



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 05:21 AM BST

PDB ID : 5MRH
Title : Crystal structure of the Vps10p domain of human sortilin/NTS3 in complex with Triazolone 1
Authors : Andersen, J.L.; Strandbygaard, D.; Thirup, S.
Deposited on : 2016-12-23
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

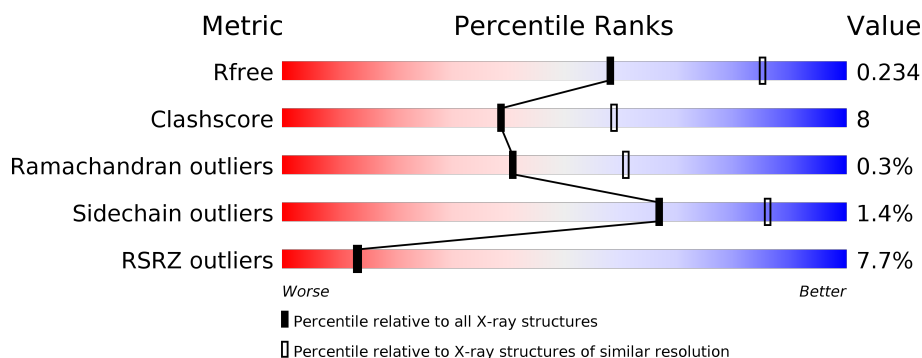
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on 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	696	<div> <div>7%</div> <div> <div></div> <div>76%</div> <div>15%</div> <div>8%</div> </div> </div>
2	B	2	<div> <div>100%</div> </div>
2	C	2	<div> <div>50%</div> <div>50%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5157 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 Sortilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	0	0
			5023	3174	845	976	28			

There are 18 discrepancies between the modelled and reference sequences:

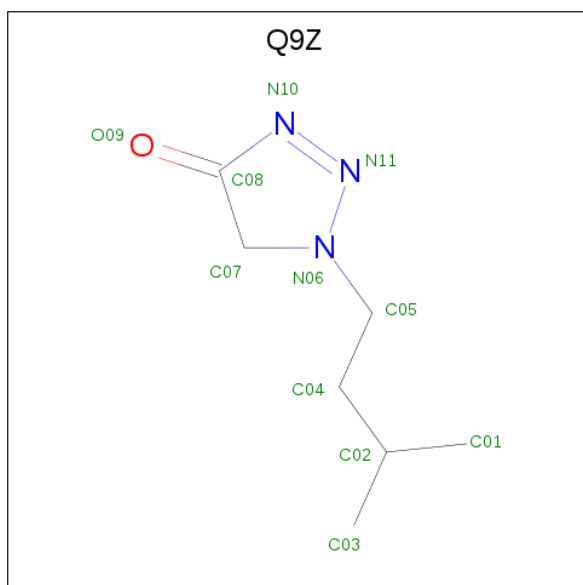
Chain	Residue	Modelled	Actual	Comment	Reference
A	617	MET	VAL	conflict	UNP Q99523
A	724	GLY	-	expression tag	UNP Q99523
A	725	SER	-	expression tag	UNP Q99523
A	726	ALA	-	expression tag	UNP Q99523
A	727	MET	-	expression tag	UNP Q99523
A	728	ILE	-	expression tag	UNP Q99523
A	729	GLU	-	expression tag	UNP Q99523
A	730	GLY	-	expression tag	UNP Q99523
A	731	ARG	-	expression tag	UNP Q99523
A	732	GLY	-	expression tag	UNP Q99523
A	733	VAL	-	expression tag	UNP Q99523
A	734	GLY	-	expression tag	UNP Q99523
A	735	HIS	-	expression tag	UNP Q99523
A	736	HIS	-	expression tag	UNP Q99523
A	737	HIS	-	expression tag	UNP Q99523
A	738	HIS	-	expression tag	UNP Q99523
A	739	HIS	-	expression tag	UNP Q99523
A	740	HIS	-	expression tag	UNP Q99523

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



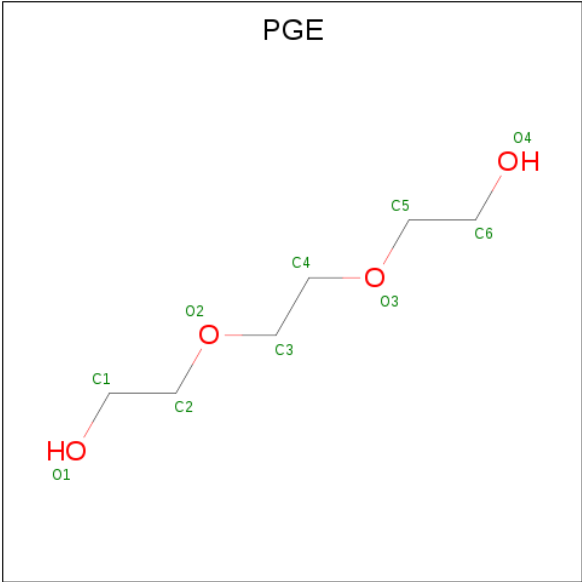
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 3-(3-methylbutyl)-4 {H}-1,2,3-triazol-5-one (three-letter code: Q9Z) (formula: C₇H₁₃N₃O).



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

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		

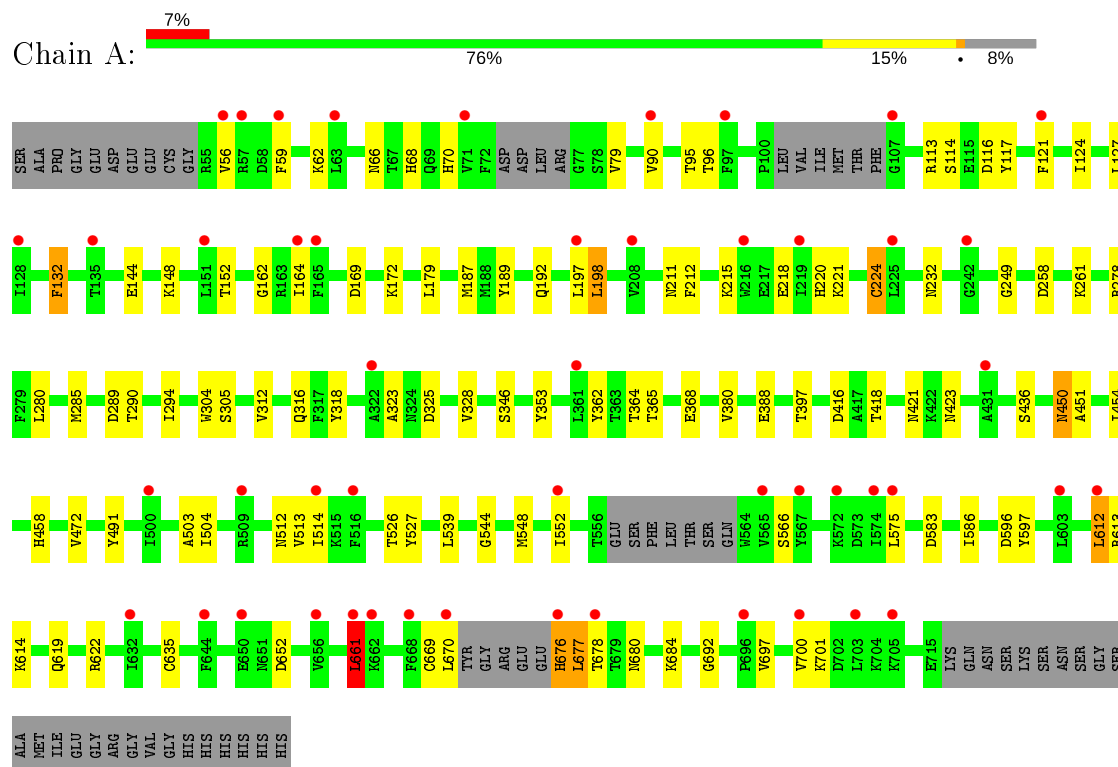
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	57	Total	O	0	0
			57	57		

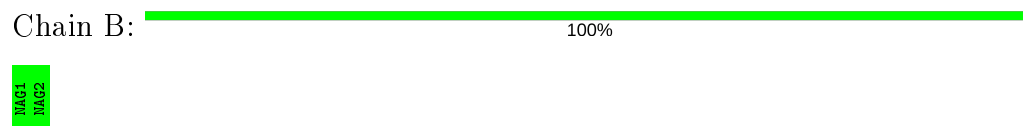
3 Residue-property plots [i](#)

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: Sortilin



- 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	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.40Å 78.23Å 111.80Å 90.00° 127.20° 90.00°	Depositor
Resolution (Å)	40.13 – 2.50 66.94 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.1 (40.13-2.50) 98.1 (66.94-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.51Å)	Xtriage
Refinement program	PHENIX (dev_2614: ???)	Depositor
R, R_{free}	0.211 , 0.233 0.211 , 0.234	Depositor DCC
R_{free} test set	1904 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	71.0	Xtriage
Anisotropy	0.524	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 76.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5157	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% 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: PGE, NAG, Q9Z

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	0.30	0/5140	0.57	4/6960 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	224	CYS	CA-CB-SG	6.14	125.05	114.00
1	A	612	LEU	CA-CB-CG	5.97	129.03	115.30
1	A	661	LEU	CA-CB-CG	5.79	128.62	115.30
1	A	198	LEU	CA-CB-CG	5.35	127.61	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5023	0	4816	75	0
2	B	28	0	25	0	0
2	C	28	0	25	3	0
3	A	11	0	0	0	0
4	A	10	0	14	0	0
5	A	57	0	0	1	0
All	All	5157	0	4880	75	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 8.

All (75) 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:364:THR:HG22	1:A:365:THR:H	1.52	0.74
1:A:278:ARG:NH1	1:A:325:ASP:O	2.27	0.67
1:A:62:LYS:O	1:A:66:ASN:ND2	2.28	0.67
1:A:152:THR:HA	1:A:164:ILE:HG22	1.75	0.67
1:A:504:ILE:HD11	1:A:514:ILE:HG13	1.78	0.65
1:A:124:ILE:HB	1:A:127:LEU:HD13	1.80	0.63
1:A:504:ILE:HG13	1:A:539:LEU:HD11	1.82	0.62
1:A:96:THR:HB	1:A:132:PHE:HE1	1.65	0.62
1:A:450:ASN:N	1:A:450:ASN:OD1	2.21	0.61
1:A:258:ASP:HB3	1:A:261:LYS:HB2	1.82	0.61
1:A:514:ILE:HD12	1:A:552:ILE:HD11	1.83	0.61
1:A:661:LEU:H	1:A:661:LEU:HD22	1.67	0.59
1:A:450:ASN:ND2	1:A:544:GLY:O	2.30	0.59
1:A:512:ASN:OD1	1:A:513:VAL:HG23	2.03	0.59
1:A:305:SER:HB3	1:A:676:HIS:CE1	2.38	0.59
1:A:312:VAL:HB	1:A:316:GLN:HB2	1.84	0.58
1:A:56:VAL:HG13	1:A:59:PHE:HB3	1.85	0.58
1:A:548:MET:H	2:C:1:NAG:H82	1.68	0.58
1:A:548:MET:N	2:C:1:NAG:H82	2.19	0.57
1:A:323:ALA:HB2	1:A:328:VAL:HG23	1.86	0.57
1:A:416:ASP:OD1	1:A:418:THR:OG1	2.22	0.56
1:A:364:THR:HG22	1:A:365:THR:N	2.20	0.55
1:A:249:GLY:HA3	1:A:285:MET:SD	2.47	0.55
1:A:612:LEU:HD23	1:A:613:ARG:O	2.08	0.53
1:A:680:ASN:HB3	1:A:697:VAL:HG13	1.91	0.53
1:A:187:MET:HE3	1:A:197:LEU:HD13	1.89	0.52
1:A:220:HIS:ND1	1:A:221:LYS:O	2.41	0.52
1:A:451:ALA:HB1	1:A:454:ILE:HD12	1.92	0.51
1:A:548:MET:HE3	1:A:575:LEU:HD12	1.92	0.51
1:A:678:THR:HA	1:A:701:LYS:O	2.10	0.51
1:A:614:LYS:NZ	5:A:912:HOH:O	2.45	0.50
1:A:346:SER:HB3	1:A:353:TYR:CE2	2.47	0.49
1:A:586:ILE:O	1:A:586:ILE:HD12	2.13	0.49
1:A:66:ASN:O	1:A:68:HIS:ND1	2.47	0.48
1:A:421:ASN:OD1	1:A:423:ASN:N	2.42	0.47
1:A:90:VAL:HA	1:A:114:SER:O	2.14	0.47
1:A:548:MET:SD	1:A:619:GLN:NE2	2.88	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:MET:CE	1:A:575:LEU:HD12	2.44	0.47
1:A:388:GLU:OE2	1:A:388:GLU:N	2.40	0.47
1:A:669:CYS:HB3	1:A:670:LEU:HD12	1.97	0.47
1:A:622:ARG:HD2	2:C:1:NAG:O7	2.15	0.47
1:A:116:ASP:O	1:A:117:TYR:HB2	2.15	0.46
1:A:169:ASP:O	1:A:172:LYS:HG3	2.16	0.46
1:A:189:TYR:CD1	1:A:197:LEU:HD21	2.50	0.46
1:A:491:TYR:O	1:A:503:ALA:HA	2.15	0.46
1:A:294:ILE:HD11	1:A:318:TYR:CG	2.51	0.46
1:A:79:VAL:HG13	1:A:95:THR:HG22	1.98	0.46
1:A:70:HIS:CD2	1:A:121:PHE:H	2.33	0.46
1:A:305:SER:HB3	1:A:676:HIS:HE1	1.79	0.45
1:A:148:LYS:HD3	1:A:212:PHE:CD1	2.52	0.45
1:A:513:VAL:HG11	1:A:526:THR:CG2	2.46	0.45
1:A:289:ASP:OD1	1:A:290:THR:N	2.48	0.45
1:A:380:VAL:HA	1:A:397:THR:O	2.17	0.44
1:A:677:LEU:H	1:A:677:LEU:HG	1.60	0.44
1:A:458:HIS:CD2	1:A:472:VAL:HG22	2.52	0.44
1:A:678:THR:HB	1:A:700:VAL:HB	1.98	0.44
1:A:635:CYS:SG	1:A:684:LYS:NZ	2.91	0.44
1:A:364:THR:CG2	1:A:365:THR:H	2.25	0.44
1:A:218:GLU:OE2	1:A:221:LYS:HG2	2.18	0.44
1:A:513:VAL:HG13	1:A:527:TYR:O	2.17	0.44
1:A:676:HIS:CD2	1:A:676:HIS:N	2.86	0.44
1:A:635:CYS:HB2	1:A:692:GLY:O	2.18	0.43
1:A:583:ASP:O	1:A:613:ARG:HG3	2.18	0.43
1:A:113:ARG:HB2	1:A:124:ILE:HD13	2.01	0.43
1:A:596:ASP:OD1	1:A:597:TYR:N	2.51	0.43
1:A:368:GLU:HG2	1:A:436:SER:HB3	2.01	0.42
1:A:189:TYR:HA	1:A:197:LEU:HD23	2.00	0.42
1:A:491:TYR:HB2	1:A:504:ILE:HB	2.01	0.42
1:A:211:ASN:HD21	1:A:215:LYS:N	2.17	0.41
1:A:212:PHE:O	1:A:212:PHE:CG	2.73	0.41
1:A:552:ILE:O	1:A:566:SER:HA	2.20	0.41
1:A:504:ILE:CD1	1:A:514:ILE:HG13	2.47	0.41
1:A:280:LEU:HD22	1:A:304:TRP:CH2	2.55	0.41
1:A:162:GLY:O	1:A:179:LEU:N	2.52	0.40
1:A:192:GLN:HB3	1:A:232:ASN:OD1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	629/696 (90%)	614 (98%)	13 (2%)	2 (0%)	41	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	144	GLU
1	A	652	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	557/605 (92%)	549 (99%)	8 (1%)	67	86

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	132	PHE
1	A	198	LEU
1	A	224	CYS
1	A	362	TYR
1	A	450	ASN
1	A	661	LEU
1	A	676	HIS
1	A	677	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	676	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	1,2	14,14,15	0.52	0	17,19,21	0.51	0
2	NAG	B	2	2	14,14,15	0.36	0	17,19,21	0.51	0
2	NAG	C	1	1,2	14,14,15	0.37	0	17,19,21	0.68	1 (5%)
2	NAG	C	2	2	14,14,15	0.31	0	17,19,21	0.51	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	C	1	NAG	C1-O5-C5	2.18	115.14	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

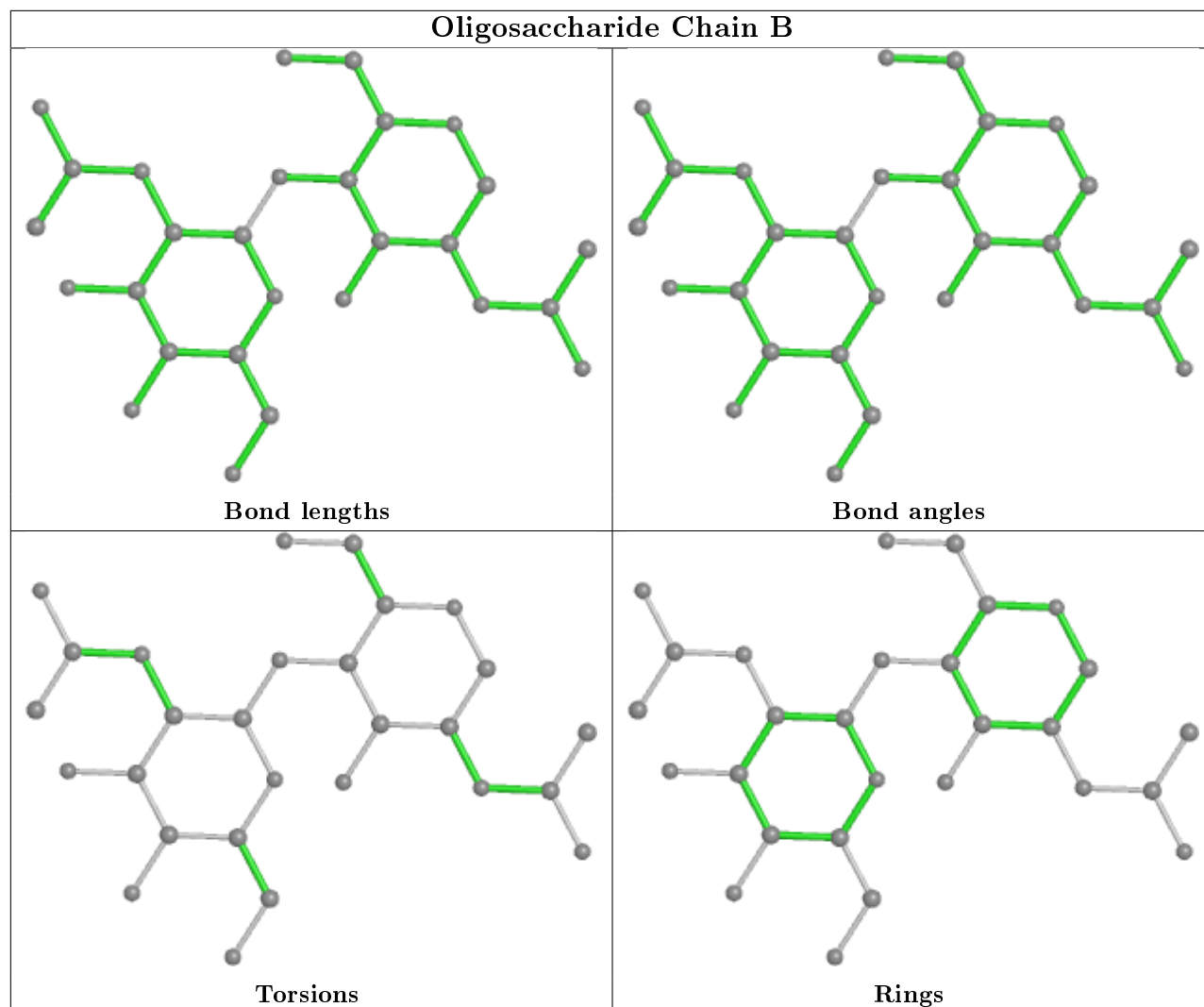
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6

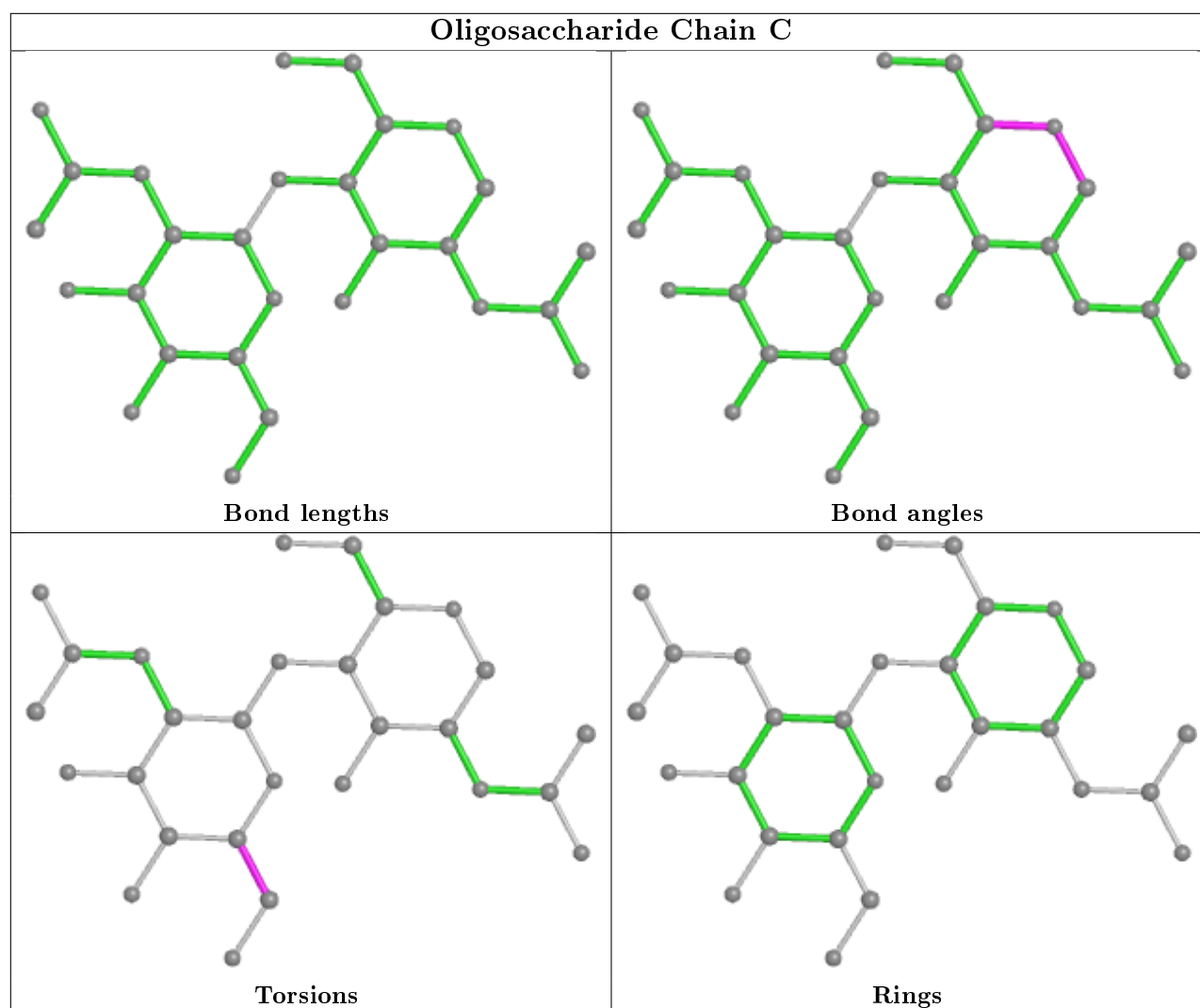
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	3	0

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





5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PGE	A	806	-	9,9,9	0.31	0	8,8,8	0.25	0
3	Q9Z	A	805	-	10,11,11	2.47	3 (30%)	9,14,14	2.60	3 (33%)

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	PGE	A	806	-	-	1/7/7/7	-
3	Q9Z	A	805	-	-	3/5/14/14	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	805	Q9Z	N10-N11	5.10	1.40	1.32
3	A	805	Q9Z	C07-N06	-4.07	1.41	1.45
3	A	805	Q9Z	C05-N06	-3.34	1.43	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	805	Q9Z	C05-N06-N11	-6.78	107.34	120.27
3	A	805	Q9Z	C07-C08-N10	-2.49	107.97	113.44
3	A	805	Q9Z	O09-C08-C07	2.26	125.72	122.06

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	805	Q9Z	C01-C02-C04-C05
3	A	805	Q9Z	C03-C02-C04-C05
3	A	805	Q9Z	C02-C04-C05-N06
4	A	806	PGE	C3-C4-O3-C5

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 [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	639/696 (91%)	0.85	49 (7%) 13 13	52, 100, 159, 207	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	97	PHE	5.1
1	A	678	THR	5.0
1	A	676	HIS	4.8
1	A	575	LEU	4.2
1	A	219	ILE	4.2
1	A	668	PHE	3.8
1	A	696	PRO	3.8
1	A	107	GLY	3.7
1	A	670	LEU	3.7
1	A	662	LYS	3.7
1	A	57	ARG	3.7
1	A	56	VAL	3.4
1	A	165	PHE	3.3
1	A	700	VAL	3.2
1	A	565	VAL	3.2
1	A	703	LEU	3.1
1	A	661	LEU	3.0
1	A	59	PHE	2.8
1	A	164	ILE	2.8
1	A	71	VAL	2.7
1	A	500	ILE	2.7
1	A	63	LEU	2.5
1	A	572	LYS	2.5
1	A	574	ILE	2.5
1	A	656	VAL	2.5
1	A	121	PHE	2.4
1	A	567	TYR	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	361	LEU	2.4
1	A	644	PHE	2.3
1	A	705	LYS	2.3
1	A	225	LEU	2.3
1	A	509	ARG	2.3
1	A	431	ALA	2.3
1	A	128	ILE	2.3
1	A	514	ILE	2.3
1	A	516	PHE	2.2
1	A	552	ILE	2.2
1	A	242	GLY	2.2
1	A	216	TRP	2.2
1	A	197	LEU	2.2
1	A	612	LEU	2.2
1	A	322	ALA	2.1
1	A	151	LEU	2.1
1	A	603	LEU	2.1
1	A	208	VAL	2.1
1	A	90	VAL	2.1
1	A	135	THR	2.1
1	A	632	ILE	2.0
1	A	650	GLU	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.

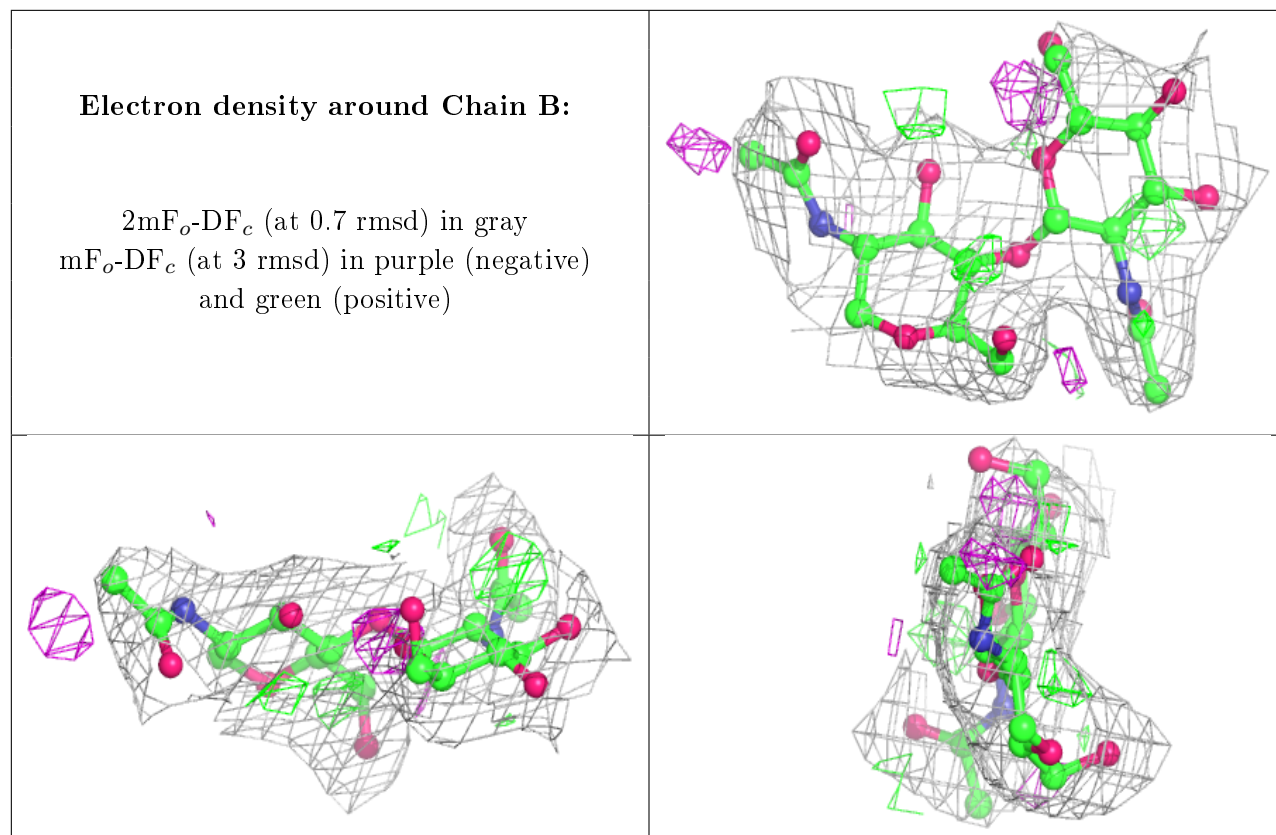
6.3 Carbohydrates [i](#)

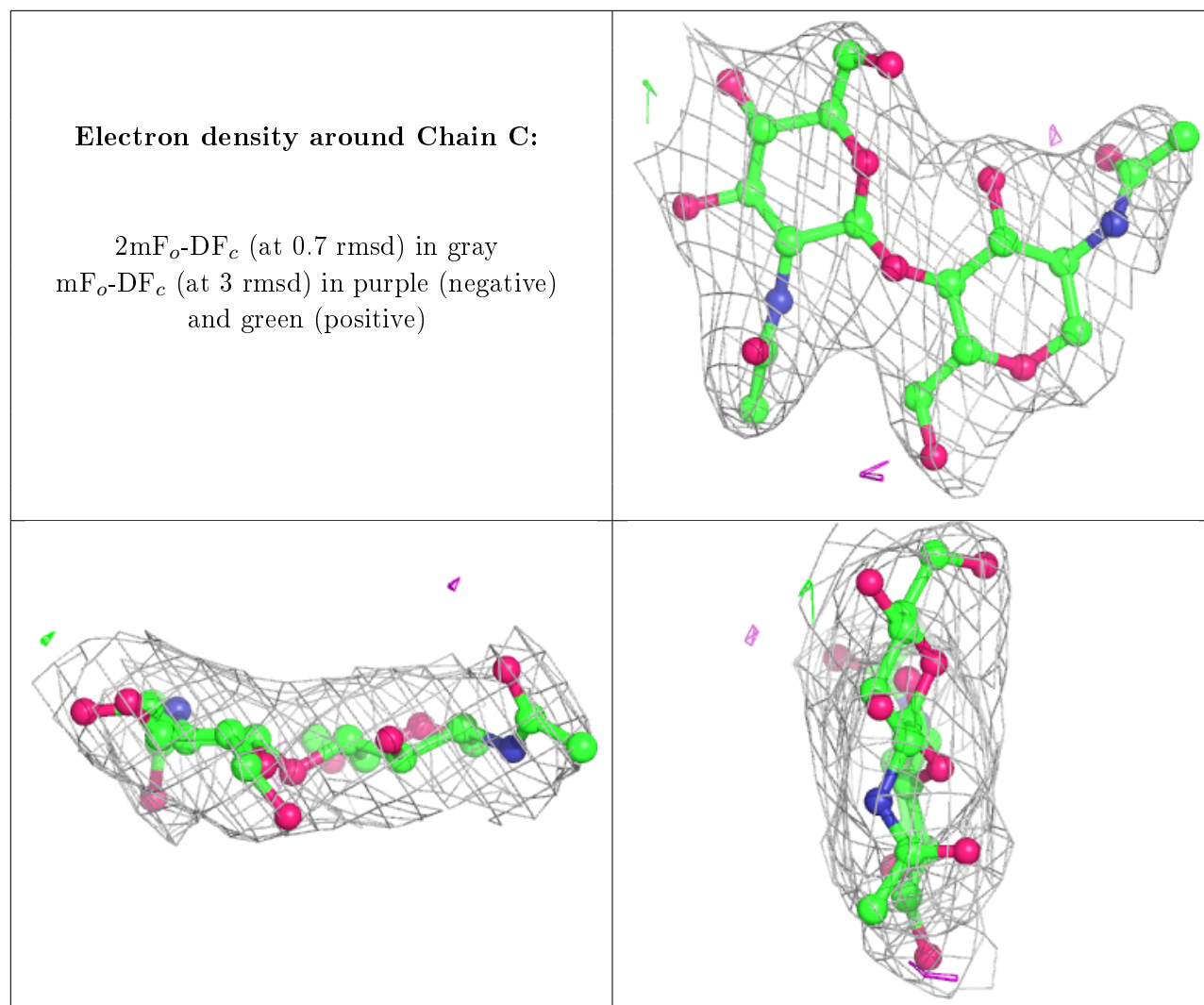
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	B	2	14/15	0.92	0.18	61,96,112,120	0
2	NAG	C	2	14/15	0.94	0.14	86,112,124,125	0
2	NAG	C	1	14/15	0.94	0.16	80,92,101,101	0
2	NAG	B	1	14/15	0.97	0.17	45,56,69,79	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. 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(\AA^2)	Q<0.9
4	PGE	A	806	10/10	0.88	0.20	87,103,118,122	0
3	Q9Z	A	805	11/11	0.97	0.23	71,74,80,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.