Full wwPDB X-ray Structure Validation Report (i)

Mar 3, 2024 – 12:48 AM EST

PDB ID	:	5WD7
Title	:	Structure of a bacterial polysialyltransferase in complex with fondaparinux
Authors	:	Worrall, L.J.; Lizak, C.; Strynadka, N.C.J.
Deposited on	:	2017-07-04
Resolution	:	3.10 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

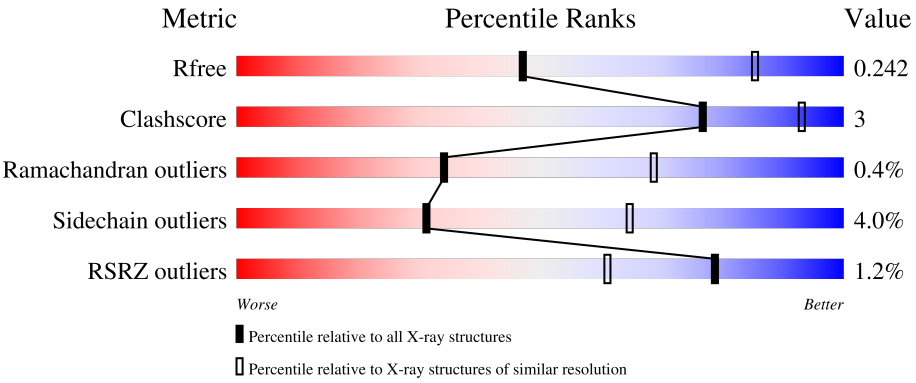
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

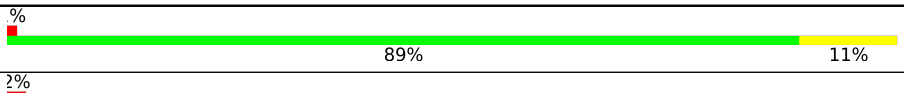
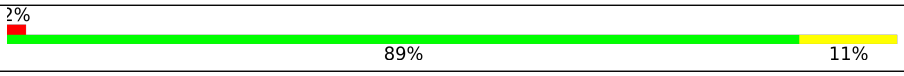
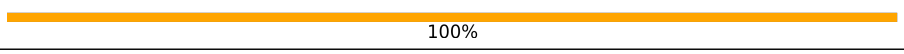
The reported resolution of this entry is 3.10 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	382	
1	M	382	
2	B	5	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6514 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.

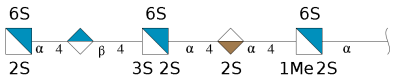
- Molecule 1 is a protein called SiaD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	0	0
			3190	2080	524	570	16			
1	M	382	Total	C	N	O	S	0	2	0
			3223	2098	536	573	16			

There are 6 discrepancies between the modelled and reference sequences:

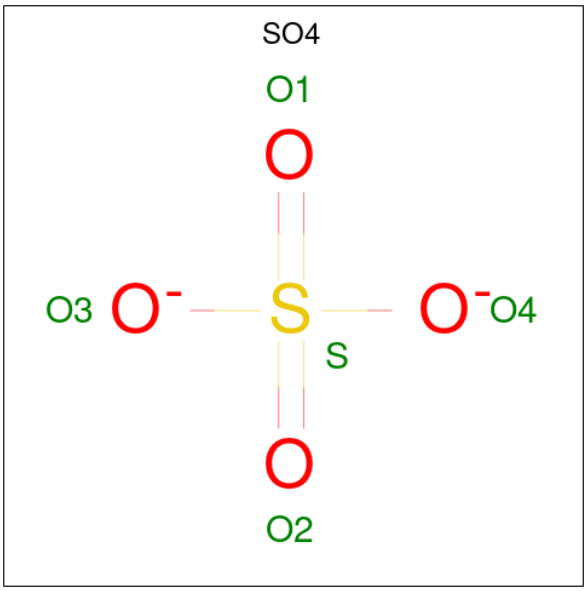
Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	initiating methionine	UNP G4RIN4
A	68	ALA	LYS	conflict	UNP G4RIN4
A	69	ALA	LYS	conflict	UNP G4RIN4
M	20	MET	-	initiating methionine	UNP G4RIN4
M	68	ALA	LYS	conflict	UNP G4RIN4
M	69	ALA	LYS	conflict	UNP G4RIN4

- Molecule 2 is an oligosaccharide called 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-3,6-di-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-methyl 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranoside.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	S	0	0	0
			91	31	3	49	8			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



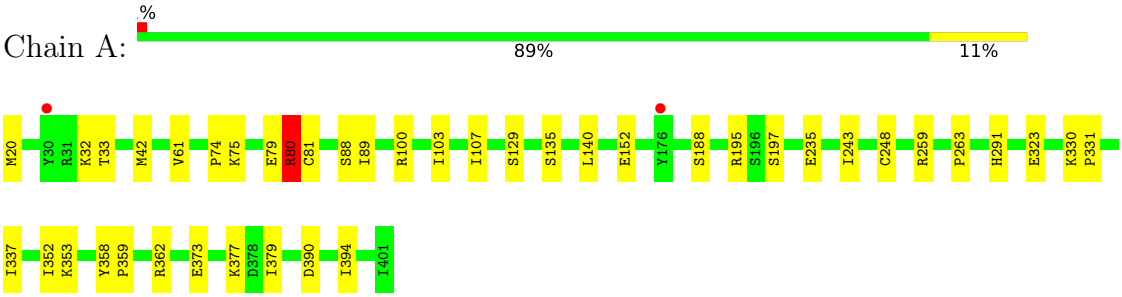
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	M	1	Total O S 5 4 1	0	0
3	M	1	Total O S 5 4 1	0	0

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.

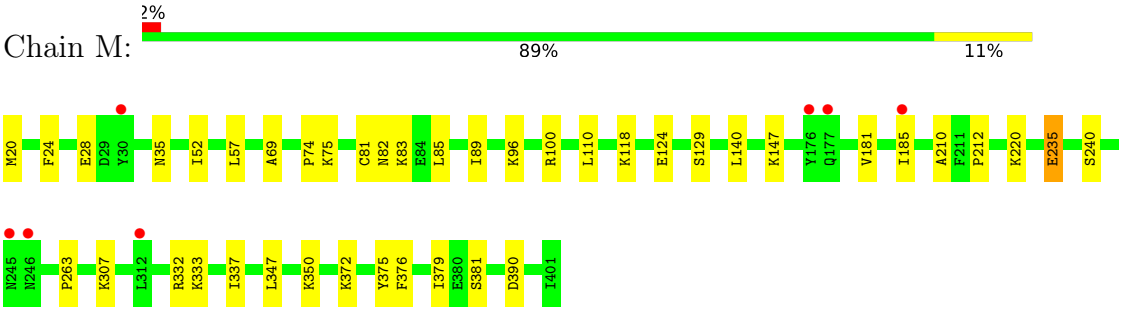
• Molecule 1: SiaD

Chain A:



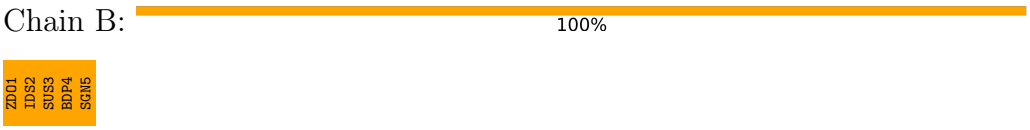
• Molecule 1: SiaD

Chain M:



• Molecule 2: 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-4)-2-deoxy-3,6-di-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-methyl 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranoside

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	78.31Å 78.31Å 299.89Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.93 – 3.10 44.93 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.93-3.10) 100.0 (44.93-3.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.50 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.187 , 0.238 0.195 , 0.242	Depositor DCC
R_{free} test set	938 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å^2)	87.6	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 58.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6514	wwPDB-VP
Average B, all atoms (Å^2)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IDS, SO4, BDP, ZDO, SUS, SGN

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	0.60	0/3255	0.83	2/4363 (0.0%)
1	M	0.57	0/3288	0.79	0/4405
All	All	0.59	0/6543	0.81	2/8768 (0.0%)

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
1	A	0	1
1	M	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	80	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	362	ARG	NE-CZ-NH2	-5.46	117.57	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	20	MET	Peptide
1	M	20	MET	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	3190	0	3317	14	0
1	M	3223	0	3354	20	0
2	B	91	0	27	5	0
3	M	10	0	0	1	0
All	All	6514	0	6698	39	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 3.

All (39) 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:M:100[B]:ARG:HH21	1:M:100[B]:ARG:HG2	1.56	0.69
1:M:337:ILE:HG21	1:M:379:ILE:HD11	1.77	0.67
1:M:124:GLU:HG2	1:M:147:LYS:HB2	1.78	0.64
1:M:100[B]:ARG:HH21	1:M:100[B]:ARG:CG	2.12	0.61
1:A:337:ILE:HG21	1:A:379:ILE:HD11	1.84	0.58
2:B:2:IDS:O1S	2:B:2:IDS:O3	2.13	0.56
1:M:74:PRO:HB2	1:M:75:LYS:HE3	1.87	0.56
1:M:372:LYS:HD2	1:M:375:TYR:CE2	2.42	0.55
1:M:332:ARG:HG2	1:M:333:LYS:HG3	1.89	0.54
1:M:118:LYS:NZ	3:M:502:SO4:O4	2.41	0.53
1:A:42:MET:HG2	1:A:80:ARG:NH1	2.23	0.53
1:M:235:GLU:OE2	1:M:240:SER:OG	2.28	0.52
1:M:100[B]:ARG:CG	1:M:100[B]:ARG:NH2	2.71	0.52
1:A:331:PRO:O	1:A:352:ILE:HG23	2.10	0.51
2:B:4:BDP:C6	2:B:5:SGN:H5	2.43	0.49
1:M:35:ASN:ND2	1:M:57:LEU:HD22	2.26	0.49
1:A:81:CYS:SG	1:A:89:ILE:HD11	2.53	0.49
1:A:243:ILE:HG23	1:A:248:CYS:HB2	1.97	0.47
1:M:332:ARG:CG	1:M:333:LYS:HG3	2.45	0.46
2:B:3:SUS:O9S	2:B:3:SUS:N2	2.49	0.46
1:A:259:ARG:CZ	1:A:291:HIS:CE1	2.99	0.45
1:A:359:PRO:HG2	1:A:394:ILE:CG2	2.46	0.45
1:A:103:ILE:HG22	1:A:107:ILE:HD12	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:210:ALA:HB1	1:M:347:LEU:HD11	2.00	0.44
1:M:81:CYS:SG	1:M:82:ASN:N	2.91	0.44
1:M:81:CYS:SG	1:M:89:ILE:HD11	2.59	0.43
1:M:52:ILE:CG2	1:M:85:LEU:HD22	2.48	0.43
1:A:75:LYS:NZ	1:A:79:GLU:OE1	2.39	0.42
1:M:376:PHE:HA	1:M:379:ILE:HG22	1.99	0.42
1:M:181:VAL:O	1:M:185:ILE:HD12	2.20	0.42
1:A:42:MET:HE1	1:A:80:ARG:HD3	2.02	0.41
2:B:1:ZDO:OS1	2:B:2:IDS:C1	2.68	0.41
2:B:2:IDS:C6	2:B:3:SUS:H5	2.50	0.41
1:M:181:VAL:HG12	1:M:185:ILE:HD12	2.02	0.41
1:A:358:TYR:HB3	1:A:359:PRO:HD3	2.02	0.41
1:M:110:LEU:HD12	1:M:140:LEU:HD11	2.01	0.41
1:A:42:MET:HG2	1:A:80:ARG:HH12	1.85	0.40
1:A:188:SER:O	1:A:195:ARG:NH2	2.53	0.40
1:A:33:THR:HG23	1:A:61:VAL:HG23	2.04	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	380/382 (100%)	361 (95%)	18 (5%)	1 (0%)	41	73
1	M	383/382 (100%)	361 (94%)	20 (5%)	2 (0%)	29	64
All	All	763/764 (100%)	722 (95%)	38 (5%)	3 (0%)	34	69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	PRO
1	M	69	ALA

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Mol	Chain	Res	Type
1	M	263	PRO

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	362/362 (100%)	345 (95%)	17 (5%)	26	59
1	M	365/362 (101%)	353 (97%)	12 (3%)	38	69
All	All	727/724 (100%)	698 (96%)	29 (4%)	31	65

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LYS
1	A	74	PRO
1	A	80	ARG
1	A	88	SER
1	A	100	ARG
1	A	129	SER
1	A	135	SER
1	A	140	LEU
1	A	152	GLU
1	A	197	SER
1	A	235	GLU
1	A	323	GLU
1	A	330	LYS
1	A	353	LYS
1	A	373	GLU
1	A	377	LYS
1	A	390	ASP
1	M	24	PHE
1	M	28	GLU
1	M	83	LYS
1	M	96	LYS
1	M	129	SER

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Mol	Chain	Res	Type
1	M	212	PRO
1	M	220	LYS
1	M	235	GLU
1	M	307	LYS
1	M	350	LYS
1	M	381	SER
1	M	390	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	ASN
1	A	258	GLN
1	M	291	HIS
1	M	310	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)

5 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	ZDO	B	1	2	20,21,21	0.65	1 (5%)	25,32,32	1.28	2 (8%)
2	IDS	B	2	2	16,16,17	0.77	1 (6%)	17,24,26	1.36	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	SUS	B	3	2	22,23,24	0.75	1 (4%)	24,36,38	1.29	2 (8%)
2	BDP	B	4	2	12,12,13	0.45	0	14,17,19	0.88	1 (7%)
2	SGN	B	5	2	18,19,20	0.80	1 (5%)	22,29,31	1.32	1 (4%)

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	ZDO	B	1	2	-	8/13/33/33	0/1/1/1
2	IDS	B	2	2	-	4/9/26/29	0/1/1/1
2	SUS	B	3	2	-	3/16/33/36	0/1/1/1
2	BDP	B	4	2	-	1/4/21/24	0/1/1/1
2	SGN	B	5	2	-	1/11/28/31	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	ZDO	OS2-S6	2.69	1.56	1.45
2	B	5	SGN	O5S-S2	2.59	1.56	1.45
2	B	2	IDS	O1S-S	2.27	1.55	1.45
2	B	3	SUS	O5S-S2	2.24	1.54	1.45

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1	ZDO	C4-C3-C2	-4.46	103.81	110.34
2	B	3	SUS	O8S-S3-O7S	4.20	129.08	112.22
2	B	2	IDS	C2-O2-S	4.11	123.27	117.91
2	B	5	SGN	O5S-S2-O4S	4.07	128.55	112.22
2	B	3	SUS	O5S-S2-O4S	3.35	125.67	112.22
2	B	1	ZDO	OS2-S6-OS1	2.91	123.89	112.22
2	B	2	IDS	O2S-S-O1S	2.24	121.22	112.22
2	B	2	IDS	O4-C4-C3	-2.20	105.27	110.35
2	B	4	BDP	C2-C3-C4	-2.14	107.19	110.89

There are no chirality outliers.

All (17) torsion outliers are listed below:

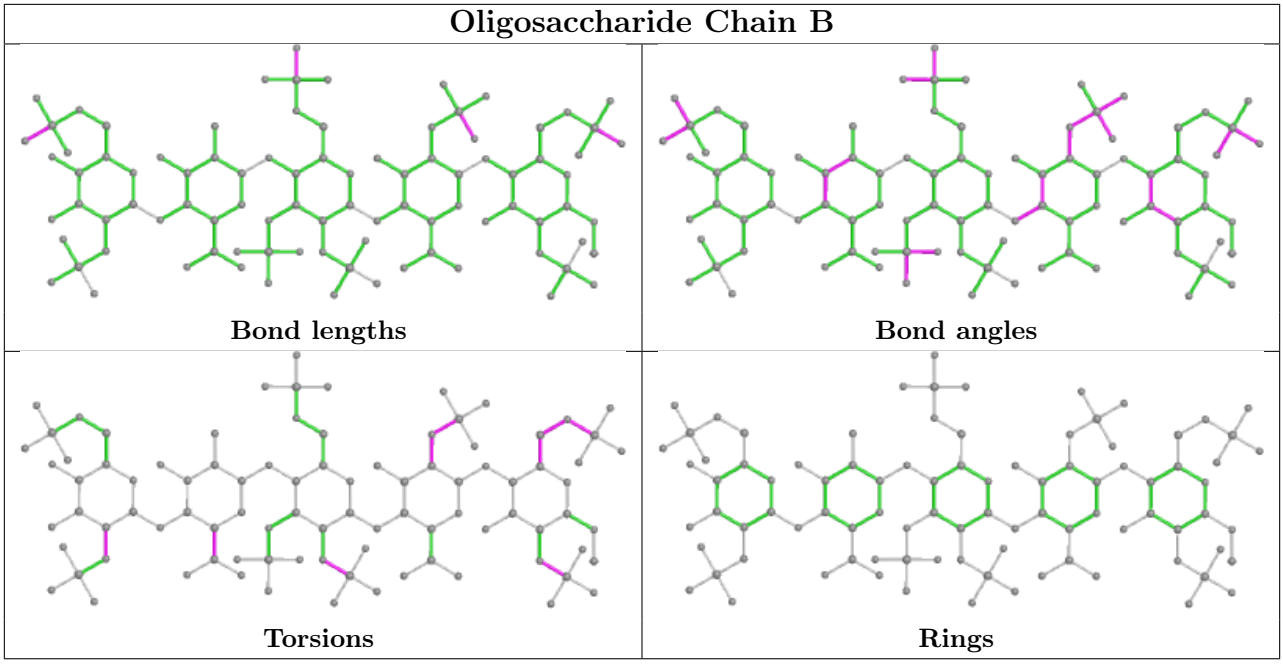
Mol	Chain	Res	Type	Atoms
2	B	1	ZDO	C4-C5-C6-O6
2	B	1	ZDO	O5-C5-C6-O6
2	B	1	ZDO	C2-N2-S2-OSB
2	B	1	ZDO	C2-N2-S2-OSC
2	B	2	IDS	C1-C2-O2-S
2	B	2	IDS	C2-O2-S-O3S
2	B	3	SUS	C2-N2-S1-O2S
2	B	1	ZDO	C6-O6-S6-OS2
2	B	3	SUS	C2-N2-S1-O1S
2	B	2	IDS	C3-C2-O2-S
2	B	1	ZDO	C5-C6-O6-S6
2	B	1	ZDO	C6-O6-S6-OS3
2	B	1	ZDO	C2-N2-S2-OSA
2	B	2	IDS	C2-O2-S-O1S
2	B	3	SUS	C2-N2-S1-O3S
2	B	5	SGN	C1-C2-N2-S1
2	B	4	BDP	O5-C5-C6-O6B

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	IDS	3	0
2	B	3	SUS	2	0
2	B	1	ZDO	1	0
2	B	4	BDP	1	0
2	B	5	SGN	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)

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	$\#\|Z\| > 2$
3	SO4	M	502	-	4,4,4	0.42	0	6,6,6	0.29	0
3	SO4	M	501	-	4,4,4	0.39	0	6,6,6	0.43	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	502	SO4	1	0

5.7 Other polymers

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, 95^{th} 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	382/382 (100%)	-0.26	2 (0%)	91	81	50, 80, 113, 196	0
1	M	382/382 (100%)	-0.05	7 (1%)	68	47	55, 93, 132, 159	0
All	All	764/764 (100%)	-0.16	9 (1%)	79	61	50, 85, 125, 196	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	30	TYR	5.7
1	M	177	GLN	2.9
1	M	30	TYR	2.4
1	M	176	TYR	2.3
1	M	245	ASN	2.2
1	M	246	ASN	2.1
1	M	185	ILE	2.1
1	A	176	TYR	2.0
1	M	312	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, 95^{th} 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	SGN	B	5	19/20	0.84	0.16	88,119,152,157	0

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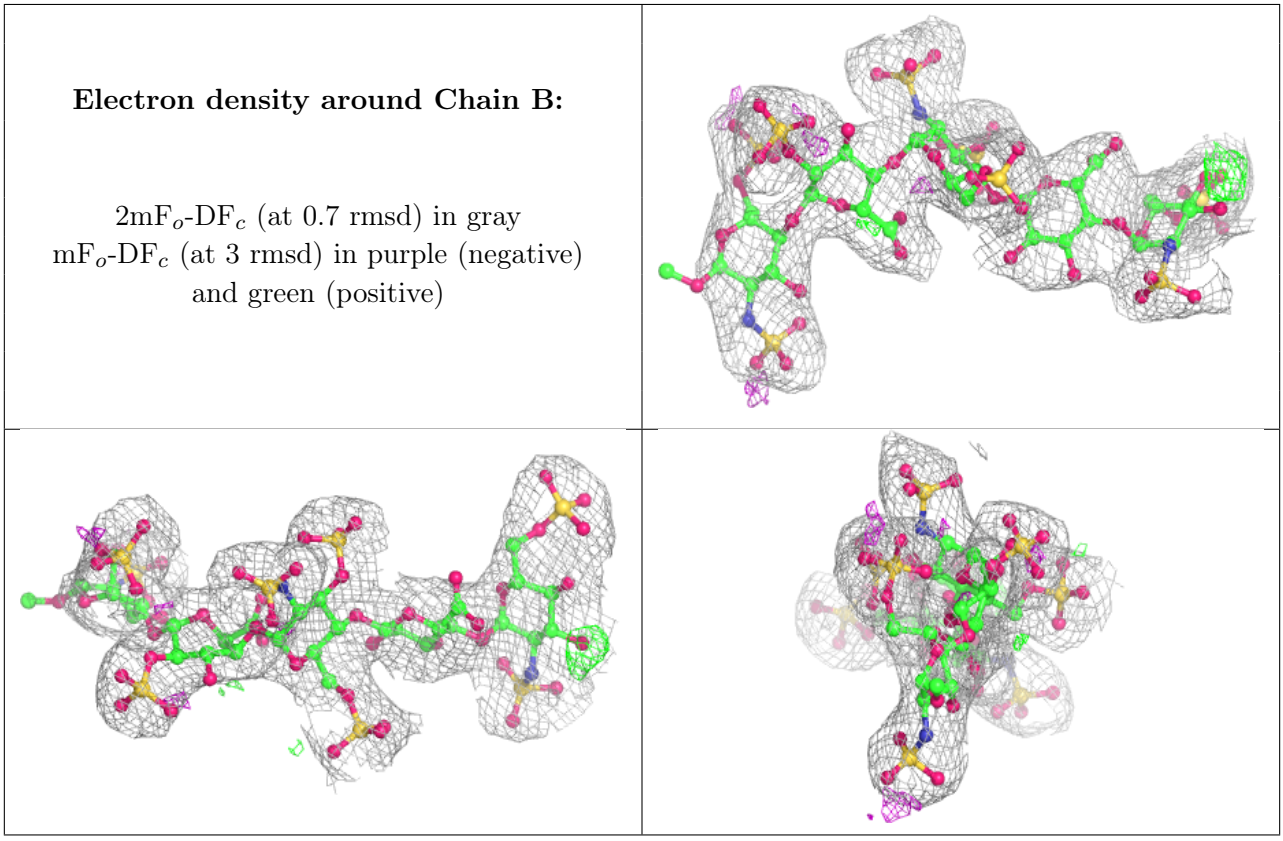
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZDO	B	1	21/21	0.90	0.17	113,128,140,159	0
2	IDS	B	2	16/17	0.92	0.12	99,118,136,138	0
2	BDP	B	4	12/13	0.93	0.16	88,98,110,127	0
2	SUS	B	3	23/24	0.95	0.15	84,94,105,123	0

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
3	SO4	M	502	5/5	0.93	0.14	92,98,104,123	0
3	SO4	M	501	5/5	0.95	0.13	85,99,108,117	0

6.5 Other polymers (i)

There are no such residues in this entry.