



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 12:35 PM BST

PDB ID : 1FSU
Title : Crystal Structure of 4-Sulfatase (human)
Authors : Bond, C.; Guss, M.
Deposited on : 1996-07-29
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

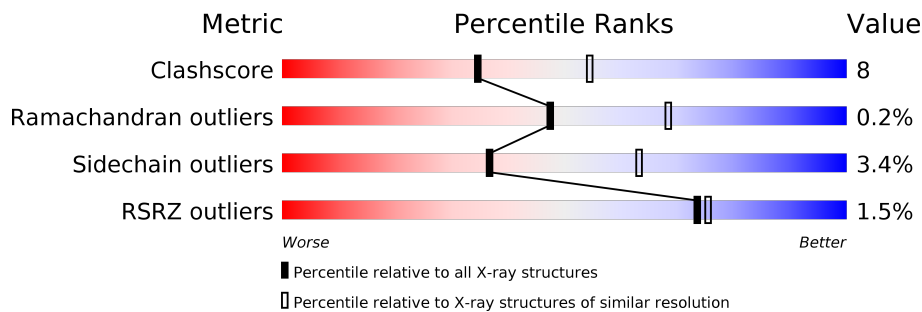
1 Overall quality at a glance

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))
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	492	 75% 20% ..
2	B	2	 100%
2	C	2	 100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4139 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 N-ACETYL GALACTOSAMINE-4-SULFATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	475	3811	2442	664	690	15	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	91	ALS	CYS	conflict	UNP P15848

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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0

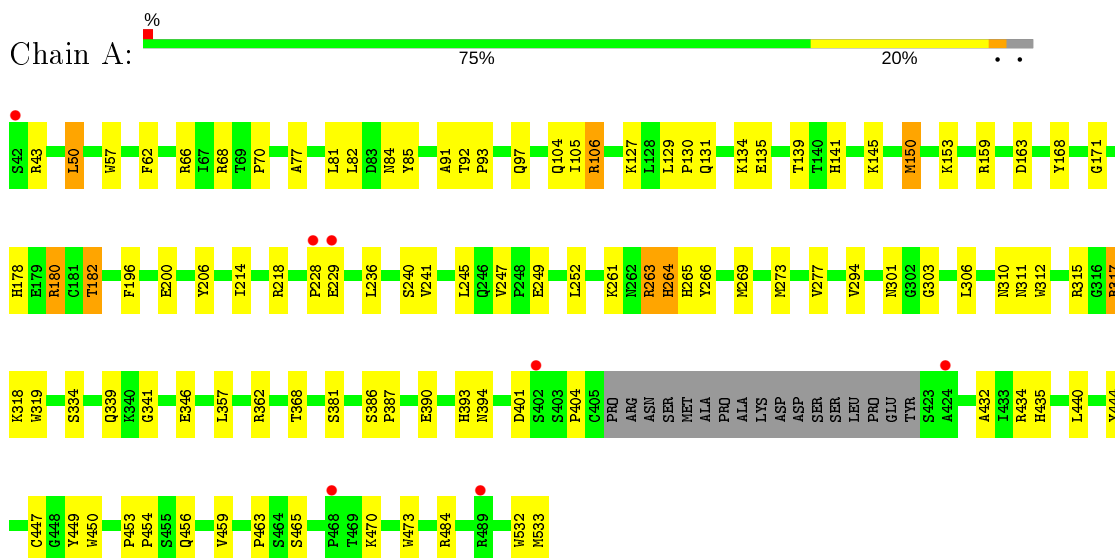
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	270	Total O 270 270	0	0

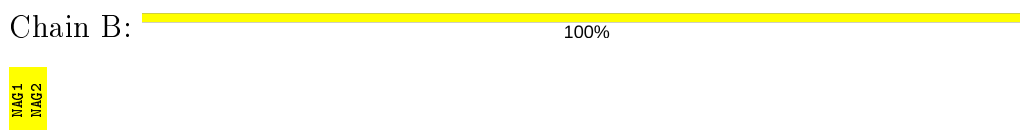
3 Residue-property plots

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.

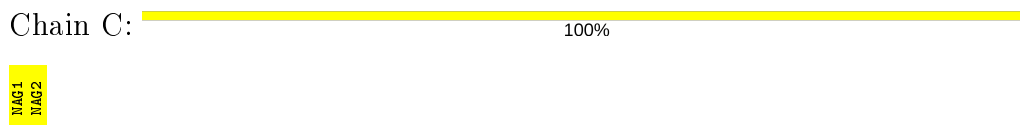
- Molecule 1: N-ACETYL GALACTOSAMINE-4-SULFATASE



- 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 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.00Å 107.00Å 144.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50 19.99 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.0 ((Not available)-2.50) 91.9 (19.99-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.30 (at 2.50Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.188 , 0.243 0.179 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtrriage
Anisotropy	0.112	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4139	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% 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: CA, CL, ALS, NAG

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	0.34	0/3922	0.64	0/5346

There are no bond length outliers.

There are no bond angle outliers.

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	3811	0	3693	63	0
2	B	28	0	25	0	0
2	C	28	0	25	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	270	0	0	5	0
All	All	4139	0	3743	63	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 (63) 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:245:LEU:HD22	1:A:269:MET:HE1	1.65	0.78
1:A:303:GLY:H	1:A:311:ASN:HD21	1.31	0.77
1:A:97:GLN:HE21	1:A:104:GLN:HE21	1.35	0.74
1:A:159:ARG:HD3	1:A:200:GLU:OE2	1.90	0.71
1:A:50:LEU:HD22	1:A:236:LEU:HD23	1.77	0.66
1:A:318:LYS:O	1:A:319:TRP:HB2	1.95	0.66
1:A:401:ASP:HB2	1:A:444:TYR:HE2	1.62	0.63
1:A:245:LEU:CD2	1:A:269:MET:HE1	2.29	0.61
1:A:178:HIS:HD2	5:A:625:HOH:O	1.85	0.59
1:A:266:TYR:HA	1:A:269:MET:HE3	1.85	0.58
1:A:228:PRO:O	1:A:229:GLU:HB2	2.04	0.58
1:A:261:LYS:HE3	1:A:265:HIS:HE1	1.70	0.57
1:A:134:LYS:HE2	1:A:163:ASP:OD1	2.05	0.56
1:A:266:TYR:HA	1:A:269:MET:CE	2.36	0.55
1:A:153:LYS:O	1:A:159:ARG:HG3	2.07	0.55
1:A:145:LYS:HE3	1:A:171:GLY:HA2	1.88	0.54
1:A:129:LEU:HD21	1:A:357:LEU:HD12	1.90	0.54
1:A:105:ILE:HG13	1:A:106:ARG:HD3	1.90	0.53
1:A:339:GLN:HG2	1:A:381:SER:O	2.09	0.52
1:A:294:VAL:HA	1:A:334:SER:HB2	1.94	0.50
1:A:180:ARG:NH1	1:A:182:THR:HG23	2.27	0.50
1:A:306:LEU:HD13	1:A:404:PRO:HD2	1.94	0.50
1:A:127:LYS:HB3	1:A:368:THR:HG22	1.94	0.49
1:A:131:GLN:O	1:A:135:GLU:HG3	2.13	0.49
1:A:303:GLY:N	1:A:311:ASN:HD21	2.06	0.49
1:A:263:ARG:HD2	1:A:454:PRO:HD2	1.95	0.49
1:A:245:LEU:HA	1:A:269:MET:HE1	1.95	0.48
1:A:43:ARG:NH2	1:A:229:GLU:HG2	2.28	0.48
1:A:92:THR:HB	1:A:93:PRO:HD3	1.95	0.48
1:A:57:TRP:H	1:A:311:ASN:HD22	1.61	0.48
1:A:532:TRP:CD1	1:A:533:MET:HG2	2.49	0.47
1:A:261:LYS:HE3	1:A:265:HIS:CE1	2.50	0.47
1:A:317:ARG:HD2	5:A:776:HOH:O	2.13	0.47
1:A:240:SER:HA	1:A:241:VAL:HA	1.75	0.46
1:A:390:GLU:HB2	1:A:435:HIS:HD2	1.81	0.46
1:A:84:ASN:HD21	1:A:434:ARG:HH22	1.64	0.45
1:A:62:PHE:CE1	1:A:81:LEU:HD11	2.51	0.45
1:A:393:HIS:O	1:A:394:ASN:HB2	2.16	0.45
1:A:465:SER:HB3	5:A:778:HOH:O	2.17	0.45
1:A:252:LEU:O	1:A:264:HIS:HE1	2.01	0.44
1:A:245:LEU:HD22	1:A:269:MET:CE	2.42	0.44
1:A:346:GLU:HB3	1:A:386:SER:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:CYS:SG	1:A:449:TYR:HD2	2.41	0.44
1:A:178:HIS:HE1	1:A:206:TYR:O	2.02	0.43
1:A:218:ARG:HD3	5:A:870:HOH:O	2.16	0.43
1:A:214:ILE:O	1:A:218:ARG:HG2	2.19	0.43
1:A:432:ALA:HA	1:A:440:LEU:O	2.19	0.43
1:A:449:TYR:CE1	1:A:463:PRO:HB3	2.54	0.42
1:A:484:ARG:NH1	5:A:875:HOH:O	2.44	0.42
1:A:68:ARG:C	1:A:70:PRO:HD3	2.40	0.42
1:A:130:PRO:O	1:A:134:LYS:HB2	2.20	0.42
1:A:470:LYS:HE3	1:A:473:TRP:NE1	2.35	0.42
1:A:315:ARG:HD2	1:A:450:TRP:CZ2	2.54	0.42
1:A:77:ALA:O	1:A:341:GLY:HA2	2.19	0.42
1:A:453:PRO:CG	1:A:456:GLN:HE21	2.32	0.42
1:A:310:ASN:HD21	1:A:312:TRP:HB2	1.86	0.41
1:A:139:THR:HG22	1:A:141:HIS:NE2	2.36	0.41
1:A:82:LEU:HB3	1:A:85:TYR:HB2	2.03	0.41
1:A:386:SER:HA	1:A:387:PRO:HD3	1.87	0.41
1:A:150:MET:SD	1:A:150:MET:C	2.99	0.41
1:A:273:MET:O	1:A:277:VAL:HG23	2.21	0.40
1:A:145:LYS:HB2	1:A:168:TYR:CZ	2.57	0.40
1:A:129:LEU:N	1:A:130:PRO:HD2	2.37	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	470/492 (96%)	450 (96%)	19 (4%)	1 (0%)	47 68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	GLU

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	413/428 (96%)	399 (97%)	14 (3%)	37 63

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LEU
1	A	66	ARG
1	A	106	ARG
1	A	150	MET
1	A	180	ARG
1	A	182	THR
1	A	196	PHE
1	A	247	VAL
1	A	263	ARG
1	A	264	HIS
1	A	301	ASN
1	A	317	ARG
1	A	362	ARG
1	A	459	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	ASN
1	A	88	GLN
1	A	104	GLN
1	A	178	HIS
1	A	264	HIS
1	A	265	HIS
1	A	310	ASN

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Mol	Chain	Res	Type
1	A	311	ASN
1	A	435	HIS
1	A	456	GLN
1	A	458	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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
1	ALS	A	91	1,3	7,10,11	1.19	1 (14%)	6,14,16	0.80	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
1	ALS	A	91	1,3	-	1/3/11/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	91	ALS	OS1-S	-2.70	1.49	1.57

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	91	ALS	N-CA-CB-OG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

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	1.04	1 (7%)	17,19,21	1.79	3 (17%)
2	NAG	B	2	2	14,14,15	1.06	1 (7%)	17,19,21	1.23	2 (11%)
2	NAG	C	1	1,2	14,14,15	1.30	2 (14%)	17,19,21	1.86	5 (29%)
2	NAG	C	2	2	14,14,15	1.12	2 (14%)	17,19,21	1.61	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	B	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
2	NAG	C	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	C	2	2	-	4/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	NAG	C1-C2	2.48	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	NAG	C4-C5	2.40	1.58	1.53
2	B	2	NAG	C1-C2	2.31	1.55	1.52
2	B	1	NAG	C3-C2	2.20	1.57	1.52
2	C	2	NAG	C3-C2	2.15	1.57	1.52
2	C	2	NAG	C4-C5	2.15	1.57	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	C1-C2-N2	5.24	119.43	110.49
2	C	2	NAG	C6-C5-C4	4.98	124.66	113.00
2	C	1	NAG	C1-C2-N2	3.76	116.91	110.49
2	B	1	NAG	O5-C1-C2	-3.55	105.69	111.29
2	B	2	NAG	C1-C2-N2	-3.42	104.65	110.49
2	C	1	NAG	O5-C1-C2	-3.38	105.94	111.29
2	C	1	NAG	C1-O5-C5	3.33	116.70	112.19
2	B	1	NAG	C2-N2-C7	3.07	127.28	122.90
2	C	1	NAG	C8-C7-N2	-2.96	111.08	116.10
2	B	2	NAG	C1-O5-C5	2.66	115.80	112.19
2	C	1	NAG	C4-C3-C2	-2.04	108.02	111.02

There are no chirality outliers.

All (12) torsion outliers are listed below:

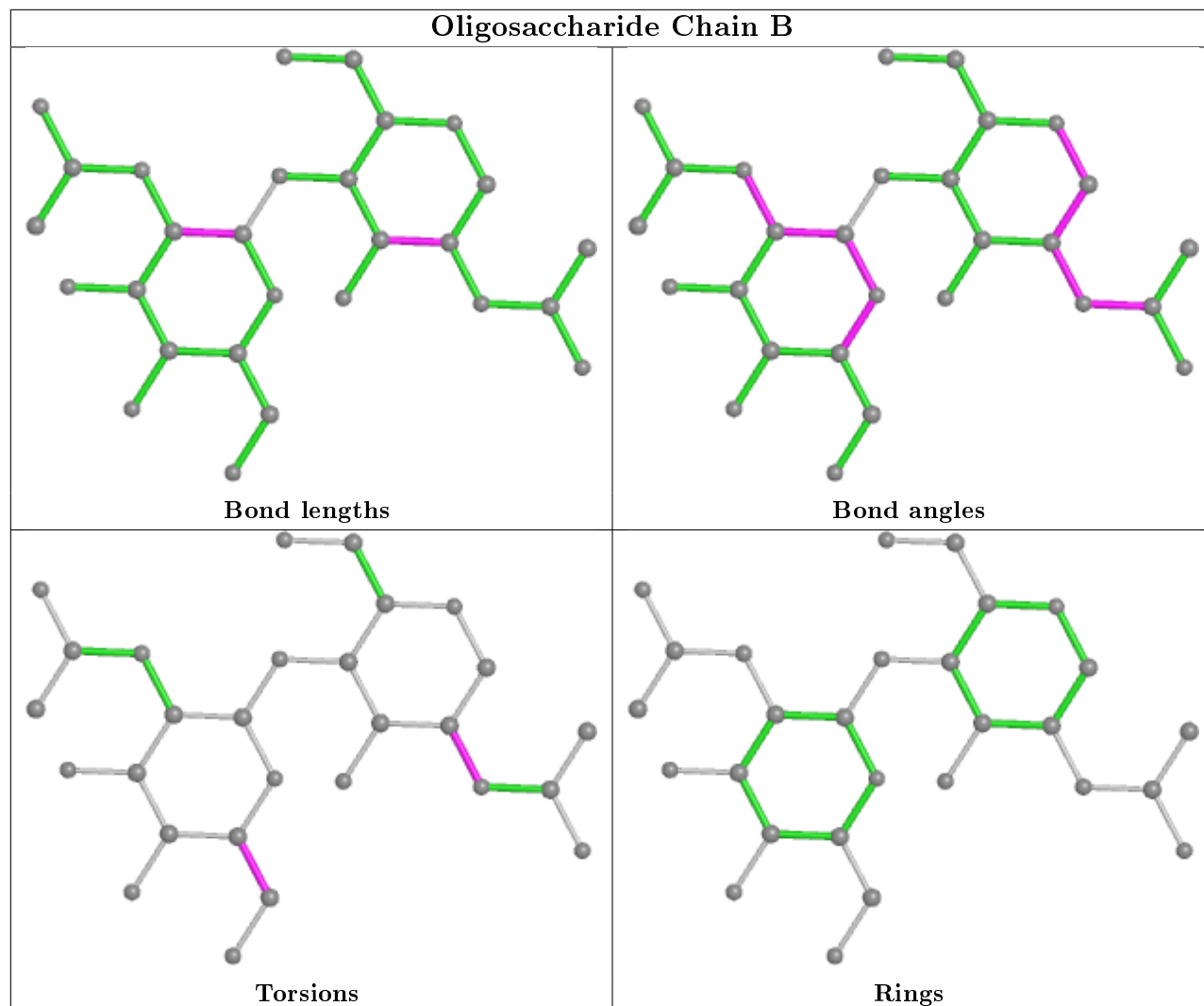
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	B	2	NAG	C4-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	B	2	NAG	O5-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C1-C2-N2-C7
2	B	1	NAG	C1-C2-N2-C7
2	C	2	NAG	C1-C2-N2-C7
2	C	1	NAG	C3-C2-N2-C7
2	C	2	NAG	C3-C2-N2-C7
2	B	1	NAG	C3-C2-N2-C7

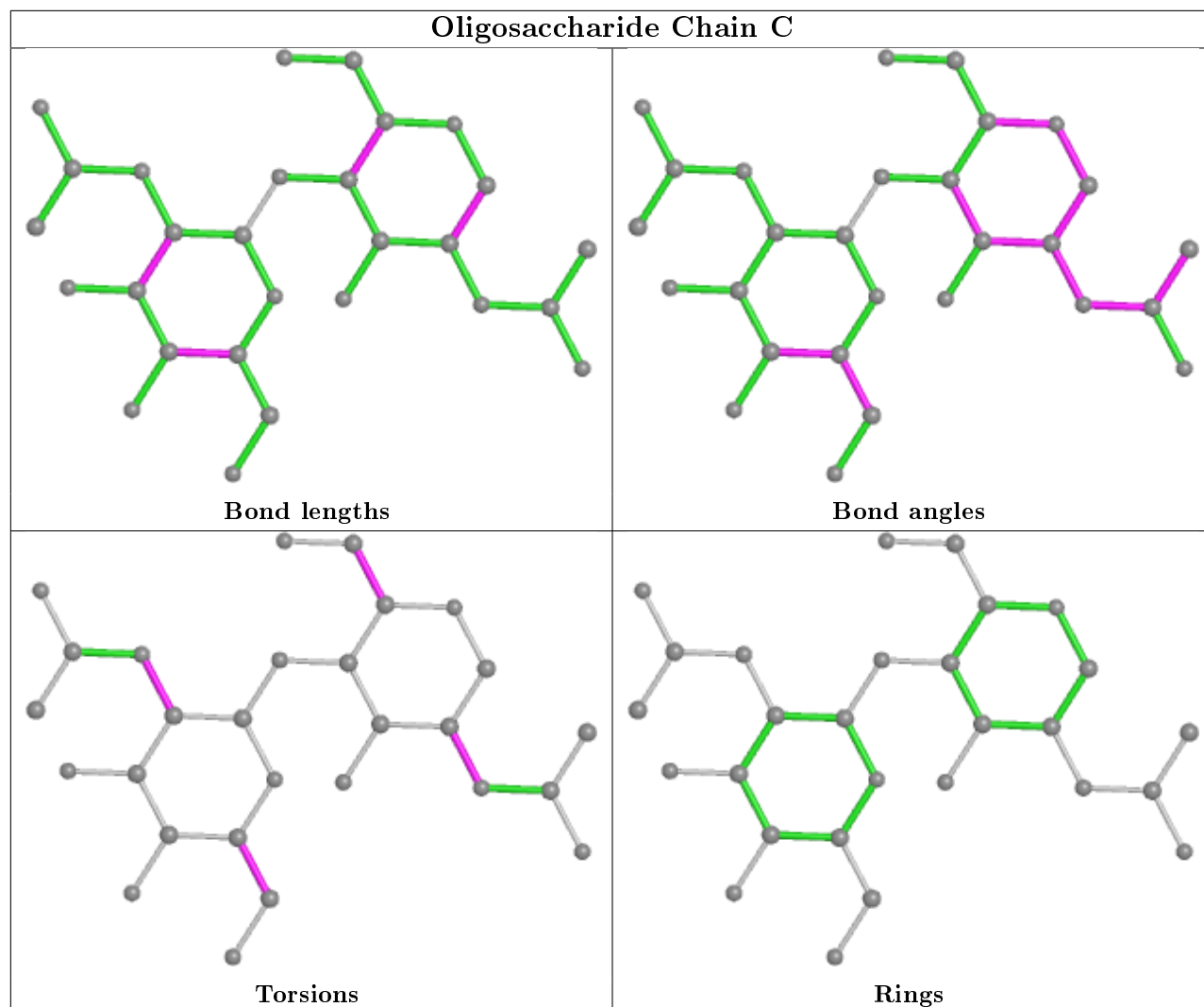
There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	474/492 (96%)	-0.72	7 (1%) 73 75	11, 20, 43, 98	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	229	GLU	4.4
1	A	402	SER	3.7
1	A	228	PRO	3.7
1	A	42	SER	3.4
1	A	489	ARG	2.3
1	A	424	ALA	2.1
1	A	468	PRO	2.1

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

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

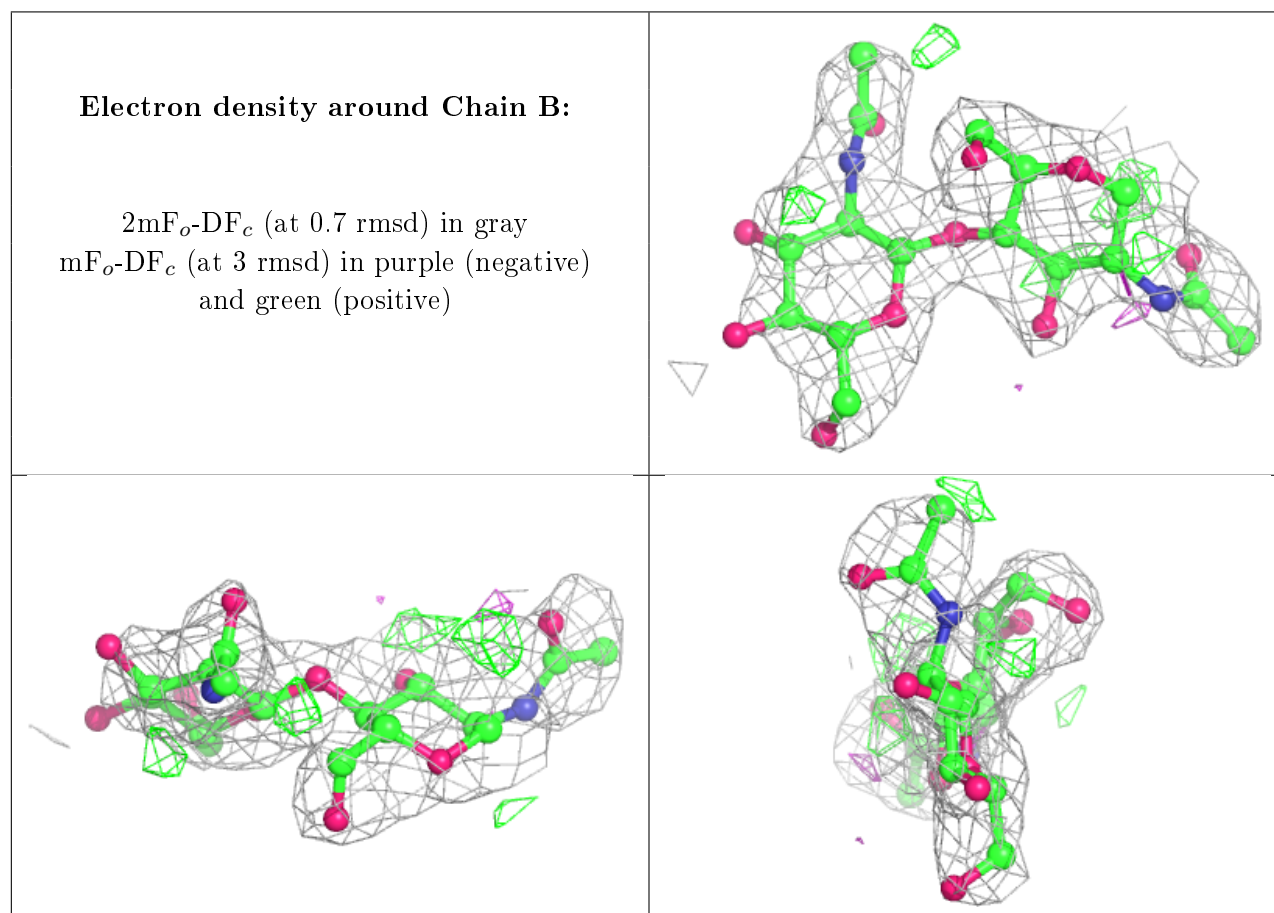
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	ALS	A	91	11/12	0.98	0.09	15,23,47,48	0

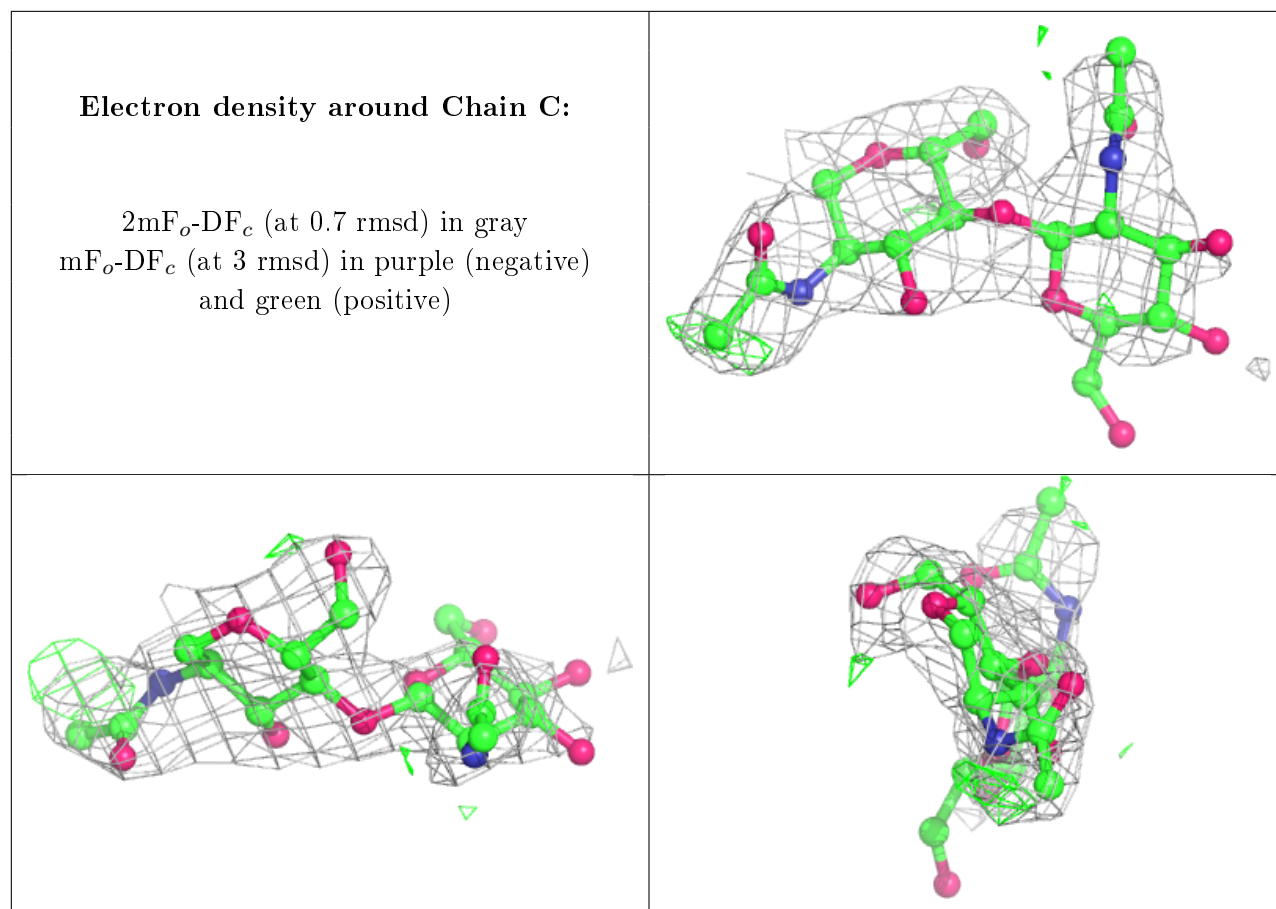
6.3 Carbohydrates [i](#)

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	0.76	0.38	69,74,79,80	0
2	NAG	C	2	14/15	0.81	0.43	72,78,82,84	0
2	NAG	C	1	14/15	0.88	0.23	42,51,59,63	0
2	NAG	B	1	14/15	0.91	0.21	43,52,56,64	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.





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
3	CA	A	604	1/1	0.97	0.04	23,23,23,23	0
4	CL	A	605	1/1	0.99	0.10	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.