAL
1	B	281	VAL
1	B	286	LEU
1	B	300	LEU
1	B	306	GLN
1	B	310	LEU
1	B	324	LEU
1	B	326	LYS
1	B	332	ARG
1	B	345	SER
1	B	366	THR
1	B	369	VAL
1	B	372	LEU
1	B	385	THR
1	B	387	LEU
1	B	388	LEU
1	B	391	LYS
1	B	420	THR
1	B	421	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	HIS
1	A	115	HIS
1	A	122	ASN
1	A	125	HIS
1	A	157	GLN
1	A	179	ASN

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Mol	Chain	Res	Type
1	A	212	GLN
1	A	221	GLN
1	A	272	ASN
1	A	379	ASN
1	A	406	HIS
1	A	408	ASN
1	A	416	GLN
1	B	46	HIS
1	B	53	ASN
1	B	57	GLN
1	B	115	HIS
1	B	125	HIS
1	B	306	GLN
1	B	406	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

11 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	2,1	14,14,15	0.52	0	17,19,21	2.01	3 (17%)
2	NAG	C	2	2	14,14,15	0.81	0	17,19,21	1.92	5 (29%)
2	MAN	C	3	2	11,11,12	0.86	0	15,15,17	1.20	1 (6%)
3	NAG	D	1	3	14,14,15	0.52	0	17,19,21	2.39	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	D	2	3	14,14,15	0.84	1 (7%)	17,19,21	1.30	2 (11%)
3	NAG	E	1	3,1	14,14,15	0.56	0	17,19,21	2.42	4 (23%)
3	NAG	E	2	3	14,14,15	0.98	1 (7%)	17,19,21	1.34	3 (17%)
3	NAG	F	1	3,1	14,14,15	0.65	0	17,19,21	1.41	3 (17%)
3	NAG	F	2	3	14,14,15	0.73	0	17,19,21	1.72	3 (17%)
3	NAG	G	1	3,1	14,14,15	0.42	0	17,19,21	1.70	4 (23%)
3	NAG	G	2	3	14,14,15	0.72	1 (7%)	17,19,21	0.97	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
2	NAG	C	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
2	MAN	C	3	2	1/1/4/5	2/2/19/22	0/1/1/1
3	NAG	D	1	3	-	5/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	0/1/1/1
3	NAG	E	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	1/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	3/6/23/26	0/1/1/1
3	NAG	G	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	2	NAG	C1-C2	3.13	1.56	1.52
3	G	2	NAG	C1-C2	2.14	1.55	1.52
3	D	2	NAG	C1-C2	2.01	1.55	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	NAG	C1-O5-C5	8.17	123.13	112.19
3	D	1	NAG	C2-N2-C7	7.02	132.31	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	6.30	120.64	112.19
3	F	2	NAG	C1-O5-C5	5.15	119.08	112.19
2	C	2	NAG	C1-O5-C5	4.57	118.31	112.19
3	D	1	NAG	C1-O5-C5	4.56	118.30	112.19
3	D	2	NAG	C4-C3-C2	3.98	116.85	111.02
3	G	1	NAG	C2-N2-C7	-3.72	117.91	122.90
3	D	1	NAG	O5-C1-C2	3.14	116.14	111.29
3	E	2	NAG	O5-C1-C2	3.06	116.03	111.29
2	C	2	NAG	C1-C2-N2	3.05	115.23	110.43
2	C	3	MAN	C1-C2-C3	3.00	114.01	109.64
3	F	2	NAG	C2-N2-C7	-2.85	119.08	122.90
2	C	2	NAG	O4-C4-C5	2.78	116.17	109.32
3	F	2	NAG	O5-C5-C6	-2.76	102.29	107.66
3	G	1	NAG	O5-C1-C2	-2.58	107.30	111.29
3	D	2	NAG	C3-C4-C5	2.50	114.76	110.23
2	C	1	NAG	C4-C3-C2	2.43	114.58	111.02
3	E	1	NAG	C2-N2-C7	2.41	126.13	122.90
3	G	1	NAG	C3-C4-C5	-2.40	105.88	110.23
3	F	1	NAG	O5-C1-C2	-2.35	107.66	111.29
2	C	1	NAG	O5-C5-C6	2.31	112.17	107.66
3	F	1	NAG	C1-O5-C5	2.27	115.23	112.19
3	E	1	NAG	O5-C5-C4	2.25	116.31	110.83
3	D	1	NAG	C8-C7-N2	2.23	119.82	116.12
2	C	2	NAG	O7-C7-C8	-2.15	118.23	122.05
3	E	2	NAG	O4-C4-C5	2.13	114.57	109.32
2	C	2	NAG	O7-C7-N2	2.12	125.72	121.98
3	E	2	NAG	O5-C5-C4	-2.09	105.75	110.83
3	F	1	NAG	C2-N2-C7	-2.08	120.11	122.90
3	G	1	NAG	C1-O5-C5	2.04	114.92	112.19
3	E	1	NAG	C1-C2-N2	-2.00	107.27	110.43

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	3	MAN	C1

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1	NAG	C1-C2-N2-C7
3	D	1	NAG	C8-C7-N2-C2
3	D	1	NAG	O7-C7-N2-C2

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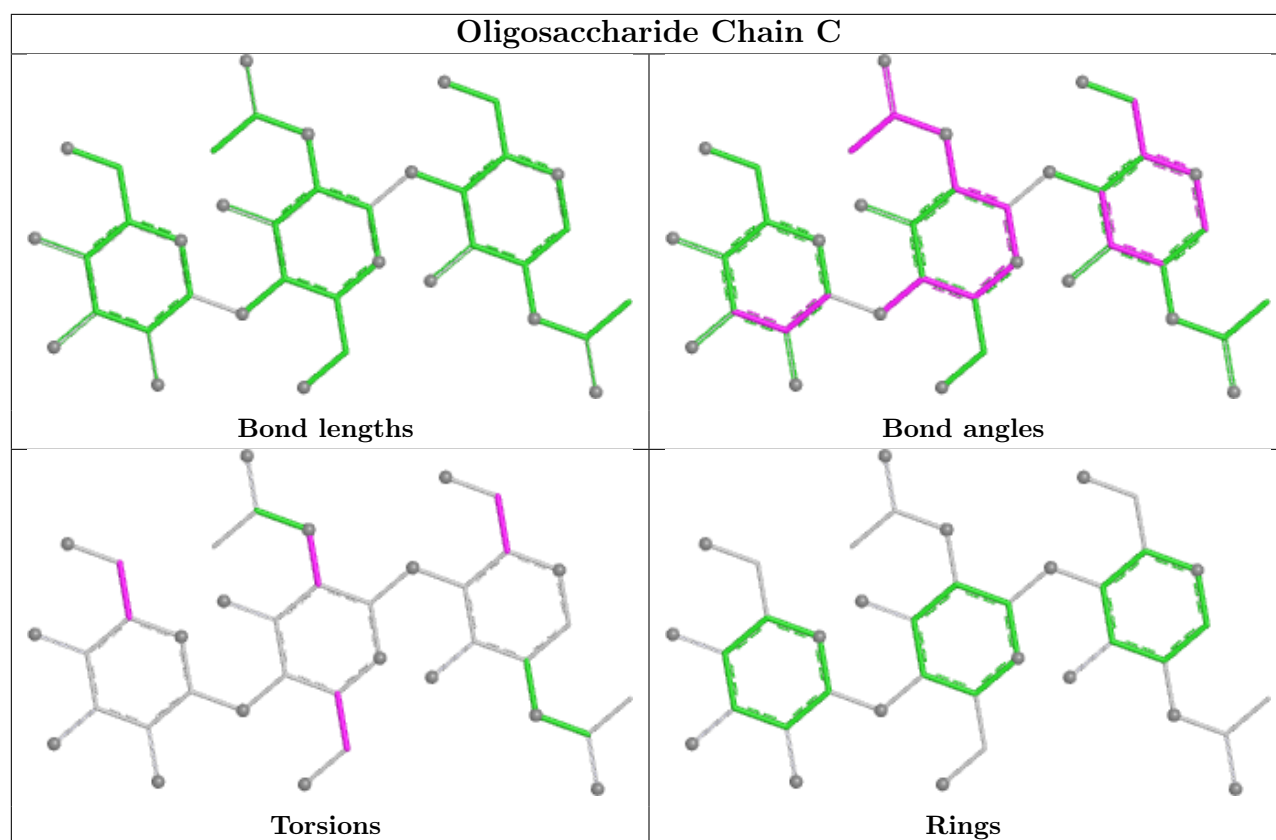
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	O5-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	F	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O5-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
2	C	3	MAN	O5-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
3	F	2	NAG	O7-C7-N2-C2
3	F	1	NAG	C4-C5-C6-O6
2	C	3	MAN	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
2	C	2	NAG	C1-C2-N2-C7
3	D	2	NAG	C8-C7-N2-C2
3	D	2	NAG	O7-C7-N2-C2
3	E	2	NAG	C4-C5-C6-O6

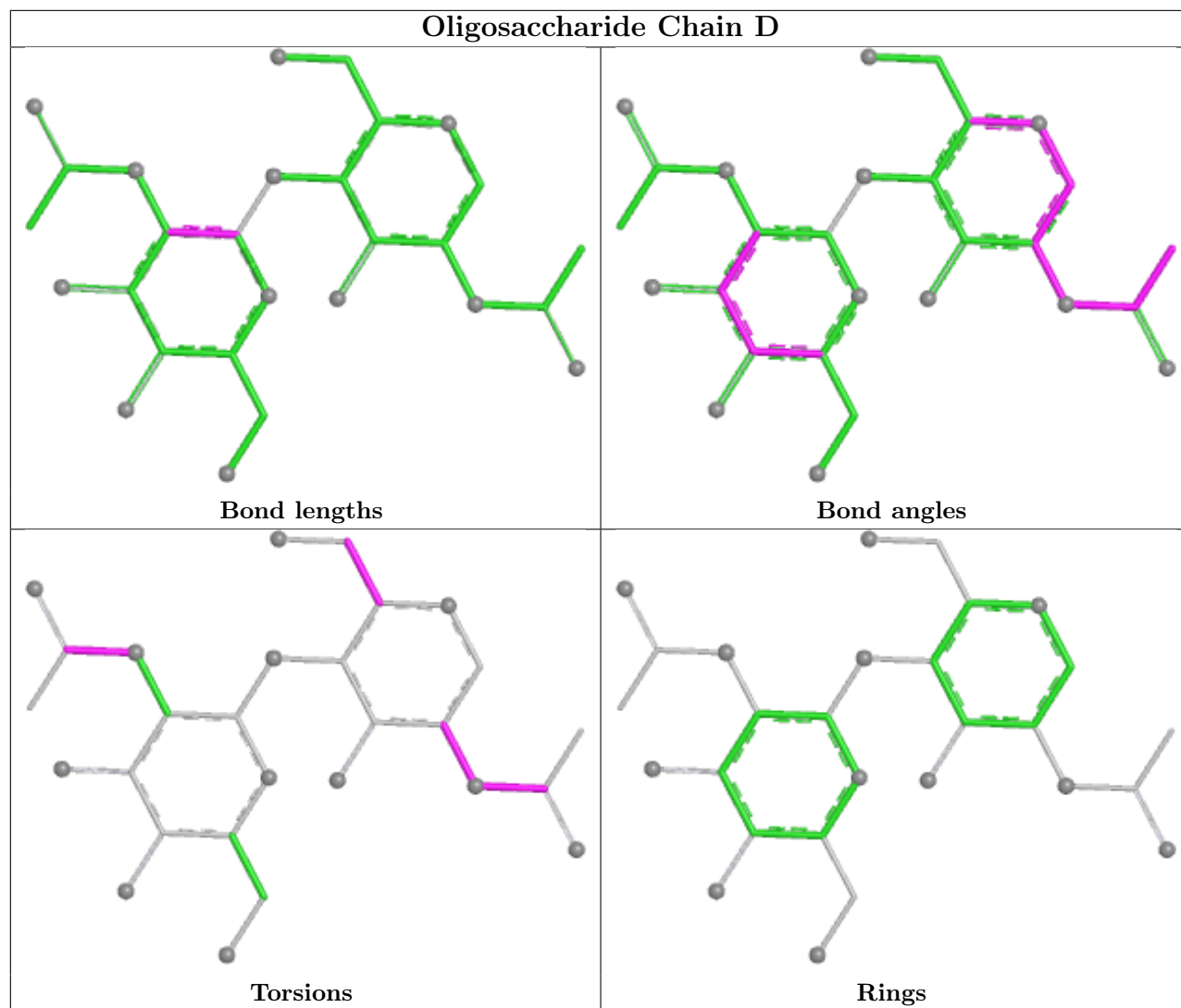
There are no ring outliers.

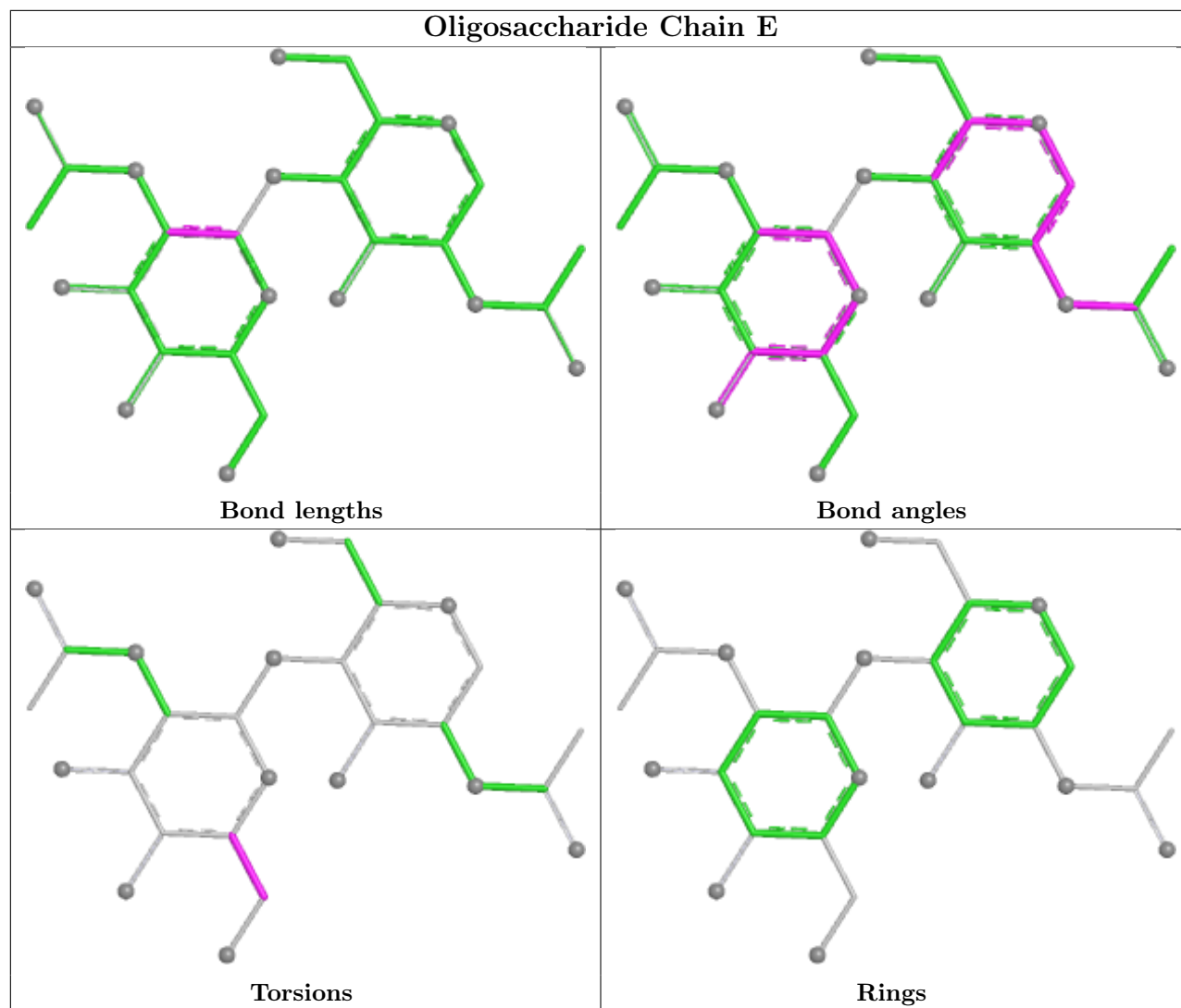
6 monomers are involved in 7 short contacts:

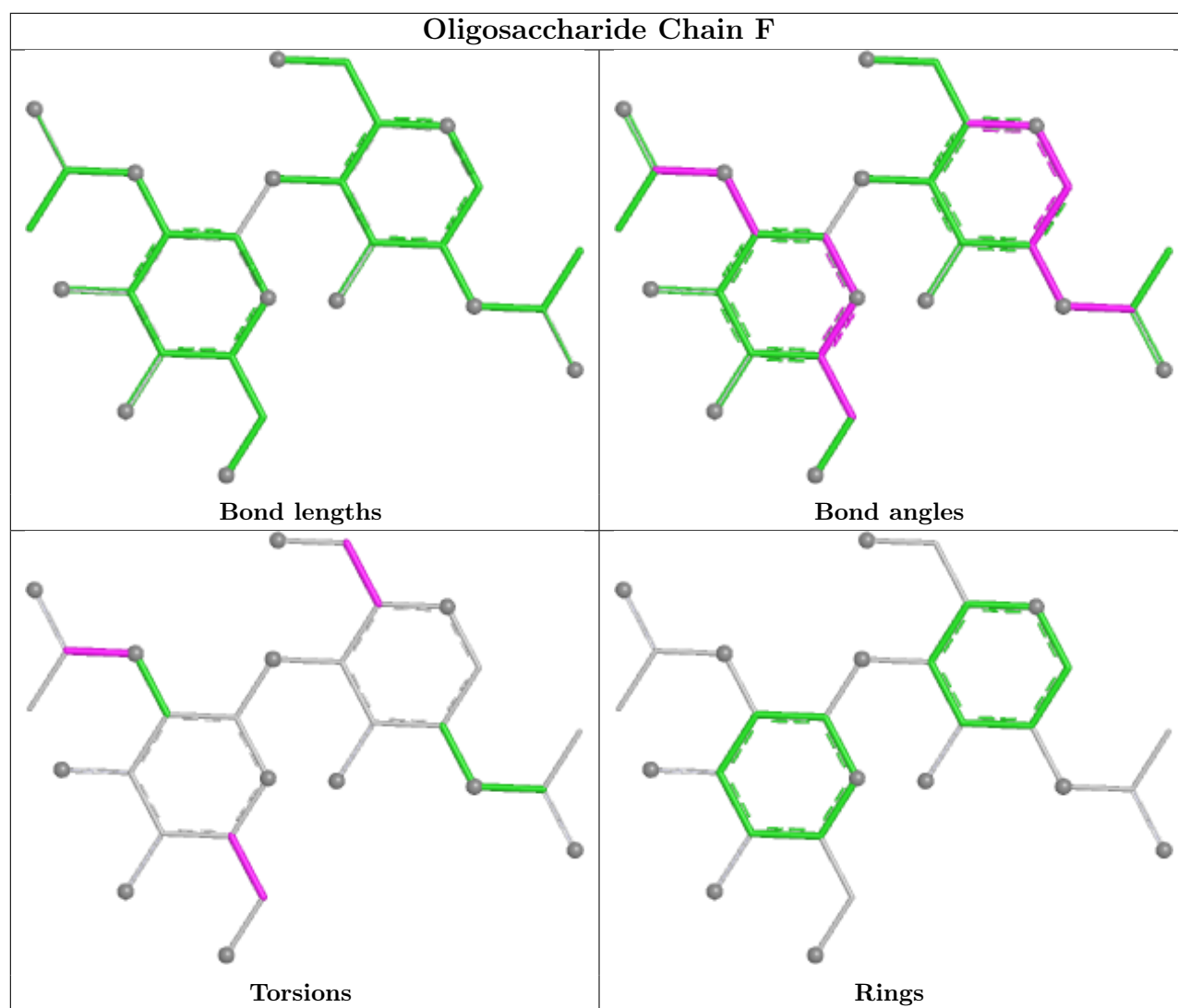
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	2	NAG	1	0
3	F	1	NAG	1	0
3	E	1	NAG	1	0
2	C	1	NAG	1	0
3	G	2	NAG	1	0
3	D	1	NAG	3	0

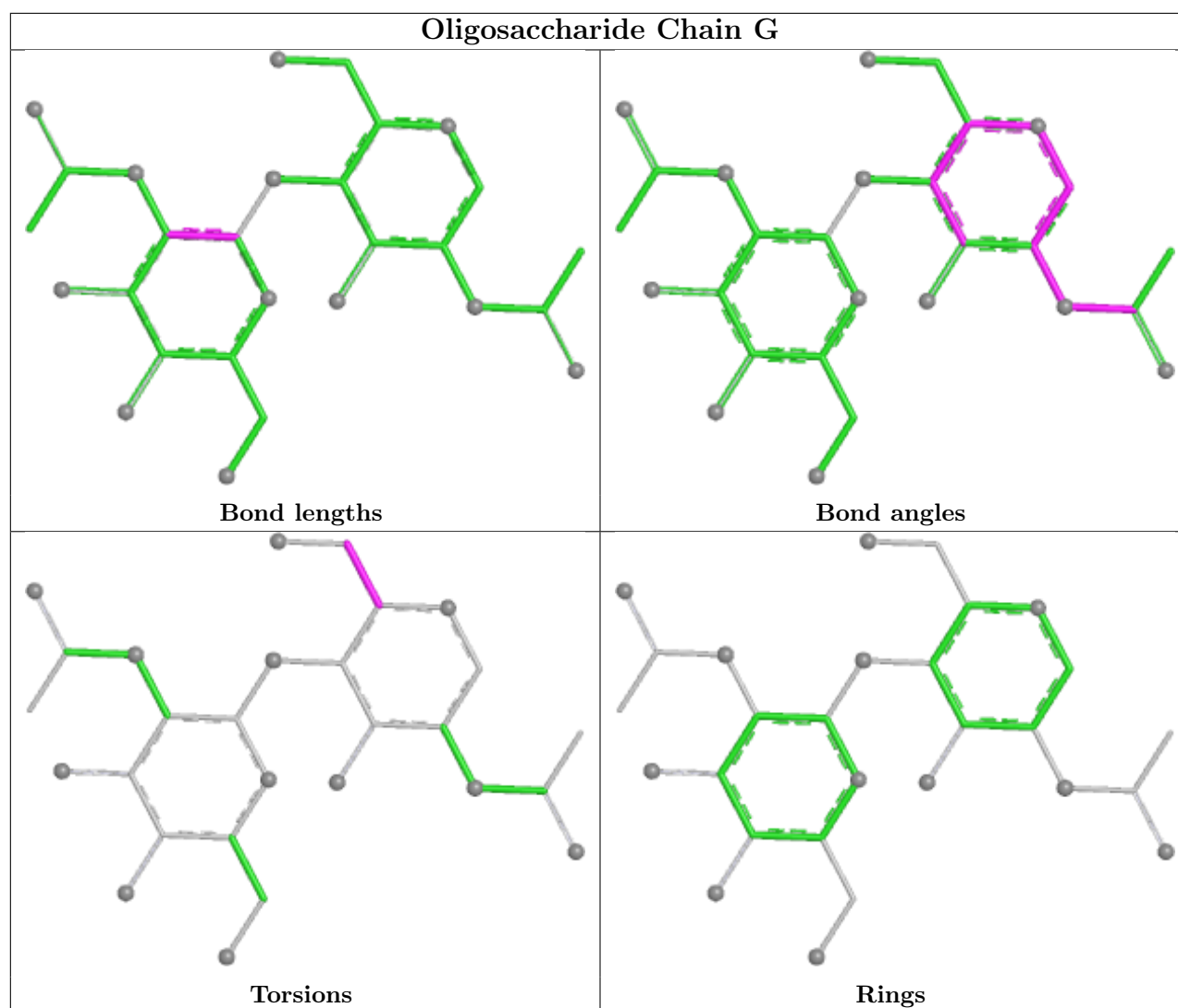
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

11 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	3	-	4,4,4	0.35	0	6,6,6	0.57	0
8	MAN	B	695	-	11,11,12	0.77	1 (9%)	15,15,17	1.56	4 (26%)
5	GLA	A	3681	-	12,12,12	1.39	1 (8%)	17,17,17	1.96	7 (41%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GLA	B	3681	-	12,12,12	0.77	0	17,17,17	1.86	5 (29%)
6	TAM	A	6744	-	10,10,10	2.90	5 (50%)	12,12,12	3.80	5 (41%)
7	NAG	A	430	-	14,14,15	0.59	0	17,19,21	1.64	5 (29%)
8	MAN	B	694	-	11,11,12	0.65	0	15,15,17	1.61	3 (20%)
7	NAG	A	498	1	14,14,15	1.04	1 (7%)	17,19,21	1.71	4 (23%)
4	SO4	A	1	-	4,4,4	0.26	0	6,6,6	1.07	0
4	SO4	B	2	-	4,4,4	0.17	0	6,6,6	0.69	0
4	SO4	B	4	-	4,4,4	0.28	0	6,6,6	0.34	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
8	MAN	B	695	-	-	2/2/19/22	0/1/1/1
5	GLA	A	3681	-	-	0/2/22/22	0/1/1/1
6	TAM	A	6744	-	-	6/12/12/12	-
5	GLA	B	3681	-	-	0/2/22/22	0/1/1/1
7	NAG	A	430	-	-	4/6/23/26	0/1/1/1
8	MAN	B	694	-	-	1/2/19/22	0/1/1/1
7	NAG	A	498	1	-	1/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	6744	TAM	C2-C	-4.62	1.47	1.53
6	A	6744	TAM	C1-C	4.27	1.59	1.53
6	A	6744	TAM	C-N	-4.00	1.36	1.49
5	A	3681	GLA	O5-C5	-3.82	1.35	1.44
6	A	6744	TAM	C3-C	-3.81	1.48	1.53
6	A	6744	TAM	C1-C4	2.94	1.57	1.52
7	A	498	NAG	C1-C2	2.22	1.55	1.52
8	B	695	MAN	C2-C3	2.11	1.55	1.52

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	6744	TAM	C2-C-C1	7.07	122.97	110.50
6	A	6744	TAM	C2-C-N	-6.39	92.19	108.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	6744	TAM	C5-C2-C	-5.65	109.31	115.97
6	A	6744	TAM	C4-C1-C	-5.25	109.78	115.97
8	B	694	MAN	C1-C2-C3	4.65	116.41	109.64
5	B	3681	GLA	O5-C1-C2	4.48	118.18	110.30
7	A	498	NAG	C3-C4-C5	4.13	117.73	110.23
6	A	6744	TAM	O6-C6-C3	-3.97	100.19	111.33
5	B	3681	GLA	C1-O5-C5	3.84	121.08	113.65
5	A	3681	GLA	O2-C2-C3	3.44	118.49	110.38
5	A	3681	GLA	O1-C1-C2	3.28	118.49	108.98
5	B	3681	GLA	O1-C1-O5	-2.92	101.72	110.41
7	A	430	NAG	C1-C2-N2	-2.90	105.87	110.43
7	A	430	NAG	O5-C1-C2	2.88	115.75	111.29
7	A	498	NAG	O4-C4-C3	-2.87	103.61	110.38
5	A	3681	GLA	C4-C3-C2	-2.85	105.83	110.83
8	B	694	MAN	C2-C3-C4	2.71	115.62	110.86
7	A	498	NAG	C1-C2-N2	2.66	114.63	110.43
5	A	3681	GLA	C1-O5-C5	2.66	118.79	113.65
7	A	430	NAG	C1-O5-C5	2.55	115.60	112.19
5	A	3681	GLA	O1-C1-O5	-2.49	103.00	110.41
8	B	695	MAN	O5-C5-C4	-2.46	104.84	110.83
8	B	695	MAN	C1-C2-C3	2.37	113.10	109.64
5	A	3681	GLA	O3-C3-C2	2.35	115.92	110.38
5	A	3681	GLA	O5-C5-C6	-2.31	100.72	106.44
8	B	695	MAN	O3-C3-C2	2.31	114.76	110.05
7	A	430	NAG	O5-C5-C4	-2.26	105.34	110.83
8	B	694	MAN	O5-C5-C6	2.15	111.86	107.66
7	A	498	NAG	C2-N2-C7	-2.14	120.04	122.90
5	B	3681	GLA	C4-C3-C2	-2.07	107.20	110.83
8	B	695	MAN	C3-C4-C5	-2.04	106.54	110.23
5	B	3681	GLA	O4-C4-C5	-2.02	104.34	109.32
7	A	430	NAG	O5-C5-C6	2.01	111.58	107.66

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	6744	TAM	C2-C-C1-C4
6	A	6744	TAM	C3-C-C1-C4
6	A	6744	TAM	N-C-C1-C4
6	A	6744	TAM	C2-C-C3-C6
6	A	6744	TAM	N-C-C3-C6
7	A	430	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
8	B	695	MAN	O5-C5-C6-O6
7	A	430	NAG	C4-C5-C6-O6
8	B	695	MAN	C4-C5-C6-O6
7	A	430	NAG	C8-C7-N2-C2
7	A	430	NAG	O7-C7-N2-C2
8	B	694	MAN	O5-C5-C6-O6
7	A	498	NAG	C4-C5-C6-O6
6	A	6744	TAM	C1-C-C3-C6

There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	3681	GLA	4	0
6	A	6744	TAM	11	0
7	A	430	NAG	2	0
7	A	498	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	390/398 (97%)	0.83	18 (4%) 37 34	56, 71, 78, 85	0
1	B	391/398 (98%)	0.94	32 (8%) 17 15	57, 70, 77, 91	0
All	All	781/796 (98%)	0.89	50 (6%) 25 22	56, 71, 78, 91	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	SER	3.2
1	A	143	ALA	3.1
1	B	309	ALA	3.1
1	B	271	GLY	2.9
1	A	368	ALA	2.9
1	B	124	VAL	2.9
1	A	415	LEU	2.7
1	B	273	PHE	2.7
1	A	233	ASP	2.7
1	B	418	GLU	2.6
1	A	421	MET	2.6
1	B	288	ALA	2.6
1	A	416	GLN	2.5
1	B	385	THR	2.5
1	A	370	ALA	2.5
1	B	415	LEU	2.4
1	B	305	PRO	2.4
1	B	352	ALA	2.4
1	A	209	TRP	2.4
1	B	116	GLY	2.4
1	A	420	THR	2.4
1	B	416	GLN	2.4
1	B	274	GLY	2.4
1	B	239	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	303	ILE	2.4
1	A	323	PRO	2.3
1	B	276	SER	2.3
1	B	245	TRP	2.3
1	A	230	ALA	2.3
1	A	291	ALA	2.3
1	B	421	MET	2.3
1	A	309	ALA	2.3
1	B	302	HIS	2.3
1	A	239	ILE	2.3
1	B	277	TRP	2.2
1	A	221	GLN	2.2
1	B	246	THR	2.2
1	A	189	LEU	2.2
1	B	286	LEU	2.2
1	B	300	LEU	2.2
1	A	372	LEU	2.1
1	B	289	ILE	2.1
1	B	269	VAL	2.1
1	B	209	TRP	2.1
1	B	232	ILE	2.1
1	B	313	ASP	2.1
1	B	152	TYR	2.1
1	B	144	GLY	2.1
1	B	270	ILE	2.0
1	A	171	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

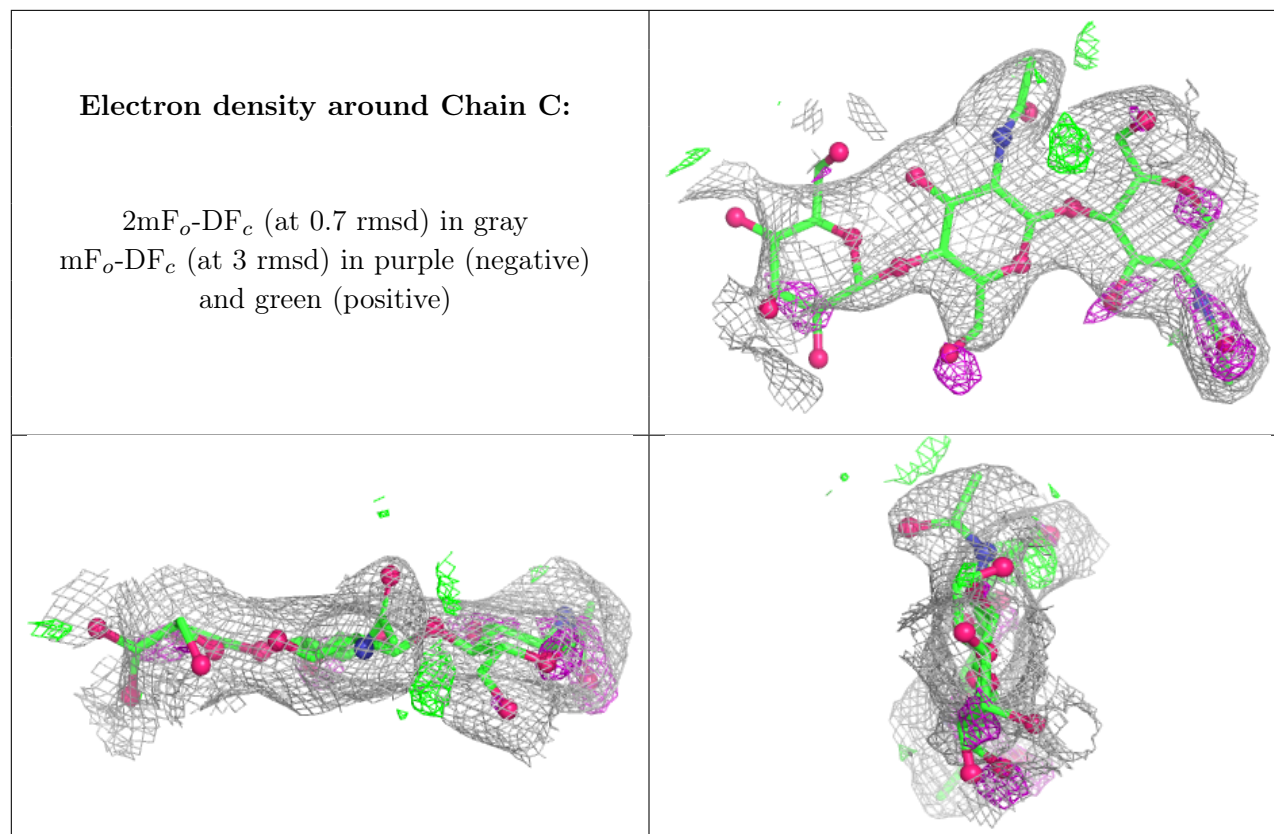
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	C	3	11/12	0.45	0.15	107,110,113,113	0
3	NAG	G	2	14/15	0.60	0.13	100,113,115,116	0

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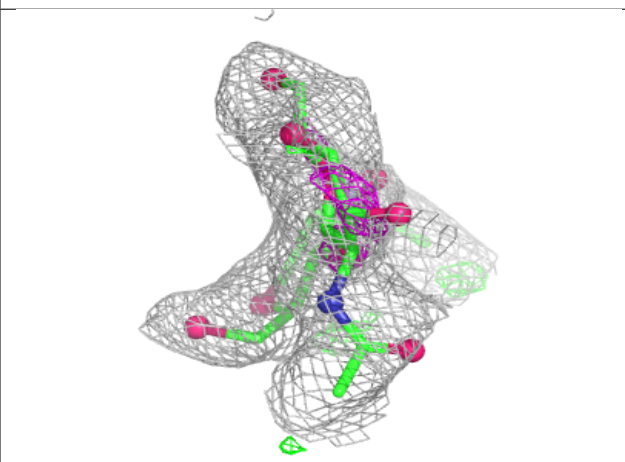
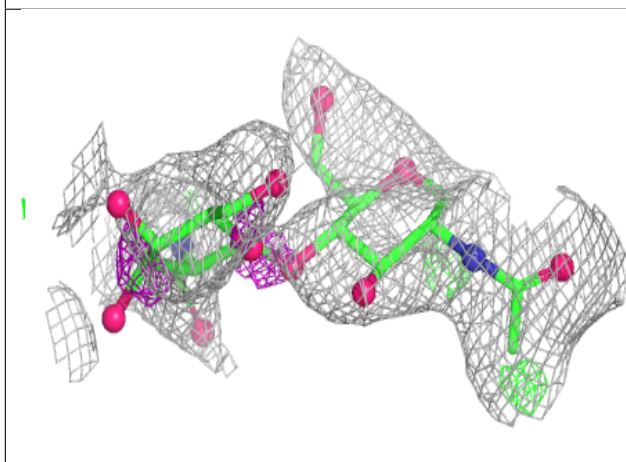
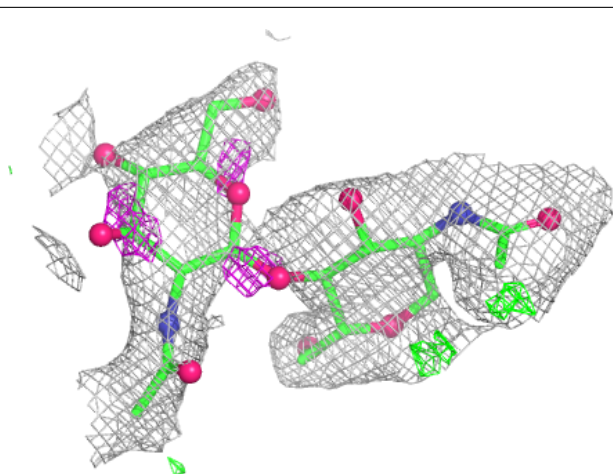
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	G	1	14/15	0.61	0.14	98,104,106,110	0
3	NAG	D	2	14/15	0.63	0.14	103,107,108,108	0
3	NAG	E	2	14/15	0.63	0.12	103,105,106,108	0
3	NAG	D	1	14/15	0.71	0.13	89,96,100,104	0
2	NAG	C	2	14/15	0.77	0.13	78,85,89,95	0
3	NAG	F	2	14/15	0.78	0.13	78,80,86,86	0
3	NAG	E	1	14/15	0.79	0.13	89,92,95,99	0
3	NAG	F	1	14/15	0.80	0.12	75,78,78,79	0
2	NAG	C	1	14/15	0.84	0.11	57,65,67,71	0

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



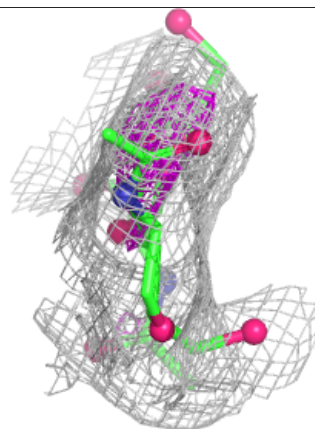
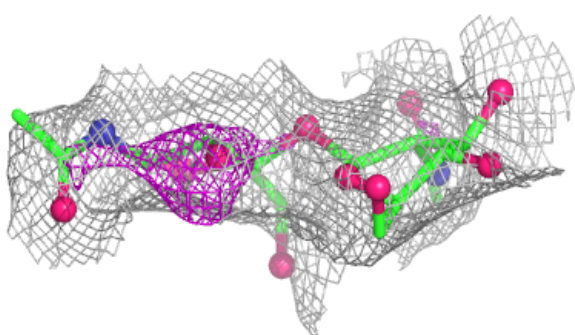
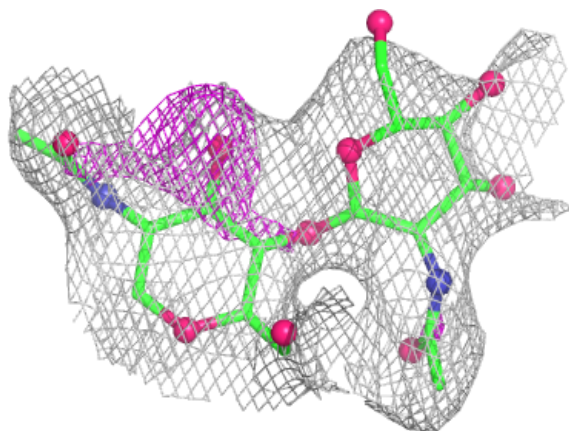
Electron density around Chain D:

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



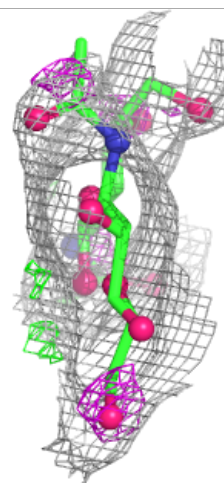
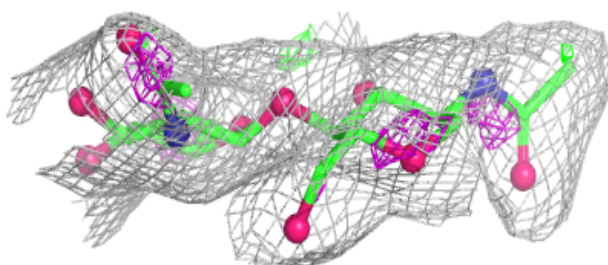
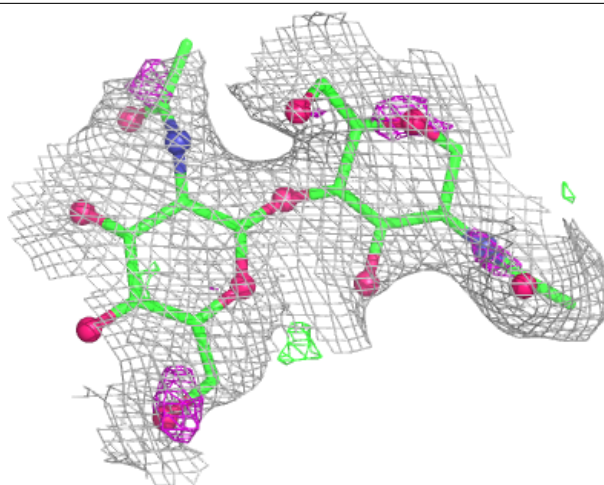
Electron density around Chain E:

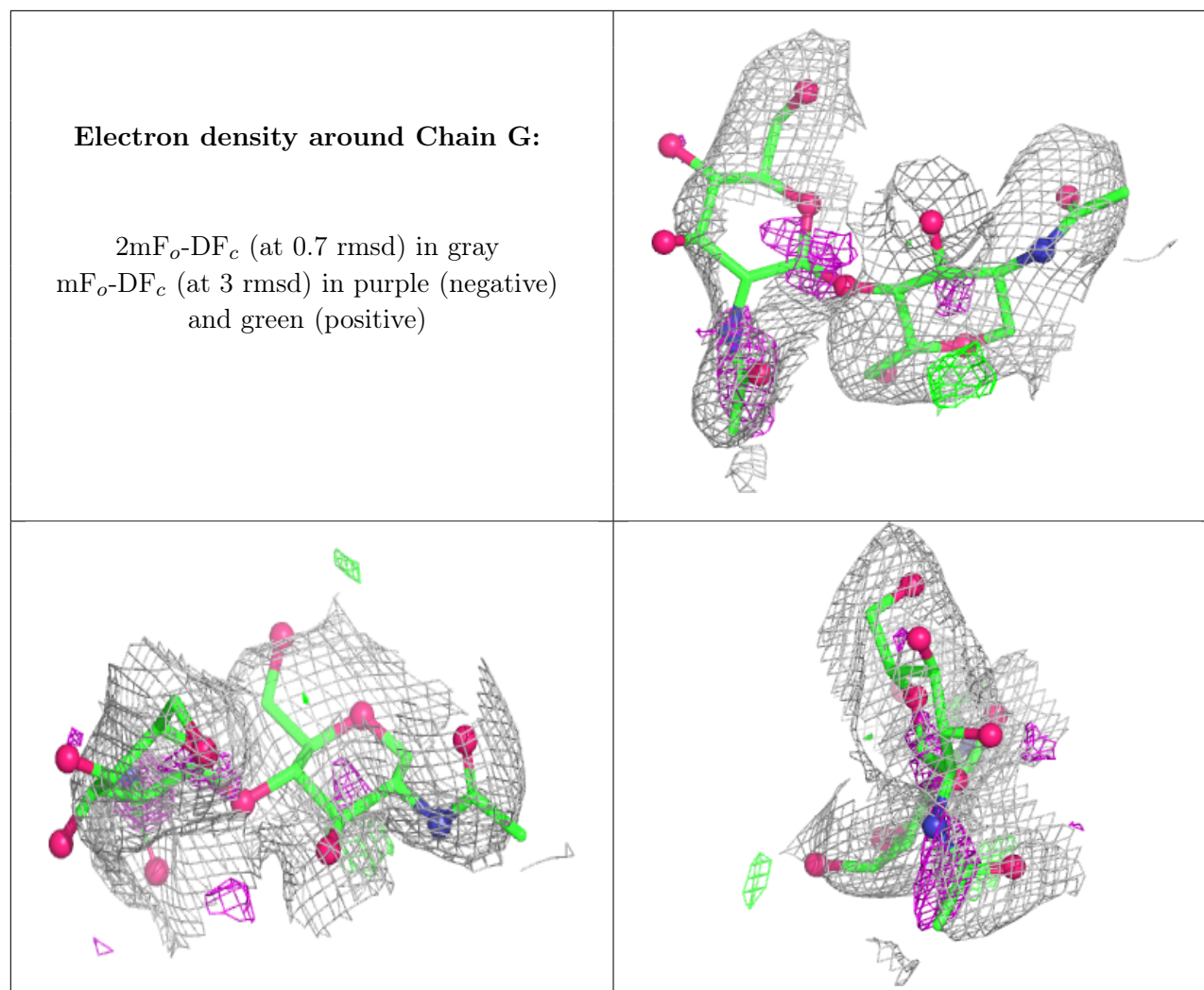
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



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
7	NAG	A	430	14/15	0.65	0.12	101,106,109,110	0
8	MAN	B	694	11/12	0.65	0.13	151,152,153,153	0
7	NAG	A	498	14/15	0.66	0.15	71,83,86,87	0
4	SO4	B	4	5/5	0.68	0.14	107,108,109,110	0
4	SO4	B	3	5/5	0.69	0.14	102,103,104,105	0
6	TAM	A	6744	11/11	0.72	0.19	50,55,64,68	0
8	MAN	B	695	11/12	0.80	0.12	90,96,97,99	0
5	GLA	B	3681	12/12	0.85	0.12	52,57,60,65	0
4	SO4	B	2	5/5	0.85	0.10	81,82,83,84	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	A	1	5/5	0.89	0.11	69,71,71,73	0
5	GLA	A	3681	12/12	0.89	0.12	47,54,56,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.