



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:16 PM UTC

PDB ID : 2XQF / pdb_00002xqf
Title : X-ray Structure of human butyrylcholinesterase inhibited by racemic VX
Authors : Wandhammer, M.; Carletti, E.; Gillon, E.; Masson, P.; Goeldner, M.; Noort, D.; Nachon, F.
Deposited on : 2010-09-02
Resolution : 2.10 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

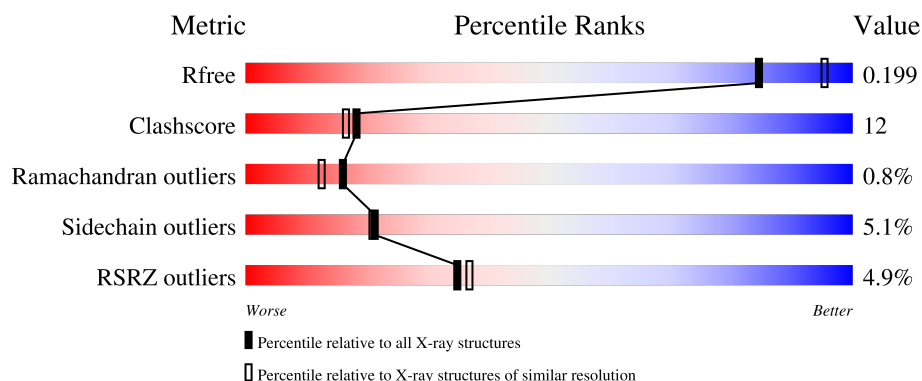
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	527	5% 77% 17% ..
2	B	3	100%
2	C	3	33% 67%

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	C	1	X	-	-	-
4	UNX	A	1554	-	-	X	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 4894 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.

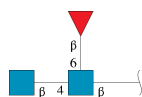
- Molecule 1 is a protein called CHOLINESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	527	4269	2759	712	782	16	0	11	0

There are 4 discrepancies between the modelled and reference sequences:

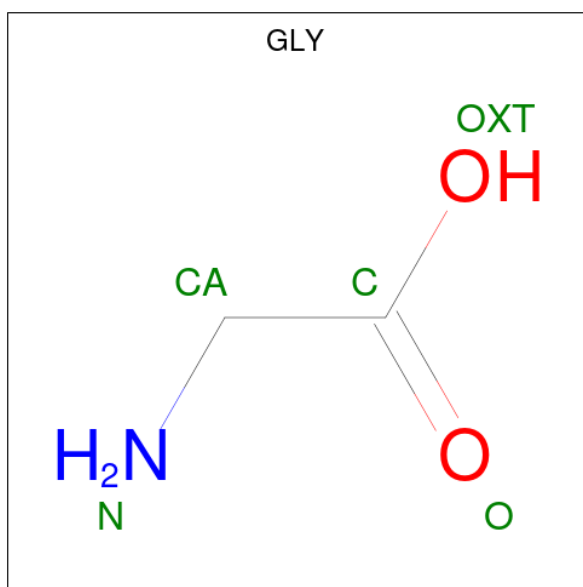
Chain	Residue	Modelled	Actual	Comment	Reference
A	17	GLN	ASN	engineered mutation	UNP P06276
A	455	GLN	ASN	engineered mutation	UNP P06276
A	481	GLN	ASN	engineered mutation	UNP P06276
A	486	GLN	ASN	engineered mutation	UNP P06276

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	3	38	22	2	14	0	0	0
2	C	3	38	22	2	14	0	0	0

- Molecule 3 is GLYCINE (CCD ID: GLY) (formula: C₂H₅NO₂).

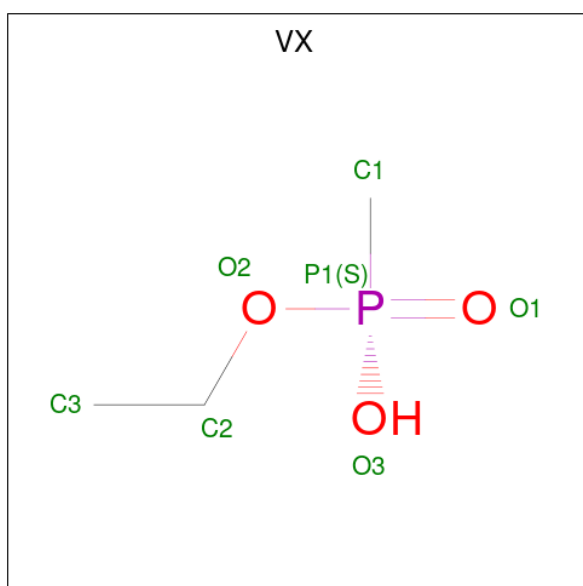


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			5	2	1	2		

- Molecule 4 is UNKNOWN ATOM OR ION (CCD ID: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	36	Total	X	0	0
			36	36		

- Molecule 5 is O-ETHYLMETHYLPHOSPHONIC ACID ESTER GROUP (CCD ID: VX) (formula: C₃H₉O₃P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	P	0	0
			6	3	2	1		

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

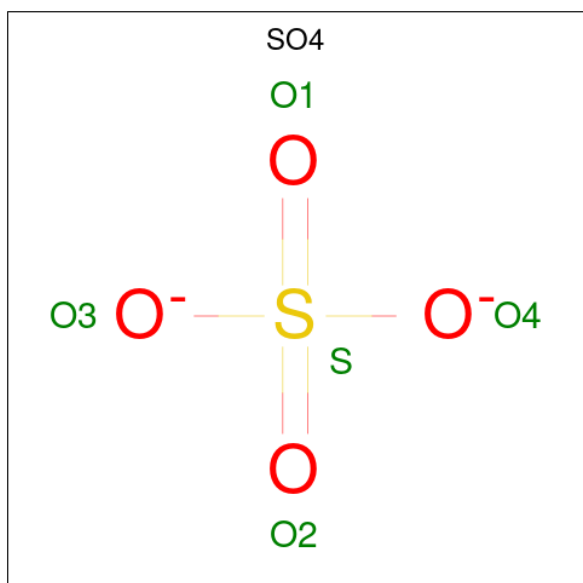
- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Cl	0	0
			2	2		

- Molecule 8 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	K	0	0
			1	1		

- Molecule 9 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



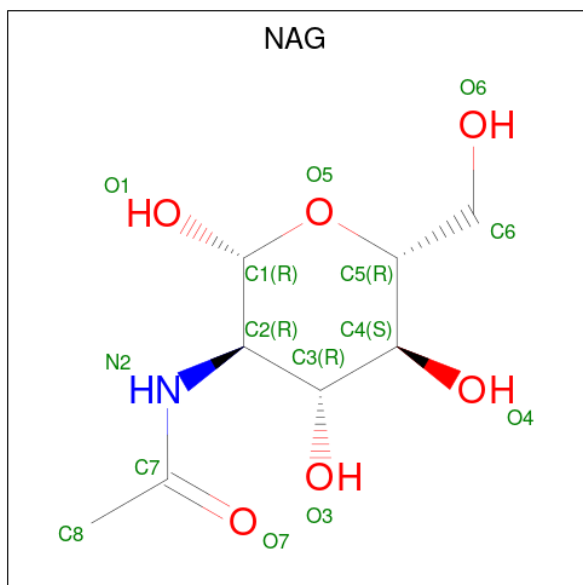
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

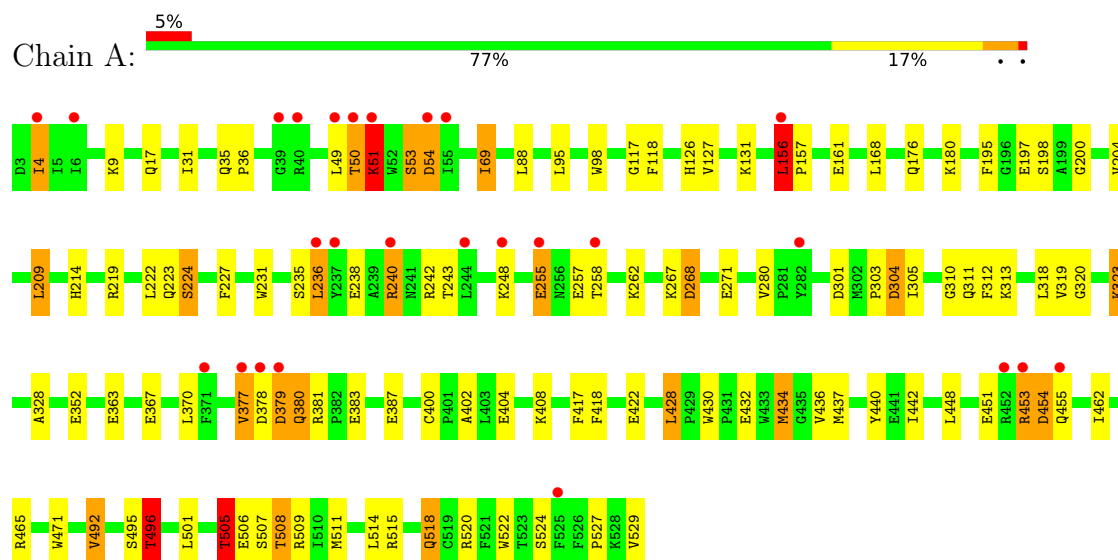
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	432	Total	O	0	0
			432	432		

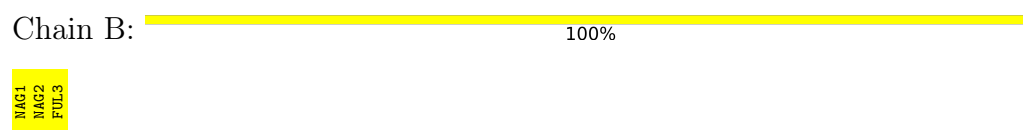
3 Residue-property plots [i](#)

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: CHOLINESTERASE



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	155.11Å 155.11Å 128.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.00 – 2.10 48.00 – 2.39	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.00-2.10) 99.5 (48.00-2.39)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.12 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.149 , 0.189 0.199 , 0.199	Depositor DCC
R_{free} test set	1242 reflections (4.00%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.762	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4894	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: FUL, K, SO4, NAG, VX, UNX, NA, CL

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	1.62	32/4419 (0.7%)	1.28	17/5997 (0.3%)

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	462	ILE	CA-CB	8.31	1.64	1.54
1	A	127	VAL	CA-CB	8.20	1.64	1.54
1	A	36	PRO	CA-C	7.84	1.59	1.52
1	A	204	VAL	CA-CB	7.24	1.62	1.54
1	A	434	MET	SD-CE	-6.99	1.62	1.79
1	A	377	VAL	CA-CB	6.69	1.60	1.53
1	A	422	GLU	CA-CB	6.66	1.62	1.53
1	A	304	ASP	CA-C	6.55	1.61	1.52
1	A	320	GLY	N-CA	6.50	1.52	1.45
1	A	126	HIS	CA-CB	6.42	1.63	1.53
1	A	219	ARG	CG-CD	6.20	1.71	1.52
1	A	492	VAL	C-O	5.79	1.32	1.24
1	A	328	ALA	CA-CB	5.79	1.62	1.53
1	A	440	TYR	CA-CB	5.77	1.61	1.53
1	A	31	ILE	N-CA	5.74	1.52	1.46
1	A	436	VAL	CA-CB	5.71	1.60	1.53
1	A	508	THR	CA-CB	-5.64	1.44	1.53
1	A	515	ARG	CZ-NH1	5.62	1.40	1.32
1	A	402	ALA	N-CA	5.50	1.53	1.46
1	A	305	ILE	N-CA	5.43	1.52	1.46
1	A	301	ASP	C-O	-5.43	1.17	1.23
1	A	514	LEU	CA-C	5.42	1.59	1.52
1	A	465	ARG	CZ-NH1	5.40	1.40	1.32
1	A	219	ARG	CD-NE	-5.36	1.38	1.46
1	A	131	LYS	CA-C	5.33	1.59	1.52
1	A	118	PHE	N-CA	5.28	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	505	THR	N-CA	-5.20	1.39	1.46
1	A	224	SER	C-O	5.18	1.30	1.23
1	A	320	GLY	C-O	5.09	1.29	1.23
1	A	524	SER	CA-C	-5.08	1.46	1.52
1	A	442	ILE	CA-CB	5.06	1.60	1.54
1	A	323	LYS	C-O	5.03	1.30	1.24

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	508	THR	N-CA-CB	-7.89	98.10	109.85
1	A	219	ARG	NE-CZ-NH1	-6.58	114.92	121.50
1	A	53	SER	N-CA-C	-6.05	100.71	110.32
1	A	505	THR	N-CA-CB	-5.99	101.07	110.30
1	A	507	SER	CA-C-N	5.90	129.21	120.95
1	A	507	SER	C-N-CA	5.90	129.21	120.95
1	A	219	ARG	CG-CD-NE	-5.79	99.27	112.00
1	A	156	LEU	CA-C-N	-5.40	114.20	119.76
1	A	156	LEU	C-N-CA	-5.40	114.20	119.76
1	A	400	CYS	CA-C-N	-5.26	113.12	118.85
1	A	400	CYS	C-N-CA	-5.26	113.12	118.85
1	A	168	LEU	N-CA-C	-5.16	105.74	111.36
1	A	524	SER	CA-C-N	-5.09	114.53	122.42
1	A	524	SER	C-N-CA	-5.09	114.53	122.42
1	A	200	GLY	N-CA-C	-5.06	106.29	112.77
1	A	471	TRP	N-CA-C	-5.02	105.89	111.36
1	A	496	THR	N-CA-CB	5.02	118.97	110.49

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	4269	0	4183	100	0
2	B	38	0	33	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	38	0	34	2	0
3	A	5	0	2	1	0
4	A	36	0	0	4	0
5	A	6	0	8	2	0
6	A	1	0	0	0	0
7	A	2	0	0	0	0
8	A	1	0	0	0	0
9	A	10	0	0	0	0
10	A	56	0	52	1	0
11	A	432	0	0	30	4
All	All	4894	0	4312	106	4

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 (106) 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:455[A]:GLN:NE2	1:A:455[A]:GLN:CD	2.07	1.13
1:A:495:SER:HA	11:A:2385:HOH:O	1.51	1.10
1:A:4:ILE:H	1:A:4:ILE:HD12	1.23	1.01
1:A:35:GLN:HG2	11:A:2025:HOH:O	1.59	0.99
1:A:496:THR:N	11:A:2385:HOH:O	1.75	0.99
1:A:495:SER:CA	11:A:2385:HOH:O	2.03	0.98
1:A:255:GLU:H	1:A:255:GLU:CD	1.75	0.95
1:A:4:ILE:HD12	1:A:4:ILE:N	1.82	0.94
1:A:268:ASP:HB2	11:A:2199:HOH:O	1.66	0.94
1:A:518:GLN:HE21	1:A:518:GLN:H	1.18	0.92
1:A:451:GLU:OE1	11:A:2345:HOH:O	1.92	0.87
1:A:176[B]:GLN:HE21	1:A:180:LYS:HG3	1.40	0.85
1:A:505:THR:HG21	11:A:2153:HOH:O	1.77	0.85
1:A:380:GLN:HA	1:A:380:GLN:NE2	1.97	0.80
1:A:262:LYS:HE3	11:A:2197:HOH:O	1.80	0.79
1:A:304:ASP:HB3	11:A:2091:HOH:O	1.84	0.78
1:A:156:LEU:HD13	1:A:243:THR:HG21	1.64	0.78
1:A:311:GLN:NE2	11:A:2238:HOH:O	1.91	0.75
1:A:380:GLN:HA	1:A:380:GLN:HE21	1.54	0.73
1:A:379:ASP:O	1:A:380:GLN:NE2	2.23	0.72
1:A:408:LYS:HE3	11:A:2311:HOH:O	1.90	0.71
1:A:495:SER:O	1:A:496:THR:OG1	2.08	0.71
1:A:451:GLU:OE2	11:A:2343:HOH:O	2.10	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:GLU:HG2	11:A:2184:HOH:O	1.93	0.69
1:A:117:GLY:HA2	5:A:1530:VX:H22	1.73	0.69
1:A:4:ILE:H	1:A:4:ILE:CD1	2.03	0.68
1:A:98:TRP:HE1	3:A:1531:GLY:N	1.92	0.68
1:A:176[B]:GLN:NE2	1:A:180:LYS:HG3	2.08	0.67
1:A:323:LYS:HE2	11:A:2250:HOH:O	1.94	0.67
4:A:1580:UNX:UNK	4:A:1581:UNX:UNK	1.38	0.67
1:A:248:LYS:HD3	2:C:3:FUL:H61	1.75	0.67
1:A:4:ILE:N	1:A:4:ILE:CD1	2.55	0.67
1:A:377:VAL:HA	1:A:378:ASP:HB2	1.77	0.65
1:A:311:GLN:HB3	11:A:2238:HOH:O	1.95	0.65
1:A:238[A]:GLU:OE1	1:A:242:ARG:NE	2.25	0.65
1:A:518:GLN:H	1:A:518:GLN:NE2	1.92	0.64
1:A:434:MET:HE2	1:A:437:MET:HE3	1.79	0.63
1:A:381:ARG:HB3	1:A:383:GLU:OE1	1.98	0.63
4:A:1554:UNX:UNK	4:A:1562:UNX:UNK	1.44	0.61
1:A:53:SER:O	1:A:54:ASP:OD1	2.18	0.61
1:A:408:LYS:HG3	11:A:2311:HOH:O	2.00	0.60
1:A:267:LYS:HE3	1:A:271:GLU:OE2	2.02	0.60
1:A:227:PHE:CD1	1:A:227:PHE:C	2.79	0.60
1:A:304:ASP:OD1	11:A:2232:HOH:O	2.16	0.59
1:A:176[B]:GLN:NE2	1:A:180:LYS:HE3	2.18	0.58
1:A:51:LYS:N	11:A:2044:HOH:O	2.36	0.58
1:A:238[A]:GLU:CD	1:A:242:ARG:HE	2.12	0.57
1:A:53:SER:O	1:A:54:ASP:CG	2.48	0.57
10:A:1542:NAG:H83	11:A:2047:HOH:O	2.05	0.56
1:A:255:GLU:CD	1:A:255:GLU:N	2.55	0.55
1:A:511[B]:MET:HE1	11:A:2320:HOH:O	2.06	0.55
1:A:209:LEU:HD12	1:A:312:PHE:HB3	1.88	0.54
1:A:522:TRP:O	1:A:527:PRO:HD3	2.07	0.54
1:A:227:PHE:CE1	1:A:303:PRO:HB2	2.42	0.54
1:A:235:SER:OG	1:A:238[B]:GLU:HG3	2.08	0.54
1:A:509:ARG:NH1	11:A:2403:HOH:O	2.24	0.53
1:A:363:GLU:O	1:A:367[B]:GLU:HG3	2.09	0.53
1:A:50:THR:O	1:A:51:LYS:HB3	2.09	0.53
1:A:434:MET:HE2	1:A:437:MET:CE	2.38	0.52
1:A:432[B]:GLU:HG2	11:A:2334:HOH:O	2.08	0.52
1:A:318:LEU:C	1:A:318:LEU:HD23	2.35	0.51
1:A:198:SER:HA	1:A:224:SER:O	2.12	0.50
1:A:267:LYS:HE3	1:A:271:GLU:CD	2.37	0.49
1:A:267:LYS:CE	1:A:271:GLU:CD	2.85	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238[A]:GLU:OE1	1:A:242:ARG:HG3	2.12	0.49
1:A:69:ILE:HD11	1:A:88:LEU:HD11	1.95	0.48
1:A:95:LEU:HD12	1:A:95:LEU:C	2.38	0.48
1:A:383:GLU:O	1:A:387:GLU:HG3	2.13	0.48
1:A:50:THR:O	1:A:51:LYS:CB	2.63	0.47
1:A:161:GLU:HG3	1:A:258:THR:HG23	1.97	0.47
1:A:231:TRP:CE3	5:A:1530:VX:H21	2.50	0.47
1:A:236:LEU:HD12	1:A:236:LEU:HA	1.81	0.46
1:A:352:GLU:HG3	11:A:2268:HOH:O	2.15	0.46
1:A:495:SER:O	1:A:496:THR:CB	2.63	0.46
1:A:313:LYS:NZ	11:A:2244:HOH:O	2.41	0.46
1:A:501:LEU:HD11	1:A:508:THR:HG22	1.97	0.45
1:A:176[B]:GLN:HE21	1:A:180:LYS:HE3	1.82	0.45
1:A:197:GLU:HA	1:A:223:GLN:O	2.17	0.44
1:A:319:VAL:O	1:A:418:PHE:HA	2.17	0.44
1:A:255:GLU:N	1:A:255:GLU:OE1	2.44	0.44
1:A:455[A]:GLN:NE2	1:A:455[A]:GLN:CG	2.80	0.44
1:A:17:GLN:HE21	1:A:17:GLN:HA	1.83	0.44
1:A:157:PRO:CD	1:A:240:ARG:HG2	2.47	0.44
4:A:1577:UNX:UNK	4:A:1578:UNX:UNK	1.60	0.44
1:A:511[B]:MET:CE	11:A:2320:HOH:O	2.63	0.43
1:A:434:MET:CE	1:A:437:MET:HE3	2.48	0.43
1:A:428:LEU:HD13	1:A:430:TRP:HB2	2.01	0.43
1:A:214:HIS:HE1	11:A:2245:HOH:O	2.00	0.43
1:A:454:ASP:O	1:A:455[A]:GLN:HB2	2.18	0.43
1:A:520:ARG:NH2	11:A:2416:HOH:O	2.52	0.42
1:A:157:PRO:HD3	1:A:240:ARG:HG2	2.01	0.42
1:A:240:ARG:NH1	1:A:257:GLU:OE2	2.53	0.42
1:A:417:PHE:CE2	1:A:492:VAL:HG12	2.54	0.42
1:A:451:GLU:OE1	11:A:2344:HOH:O	2.22	0.42
1:A:378:ASP:O	1:A:379:ASP:C	2.62	0.41
4:A:1554:UNX:UNK	4:A:1555:UNX:UNK	1.64	0.41
1:A:501:LEU:HD11	1:A:508:THR:CG2	2.50	0.41
1:A:505:THR:HG23	1:A:506[B]:GLU:OE2	2.21	0.41
1:A:383:GLU:CD	1:A:383:GLU:H	2.29	0.41
1:A:222:LEU:HD12	1:A:222:LEU:N	2.36	0.41
1:A:378:ASP:O	1:A:379:ASP:O	2.39	0.40
1:A:454:ASP:OD2	11:A:2348:HOH:O	2.22	0.40
2:C:2:NAG:H5	2:C:2:NAG:N2	2.36	0.40
1:A:156:LEU:HA	1:A:156:LEU:HD12	1.76	0.40
1:A:310:GLY:HA2	1:A:312:PHE:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:453:ARG:HG3	11:A:2345:HOH:O	2.20	0.40

All (4) 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)
11:A:2240:HOH:O	11:A:2365:HOH:O[3_555]	1.86	0.34
11:A:2239:HOH:O	11:A:2365:HOH:O[3_555]	2.05	0.15
11:A:2179:HOH:O	11:A:2239:HOH:O[4_555]	2.14	0.06
11:A:2233:HOH:O	11:A:2368:HOH:O[3_555]	2.19	0.01

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	536/527 (102%)	516 (96%)	16 (3%)	4 (1%)	18 15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	379	ASP
1	A	496	THR
1	A	51	LYS
1	A	54	ASP

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	463/452 (102%)	440 (95%)	23 (5%)	22	22

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	9	LYS
1	A	49	LEU
1	A	50	THR
1	A	51	LYS
1	A	69	ILE
1	A	156	LEU
1	A	195	PHE
1	A	209	LEU
1	A	236	LEU
1	A	240	ARG
1	A	255	GLU
1	A	268	ASP
1	A	280	VAL
1	A	370	LEU
1	A	380	GLN
1	A	428	LEU
1	A	448	LEU
1	A	453	ARG
1	A	454	ASP
1	A	505	THR
1	A	518	GLN
1	A	529	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	179	GLN
1	A	214	HIS
1	A	275	ASN
1	A	289	ASN
1	A	380	GLN
1	A	518	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 ⓘ

6 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	1,2	14,14,15	1.71	3 (21%)	17,19,21	2.11	5 (29%)
2	NAG	B	2	2	14,14,15	0.59	0	17,19,21	2.32	6 (35%)
2	FUL	B	3	2	10,10,11	1.11	1 (10%)	14,14,16	3.13	4 (28%)
2	NAG	C	1	1,2	14,14,15	0.79	0	17,19,21	2.04	5 (29%)
2	NAG	C	2	2	14,14,15	1.57	2 (14%)	17,19,21	2.50	6 (35%)
2	FUL	C	3	2	10,10,11	0.84	0	14,14,16	2.45	4 (28%)

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	1,2	-	4/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	FUL	B	3	2	-	-	0/1/1/1
2	NAG	C	1	1,2	1/1/5/7	3/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	FUL	C	3	2	-	-	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	NAG	C1-C2	4.53	1.58	1.52
2	B	1	NAG	C2-N2	-3.51	1.40	1.46
2	B	1	NAG	C1-C2	3.18	1.56	1.52
2	B	1	NAG	O5-C1	-3.13	1.38	1.43
2	C	2	NAG	C2-N2	2.91	1.51	1.46
2	B	3	FUL	C4-C5	2.10	1.57	1.52

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	FUL	C1-C2-C3	-9.40	95.96	109.64
2	C	2	NAG	C2-N2-C7	7.64	133.13	122.90
2	B	2	NAG	C3-C4-C5	-5.84	99.65	110.23
2	C	3	FUL	C1-C2-C3	-5.76	101.26	109.64
2	C	1	NAG	C1-O5-C5	5.52	119.58	112.19
2	B	1	NAG	C1-C2-N2	-4.43	103.45	110.43
2	B	3	FUL	C1-O5-C5	-4.40	102.59	112.97
2	B	1	NAG	C3-C4-C5	-4.10	102.80	110.23
2	C	3	FUL	C1-O5-C5	-4.02	103.48	112.97
2	C	3	FUL	C3-C4-C5	3.50	115.14	109.81
2	C	1	NAG	C1-C2-N2	-3.43	105.03	110.43
2	B	2	NAG	O5-C5-C6	3.34	114.16	107.66
2	B	2	NAG	O5-C5-C4	-3.19	103.06	110.83
2	B	3	FUL	O5-C5-C6	3.19	114.32	107.40
2	C	3	FUL	O2-C2-C1	3.13	116.39	109.22
2	B	1	NAG	C1-O5-C5	3.08	116.31	112.19
2	C	1	NAG	C2-N2-C7	3.03	126.95	122.90
2	B	2	NAG	C2-N2-C7	-3.00	118.87	122.90
2	C	2	NAG	O5-C5-C6	2.90	113.30	107.66
2	C	2	NAG	O3-C3-C4	2.73	116.80	110.38
2	B	3	FUL	C6-C5-C4	2.69	118.00	113.08
2	C	2	NAG	C1-C2-N2	2.53	114.42	110.43
2	B	1	NAG	C2-N2-C7	2.49	126.24	122.90
2	C	1	NAG	C8-C7-N2	2.40	120.10	116.12
2	C	2	NAG	O7-C7-N2	2.39	126.21	121.98
2	B	1	NAG	C6-C5-C4	2.25	118.54	113.02
2	B	2	NAG	O4-C4-C3	2.19	115.55	110.38
2	C	2	NAG	C4-C3-C2	-2.17	107.84	111.02
2	B	2	NAG	O3-C3-C2	2.10	113.76	109.40
2	C	1	NAG	O3-C3-C4	-2.10	105.43	110.38

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	NAG	C1

All (9) torsion outliers are listed below:

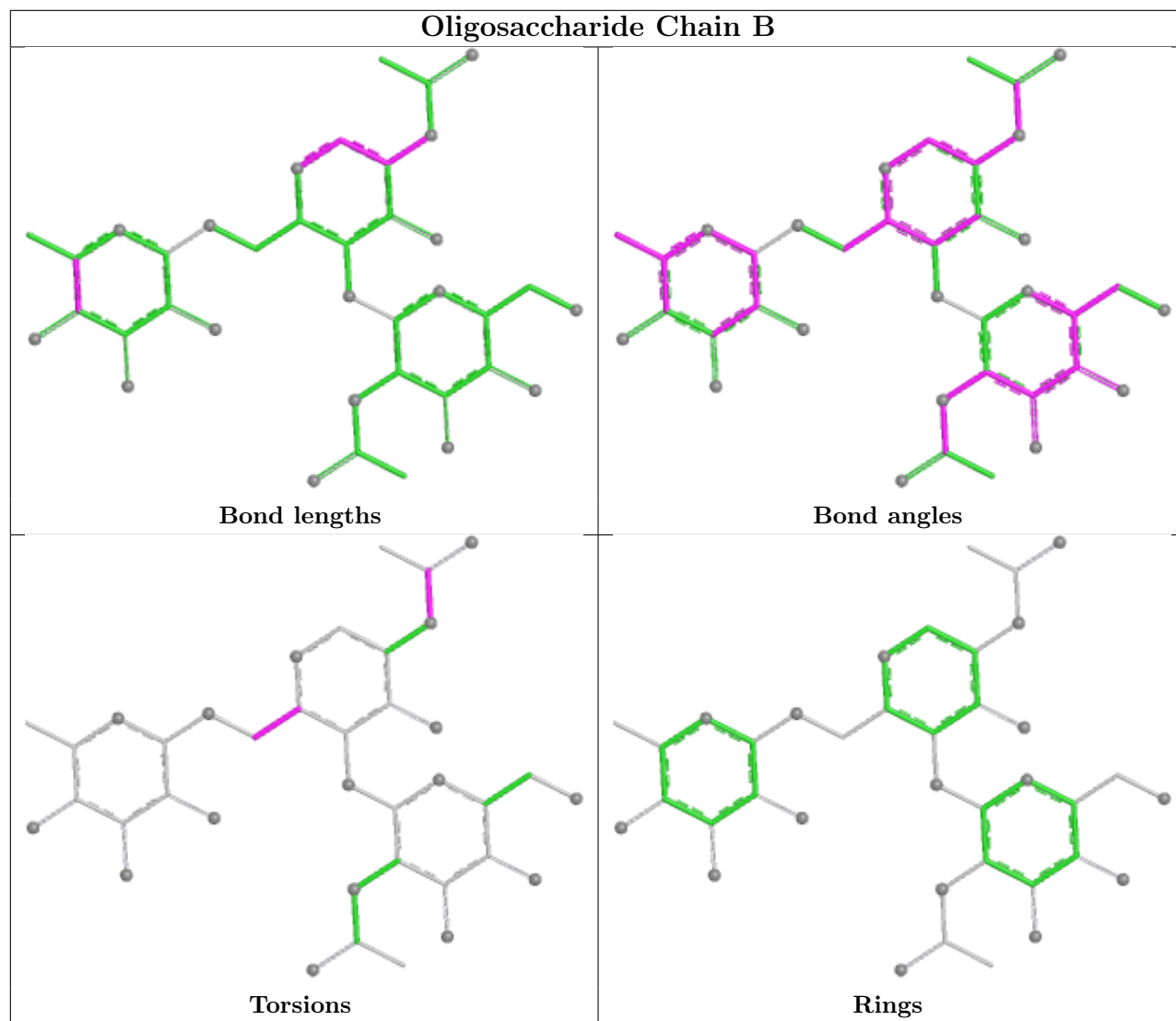
Mol	Chain	Res	Type	Atoms
2	B	1	NAG	C8-C7-N2-C2
2	B	1	NAG	O7-C7-N2-C2
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	B	1	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	C	2	NAG	C1-C2-N2-C7
2	C	2	NAG	C3-C2-N2-C7

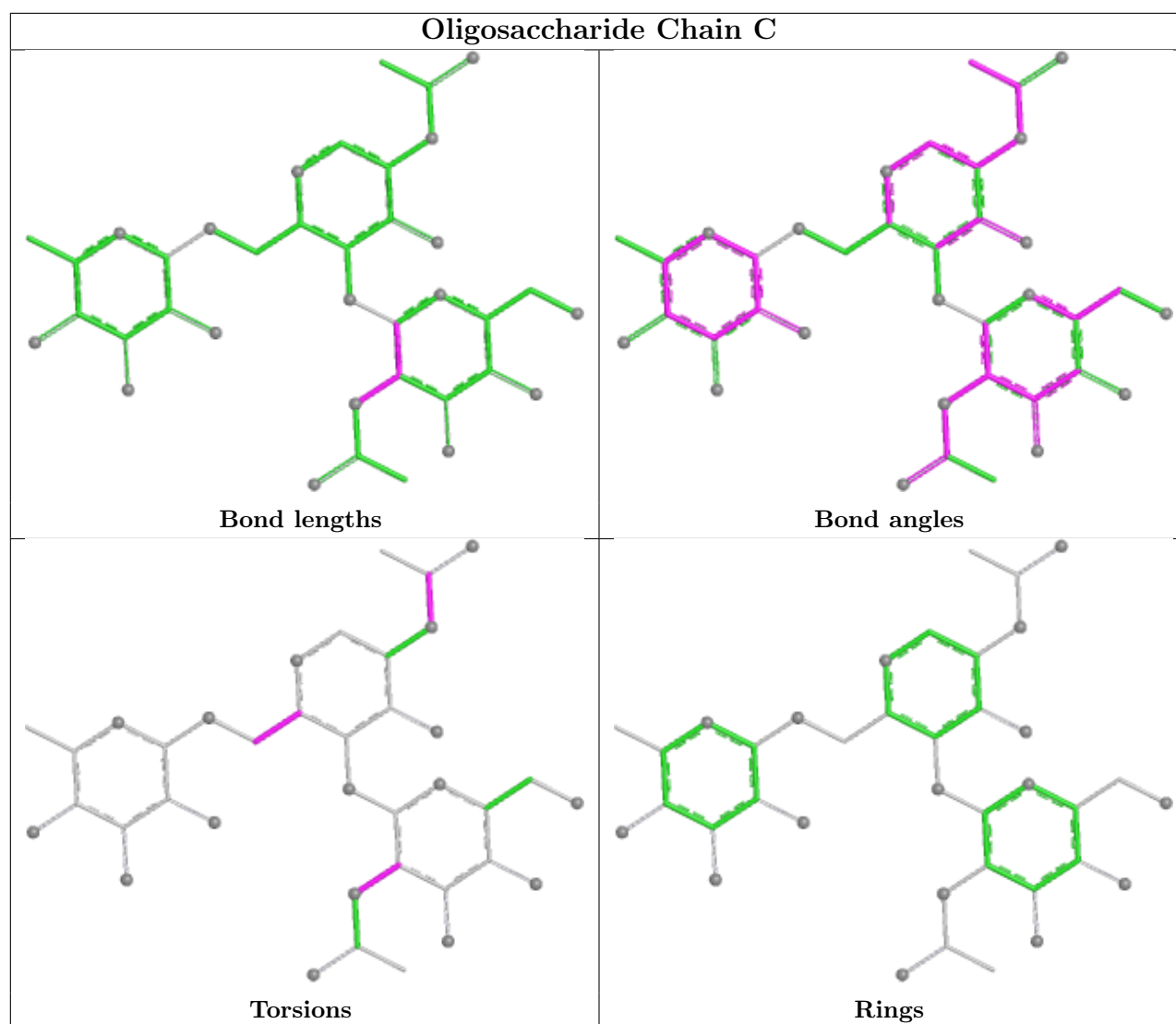
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	1	0
2	C	3	FUL	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 48 ligands modelled in this entry, 36 are unknown and 4 are monoatomic - leaving 8 for Mogul analysis.

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	$\# Z > 2$
9	SO4	A	1537	-	4,4,4	0.24	0	6,6,6	0.56	0
10	NAG	A	1542	1	14,14,15	0.92	0	17,19,21	3.05	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	NAG	A	1543	1	14,14,15	1.21	2 (14%)	17,19,21	2.20	6 (35%)
10	NAG	A	1541	1	14,14,15	1.53	2 (14%)	17,19,21	2.29	5 (29%)
5	VX	A	1530	1	2,5,6	1.40	0	1,5,8	0.11	0
3	GLY	A	1531	-	4,4,4	0.91	0	3,4,4	1.13	0
9	SO4	A	1536	-	4,4,4	0.80	0	6,6,6	0.51	0
10	NAG	A	1544	1	14,14,15	0.87	1 (7%)	17,19,21	1.81	4 (23%)

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
10	NAG	A	1542	1	-	4/6/23/26	0/1/1/1
10	NAG	A	1543	1	-	2/6/23/26	0/1/1/1
10	NAG	A	1541	1	-	2/6/23/26	0/1/1/1
5	VX	A	1530	1	-	0/1/3/4	-
3	GLY	A	1531	-	-	0/2/2/2	-
10	NAG	A	1544	1	-	2/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	A	1541	NAG	C1-C2	3.54	1.57	1.52
10	A	1541	NAG	C2-N2	2.54	1.50	1.46
10	A	1543	NAG	C1-C2	2.52	1.55	1.52
10	A	1544	NAG	C1-C2	2.28	1.55	1.52
10	A	1543	NAG	C3-C2	2.11	1.56	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	A	1542	NAG	C1-O5-C5	10.43	126.16	112.19
10	A	1541	NAG	C8-C7-N2	4.96	124.34	116.12
10	A	1543	NAG	C1-O5-C5	4.81	118.64	112.19
10	A	1541	NAG	O7-C7-C8	-4.25	114.49	122.05
10	A	1543	NAG	O5-C5-C6	3.98	115.42	107.66
10	A	1541	NAG	C1-C2-N2	3.96	116.67	110.43
10	A	1544	NAG	C2-N2-C7	3.73	127.91	122.90
10	A	1544	NAG	C1-O5-C5	3.65	117.08	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	A	1542	NAG	C8-C7-N2	3.50	121.93	116.12
10	A	1541	NAG	C2-N2-C7	3.49	127.58	122.90
10	A	1542	NAG	C1-C2-N2	3.20	115.47	110.43
10	A	1543	NAG	O7-C7-N2	3.03	127.33	121.98
10	A	1541	NAG	C6-C5-C4	2.72	119.70	113.02
10	A	1543	NAG	C4-C3-C2	2.51	114.70	111.02
10	A	1542	NAG	O7-C7-C8	-2.42	117.75	122.05
10	A	1544	NAG	C1-C2-N2	-2.35	106.73	110.43
10	A	1542	NAG	O4-C4-C5	2.20	114.74	109.32
10	A	1543	NAG	O5-C5-C4	-2.20	105.48	110.83
10	A	1543	NAG	C8-C7-N2	-2.15	112.56	116.12
10	A	1544	NAG	C4-C3-C2	2.07	114.05	111.02

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	1543	NAG	O5-C5-C6-O6
10	A	1542	NAG	C4-C5-C6-O6
10	A	1543	NAG	C4-C5-C6-O6
10	A	1542	NAG	O5-C5-C6-O6
10	A	1541	NAG	C8-C7-N2-C2
10	A	1541	NAG	O7-C7-N2-C2
10	A	1542	NAG	C8-C7-N2-C2
10	A	1542	NAG	O7-C7-N2-C2
10	A	1544	NAG	C4-C5-C6-O6
10	A	1544	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	1542	NAG	1	0
5	A	1530	VX	2	0
3	A	1531	GLY	1	0

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	527/527 (100%)	0.37	26 (4%) 35 37	15, 32, 58, 92	25 (4%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	377	VAL	5.7
1	A	258	THR	4.1
1	A	237	TYR	3.5
1	A	379	ASP	3.1
1	A	54	ASP	3.1
1	A	282	TYR	3.0
1	A	40	ARG	2.8
1	A	255	GLU	2.8
1	A	378	ASP	2.6
1	A	49	LEU	2.6
1	A	51	LYS	2.6
1	A	240	ARG	2.6
1	A	236	LEU	2.5
1	A	453	ARG	2.5
1	A	50	THR	2.5
1	A	4	ILE	2.4
1	A	455[A]	GLN	2.4
1	A	244	LEU	2.3
1	A	55	ILE	2.3
1	A	371	PHE	2.2
1	A	452	ARG	2.2
1	A	39	GLY	2.2
1	A	6	ILE	2.1
1	A	248	LYS	2.1
1	A	525	PHE	2.1
1	A	156	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

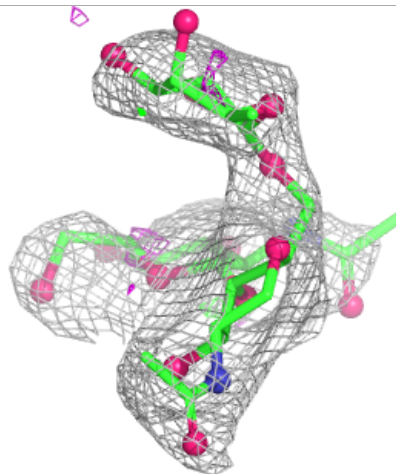
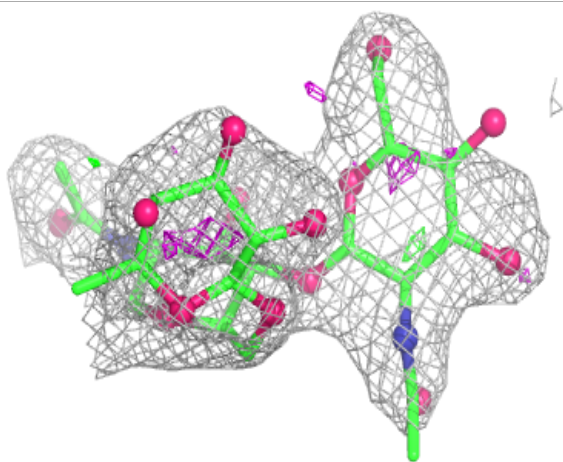
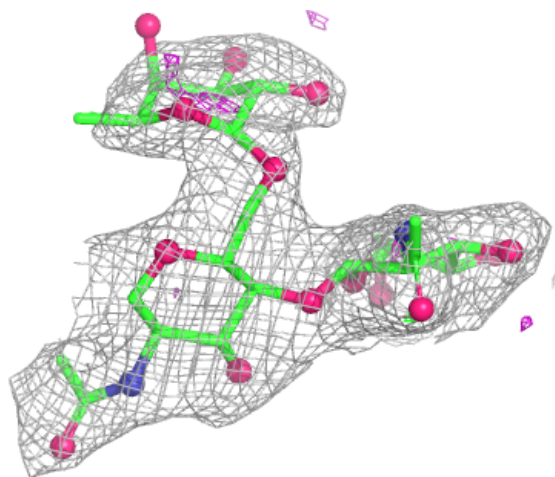
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

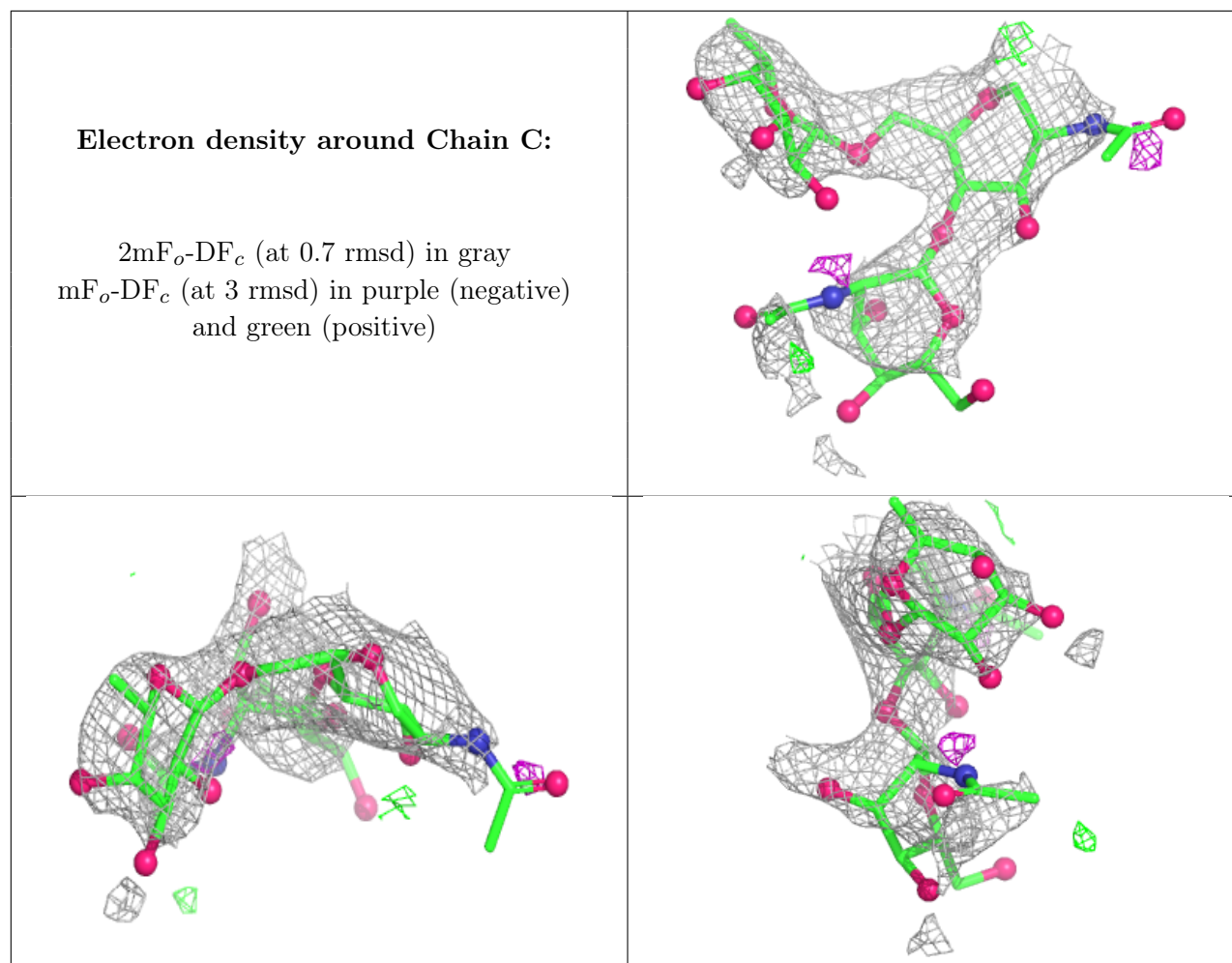
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	C	2	14/15	0.46	0.20	91,95,99,100	0
2	FUL	B	3	10/11	0.70	0.18	67,72,74,74	0
2	NAG	C	1	14/15	0.74	0.15	65,76,85,86	0
2	NAG	B	2	14/15	0.77	0.15	63,71,75,78	0
2	FUL	C	3	10/11	0.82	0.15	73,76,77,80	0
2	NAG	B	1	14/15	0.89	0.10	36,49,55,64	0

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
10	NAG	A	1542	14/15	0.53	0.17	59,69,73,75	0
4	UNX	A	1552	1/1	0.54	0.34	37,37,37,37	0
4	UNX	A	1560	1/1	0.59	0.18	56,56,56,56	0
10	NAG	A	1543	14/15	0.65	0.17	48,63,72,73	0
10	NAG	A	1544	14/15	0.65	0.15	78,87,93,94	0
4	UNX	A	1556	1/1	0.70	0.21	59,59,59,59	0
4	UNX	A	1557	1/1	0.72	0.20	61,61,61,61	0
4	UNX	A	1559	1/1	0.72	0.21	27,27,27,27	0
4	UNX	A	1577	1/1	0.74	0.15	53,53,53,53	0
4	UNX	A	1579	1/1	0.75	0.18	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	UNX	A	1578	1/1	0.79	0.10	46,46,46,46	0
4	UNX	A	1569	1/1	0.79	0.18	35,35,35,35	0
10	NAG	A	1541	14/15	0.79	0.13	57,65,69,71	0
4	UNX	A	1553	1/1	0.80	0.23	47,47,47,47	0
4	UNX	A	1573	1/1	0.81	0.39	60,60,60,60	0
4	UNX	A	1567	1/1	0.82	0.25	33,33,33,33	0
4	UNX	A	1565	1/1	0.84	0.22	30,30,30,30	0
4	UNX	A	1551	1/1	0.86	0.23	47,47,47,47	0
4	UNX	A	1566	1/1	0.86	0.14	40,40,40,40	0
4	UNX	A	1586	1/1	0.86	0.18	41,41,41,41	0
8	K	A	1535	1/1	0.86	0.21	89,89,89,89	1
4	UNX	A	1571	1/1	0.87	0.12	39,39,39,39	0
4	UNX	A	1558	1/1	0.87	0.10	47,47,47,47	0
4	UNX	A	1576	1/1	0.87	0.17	47,47,47,47	0
4	UNX	A	1561	1/1	0.87	0.17	31,31,31,31	0
5	VX	A	1530	6/7	0.88	0.21	24,30,42,43	0
3	GLY	A	1531	5/5	0.88	0.10	32,38,43,47	0
4	UNX	A	1580	1/1	0.89	0.21	26,26,26,26	0
4	UNX	A	1582	1/1	0.89	0.24	39,39,39,39	0
4	UNX	A	1568	1/1	0.89	0.15	28,28,28,28	0
4	UNX	A	1572	1/1	0.90	0.14	36,36,36,36	0
9	SO4	A	1537	5/5	0.90	0.18	67,70,73,73	5
4	UNX	A	1585	1/1	0.90	0.17	43,43,43,43	0
4	UNX	A	1564	1/1	0.91	0.35	54,54,54,54	0
6	NA	A	1532	1/1	0.91	0.17	57,57,57,57	1
4	UNX	A	1554	1/1	0.91	0.15	41,41,41,41	0
4	UNX	A	1581	1/1	0.91	0.17	35,35,35,35	0
7	CL	A	1533	1/1	0.92	0.22	72,72,72,72	0
4	UNX	A	1562	1/1	0.92	0.12	40,40,40,40	0
9	SO4	A	1536	5/5	0.92	0.13	25,29,34,37	5
4	UNX	A	1583	1/1	0.92	0.22	45,45,45,45	0
4	UNX	A	1574	1/1	0.93	0.17	53,53,53,53	0
4	UNX	A	1584	1/1	0.93	0.14	31,31,31,31	0
7	CL	A	1534	1/1	0.94	0.10	69,69,69,69	0
4	UNX	A	1563	1/1	0.94	0.36	44,44,44,44	0
4	UNX	A	1555	1/1	0.96	0.08	37,37,37,37	0
4	UNX	A	1570	1/1	0.96	0.09	37,37,37,37	0
4	UNX	A	1575	1/1	0.97	0.18	46,46,46,46	0

6.5 Other polymers ⓘ

There are no such residues in this entry.