



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:58 AM UTC

PDB ID : 3H9V / pdb_00003h9v
Title : Crystal structure of the ATP-gated P2X4 ion channel in the closed, apo state at 3.1 Angstroms
Authors : Kawate, T.; Michel, J.C.; Gouaux, E.
Deposited on : 2009-04-30
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 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

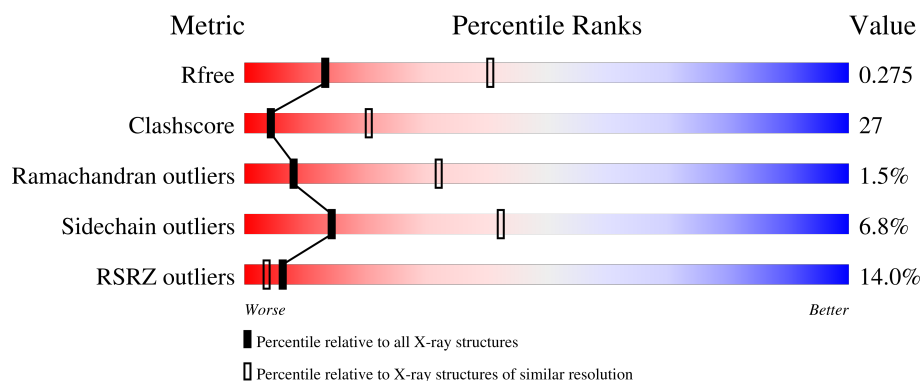
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	356	
2	B	2	
2	C	2	

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	NAG	B	1	X	-	-	-
2	NAG	B	2	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2501 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 P2X purinoceptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2443	1554	408	465	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP Q6NYR1
A	27	SER	-	expression tag	UNP Q6NYR1
A	51	PHE	CYS	engineered mutation	UNP Q6NYR1
A	78	LYS	ASN	engineered mutation	UNP Q6NYR1
A	187	ARG	ASN	engineered mutation	UNP Q6NYR1
A	252	ARG	HIS	engineered mutation	UNP Q6NYR1

- 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	B	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is GADOLINIUM ATOM (CCD ID: GD) (formula: Gd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Gd	0	0
			2	2		

4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	100.84Å 100.84Å 431.32Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.11 – 3.10 29.11 – 3.10	Depositor EDS
% Data completeness (in resolution range)	90.7 (29.11-3.10) 90.6 (29.11-3.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 3.06Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.248 , 0.278 0.251 , 0.275	Depositor DCC
R_{free} test set	732 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	85.8	Xtriage
Anisotropy	0.701	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 66.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2501	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% 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: NAG, GD

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	0.50	1/2498 (0.0%)	0.91	7/3413 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	167	TRP	C-O	-6.86	1.15	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	PRO	CA-C-N	6.11	126.14	119.90
1	A	179	PRO	C-N-CA	6.11	126.14	119.90
1	A	186	GLU	N-CA-C	-5.99	104.09	111.33
1	A	198	TYR	CA-C-N	5.40	124.85	119.24
1	A	198	TYR	C-N-CA	5.40	124.85	119.24
1	A	290	ASN	N-CA-C	5.13	115.30	108.38
1	A	176	PRO	O-C-N	5.11	123.55	121.15

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	2443	0	2263	128	0
2	B	28	0	25	1	0
2	C	28	0	25	0	0
3	A	2	0	0	0	0
All	All	2501	0	2313	129	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 27.

All (129) 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:89:VAL:HA	1:A:93:ILE:HB	1.40	1.01
1:A:89:VAL:HA	1:A:93:ILE:CB	2.03	0.87
1:A:255:VAL:HG23	1:A:282:LEU:HB2	1.56	0.86
1:A:89:VAL:HG23	1:A:93:ILE:HG21	1.58	0.85
1:A:46:VAL:HG23	1:A:47:ILE:H	1.43	0.83
1:A:73:GLY:O	1:A:89:VAL:HG12	1.78	0.83
1:A:176:PRO:HD2	1:A:252:ARG:HH21	1.45	0.82
1:A:255:VAL:CG2	1:A:282:LEU:HB2	2.12	0.79
1:A:45:TYR:O	1:A:49:TYR:HB3	1.83	0.78
1:A:105:LEU:HD21	1:A:108:MET:HE3	1.63	0.77
1:A:129:CYS:HB2	1:A:134:ASP:HB2	1.68	0.76
1:A:89:VAL:HA	1:A:93:ILE:CG1	2.17	0.75
1:A:188:PHE:HB2	1:A:235:LEU:HD12	1.71	0.72
1:A:124:THR:HG22	1:A:126:ALA:H	1.53	0.71
1:A:42:VAL:HG11	1:A:349:LEU:HB3	1.72	0.71
1:A:89:VAL:HG23	1:A:93:ILE:CG2	2.20	0.71
1:A:98:GLU:HG3	1:A:99:ASP:OD2	1.93	0.69
1:A:74:ILE:HD12	1:A:86:ILE:HD13	1.75	0.69
1:A:82:LEU:HD12	1:A:82:LEU:H	1.58	0.69
1:A:98:GLU:HG3	1:A:99:ASP:N	2.08	0.67
2:B:1:NAG:O3	2:B:2:NAG:O5	2.13	0.66
1:A:176:PRO:HD2	1:A:252:ARG:NH2	2.12	0.65
1:A:212:ILE:HG21	1:A:217:LEU:HD21	1.77	0.65
1:A:112:THR:HG22	1:A:313:THR:HG23	1.79	0.65
1:A:49:TYR:O	1:A:53:TYR:HB3	1.98	0.64
1:A:249:MET:HE2	1:A:281:ARG:HE	1.62	0.64
1:A:155:TYR:HB2	1:A:161:THR:HG22	1.79	0.63
1:A:48:GLY:O	1:A:52:VAL:HG22	1.98	0.63
1:A:79:THR:OG1	1:A:82:LEU:HB2	1.98	0.63
1:A:51:PHE:O	1:A:55:LYS:HA	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:ILE:HG23	1:A:279:PHE:CD2	2.34	0.62
1:A:89:VAL:CA	1:A:93:ILE:HD12	2.29	0.62
1:A:222:PHE:HB2	1:A:230:CYS:O	2.00	0.62
1:A:239:VAL:HG21	1:A:246:PHE:HD1	1.66	0.61
1:A:42:VAL:O	1:A:46:VAL:HG22	2.00	0.60
1:A:105:LEU:HD21	1:A:108:MET:CE	2.32	0.60
1:A:89:VAL:CA	1:A:93:ILE:HB	2.25	0.59
1:A:97:GLN:N	1:A:97:GLN:CD	2.60	0.59
1:A:115:THR:HG22	1:A:311:THR:HG22	1.84	0.58
1:A:261:ARG:HG3	1:A:327:PHE:CE2	2.36	0.58
1:A:67:VAL:HG12	1:A:194:ASN:ND2	2.18	0.58
1:A:89:VAL:HB	1:A:93:ILE:HD12	1.86	0.57
1:A:176:PRO:HG3	1:A:318:TYR:CZ	2.39	0.57
1:A:67:VAL:HG12	1:A:194:ASN:HD22	1.69	0.57
1:A:148:ARG:HA	1:A:164:VAL:HG12	1.87	0.57
1:A:89:VAL:O	1:A:94:ILE:HG13	2.04	0.57
1:A:239:VAL:HG21	1:A:246:PHE:CD1	2.40	0.57
1:A:274:VAL:HG13	1:A:275:PRO:HD2	1.88	0.56
1:A:98:GLU:HG3	1:A:99:ASP:H	1.68	0.56
1:A:281:ARG:NH1	1:A:283:ASP:OD2	2.39	0.56
1:A:262:TRP:CH2	1:A:275:PRO:HG3	2.42	0.55
1:A:34:LEU:O	1:A:38:THR:HG23	2.07	0.55
1:A:46:VAL:HG23	1:A:47:ILE:N	2.19	0.55
1:A:89:VAL:HA	1:A:93:ILE:HD12	1.89	0.55
1:A:89:VAL:CB	1:A:93:ILE:HD12	2.37	0.55
1:A:118:LYS:HA	1:A:165:LEU:HA	1.89	0.55
1:A:259:GLN:HE21	1:A:278:THR:HB	1.72	0.55
1:A:89:VAL:O	1:A:90:ALA:O	2.25	0.54
1:A:94:ILE:HG22	1:A:96:PRO:O	2.07	0.54
1:A:47:ILE:HA	1:A:51:PHE:HD2	1.73	0.54
1:A:47:ILE:HA	1:A:51:PHE:CD2	2.43	0.54
1:A:75:ALA:HB2	1:A:188:PHE:HZ	1.73	0.53
1:A:155:TYR:HB2	1:A:161:THR:CG2	2.38	0.53
1:A:249:MET:HE2	1:A:281:ARG:HH21	1.74	0.53
1:A:135:CYS:HB2	1:A:150:GLY:O	2.10	0.52
1:A:51:PHE:CE1	1:A:57:TYR:HB3	2.45	0.52
1:A:76:LEU:HD12	1:A:85:ARG:O	2.09	0.52
1:A:255:VAL:HG23	1:A:255:VAL:O	2.09	0.52
1:A:89:VAL:HA	1:A:93:ILE:CD1	2.41	0.51
1:A:186:GLU:HG3	1:A:246:PHE:HB3	1.91	0.51
1:A:268:MET:HB3	1:A:269:PRO:HD2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ALA:N	1:A:293:PRO:HD3	2.26	0.51
1:A:98:GLU:HA	1:A:98:GLU:OE1	2.11	0.51
1:A:105:LEU:CD2	1:A:108:MET:HE3	2.38	0.50
1:A:268:MET:HE3	1:A:269:PRO:HD2	1.93	0.50
1:A:351:LEU:O	1:A:351:LEU:HD13	2.11	0.50
1:A:87:TRP:CG	1:A:182:LEU:HD11	2.46	0.50
1:A:97:GLN:N	1:A:97:GLN:OE1	2.45	0.50
1:A:46:VAL:HG23	1:A:47:ILE:HG13	1.94	0.49
1:A:226:THR:O	1:A:228:PRO:HD3	2.12	0.49
1:A:249:MET:CE	1:A:281:ARG:HE	2.23	0.49
1:A:63:VAL:HG23	1:A:328:GLY:O	2.13	0.49
1:A:168:CYS:HA	1:A:169:PRO:C	2.38	0.48
1:A:51:PHE:CD1	1:A:57:TYR:HB3	2.49	0.48
1:A:61:ASP:OD2	1:A:198:TYR:HB3	2.13	0.48
1:A:88:ASP:OD1	1:A:92:TYR:HB2	2.14	0.48
1:A:201:PHE:O	1:A:202:ASN:C	2.56	0.48
1:A:42:VAL:HA	1:A:45:TYR:HB3	1.95	0.48
1:A:249:MET:HE2	1:A:281:ARG:NE	2.27	0.47
1:A:75:ALA:HB2	1:A:188:PHE:CZ	2.50	0.47
1:A:219:HIS:O	1:A:219:HIS:CG	2.67	0.47
1:A:239:VAL:O	1:A:242:ALA:HB3	2.15	0.47
1:A:222:PHE:HA	1:A:230:CYS:HB3	1.97	0.47
1:A:95:PRO:HA	1:A:96:PRO:HA	1.69	0.46
1:A:85:ARG:HD3	1:A:87:TRP:CE2	2.51	0.46
1:A:281:ARG:HD2	1:A:283:ASP:OD1	2.15	0.46
1:A:55:LYS:HE2	1:A:58:GLN:HE22	1.79	0.46
1:A:268:MET:HE2	1:A:272:TRP:CZ3	2.51	0.46
1:A:269:PRO:HG2	1:A:272:TRP:CD1	2.51	0.46
1:A:239:VAL:CG2	1:A:246:PHE:HD1	2.29	0.46
1:A:86:ILE:CG2	1:A:87:TRP:N	2.79	0.45
1:A:154:SER:HA	1:A:160:LYS:HA	1.98	0.45
1:A:213:ASN:ND2	1:A:213:ASN:H	2.15	0.44
1:A:85:ARG:HD3	1:A:87:TRP:NE1	2.32	0.44
1:A:166:SER:OG	1:A:167:TRP:N	2.51	0.44
1:A:238:ILE:HG23	1:A:279:PHE:CE2	2.53	0.44
1:A:89:VAL:O	1:A:90:ALA:C	2.58	0.43
1:A:336:ILE:N	1:A:337:PRO:HD2	2.33	0.43
1:A:82:LEU:H	1:A:82:LEU:CD1	2.30	0.43
1:A:349:LEU:HD12	1:A:350:GLY:N	2.33	0.43
1:A:191:LEU:HD12	1:A:192:ILE:N	2.34	0.43
1:A:286:ASP:OD1	1:A:287:PRO:HD2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LYS:HE2	1:A:58:GLN:NE2	2.34	0.43
1:A:119:CYS:SG	1:A:170:LEU:CD2	3.07	0.43
1:A:274:VAL:HA	1:A:275:PRO:HD3	1.81	0.43
1:A:209:LEU:HA	1:A:210:PRO:HD3	1.90	0.42
1:A:98:GLU:HB3	1:A:101:SER:O	2.19	0.42
1:A:145:ASP:HB3	1:A:170:LEU:HD11	2.02	0.42
1:A:93:ILE:O	1:A:95:PRO:HD3	2.18	0.42
1:A:254:GLY:O	1:A:320:ILE:HA	2.19	0.42
1:A:64:LEU:HD12	1:A:64:LEU:HA	1.89	0.42
1:A:230:CYS:HA	1:A:231:PRO:HD3	1.75	0.42
1:A:269:PRO:C	1:A:271:SER:H	2.28	0.42
1:A:93:ILE:O	1:A:93:ILE:HG22	2.19	0.41
1:A:102:PHE:CE2	1:A:322:PHE:HB2	2.55	0.41
1:A:300:ALA:HB1	1:A:312:ARG:HD2	2.02	0.41
1:A:74:ILE:CD1	1:A:86:ILE:HD13	2.49	0.41
1:A:42:VAL:HG21	1:A:349:LEU:HD23	2.02	0.40
1:A:88:ASP:O	1:A:89:VAL:C	2.64	0.40

There are no symmetry-related clashes.

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	325/356 (91%)	300 (92%)	20 (6%)	5 (2%)	8 32

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	ALA
1	A	101	SER
1	A	95	PRO
1	A	46	VAL

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Mol	Chain	Res	Type
1	A	96	PRO

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	250/314 (80%)	233 (93%)	17 (7%)	14	42

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	VAL
1	A	74	ILE
1	A	86	ILE
1	A	94	ILE
1	A	111	THR
1	A	174	VAL
1	A	181	LEU
1	A	189	THR
1	A	191	LEU
1	A	204	ASN
1	A	213	ASN
1	A	218	THR
1	A	248	ILE
1	A	278	THR
1	A	291	VAL
1	A	298	ARG
1	A	310	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	219	HIS
1	A	259	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	B	1	2,1	14,14,15	0.86	1 (7%)	17,19,21	2.82	6 (35%)
2	NAG	B	2	2	14,14,15	1.02	1 (7%)	17,19,21	3.19	10 (58%)
2	NAG	C	1	2,1	14,14,15	0.49	0	17,19,21	2.32	4 (23%)
2	NAG	C	2	2	14,14,15	0.53	0	17,19,21	2.38	6 (35%)

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	B	1	2,1	1/1/5/7	2/6/23/26	0/1/1/1
2	NAG	B	2	2	1/1/5/7	4/6/23/26	0/1/1/1
2	NAG	C	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	C	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	B	2	NAG	C1-C2	3.28	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAG	O4-C4	2.01	1.47	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	O4-C4-C5	-6.57	93.15	109.32
2	B	2	NAG	C1-O5-C5	6.56	120.98	112.19
2	C	1	NAG	C1-O5-C5	6.27	120.59	112.19
2	B	1	NAG	O3-C3-C2	-6.10	96.73	109.40
2	B	1	NAG	C1-O5-C5	5.60	119.69	112.19
2	B	1	NAG	O5-C1-C2	4.65	118.49	111.29
2	C	2	NAG	O5-C1-C2	4.64	118.47	111.29
2	C	2	NAG	C1-O5-C5	4.46	118.17	112.19
2	C	1	NAG	C2-N2-C7	-4.16	117.33	122.90
2	B	2	NAG	O5-C5-C6	4.04	115.53	107.66
2	B	2	NAG	O5-C5-C4	3.95	120.45	110.83
2	C	2	NAG	C2-N2-C7	-3.93	117.63	122.90
2	B	2	NAG	O3-C3-C2	3.72	117.12	109.40
2	C	1	NAG	O5-C1-C2	3.67	116.96	111.29
2	B	1	NAG	O5-C5-C4	3.51	119.36	110.83
2	B	1	NAG	C2-N2-C7	3.49	127.58	122.90
2	C	2	NAG	O5-C5-C4	3.48	119.29	110.83
2	C	1	NAG	O5-C5-C4	3.36	118.99	110.83
2	B	2	NAG	C2-N2-C7	3.18	127.16	122.90
2	B	1	NAG	O4-C4-C3	2.89	117.20	110.38
2	C	2	NAG	O5-C5-C6	2.82	113.14	107.66
2	B	2	NAG	C4-C3-C2	2.60	114.83	111.02
2	B	2	NAG	C1-C2-N2	-2.44	106.58	110.43
2	B	2	NAG	O7-C7-N2	2.20	125.86	121.98
2	C	2	NAG	O3-C3-C2	-2.17	104.89	109.40
2	B	2	NAG	O5-C1-C2	2.10	114.54	111.29

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1	NAG	C1
2	B	2	NAG	C1

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	NAG	C8-C7-N2-C2

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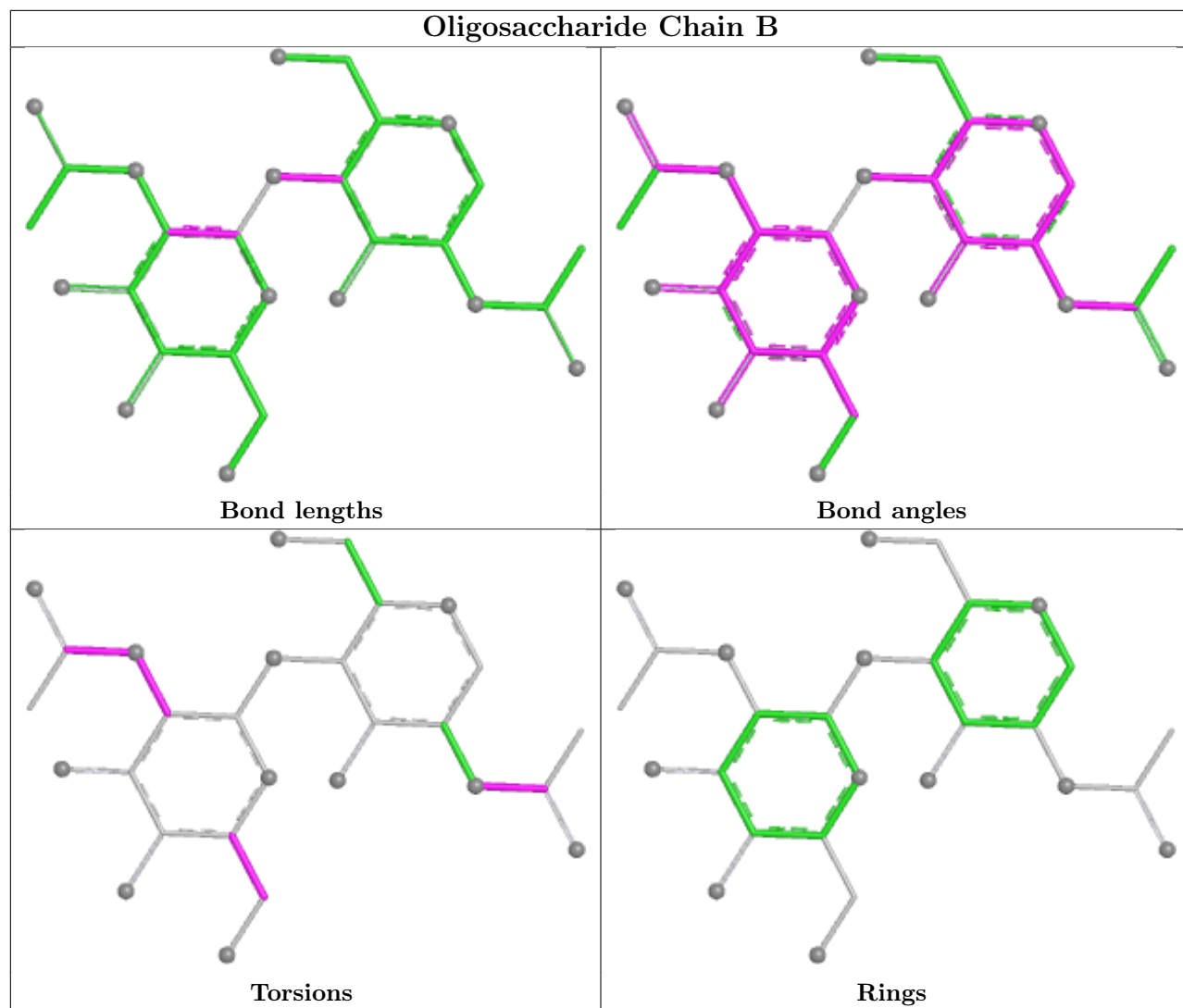
Mol	Chain	Res	Type	Atoms
2	B	1	NAG	O7-C7-N2-C2
2	B	2	NAG	C3-C2-N2-C7
2	B	2	NAG	C8-C7-N2-C2
2	C	1	NAG	C8-C7-N2-C2
2	B	2	NAG	O7-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	B	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2

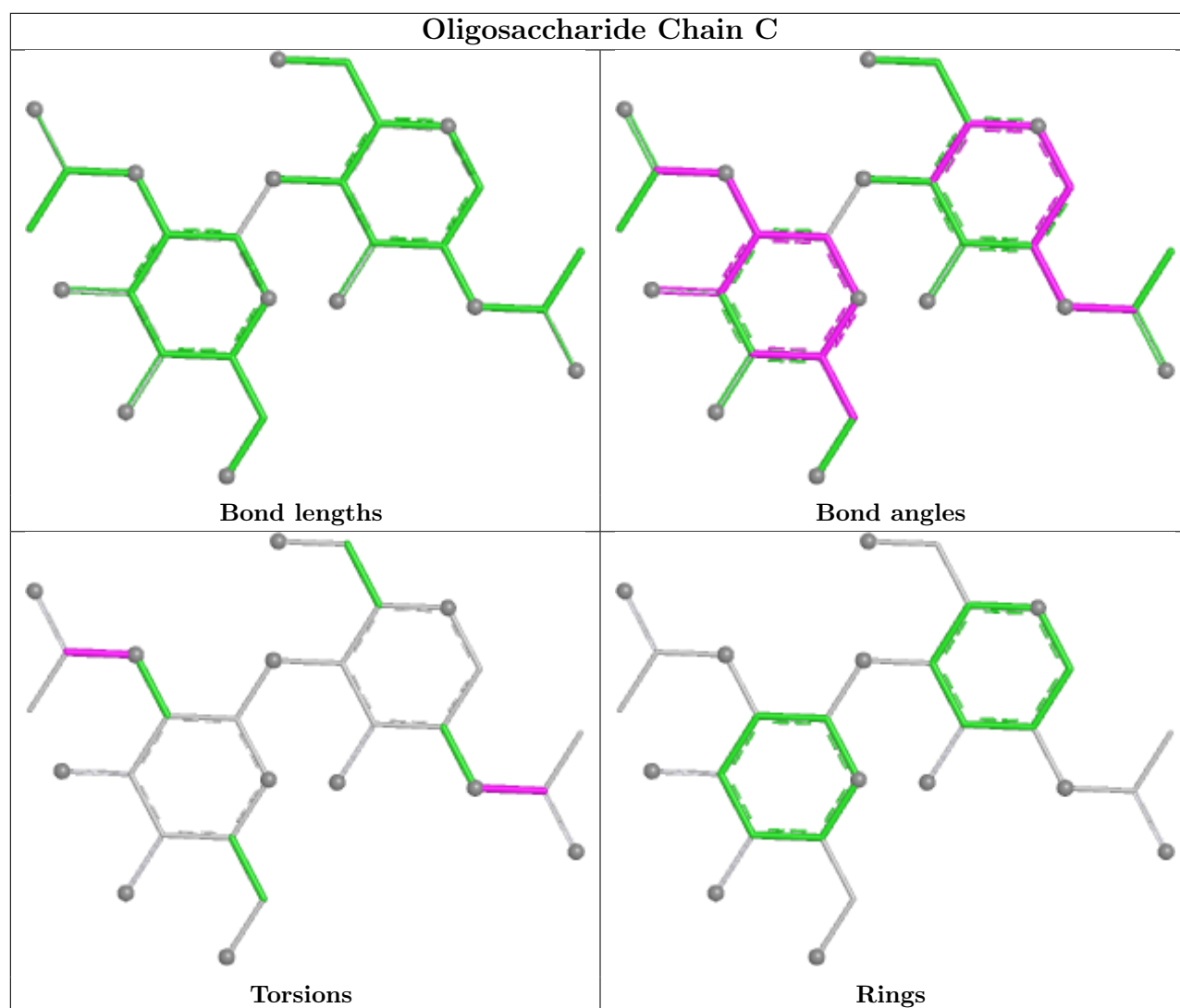
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	NAG	1	0
2	B	1	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](#)

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(Å ²)	Q<0.9
1	A	329/356 (92%)	0.86	46 (13%) 6 4	64, 99, 171, 191	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	142	ALA	8.0
1	A	53	TYR	4.6
1	A	344	ALA	4.1
1	A	138	GLY	3.5
1	A	90	ALA	3.4
1	A	263	ASP	3.4
1	A	139	PHE	3.3
1	A	165	LEU	3.0
1	A	148	ARG	3.0
1	A	336	ILE	2.8
1	A	268	MET	2.8
1	A	92	TYR	2.7
1	A	141	ASP	2.6
1	A	146	GLY	2.6
1	A	130	THR	2.6
1	A	52	VAL	2.6
1	A	74	ILE	2.5
1	A	347	ALA	2.5
1	A	343	GLY	2.5
1	A	306	SER	2.5
1	A	150	GLY	2.5
1	A	158	SER	2.4
1	A	350	GLY	2.4
1	A	305	ASN	2.4
1	A	349	LEU	2.4
1	A	310	GLU	2.3
1	A	151	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	348	LEU	2.3
1	A	136	LYS	2.3
1	A	32	GLY	2.3
1	A	117	SER	2.3
1	A	277	TYR	2.3
1	A	50	VAL	2.3
1	A	346	LEU	2.2
1	A	331	GLY	2.2
1	A	149	THR	2.2
1	A	49	TYR	2.2
1	A	329	GLN	2.1
1	A	272	TRP	2.1
1	A	243	GLU	2.1
1	A	116	GLN	2.1
1	A	270	GLN	2.1
1	A	42	VAL	2.0
1	A	86	ILE	2.0
1	A	342	ILE	2.0
1	A	288	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](#)

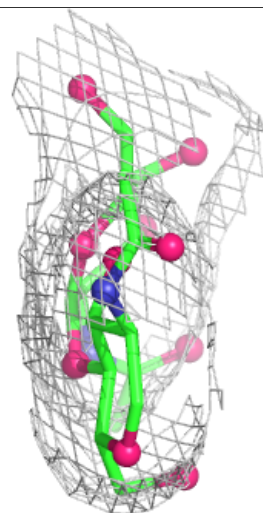
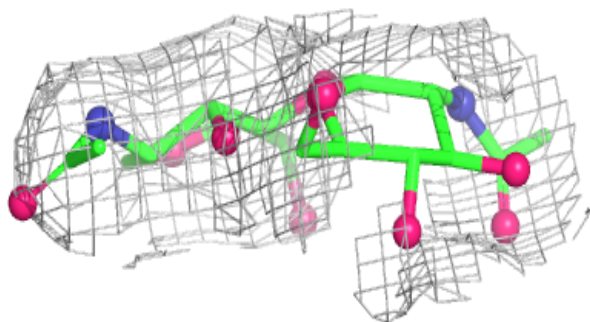
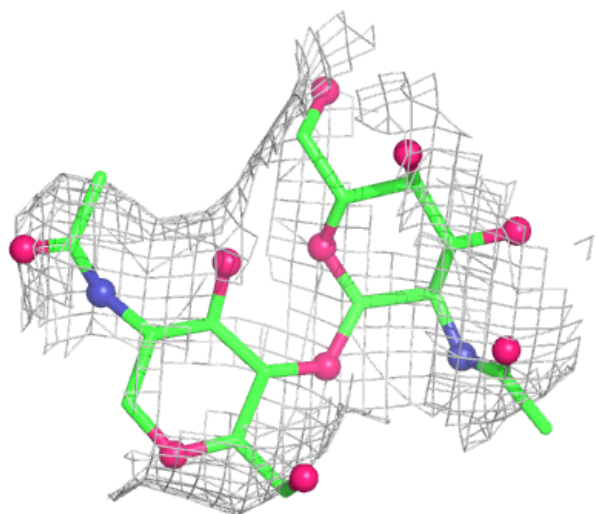
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

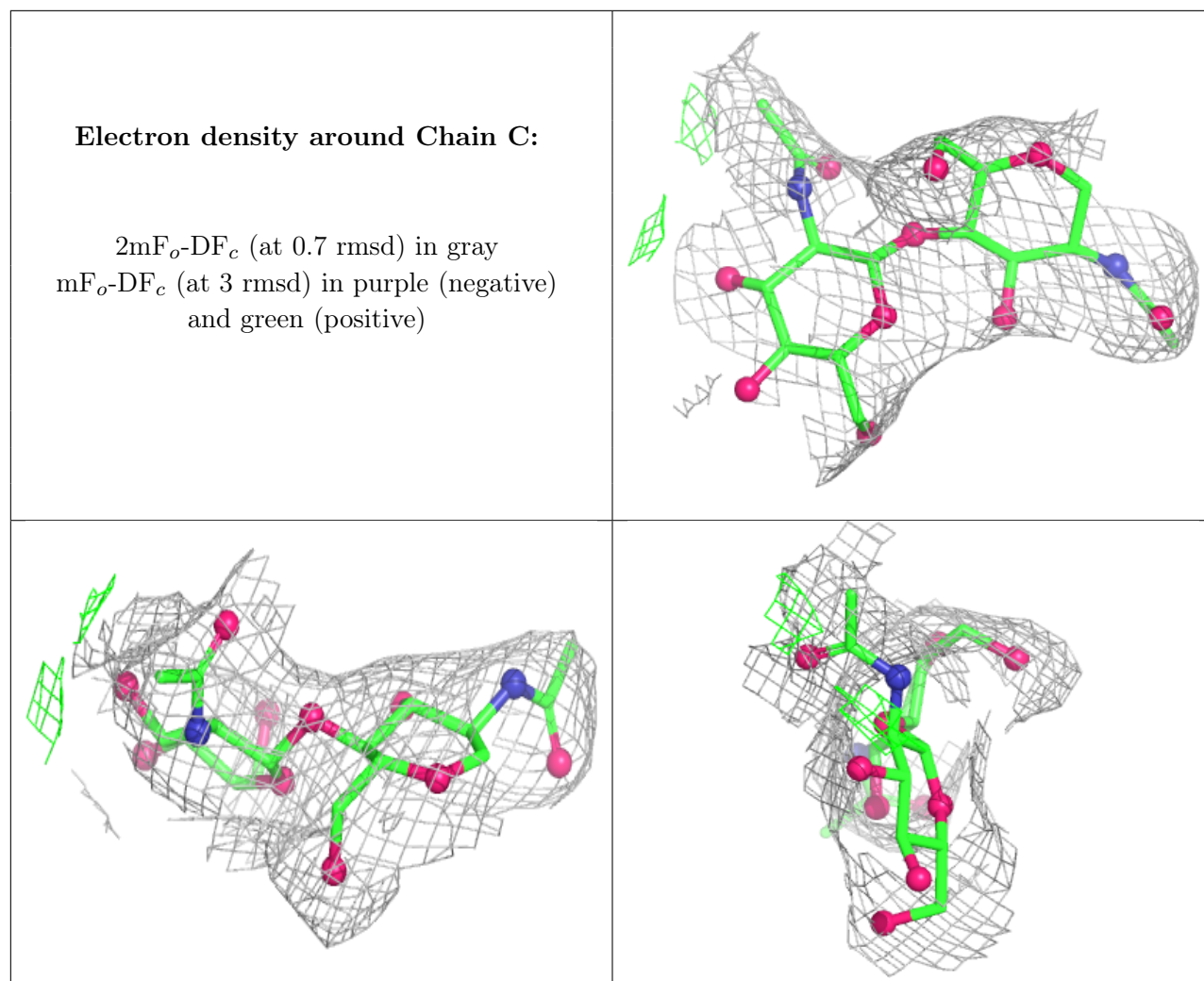
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	C	2	14/15	0.50	0.16	114,135,140,143	0
2	NAG	B	2	14/15	0.53	0.15	159,168,171,171	0
2	NAG	B	1	14/15	0.82	0.13	121,152,161,164	0
2	NAG	C	1	14/15	0.89	0.14	114,121,127,131	0

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 B:

$2mF_o-DF_c$ (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
3	GD	A	383	1/1	0.92	0.10	168,168,168,168	1
3	GD	A	382	1/1	0.96	0.26	152,152,152,152	1

6.5 Other polymers [i](#)

There are no such residues in this entry.