



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:29 AM UTC

PDB ID : 5ZSJ / pdb_00005zsj
Title : Crystal structure of monkey TLR7 in complex with GS9620
Authors : Zhang, Z.; Ohto, U.; Shimizu, T.
Deposited on : 2018-04-28
Resolution : 2.40 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

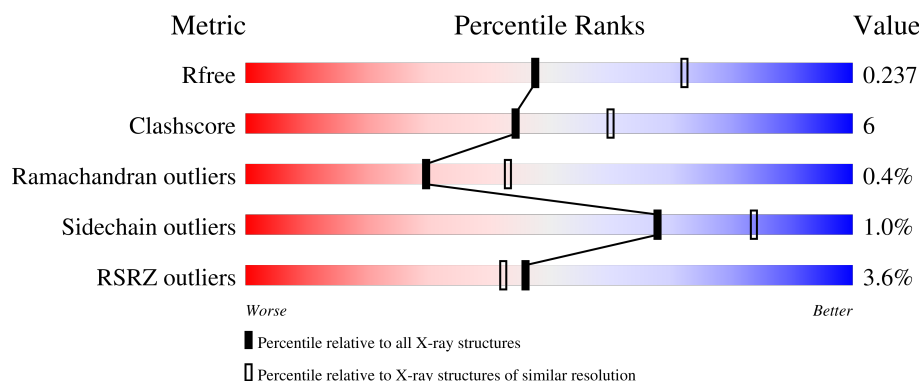
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	823	3% 82% 11% 6%
1	B	823	4% 80% 12% 6%
2	C	2	50% 50%
2	D	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
5	SO4	A	1017	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13545 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 Toll-like receptor 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	772	Total	C	N	O	S	0	0	0
			6236	3996	1062	1148	30			
1	B	772	Total	C	N	O	S	0	0	0
			6241	4000	1063	1148	30			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	ARG	-	expression tag	UNP B3Y653
A	24	SER	-	expression tag	UNP B3Y653
A	25	PRO	-	expression tag	UNP B3Y653
A	26	TRP	-	expression tag	UNP B3Y653
A	167	GLN	ASN	engineered mutation	UNP B3Y653
A	389	GLN	ASN	engineered mutation	UNP B3Y653
A	440	LEU	SER	see sequence details	UNP B3Y653
A	441	VAL	GLU	see sequence details	UNP B3Y653
A	442	PRO	VAL	see sequence details	UNP B3Y653
A	443	ARG	GLY	see sequence details	UNP B3Y653
A	444	GLY	PHE	see sequence details	UNP B3Y653
A	445	SER	CYS	see sequence details	UNP B3Y653
A	488	GLN	ASN	engineered mutation	UNP B3Y653
A	799	GLN	ASN	engineered mutation	UNP B3Y653
A	840	GLU	-	expression tag	UNP B3Y653
A	841	PHE	-	expression tag	UNP B3Y653
A	842	LEU	-	expression tag	UNP B3Y653
A	843	VAL	-	expression tag	UNP B3Y653
A	844	PRO	-	expression tag	UNP B3Y653
A	845	ARG	-	expression tag	UNP B3Y653
B	23	ARG	-	expression tag	UNP B3Y653
B	24	SER	-	expression tag	UNP B3Y653
B	25	PRO	-	expression tag	UNP B3Y653
B	26	TRP	-	expression tag	UNP B3Y653
B	167	GLN	ASN	engineered mutation	UNP B3Y653

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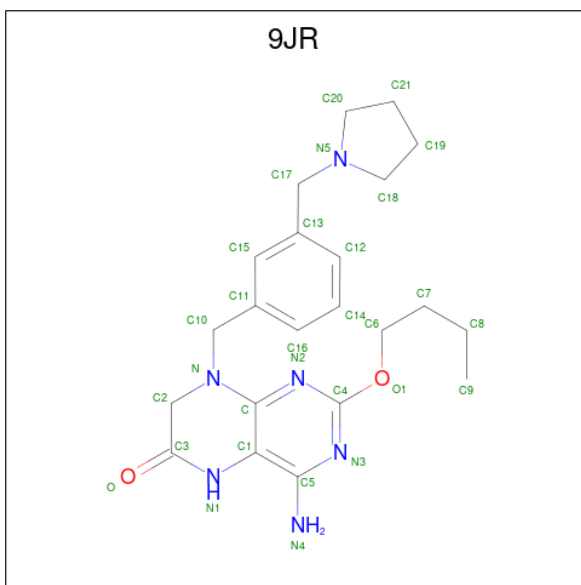
Chain	Residue	Modelled	Actual	Comment	Reference
B	389	GLN	ASN	engineered mutation	UNP B3Y653
B	440	LEU	SER	see sequence details	UNP B3Y653
B	441	VAL	GLU	see sequence details	UNP B3Y653
B	442	PRO	VAL	see sequence details	UNP B3Y653
B	443	ARG	GLY	see sequence details	UNP B3Y653
B	444	GLY	PHE	see sequence details	UNP B3Y653
B	445	SER	CYS	see sequence details	UNP B3Y653
B	488	GLN	ASN	engineered mutation	UNP B3Y653
B	799	GLN	ASN	engineered mutation	UNP B3Y653
B	840	GLU	-	expression tag	UNP B3Y653
B	841	PHE	-	expression tag	UNP B3Y653
B	842	LEU	-	expression tag	UNP B3Y653
B	843	VAL	-	expression tag	UNP B3Y653
B	844	PRO	-	expression tag	UNP B3Y653
B	845	ARG	-	expression tag	UNP B3Y653

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



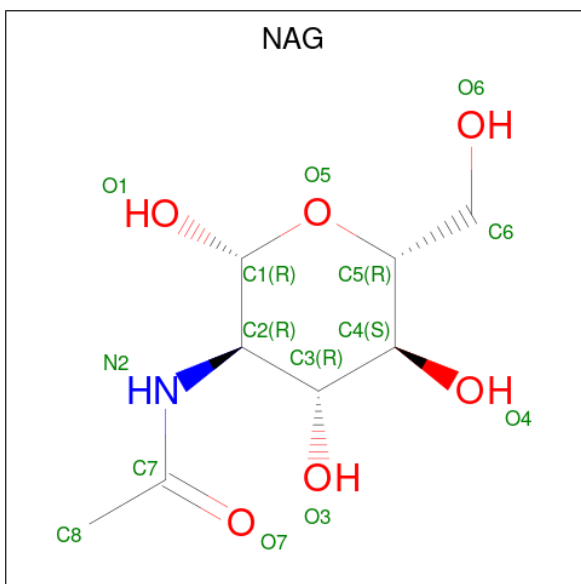
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 4-azanyl-2-butoxy-8-[[3-(pyrrolidin-1-ylmethyl)phenyl]methyl]-5,7-dihydro pteridin-6-one (CCD ID: 9JR) (formula: C₂₂H₃₀N₆O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			30	22	6	2		
3	B	1	Total	C	N	O	0	0
			30	22	6	2		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

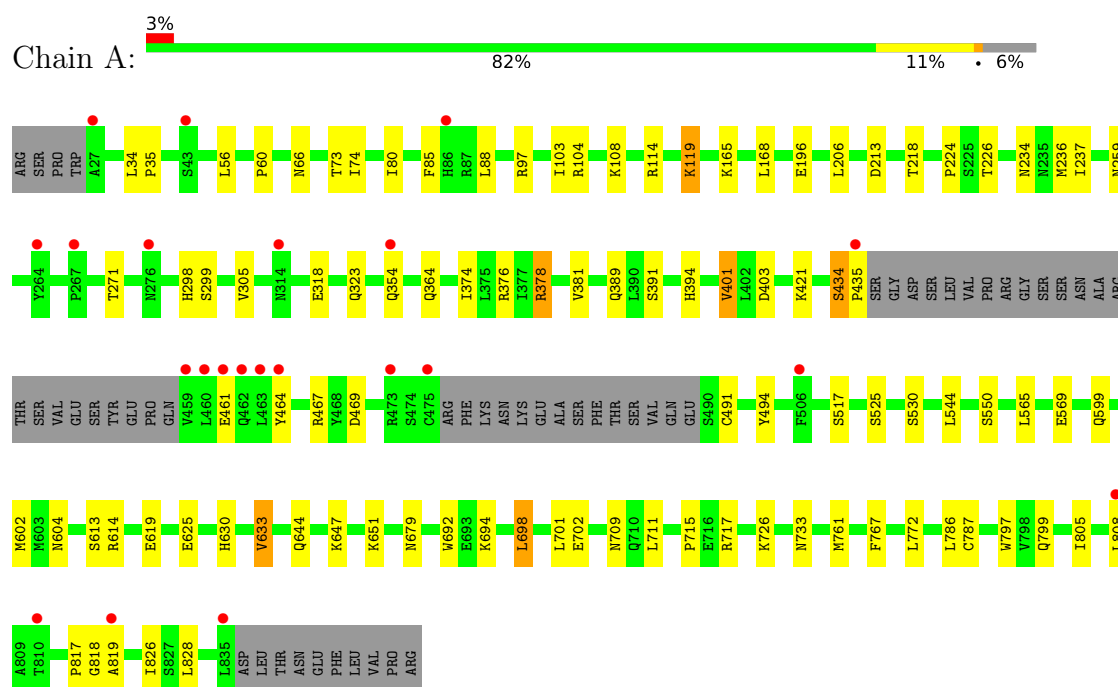
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	415	Total O 415 415	0	0
6	B	314	Total O 314 314	0	0

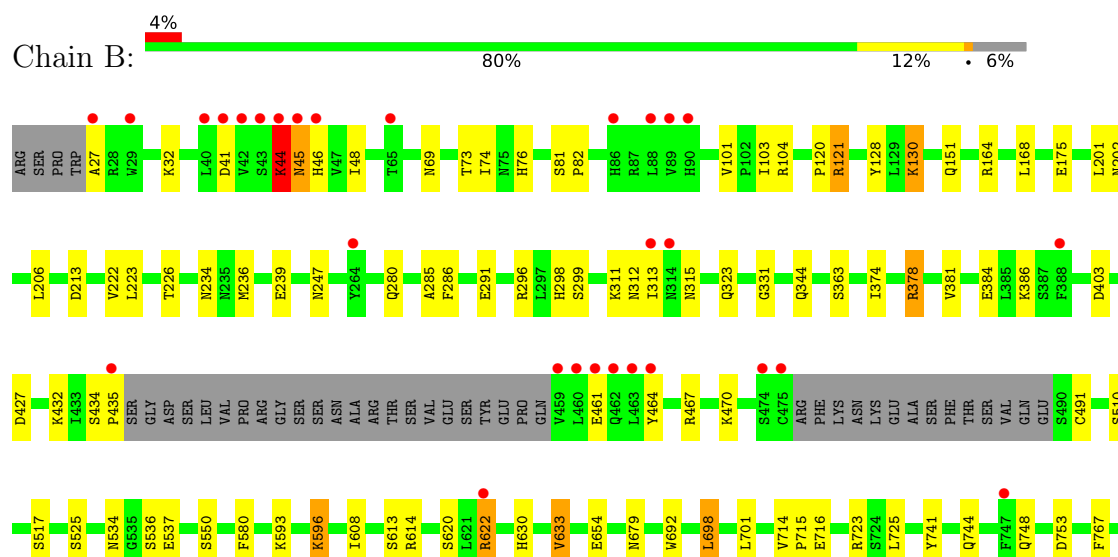
3 Residue-property plots [i](#)

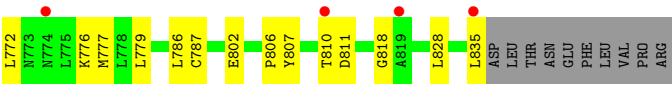
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Toll-like receptor 7



• Molecule 1: Toll-like receptor 7





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



• Molecule 2: 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	99.07Å 138.64Å 149.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.85 – 2.40 46.85 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.85-2.40) 99.8 (46.85-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.195 , 0.236 0.197 , 0.237	Depositor DCC
R_{free} test set	3976 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	31.1	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13545	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.29% 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: NAG, SO4, 9JR

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	0.39	0/6367	0.63	3/8627 (0.0%)
1	B	0.36	1/6372 (0.0%)	0.63	5/8633 (0.1%)
All	All	0.38	1/12739 (0.0%)	0.63	8/17260 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	130	LYS	CD-CE	6.18	1.71	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	614	ARG	NE-CZ-NH2	8.68	127.01	119.20
1	B	776	LYS	CA-CB-CG	-7.71	98.69	114.10
1	A	614	ARG	NE-CZ-NH1	-5.97	115.53	121.50
1	B	622	ARG	CG-CD-NE	5.62	124.35	112.00
1	B	596	LYS	CA-CB-CG	-5.60	102.91	114.10
1	B	121	ARG	CA-CB-CG	5.29	124.67	114.10
1	B	130	LYS	CG-CD-CE	-5.18	99.38	111.30
1	A	119	LYS	CA-CB-CG	5.15	124.40	114.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	434	SER	Peptide
1	B	44	LYS	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6236	0	6267	73	0
1	B	6241	0	6283	79	0
2	C	28	0	25	1	0
2	D	28	0	25	0	0
3	A	30	0	0	0	0
3	B	30	0	0	0	0
4	A	112	0	104	1	0
4	B	56	0	52	2	0
5	A	40	0	0	3	0
5	B	15	0	0	0	0
6	A	415	0	0	28	0
6	B	314	0	0	25	1
All	All	13545	0	12756	154	1

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 6.

All (154) 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:344:GLN:NE2	6:B:1103:HOH:O	1.95	0.97
1:A:271:THR:O	6:A:1101:HOH:O	1.83	0.95
1:A:196:GLU:OE2	6:A:1102:HOH:O	1.88	0.91
1:A:711:LEU:H	1:A:733:ASN:HD22	1.19	0.91
1:A:236:MET:SD	6:A:1447:HOH:O	2.31	0.87
1:B:164:ARG:NH2	6:B:1108:HOH:O	2.08	0.87
1:A:218:THR:OG1	6:A:1103:HOH:O	1.91	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:TYR:OH	6:B:1101:HOH:O	1.93	0.86
1:A:494:TYR:O	6:A:1104:HOH:O	1.92	0.85
1:B:236:MET:SD	6:B:1348:HOH:O	2.34	0.84
1:B:716:GLU:O	6:B:1102:HOH:O	1.94	0.84
1:A:165:LYS:NZ	6:A:1112:HOH:O	2.08	0.83
6:A:1105:HOH:O	2:C:2:NAG:O4	1.96	0.83
1:B:517:SER:OG	6:B:1104:HOH:O	1.96	0.82
1:A:114:ARG:NH2	6:A:1107:HOH:O	2.01	0.82
1:B:620:SER:O	1:B:622:ARG:NH1	2.13	0.81
1:B:777:MET:SD	6:B:1412:HOH:O	2.41	0.79
5:A:1012:SO4:O4	6:A:1106:HOH:O	2.00	0.78
5:A:1017:SO4:O3	6:A:1108:HOH:O	2.02	0.78
1:A:602:MET:SD	6:A:1449:HOH:O	2.44	0.76
1:B:73:THR:HG22	1:B:74:ILE:HG13	1.67	0.76
1:B:120:PRO:O	1:B:121:ARG:HB3	1.87	0.75
1:A:679:ASN:OD1	6:A:1109:HOH:O	2.03	0.75
1:A:389:GLN:O	6:A:1110:HOH:O	2.05	0.74
1:A:206:LEU:O	1:A:226:THR:HG23	1.87	0.74
1:B:698:LEU:O	6:B:1105:HOH:O	2.05	0.73
1:B:285:ALA:O	6:B:1106:HOH:O	2.06	0.73
1:B:291:GLU:OE2	6:B:1107:HOH:O	2.06	0.73
1:A:702:GLU:OE1	6:A:1111:HOH:O	2.06	0.73
1:B:470:LYS:NZ	6:B:1116:HOH:O	2.20	0.72
1:B:835:LEU:O	6:B:1109:HOH:O	2.08	0.72
1:B:767:PHE:HB3	1:B:772:LEU:HD11	1.73	0.71
1:A:66:ASN:ND2	6:A:1115:HOH:O	2.17	0.69
1:A:544:LEU:HB3	1:A:565:LEU:HD22	1.75	0.68
1:A:517:SER:OG	6:A:1114:HOH:O	2.12	0.66
1:B:121:ARG:HH11	1:B:121:ARG:HG2	1.59	0.66
1:A:298:HIS:CE1	1:A:323:GLN:HG3	2.31	0.66
1:B:384:GLU:HG2	1:B:386:LYS:HD2	1.78	0.66
1:B:76:HIS:ND1	6:B:1123:HOH:O	2.29	0.66
1:A:73:THR:HG22	1:A:74:ILE:HG13	1.78	0.65
1:B:537:GLU:OE2	6:B:1112:HOH:O	2.14	0.65
1:A:767:PHE:HB3	1:A:772:LEU:HD11	1.79	0.64
1:B:206:LEU:O	1:B:226:THR:HG23	1.97	0.64
1:B:654:GLU:HB2	6:B:1140:HOH:O	1.96	0.64
1:B:41:ASP:OD2	1:B:44:LYS:HG2	1.98	0.63
1:A:694:LYS:NZ	6:A:1113:HOH:O	2.11	0.61
1:B:741:TYR:HB3	1:B:744:GLN:HG3	1.83	0.61
1:A:711:LEU:H	1:A:733:ASN:ND2	1.95	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:HIS:CG	6:B:1123:HOH:O	2.54	0.60
1:B:298:HIS:CE1	1:B:323:GLN:HG3	2.38	0.58
1:B:593:LYS:O	1:B:596:LYS:HE2	2.03	0.58
1:A:56:LEU:HD13	1:A:60:PRO:HD3	1.84	0.58
1:B:725:LEU:O	1:B:748:GLN:HG2	2.04	0.58
1:A:709:ASN:HB2	1:A:733:ASN:HD21	1.69	0.57
1:B:787:CYS:HB3	1:B:828:LEU:HD21	1.85	0.57
1:A:772:LEU:HD13	1:A:797:TRP:CZ2	2.39	0.57
1:B:103:ILE:HG23	1:B:104:ARG:HG3	1.86	0.57
1:B:202:ASN:OD1	6:B:1113:HOH:O	2.18	0.57
1:B:311:LYS:HD3	1:B:312:ASN:OD1	2.04	0.57
1:A:318:GLU:OE1	6:A:1116:HOH:O	2.18	0.57
1:A:619:GLU:O	1:A:651:LYS:HD2	2.06	0.56
1:B:698:LEU:HB3	1:B:701:LEU:HB2	1.87	0.56
1:B:384:GLU:OE2	1:B:386:LYS:NZ	2.37	0.55
1:A:544:LEU:HB3	1:A:565:LEU:CD2	2.36	0.55
1:B:81:SER:HB2	1:B:82:PRO:HD2	1.88	0.55
1:A:717:ARG:HD2	6:A:1138:HOH:O	2.06	0.54
1:A:698:LEU:HB3	1:A:701:LEU:HB2	1.89	0.54
1:A:644:GLN:HG2	1:A:647:LYS:HD2	1.90	0.54
1:A:354:GLN:HA	1:A:354:GLN:OE1	2.06	0.54
1:B:630:HIS:HB3	1:B:633:VAL:HG13	1.89	0.53
1:B:299:SER:HA	1:B:323:GLN:O	2.08	0.53
1:A:787:CYS:HB3	1:A:828:LEU:HD21	1.90	0.53
1:A:364:GLN:HG2	6:A:1137:HOH:O	2.09	0.53
1:B:201:LEU:O	6:B:1114:HOH:O	2.19	0.53
1:A:805:ILE:HB	1:A:808:LEU:HD21	1.91	0.51
1:B:491:CYS:HB3	4:B:1005:NAG:H81	1.93	0.51
1:A:299:SER:HA	1:A:323:GLN:O	2.10	0.51
1:A:530:SER:O	1:B:432:LYS:NZ	2.44	0.51
1:A:692:TRP:CE2	1:A:715:PRO:HD3	2.46	0.50
1:A:80:ILE:O	1:A:119:LYS:HD3	2.12	0.50
1:A:569:GLU:OE2	6:A:1117:HOH:O	2.20	0.50
1:B:121:ARG:HB2	1:B:121:ARG:CZ	2.41	0.50
1:A:237:ILE:O	1:A:259:ASN:HB3	2.12	0.50
1:A:711:LEU:N	1:A:733:ASN:HD22	1.99	0.50
1:B:76:HIS:CE1	6:B:1123:HOH:O	2.63	0.50
1:A:726:LYS:NZ	6:A:1145:HOH:O	2.45	0.49
1:A:224:PRO:HB2	1:A:226:THR:HG22	1.93	0.49
1:B:151:GLN:NE2	1:B:175:GLU:OE2	2.42	0.49
1:A:378:ARG:HD2	1:A:403:ASP:OD2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:434:SER:HB2	1:B:435:PRO:HD2	1.93	0.49
1:A:85:PHE:HA	1:A:88:LEU:HD12	1.94	0.49
1:B:213:ASP:OD2	1:B:234:ASN:HB2	2.13	0.49
1:B:32:LYS:N	6:B:1110:HOH:O	2.10	0.49
1:B:239:GLU:HG2	1:B:280:GLN:HB2	1.94	0.49
1:A:108:LYS:N	5:A:1017:SO4:O4	2.46	0.49
1:A:374:ILE:CD1	1:A:401:VAL:HG13	2.43	0.49
1:A:467:ARG:HD2	6:A:1335:HOH:O	2.12	0.48
1:A:692:TRP:CD2	1:A:715:PRO:HD3	2.48	0.48
1:B:510:SER:OG	1:B:536:SER:O	2.29	0.48
1:B:802:GLU:O	6:B:1115:HOH:O	2.20	0.48
1:A:103:ILE:HG23	1:A:104:ARG:HG3	1.95	0.48
1:A:461:GLU:O	1:A:464:TYR:HB2	2.12	0.48
1:A:434:SER:HB2	1:A:435:PRO:HD2	1.94	0.48
1:A:374:ILE:HD13	1:A:401:VAL:HG13	1.95	0.47
1:A:491:CYS:HB3	4:A:1002:NAG:H81	1.96	0.47
1:B:286:PHE:CB	1:B:313:ILE:HD11	2.43	0.47
1:B:806:PRO:HB2	1:B:807:TYR:CD2	2.49	0.47
1:A:467:ARG:HE	1:A:469:ASP:H	1.61	0.47
1:B:467:ARG:HD3	1:B:470:LYS:HA	1.96	0.47
1:B:654:GLU:HG3	1:B:679:ASN:HB3	1.96	0.47
1:A:761:MET:HE2	1:A:786:LEU:HB2	1.97	0.46
1:B:315:ASN:OD1	6:B:1117:HOH:O	2.21	0.46
1:A:421:LYS:HB3	1:A:421:LYS:NZ	2.30	0.46
1:B:82:PRO:HG3	1:B:121:ARG:HG2	1.97	0.46
1:B:692:TRP:CE2	1:B:715:PRO:HD3	2.50	0.46
1:B:810:THR:HG22	1:B:811:ASP:HB2	1.98	0.46
1:B:46:HIS:HB3	1:B:48:ILE:HD11	1.99	0.45
1:B:714:VAL:HB	1:B:715:PRO:HD2	1.98	0.45
1:B:27:ALA:HB1	1:B:802:GLU:OE2	2.17	0.45
1:A:103:ILE:HD12	6:A:1194:HOH:O	2.17	0.45
1:B:121:ARG:HH11	1:B:121:ARG:CG	2.27	0.44
1:B:130:LYS:HE3	1:B:130:LYS:HB2	1.83	0.44
1:B:461:GLU:O	1:B:464:TYR:HB2	2.17	0.44
1:B:378:ARG:NH1	6:B:1154:HOH:O	2.50	0.44
1:B:201:LEU:HD12	1:B:222:VAL:HG12	2.00	0.44
1:B:331:GLY:C	1:B:363:SER:HB2	2.42	0.44
1:B:534:ASN:ND2	6:B:1133:HOH:O	2.39	0.44
1:A:165:LYS:HD2	6:A:1147:HOH:O	2.17	0.44
1:B:69:ASN:HD22	4:B:1004:NAG:H83	1.82	0.43
1:B:525:SER:HA	1:B:550:SER:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:LEU:HD12	1:A:35:PRO:HD2	1.99	0.43
1:B:121:ARG:CG	1:B:121:ARG:NH1	2.81	0.43
1:A:213:ASP:OD2	1:A:234:ASN:HB2	2.19	0.43
1:B:344:GLN:HG3	1:B:374:ILE:HB	1.99	0.43
1:A:73:THR:HA	1:A:97:ARG:O	2.18	0.43
1:B:753:ASP:HA	1:B:779:LEU:HB2	2.00	0.43
1:B:286:PHE:HB3	1:B:313:ILE:HD11	2.01	0.42
1:A:799:GLN:HA	1:A:826:ILE:HD12	2.02	0.42
1:B:296:ARG:NH1	6:B:1157:HOH:O	2.51	0.42
1:A:467:ARG:NH1	6:A:1154:HOH:O	2.49	0.42
1:A:525:SER:HA	1:A:550:SER:O	2.20	0.42
1:B:580:PHE:CD1	1:B:608:ILE:HG12	2.54	0.42
1:A:376:ARG:HA	1:A:403:ASP:HB3	2.02	0.42
1:A:817:PRO:O	1:A:819:ALA:N	2.53	0.42
1:B:403:ASP:HA	1:B:427:ASP:HB3	2.01	0.41
1:B:786:LEU:HA	1:B:786:LEU:HD23	1.77	0.41
1:A:599:GLN:CD	6:A:1144:HOH:O	2.64	0.41
1:A:604:ASN:OD1	1:A:625:GLU:HB3	2.21	0.41
1:A:630:HIS:HB3	1:A:633:VAL:HG13	2.03	0.41
1:A:104:ARG:NH1	6:A:1133:HOH:O	2.36	0.40
1:B:44:LYS:O	1:B:45:ASN:HB2	2.21	0.40
1:A:298:HIS:HE1	1:A:323:GLN:HG3	1.82	0.40
1:B:223:LEU:O	1:B:247:ASN:ND2	2.51	0.40
1:A:391:SER:HA	1:A:394:HIS:CE1	2.57	0.40

All (1) 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)
6:B:1409:HOH:O	6:B:1414:HOH:O[2_554]	2.06	0.14

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	766/823 (93%)	714 (93%)	50 (6%)	2 (0%)	36	50
1	B	766/823 (93%)	714 (93%)	48 (6%)	4 (0%)	24	37
All	All	1532/1646 (93%)	1428 (93%)	98 (6%)	6 (0%)	30	43

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	818	GLY
1	B	45	ASN
1	B	818	GLY
1	B	44	LYS
1	A	381	VAL
1	B	381	VAL

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	722/774 (93%)	715 (99%)	7 (1%)	68	84
1	B	723/774 (93%)	715 (99%)	8 (1%)	65	82
All	All	1445/1548 (93%)	1430 (99%)	15 (1%)	68	84

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

Mol	Chain	Res	Type
1	A	168	LEU
1	A	305	VAL
1	A	378	ARG
1	A	401	VAL
1	A	613	SER
1	A	633	VAL
1	A	698	LEU
1	B	101	VAL

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Mol	Chain	Res	Type
1	B	168	LEU
1	B	378	ARG
1	B	613	SER
1	B	614	ARG
1	B	633	VAL
1	B	698	LEU
1	B	723	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	ASN
1	A	252	GLN
1	A	280	GLN
1	A	300	ASN
1	A	344	GLN
1	A	398	ASN
1	A	539	GLN
1	A	578	HIS
1	A	594	ASN
1	A	599	GLN
1	A	732	ASN
1	A	733	ASN
1	B	55	HIS
1	B	110	ASN
1	B	181	GLN
1	B	300	ASN
1	B	344	GLN
1	B	389	GLN
1	B	394	HIS
1	B	398	ASN
1	B	514	GLN
1	B	581	GLN
1	B	594	ASN
1	B	599	GLN
1	B	732	ASN
1	B	760	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 ⓘ

4 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	0.58	0	17,19,21	0.44	0
2	NAG	C	2	2	14,14,15	0.68	1 (7%)	17,19,21	0.53	0
2	NAG	D	1	2,1	14,14,15	0.23	0	17,19,21	0.59	0
2	NAG	D	2	2	14,14,15	0.94	1 (7%)	17,19,21	1.18	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	2,1	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAG	C1-C2	-2.68	1.48	1.52
2	C	2	NAG	O5-C1	-2.25	1.39	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	2	NAG	C1-O5-C5	4.00	117.54	112.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

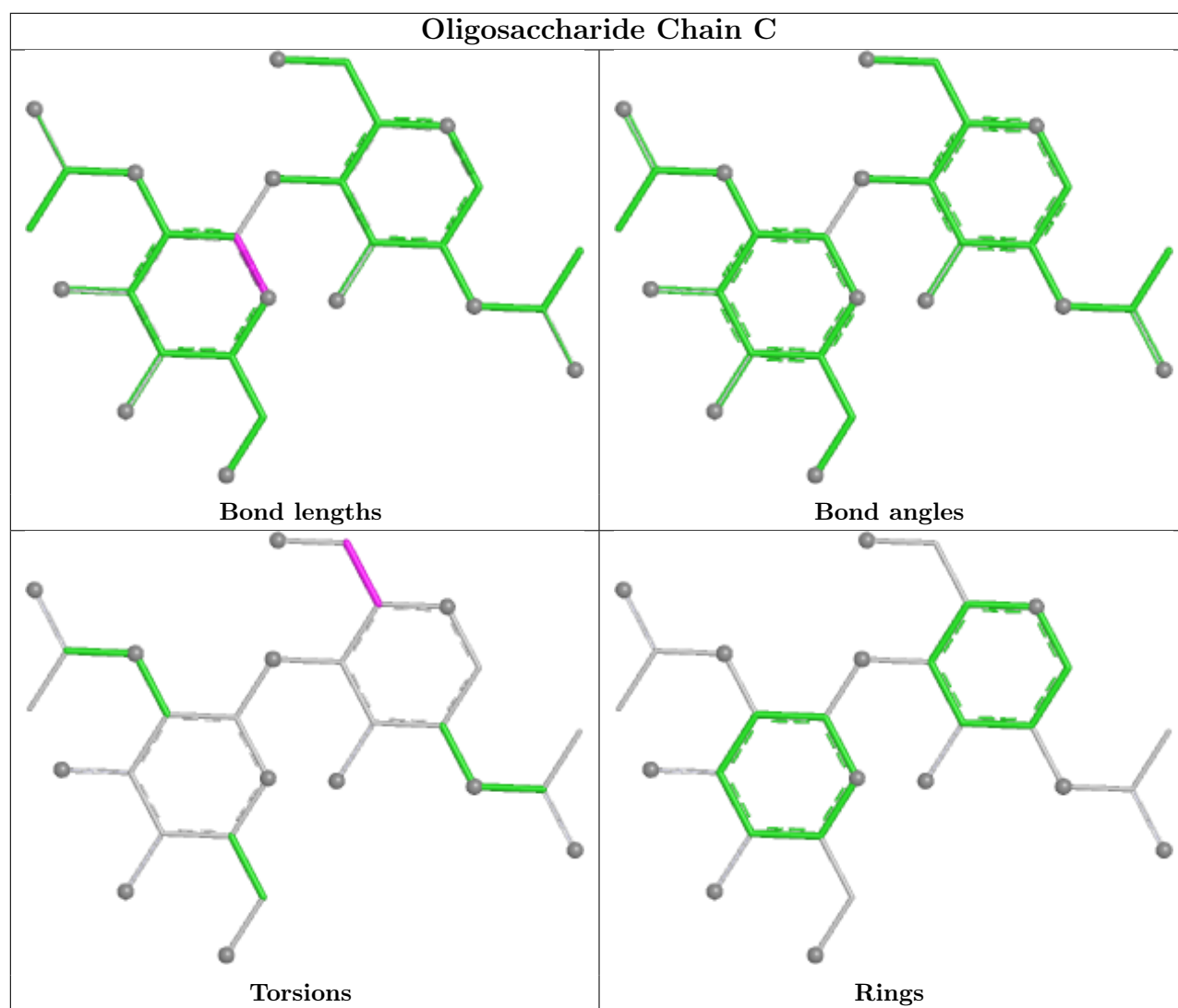
Mol	Chain	Res	Type	Atoms
2	D	2	NAG	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6

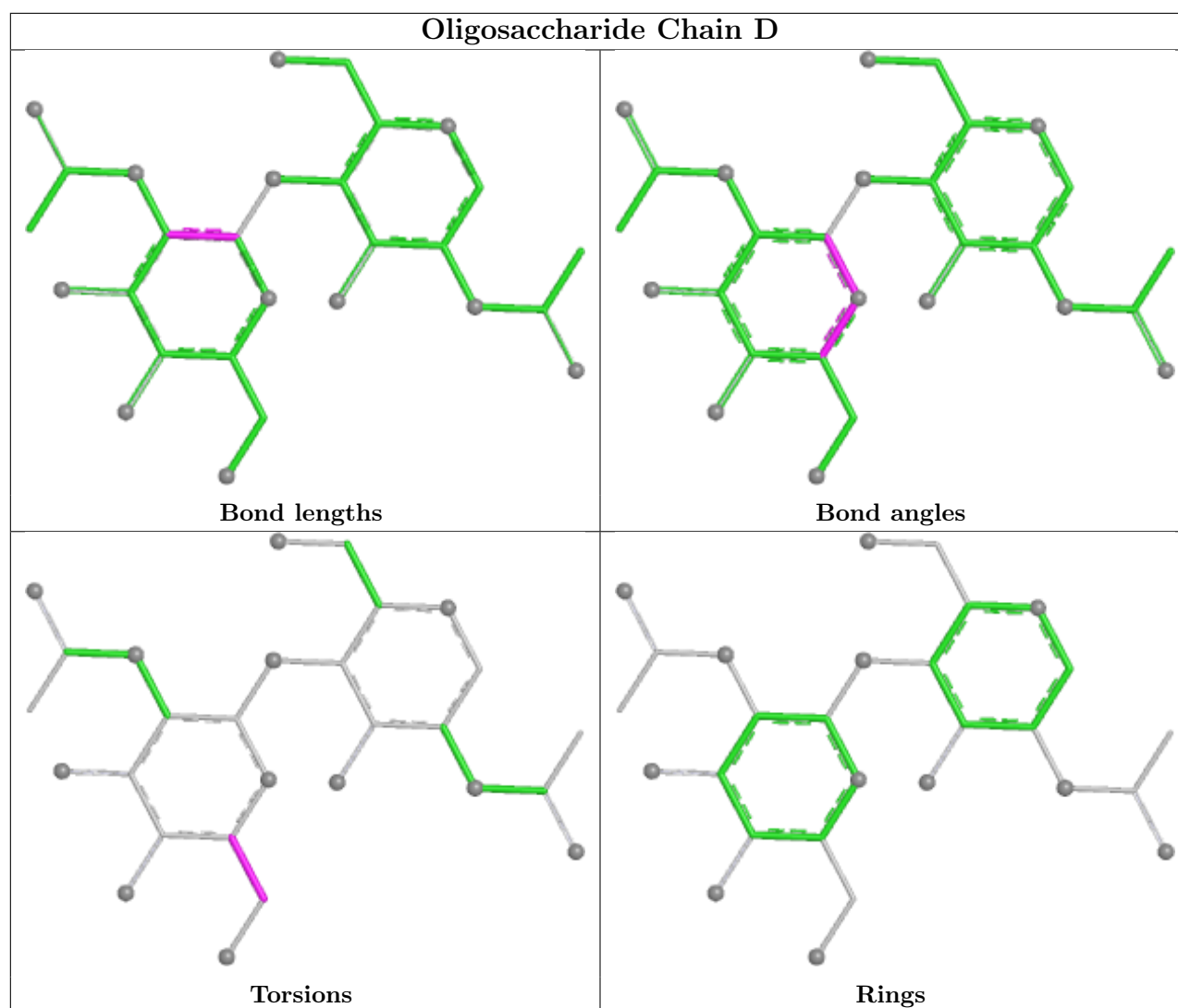
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
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.





5.6 Ligand geometry [i](#)

25 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	$\# Z > 2$
5	SO4	A	1014	-	4,4,4	0.29	0	6,6,6	0.10	0
5	SO4	B	1008	-	4,4,4	0.29	0	6,6,6	0.24	0
4	NAG	A	1002	1	14,14,15	0.23	0	17,19,21	0.78	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	A	1018	-	4,4,4	0.25	0	6,6,6	0.17	0
4	NAG	A	1008	1	14,14,15	0.49	0	17,19,21	0.85	1 (5%)
5	SO4	B	1009	-	4,4,4	0.27	0	6,6,6	0.33	0
5	SO4	A	1013	-	4,4,4	0.33	0	6,6,6	0.25	0
4	NAG	B	1006	1	14,14,15	1.22	1 (7%)	17,19,21	1.56	1 (5%)
4	NAG	A	1010	1	14,14,15	0.55	0	17,19,21	0.77	1 (5%)
5	SO4	A	1017	-	4,4,4	0.26	0	6,6,6	0.25	0
5	SO4	A	1015	-	4,4,4	0.39	0	6,6,6	0.29	0
4	NAG	A	1006	1	14,14,15	0.37	0	17,19,21	0.51	0
5	SO4	A	1016	-	4,4,4	0.31	0	6,6,6	0.27	0
5	SO4	A	1012	-	4,4,4	0.29	0	6,6,6	0.40	0
3	9JR	A	1001	-	33,33,33	1.34	4 (12%)	35,45,45	2.17	6 (17%)
4	NAG	A	1003	1	14,14,15	0.56	0	17,19,21	0.62	0
4	NAG	B	1005	1	14,14,15	0.44	0	17,19,21	0.67	0
4	NAG	A	1009	1	14,14,15	0.32	0	17,19,21	0.64	0
5	SO4	B	1010	-	4,4,4	0.33	0	6,6,6	0.21	0
4	NAG	A	1011	1	14,14,15	0.56	0	17,19,21	0.78	1 (5%)
4	NAG	A	1007	1	14,14,15	0.52	0	17,19,21	0.38	0
3	9JR	B	1001	-	33,33,33	1.24	4 (12%)	35,45,45	2.08	9 (25%)
5	SO4	A	1019	-	4,4,4	0.41	0	6,6,6	0.26	0
4	NAG	B	1004	1	14,14,15	0.24	0	17,19,21	0.66	0
4	NAG	B	1007	1	14,14,15	0.25	0	17,19,21	0.59	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
4	NAG	B	1006	1	-	1/6/23/26	0/1/1/1
4	NAG	A	1010	1	-	0/6/23/26	0/1/1/1
3	9JR	B	1001	-	-	0/13/32/32	0/4/4/4
4	NAG	A	1011	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1006	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1003	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1007	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1002	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1008	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1004	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1005	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1009	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1007	1	-	2/6/23/26	0/1/1/1
3	9JR	A	1001	-	-	2/13/32/32	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1006	NAG	O5-C1	4.47	1.51	1.43
3	A	1001	9JR	O1-C4	4.07	1.38	1.34
3	A	1001	9JR	C1-C	3.76	1.48	1.40
3	B	1001	9JR	O1-C4	3.50	1.38	1.34
3	B	1001	9JR	C1-C	3.24	1.47	1.40
3	B	1001	9JR	C1-N1	-3.08	1.33	1.39
3	A	1001	9JR	C1-N1	-2.92	1.33	1.39
3	B	1001	9JR	C1-C5	2.43	1.49	1.40
3	A	1001	9JR	C1-C5	2.15	1.48	1.40

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1001	9JR	C2-C3-N1	7.27	125.17	116.13
3	B	1001	9JR	C2-C3-N1	6.72	124.49	116.13
4	B	1006	NAG	C1-O5-C5	6.01	120.24	112.19
3	A	1001	9JR	O-C3-N1	-5.25	115.83	121.39
3	A	1001	9JR	C4-N3-C5	4.55	122.65	115.42
3	B	1001	9JR	C4-N3-C5	4.23	122.13	115.42
3	A	1001	9JR	N3-C4-N2	-3.63	121.05	127.65
3	B	1001	9JR	O-C3-N1	-3.62	117.56	121.39
3	B	1001	9JR	N3-C4-N2	-3.22	121.80	127.65
3	A	1001	9JR	C20-N5-C18	3.19	107.81	104.04
3	A	1001	9JR	C10-N-C2	3.10	120.24	116.08
4	A	1008	NAG	C1-O5-C5	3.06	116.29	112.19
3	B	1001	9JR	C20-N5-C18	2.96	107.54	104.04
3	B	1001	9JR	C11-C10-N	-2.76	109.66	114.13
4	A	1011	NAG	C1-O5-C5	2.66	115.76	112.19
4	A	1002	NAG	C1-O5-C5	2.62	115.69	112.19
4	A	1010	NAG	C1-O5-C5	2.49	115.52	112.19
3	B	1001	9JR	C16-C11-C15	2.14	121.50	118.55
3	B	1001	9JR	C16-C14-C12	-2.07	117.58	120.24
3	B	1001	9JR	C14-C12-C13	2.04	123.48	120.61

There are no chirality outliers.

All (13) torsion outliers are listed below:

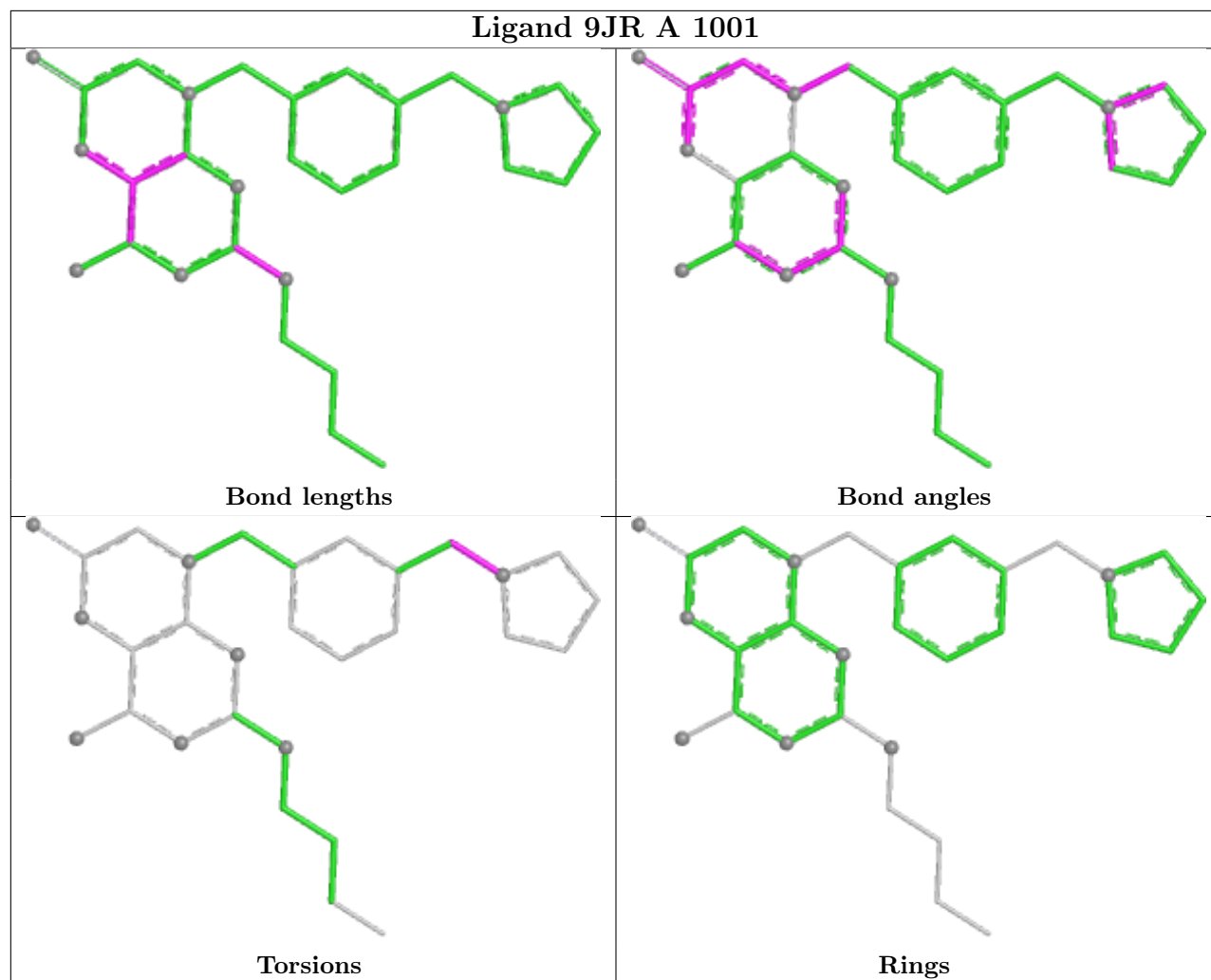
Mol	Chain	Res	Type	Atoms
4	A	1009	NAG	O5-C5-C6-O6
4	A	1007	NAG	O5-C5-C6-O6
4	A	1009	NAG	C4-C5-C6-O6
4	A	1007	NAG	C4-C5-C6-O6
4	B	1004	NAG	C8-C7-N2-C2
4	B	1004	NAG	O7-C7-N2-C2
4	A	1003	NAG	C4-C5-C6-O6
3	A	1001	9JR	C13-C17-N5-C20
4	A	1003	NAG	O5-C5-C6-O6
4	B	1007	NAG	C4-C5-C6-O6
4	B	1006	NAG	O5-C5-C6-O6
3	A	1001	9JR	C13-C17-N5-C18
4	B	1007	NAG	O5-C5-C6-O6

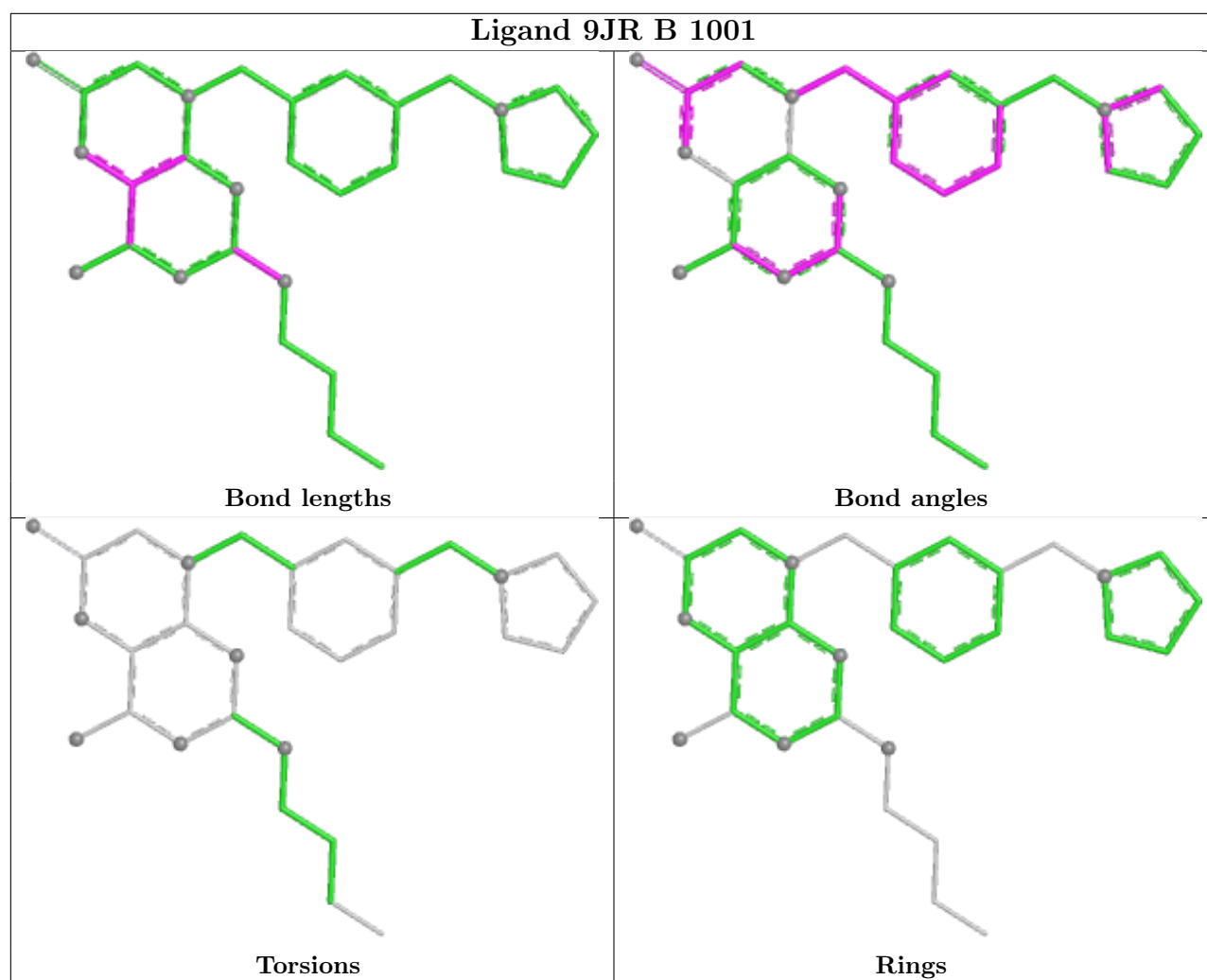
There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	NAG	1	0
5	A	1017	SO4	2	0
5	A	1012	SO4	1	0
4	B	1005	NAG	1	0
4	B	1004	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	772/823 (93%)	-0.21	22 (2%)	55 51	16, 30, 51, 93	0
1	B	772/823 (93%)	0.15	33 (4%)	40 36	21, 38, 57, 98	0
All	All	1544/1646 (93%)	-0.03	55 (3%)	46 42	16, 34, 55, 98	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	459	VAL	4.3
1	A	435	PRO	4.2
1	B	42	VAL	4.0
1	B	435	PRO	3.9
1	B	461	GLU	3.8
1	B	463	LEU	3.6
1	B	46	HIS	3.6
1	B	43	SER	3.6
1	A	462	GLN	3.5
1	A	459	VAL	3.5
1	A	460	LEU	3.4
1	B	462	GLN	3.4
1	B	313	ILE	3.4
1	B	474	SER	3.2
1	A	43	SER	3.2
1	A	819	ALA	3.1
1	A	473	ARG	3.1
1	A	810	THR	3.0
1	B	86	HIS	3.0
1	B	622	ARG	3.0
1	A	354	GLN	3.0
1	B	41	ASP	2.9
1	B	388	PHE	2.9
1	A	463	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	774	ASN	2.9
1	B	65	THR	2.8
1	B	29	TRP	2.8
1	B	475	CYS	2.8
1	B	27	ALA	2.8
1	B	810	THR	2.8
1	B	464	TYR	2.7
1	A	27	ALA	2.7
1	B	460	LEU	2.6
1	B	90	HIS	2.6
1	B	819	ALA	2.6
1	B	264	TYR	2.5
1	A	835	LEU	2.4
1	B	40	LEU	2.4
1	B	88	LEU	2.4
1	A	461	GLU	2.4
1	A	267	PRO	2.3
1	B	44	LYS	2.3
1	B	45	ASN	2.3
1	B	835	LEU	2.3
1	B	89	VAL	2.3
1	A	86	HIS	2.2
1	B	747	PHE	2.2
1	A	506	PHE	2.2
1	A	464	TYR	2.2
1	A	314	ASN	2.1
1	A	808	LEU	2.1
1	A	475	CYS	2.1
1	A	276	ASN	2.0
1	B	314	ASN	2.0
1	A	264	TYR	2.0

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

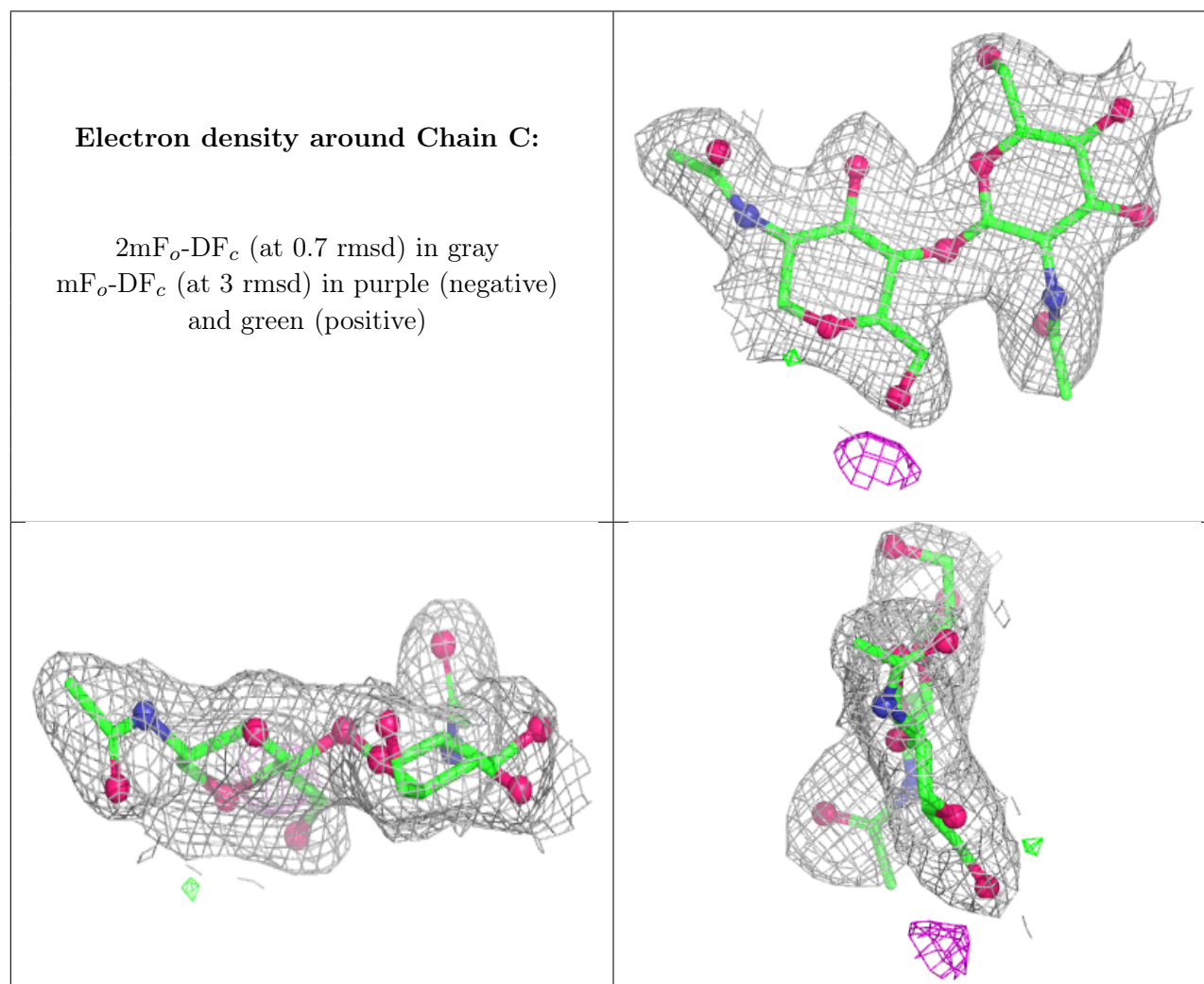
There are no non-standard protein/DNA/RNA residues in this entry.

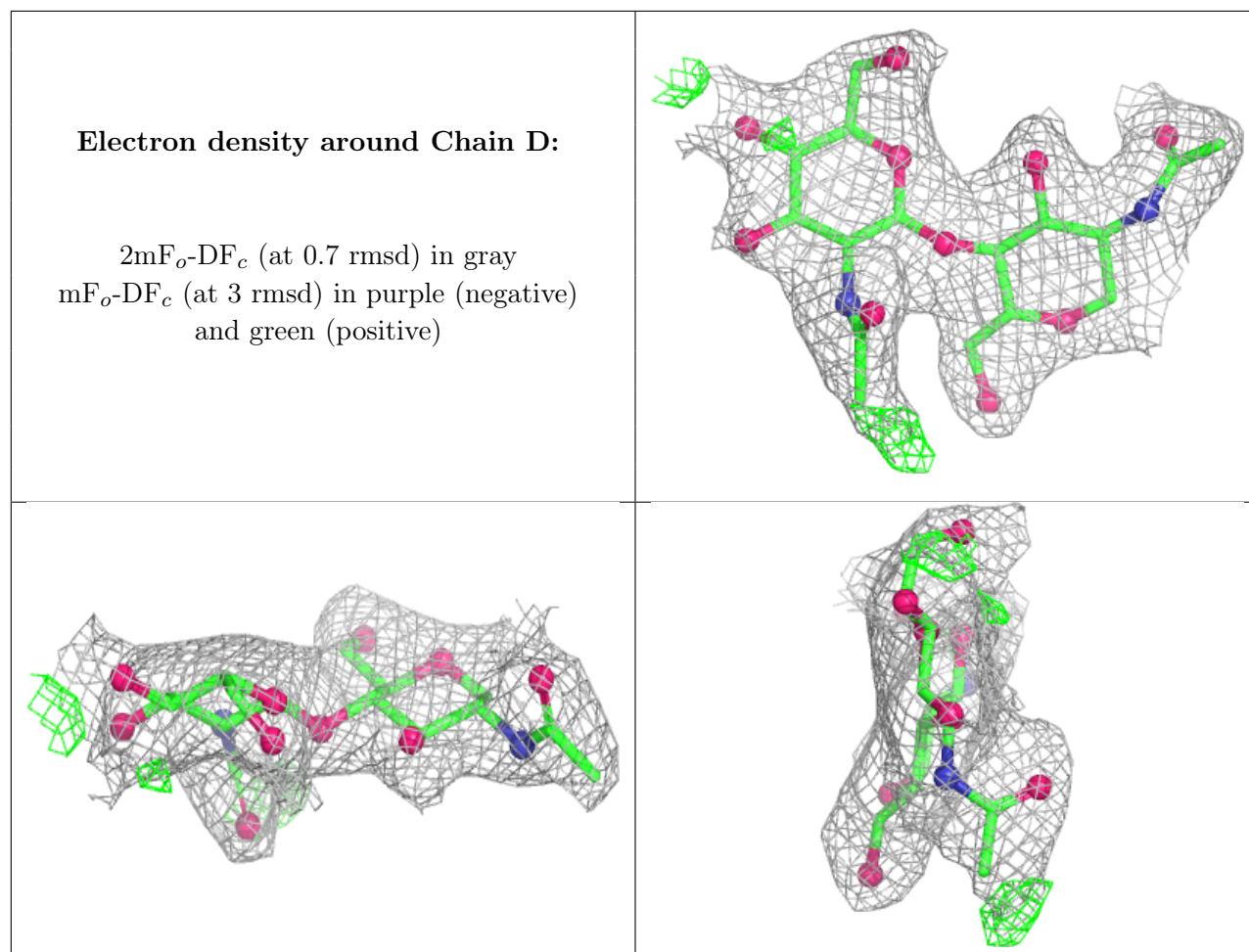
6.3 Carbohydrates [i](#)

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
2	NAG	D	2	14/15	0.80	0.12	45,53,58,59	0
2	NAG	C	2	14/15	0.87	0.11	44,53,61,62	0
2	NAG	C	1	14/15	0.93	0.08	27,38,43,45	0
2	NAG	D	1	14/15	0.94	0.08	31,37,42,43	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.





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.

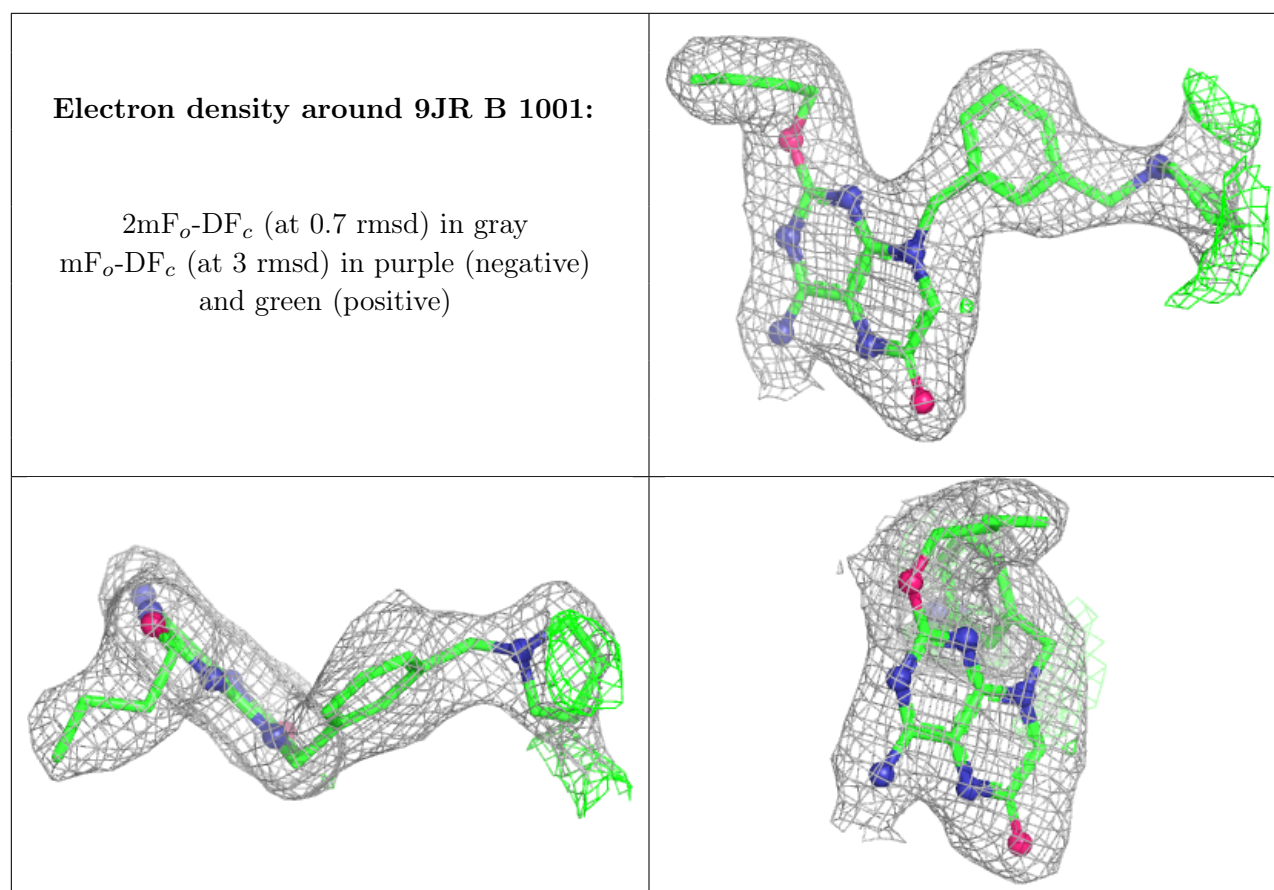
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	A	1007	14/15	0.77	0.16	57,65,76,78	0
5	SO4	A	1019	5/5	0.77	0.20	45,47,63,75	0
5	SO4	A	1016	5/5	0.79	0.21	59,60,73,78	0
4	NAG	B	1006	14/15	0.80	0.13	41,49,63,65	0
4	NAG	A	1009	14/15	0.81	0.13	37,47,50,53	0
4	NAG	A	1003	14/15	0.83	0.12	44,50,55,58	0
5	SO4	A	1018	5/5	0.84	0.14	60,67,75,84	0
5	SO4	B	1010	5/5	0.84	0.18	50,62,79,81	0
4	NAG	A	1008	14/15	0.86	0.11	36,44,53,55	0
4	NAG	B	1004	14/15	0.87	0.11	46,51,57,64	0
4	NAG	A	1011	14/15	0.87	0.10	31,40,46,47	0

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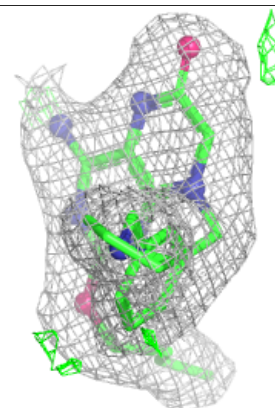
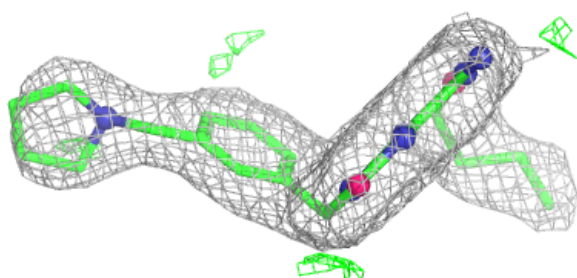
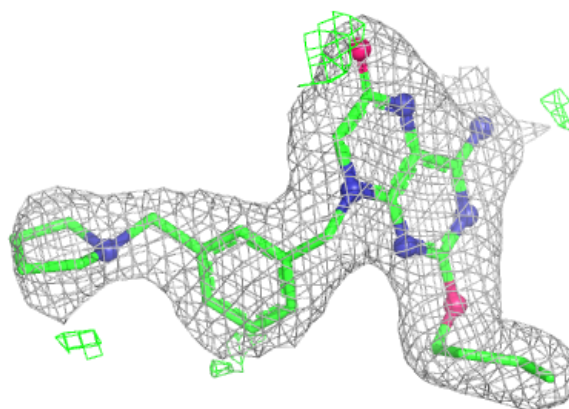
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	A	1014	5/5	0.87	0.14	60,62,74,74	0
5	SO4	A	1015	5/5	0.87	0.18	45,48,64,67	0
5	SO4	A	1017	5/5	0.89	0.18	57,60,65,68	0
5	SO4	B	1009	5/5	0.90	0.11	46,47,61,64	0
4	NAG	B	1007	14/15	0.90	0.10	42,51,56,56	0
4	NAG	B	1005	14/15	0.91	0.10	28,35,38,40	0
4	NAG	A	1010	14/15	0.92	0.08	38,43,50,52	0
4	NAG	A	1006	14/15	0.92	0.09	38,46,53,55	0
3	9JR	B	1001	30/30	0.92	0.10	24,29,54,62	0
5	SO4	A	1012	5/5	0.92	0.18	36,38,53,61	0
5	SO4	A	1013	5/5	0.93	0.13	49,54,62,63	0
5	SO4	B	1008	5/5	0.93	0.15	53,53,57,60	0
4	NAG	A	1002	14/15	0.94	0.08	23,27,31,40	0
3	9JR	A	1001	30/30	0.95	0.08	21,29,48,54	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around 9JR A 1001:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.