



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 21, 2020 – 04:54 PM BST

PDB ID : 3T77
Title : S25-2- A(2-4)KDO disaccharide complex
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Evans, S.V.
Deposited on : 2011-07-29
Resolution : 1.74 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 : 2.14.6
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) : 2.14.6

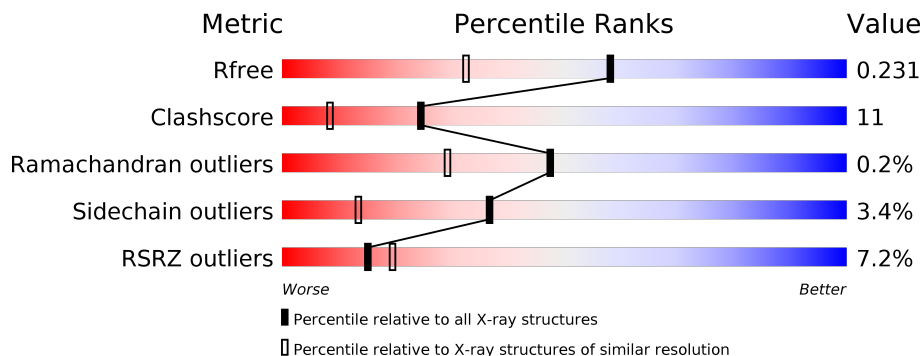
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.74 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	3764 (1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)
RSRZ outliers	127900	3705 (1.76-1.72)

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	219	
2	B	224	
3	C	2	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3856 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 S25-2 FAB (IGG1K) light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	219	1701	1056	290	347	8	0	0	0

- Molecule 2 is a protein called S25-2 FAB (IGG1K) heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	224	1712	1082	285	337	8	0	0	0

- Molecule 3 is an oligosaccharide called 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-prop-2-en-1-yl 3-deoxy-alpha-D-manno-oct-2-ulopyranosidonic acid.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
3	C	2	34	19	15	0	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		

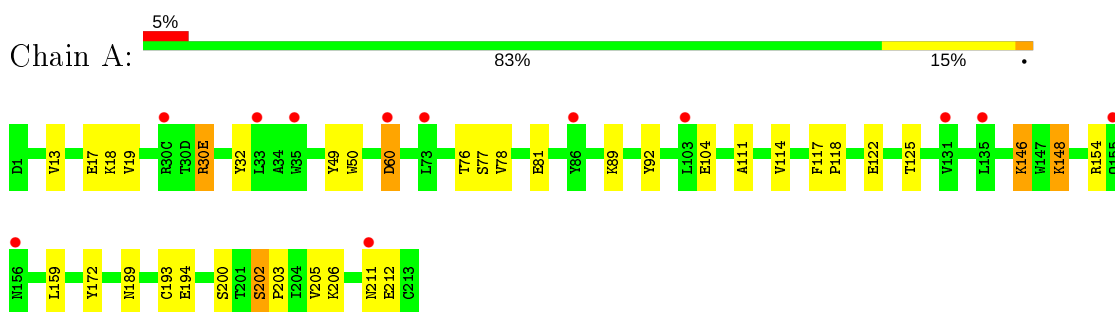
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	215	Total 215	O 215	0	0
6	B	191	Total 191	O 191	0	0

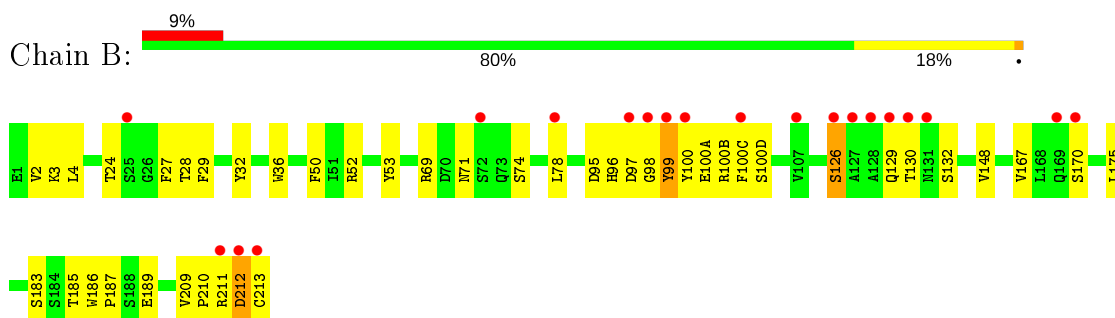
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: S25-2 FAB (IGG1K) light chain



- Molecule 2: S25-2 FAB (IGG1K) heavy chain



- Molecule 3: 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-prop-2-en-1-yl 3-deoxy-alpha-D-manno-oct-2-ulopyranosidonic acid



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.90Å 81.80Å 130.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.82 – 1.74 19.82 – 1.74	Depositor EDS
% Data completeness (in resolution range)	92.8 (19.82-1.74) 92.8 (19.82-1.74)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.90 (at 1.73Å)	Xtrriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.205 , 0.236 0.199 , 0.231	Depositor DCC
R_{free} test set	4818 reflections (10.12%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtrriage
Anisotropy	0.537	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3856	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% 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: ZN, KDO, MG, KDA

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	0.38	0/1736	0.51	0/2351
2	B	0.38	0/1758	0.57	0/2400
All	All	0.38	0/3494	0.55	0/4751

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

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	212	ASP	Peptide

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	1701	0	1648	37	1
2	B	1712	0	1657	45	0
3	C	34	0	28	1	0
4	A	2	0	0	0	0
5	A	1	0	0	0	0
6	A	215	0	0	8	1
6	B	191	0	0	7	0
All	All	3856	0	3333	74	2

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 11.

All (74) 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)
2:B:211:ARG:NH2	6:B:217:HOH:O	1.82	1.12
2:B:100(D):SER:OG	6:B:277:HOH:O	1.87	0.91
6:B:267:HOH:O	3:C:2:KDO:O8	1.91	0.88
1:A:146:LYS:HD3	1:A:148:LYS:HZ1	1.37	0.87
1:A:104:GLU:CG	6:A:317:HOH:O	2.27	0.81
2:B:69:ARG:HE	2:B:71:ASN:HD21	1.29	0.79
1:A:104:GLU:HG3	6:A:317:HOH:O	1.82	0.79
1:A:159:LEU:HG	2:B:167:VAL:HG21	1.69	0.72
1:A:146:LYS:HD3	1:A:148:LYS:NZ	2.04	0.71
2:B:96:HIS:CE1	2:B:98:GLY:HA2	2.28	0.69
1:A:19:VAL:HG23	1:A:78:VAL:HG21	1.76	0.68
1:A:159:LEU:CG	2:B:167:VAL:HG21	2.23	0.67
1:A:148:LYS:NZ	1:A:194:GLU:OE2	2.31	0.64
1:A:159:LEU:HD11	2:B:167:VAL:CG2	2.28	0.63
1:A:122:GLU:O	1:A:125:THR:HG22	2.00	0.61
2:B:69:ARG:NE	2:B:71:ASN:HD21	1.99	0.61
2:B:210:PRO:HB2	2:B:213:CYS:SG	2.41	0.60
2:B:99:TYR:N	2:B:99:TYR:HD1	1.99	0.60
2:B:126:SER:OG	2:B:126:SER:O	2.20	0.60
1:A:104:GLU:HG2	6:A:317:HOH:O	1.98	0.60
1:A:104:GLU:HG3	6:A:426:HOH:O	2.02	0.57
1:A:18:LYS:HE2	1:A:76:THR:HG22	1.85	0.57
2:B:99:TYR:N	2:B:99:TYR:CD1	2.69	0.57
2:B:32:TYR:CE1	2:B:97:ASP:HB3	2.43	0.54
2:B:96:HIS:NE2	2:B:100(A):GLU:OE1	2.42	0.53
1:A:104:GLU:OE1	1:A:172:TYR:OH	2.21	0.53
2:B:96:HIS:HD2	2:B:100:TYR:O	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:98:GLY:C	2:B:99:TYR:HD1	2.14	0.51
2:B:96:HIS:HA	6:B:276:HOH:O	2.12	0.50
2:B:4:LEU:HD22	2:B:24:THR:HG22	1.94	0.50
1:A:19:VAL:CG2	1:A:78:VAL:HG21	2.40	0.50
1:A:159:LEU:HD21	2:B:167:VAL:CG2	2.42	0.50
2:B:210:PRO:CB	2:B:213:CYS:SG	3.00	0.50
2:B:212:ASP:O	2:B:213:CYS:HB2	2.11	0.49
1:A:194:GLU:HG2	1:A:205:VAL:HG22	1.95	0.49
2:B:210:PRO:HG2	2:B:213:CYS:CB	2.43	0.49
2:B:100(A):GLU:HA	6:B:316:HOH:O	2.13	0.49
2:B:100(B):ARG:NH1	6:B:264:HOH:O	2.12	0.48
1:A:154:ARG:NH1	6:A:398:HOH:O	2.27	0.48
2:B:132:SER:O	2:B:183:SER:OG	2.32	0.48
1:A:200:SER:HB3	6:A:417:HOH:O	2.13	0.48
1:A:111:ALA:O	6:A:336:HOH:O	2.20	0.48
1:A:32:TYR:HD1	1:A:92:TYR:HA	1.80	0.47
1:A:89:LYS:NZ	2:B:100(C):PHE:CD1	2.82	0.47
2:B:210:PRO:HG2	2:B:213:CYS:HB3	1.97	0.47
2:B:185:THR:O	2:B:189:GLU:HB2	2.16	0.46
1:A:159:LEU:CD2	2:B:167:VAL:HG21	2.46	0.46
1:A:189:ASN:OD1	1:A:211:ASN:OD1	2.34	0.46
2:B:175:LEU:HD12	2:B:175:LEU:C	2.36	0.46
2:B:95:ASP:OD2	2:B:96:HIS:N	2.46	0.45
2:B:69:ARG:HE	2:B:71:ASN:ND2	2.05	0.45
2:B:186:TRP:CG	2:B:187:PRO:HA	2.51	0.45
2:B:3:LYS:O	2:B:24:THR:HA	2.16	0.45
2:B:2:VAL:HG13	2:B:27:PHE:CD1	2.52	0.44
2:B:148:VAL:CG2	2:B:175:LEU:HD21	2.47	0.44
2:B:36:TRP:CE2	2:B:78:LEU:HB2	2.53	0.44
1:A:146:LYS:HE3	1:A:146:LYS:HB3	1.49	0.43
1:A:81:GLU:HG2	6:A:283:HOH:O	2.17	0.43
2:B:2:VAL:HG13	2:B:27:PHE:HD1	1.83	0.43
2:B:100(B):ARG:HA	6:B:401:HOH:O	2.17	0.43
1:A:159:LEU:CD1	2:B:167:VAL:CG2	2.97	0.43
2:B:52:ARG:O	2:B:53:TYR:HA	2.19	0.42
1:A:202:SER:HA	1:A:203:PRO:HD3	1.92	0.42
1:A:114:VAL:O	1:A:206:LYS:HE3	2.20	0.42
1:A:60:ASP:OD1	1:A:60:ASP:N	2.53	0.42
1:A:117:PHE:HA	1:A:118:PRO:HD3	1.95	0.41
1:A:30(E):ARG:HB2	1:A:30(E):ARG:HE	1.61	0.41
1:A:49:TYR:HE2	1:A:50:TRP:CZ3	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LEU:HD21	2:B:167:VAL:HG22	2.02	0.41
2:B:29:PHE:CD2	2:B:74:SER:HA	2.55	0.41
2:B:187:PRO:HG3	2:B:210:PRO:HG3	2.03	0.41
1:A:159:LEU:CG	2:B:167:VAL:CG2	2.98	0.40
1:A:13:VAL:HG22	1:A:17:GLU:HB2	2.02	0.40
1:A:76:THR:O	1:A:77:SER:C	2.59	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:308:HOH:O	6:A:384:HOH:O[3_556]	1.91	0.29
1:A:49:TYR:OH	1:A:212:GLU:OE2[2_465]	2.01	0.19

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	217/219 (99%)	210 (97%)	7 (3%)	0	100	100
2	B	222/224 (99%)	212 (96%)	9 (4%)	1 (0%)	29	12
All	All	439/443 (99%)	422 (96%)	16 (4%)	1 (0%)	47	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	170	SER

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	195/195 (100%)	189 (97%)	6 (3%)	40	16
2	B	192/192 (100%)	185 (96%)	7 (4%)	35	12
All	All	387/387 (100%)	374 (97%)	13 (3%)	37	13

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30(E)	ARG
1	A	60	ASP
1	A	146	LYS
1	A	148	LYS
1	A	193	CYS
1	A	202	SER
2	B	28	THR
2	B	50	PHE
2	B	99	TYR
2	B	126	SER
2	B	129	GLN
2	B	130	THR
2	B	209	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	71	ASN
2	B	96	HIS
2	B	169	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates i

2 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
3	KDA	C	1	3	15,19,19	1.90	5 (33%)	19,27,27	1.44	2 (10%)
3	KDO	C	2	3	12,15,16	1.91	3 (25%)	16,21,24	1.62	5 (31%)

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
3	KDA	C	1	3	-	4/11/35/35	0/1/1/1
3	KDO	C	2	3	-	0/6/26/30	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2	KDO	C4-C5	-4.47	1.45	1.52
3	C	1	KDA	C4-C5	-3.64	1.47	1.52
3	C	1	KDA	O6-C2	2.89	1.46	1.42
3	C	1	KDA	O6-C6	2.84	1.48	1.44
3	C	2	KDO	O6-C6	2.76	1.48	1.44
3	C	1	KDA	O7-C7	-2.65	1.37	1.43
3	C	1	KDA	C3-C4	-2.54	1.49	1.53
3	C	2	KDO	O7-C7	-2.51	1.38	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	KDA	O8-C8-C7	3.39	118.45	111.07
3	C	1	KDA	C3-C2-C1	-3.17	108.62	114.12
3	C	2	KDO	O8-C8-C7	3.12	117.87	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	KDO	O6-C6-C5	2.30	111.10	107.87
3	C	2	KDO	O4-C4-C5	-2.17	105.78	110.14
3	C	2	KDO	C3-C4-C5	2.16	113.82	110.69
3	C	2	KDO	C8-C7-C6	-2.08	107.95	112.17

There are no chirality outliers.

All (4) torsion outliers are listed below:

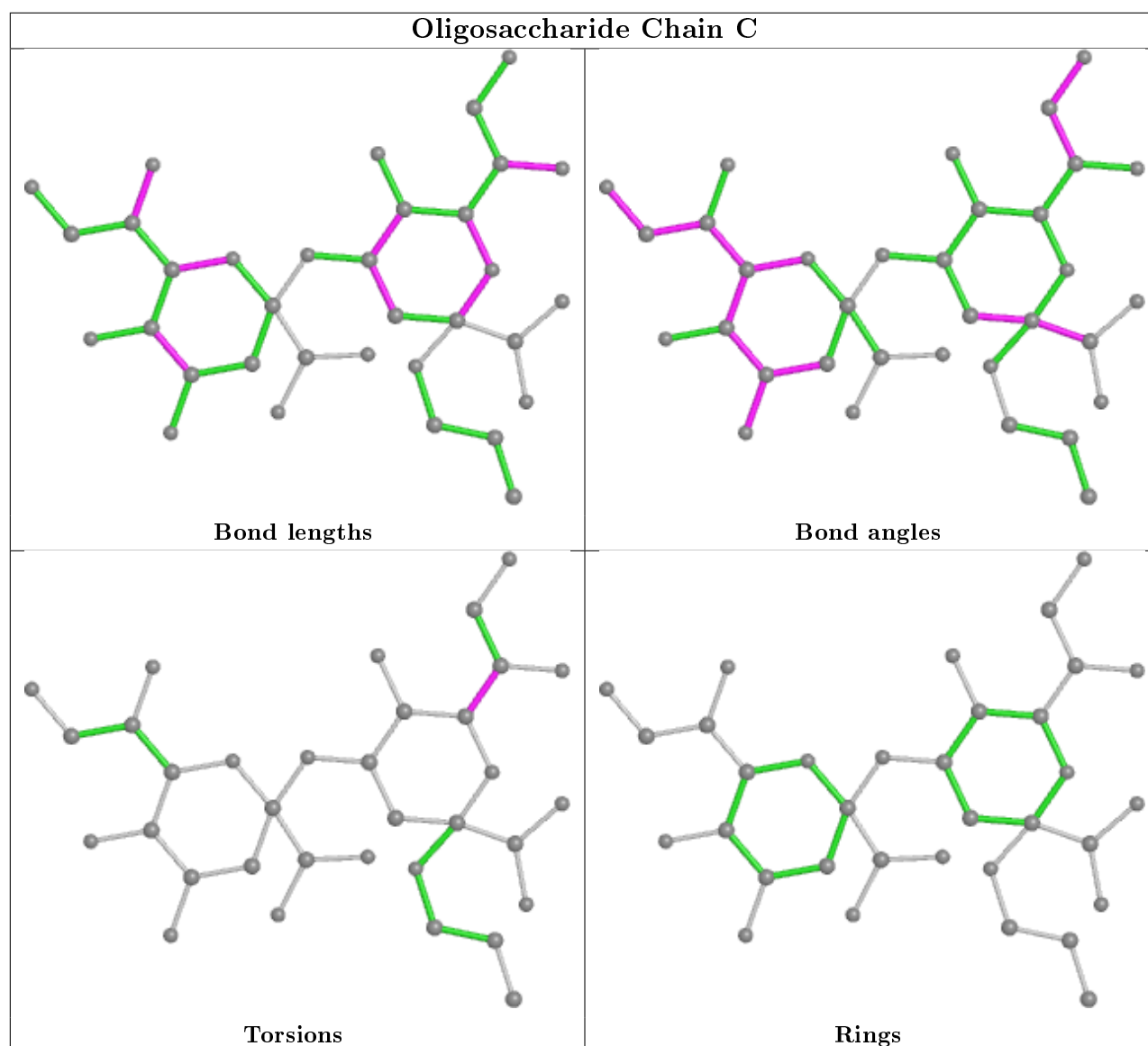
Mol	Chain	Res	Type	Atoms
3	C	1	KDA	C5-C6-C7-O7
3	C	1	KDA	O6-C6-C7-O7
3	C	1	KDA	O6-C6-C7-C8
3	C	1	KDA	C5-C6-C7-C8

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2	KDO	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 3 ligands modelled in this entry, 3 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 [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	219/219 (100%)	0.33	12 (5%) 25 30	20, 29, 42, 56	0
2	B	224/224 (100%)	0.65	20 (8%) 9 12	19, 29, 55, 73	0
All	All	443/443 (100%)	0.49	32 (7%) 15 19	19, 29, 49, 73	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	213	CYS	12.4
2	B	98	GLY	12.1
2	B	99	TYR	10.5
2	B	100	TYR	7.1
2	B	129	GLN	6.0
2	B	128	ALA	5.7
2	B	126	SER	5.7
2	B	97	ASP	4.1
1	A	211	ASN	4.0
1	A	33	LEU	3.9
1	A	155	GLN	3.4
2	B	130	THR	3.2
1	A	103	LEU	3.2
2	B	100(C)	PHE	3.1
1	A	30(C)	ARG	3.1
2	B	131	ASN	2.8
1	A	60	ASP	2.7
2	B	212	ASP	2.7
2	B	72	SER	2.6
1	A	156	ASN	2.5
2	B	107	VAL	2.5
2	B	211	ARG	2.4
2	B	127	ALA	2.4
1	A	131	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	135	LEU	2.4
2	B	25	SER	2.3
1	A	86	TYR	2.3
1	A	73	LEU	2.2
1	A	35	TRP	2.1
2	B	169	GLN	2.1
2	B	170	SER	2.1
2	B	78	LEU	2.0

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

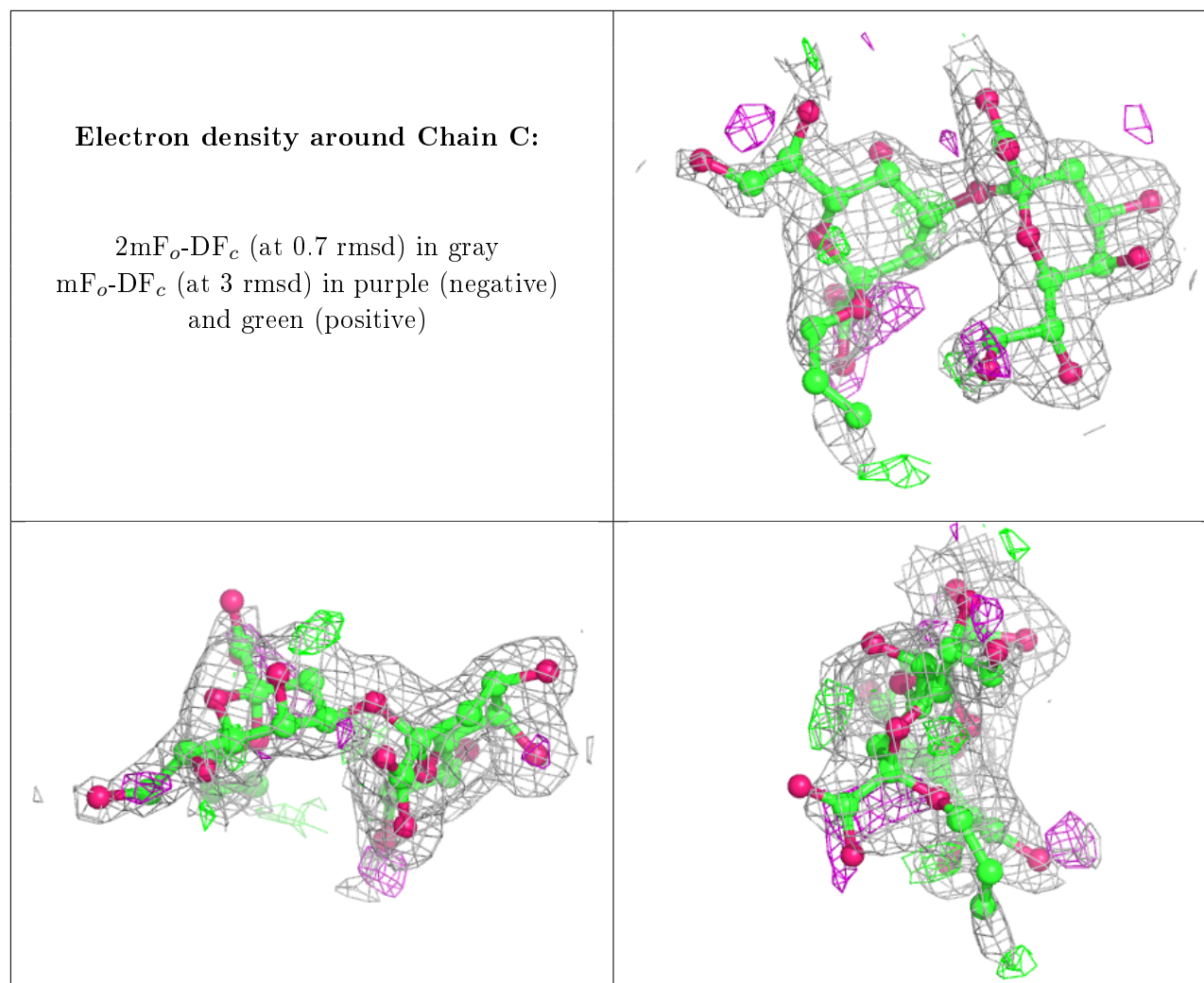
There are no non-standard protein/DNA/RNA residues in this entry.

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(Å ²)	Q<0.9
3	KDA	C	1	19/19	0.82	0.28	35,48,57,58	0
3	KDO	C	2	15/16	0.95	0.08	28,32,39,47	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(Å ²)	Q < 0.9
4	MG	A	216	1/1	0.67	0.16	39,39,39,39	0
4	MG	A	217	1/1	0.92	0.05	38,38,38,38	0
5	ZN	A	218	1/1	0.98	0.03	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.