



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2024 – 01:44 pm BST

PDB ID : 1O75
Title : Tp47, the 47-Kilodalton Lipoprotein of *Treponema pallidum*
Authors : Deka, R.K.; Machius, M.; Norgard, M.V.; Tomchick, D.R.
Deposited on : 2002-10-23
Resolution : 1.95 Å(reported)

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

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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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
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)	:	2.36.2

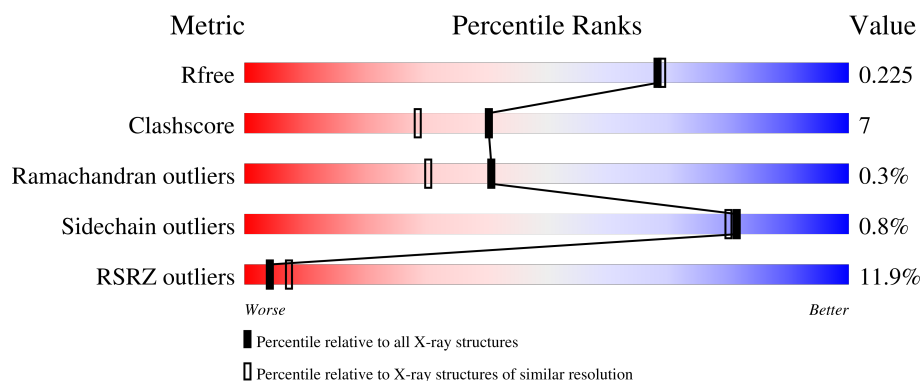
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	415	<div> <div>9%</div> <div>87%</div> <div>9%</div> <div>.</div> </div>
1	B	415	<div> <div>13%</div> <div>80%</div> <div>16%</div> <div>.</div> </div>
2	C	2	<div> <div>50%</div> <div>50%</div> </div>

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
3	XE	B	1415	-	-	X	-
3	XE	B	1417	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7017 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 47 KDA MEMBRANE ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	9	0
			3175	1993	539	630	13			
1	B	402	Total	C	N	O	S	0	5	0
			3179	1998	542	626	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	SER	HIS	engineered mutation	UNP P29723
A	9	SER	HIS	engineered mutation	UNP P29273
B	5	SER	HIS	engineered mutation	UNP P29723
B	9	SER	HIS	engineered mutation	UNP P29273

- Molecule 2 is an oligosaccharide called 2,3-di-O-sulfo-alpha-D-glucopyranose-(1-6)-2,3-di-O-sulfo-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	S	0	0	0
			39	12	23	4			

- Molecule 3 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Xe	0	0
			2	2		
3	B	3	Total	Xe	0	0
			3	3		

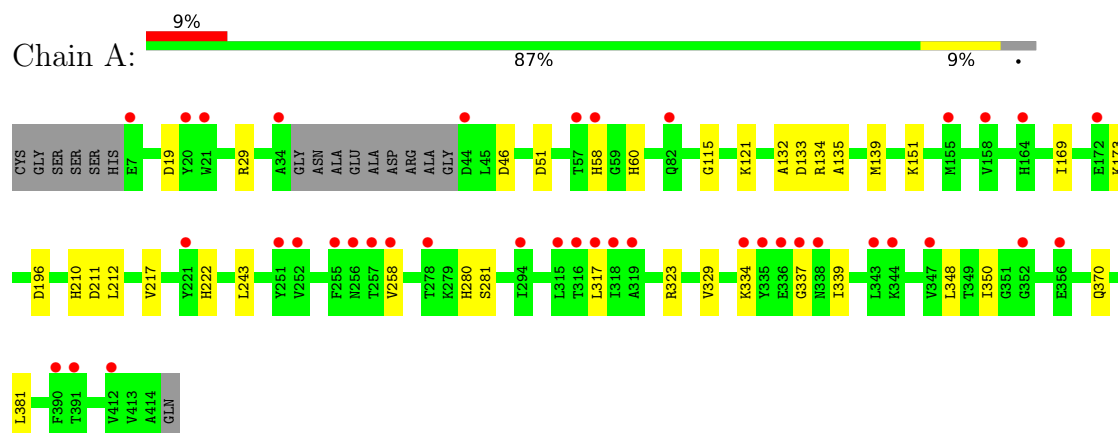
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	311	Total 311	O 311	0	0
4	B	308	Total 308	O 308	0	0

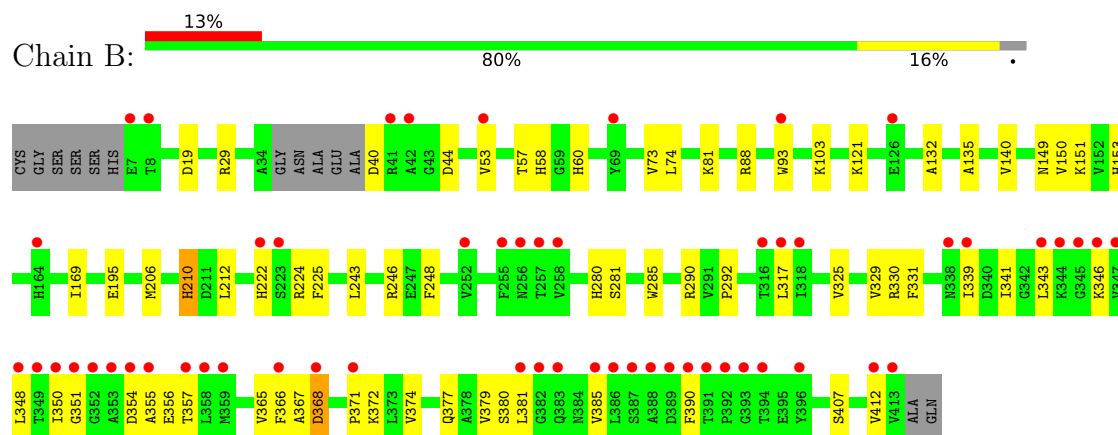
3 Residue-property plots

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: 47 KDA MEMBRANE ANTIGEN



• Molecule 1: 47 KDA MEMBRANE ANTIGEN



• Molecule 2: 2,3-di-O-sulfo-alpha-D-glucopyranose-(1-6)-2,3-di-O-sulfo-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.93Å 128.93Å 151.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.20 – 1.95 28.15 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (28.20-1.95) 94.2 (28.15-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.205 , 0.228 0.200 , 0.225	Depositor DCC
R_{free} test set	5036 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	33.7	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7017	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PDX, XE

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	0.51	0/3239	0.68	0/4372
1	B	0.51	0/3245	0.68	0/4380
All	All	0.51	0/6484	0.68	0/8752

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

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	3175	0	3036	26	0
1	B	3179	0	3042	67	0
2	C	39	0	19	1	0
3	A	2	0	0	1	0
3	B	3	0	0	7	0
4	A	311	0	0	2	0
4	B	308	0	0	1	0
All	All	7017	0	6097	90	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 7.

All (90) 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:B:339:ILE:HD11	1:B:350:ILE:HG23	1.44	0.98
1:B:151:LYS:HE2	1:B:153:HIS:CE1	2.06	0.91
1:B:169:ILE:HD13	1:B:195:GLU:HB2	1.54	0.88
1:B:350:ILE:HD12	1:B:379:VAL:HG11	1.57	0.85
1:B:19:ASP:OD1	1:B:29:ARG:HD3	1.79	0.83
1:A:211:ASP:OD2	1:A:323[A]:ARG:HD3	1.80	0.79
1:B:93[B]:TRP:CZ3	3:B:1417:XE:XE	3.20	0.72
1:B:350:ILE:HD12	1:B:379:VAL:CG1	2.20	0.70
1:B:350:ILE:CD1	1:B:379:VAL:HG11	2.27	0.64
1:B:169:ILE:CD1	1:B:195:GLU:HB2	2.25	0.64
1:A:222:HIS:O	1:B:371:PRO:HG3	1.97	0.63
1:B:339:ILE:HD12	1:B:351:GLY:O	1.98	0.63
1:B:53:VAL:O	1:B:151:LYS:HA	2.00	0.62
1:B:58:HIS:CD2	1:B:60:HIS:H	2.18	0.61
1:B:212:LEU:HD22	1:B:243:LEU:HD21	1.83	0.60
1:A:212:LEU:HD22	1:A:243:LEU:HD21	1.84	0.60
1:B:57:THR:CG2	1:B:150:VAL:HG23	2.32	0.60
1:A:370:GLN:HG2	1:B:225:PHE:CE1	2.37	0.59
1:A:217:VAL:HG21	1:A:329[A]:VAL:CG1	2.35	0.57
1:B:29:ARG:HH11	1:B:29:ARG:HG2	1.71	0.56
1:A:258:VAL:HG13	1:A:317:LEU:HD23	1.88	0.55
1:A:173:LYS:NZ	1:A:196:ASP:HB3	2.21	0.55
1:B:81:LYS:HE3	3:B:1417:XE:XE	2.85	0.54
1:A:169:ILE:HG13	4:A:2139:HOH:O	2.07	0.54
1:A:58:HIS:HD2	1:A:60:HIS:CG	2.25	0.54
1:B:331:PHE:CZ	3:B:1415:XE:XE	3.39	0.53
1:B:88:ARG:NH2	2:C:2:PDX:H4	2.22	0.53
1:B:103:LYS:HG2	1:B:140:VAL:HG21	1.91	0.53
1:B:246:ARG:CZ	4:B:2209:HOH:O	2.57	0.53
1:B:350:ILE:O	1:B:381:LEU:HD12	2.08	0.53
1:A:173:LYS:HZ1	1:A:196:ASP:HB3	1.73	0.52
1:A:19:ASP:OD1	1:A:29:ARG:HD3	2.09	0.51
1:B:379:VAL:CG1	1:B:380:SER:N	2.74	0.51
1:B:341:ILE:HG23	1:B:348:LEU:HD11	1.92	0.51
1:B:222:HIS:HE1	1:B:407:SER:C	2.14	0.50
1:B:73:VAL:HG12	1:B:150:VAL:HG22	1.93	0.50
1:B:354:ASP:OD1	1:B:355:ALA:N	2.44	0.49
1:B:354:ASP:O	1:B:357:THR:HG22	2.12	0.49
1:B:224:ARG:NH1	1:B:285:TRP:CG	2.81	0.49
1:B:343:LEU:HD22	1:B:390:PHE:CE2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:LYS:HE3	4:A:2097:HOH:O	2.12	0.48
1:A:350:ILE:O	1:A:381:LEU:HD12	2.13	0.48
1:B:354:ASP:OD1	1:B:356:GLU:N	2.42	0.48
1:B:93[A]:TRP:CD1	3:B:1417:XE:XE	3.44	0.48
1:B:40:ASP:HB3	1:B:44:ASP:HB2	1.96	0.48
1:B:151:LYS:HE2	1:B:153:HIS:HE1	1.72	0.48
1:A:46[B]:ASP:OD2	1:A:46[B]:ASP:N	2.46	0.47
1:B:93[B]:TRP:CE3	3:B:1417:XE:XE	3.45	0.47
1:B:331:PHE:HZ	3:B:1415:XE:XE	2.75	0.47
1:B:331:PHE:CE2	3:B:1415:XE:XE	3.45	0.47
1:B:58:HIS:CD2	1:B:60:HIS:HB2	2.50	0.47
1:B:248:PHE:CD1	1:B:292:PRO:HA	2.50	0.47
1:A:115:GLY:HA3	1:A:121:LYS:HE3	1.96	0.47
1:A:370:GLN:HG2	1:B:225:PHE:CZ	2.50	0.46
1:A:339:ILE:HG23	1:A:339:ILE:O	2.16	0.46
1:B:224:ARG:NH2	1:B:281:SER:O	2.48	0.46
1:B:374:VAL:HG13	1:B:377:GLN:CG	2.46	0.46
1:B:57:THR:HG23	1:B:150:VAL:HG23	1.98	0.46
1:A:280:HIS:O	1:A:281:SER:HB3	2.15	0.45
1:B:280:HIS:O	1:B:281:SER:HB3	2.17	0.45
1:B:248:PHE:CE1	1:B:292:PRO:HA	2.52	0.45
1:B:74:LEU:HD23	1:B:149:ASN:HD22	1.82	0.44
1:B:132:ALA:O	1:B:135:ALA:HB2	2.17	0.44
1:A:151:LYS:HB2	3:A:1416:XE:XE	2.96	0.44
1:A:339:ILE:HD11	1:A:350:ILE:CG2	2.48	0.44
1:B:354:ASP:C	1:B:357:THR:HG22	2.38	0.44
1:A:58:HIS:CD2	1:A:60:HIS:CG	3.05	0.44
1:B:74:LEU:HD23	1:B:149:ASN:ND2	2.33	0.43
1:B:341:ILE:HG23	1:B:348:LEU:CD1	2.48	0.43
1:A:132:ALA:O	1:A:135:ALA:HB2	2.18	0.43
1:B:212:LEU:CD2	1:B:243:LEU:HD21	2.48	0.43
1:B:366:PHE:HB3	1:B:372:LYS:HA	2.01	0.43
1:B:317:LEU:HB2	1:B:325:VAL:HB	2.00	0.42
1:A:133:ASP:O	1:A:134:ARG:HB2	2.19	0.42
1:B:365:VAL:HG11	1:B:390:PHE:CE1	2.55	0.42
1:A:217:VAL:HG21	1:A:329[B]:VAL:HG22	2.01	0.42
1:B:121:LYS:HA	1:B:121:LYS:HD3	1.88	0.42
1:B:222:HIS:CE1	1:B:407:SER:O	2.73	0.41
1:B:367:ALA:O	1:B:368:ASP:C	2.59	0.41
1:B:29:ARG:HG2	1:B:29:ARG:NH1	2.34	0.41
1:B:290:ARG:NH1	1:B:290:ARG:HG2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:VAL:HG22	1:B:330:ARG:N	2.35	0.41
1:A:217:VAL:HG21	1:A:329[B]:VAL:CG2	2.50	0.41
1:B:88:ARG:HG2	1:B:88:ARG:O	2.20	0.41
1:B:346:LYS:O	1:B:385:VAL:HA	2.21	0.41
1:A:348:LEU:C	1:A:348:LEU:HD23	2.41	0.40
1:B:210:HIS:CD2	1:B:210:HIS:C	2.94	0.40
1:B:343:LEU:HG	1:B:412:VAL:HG11	2.03	0.40
1:B:290:ARG:HG2	1:B:290:ARG:HH11	1.86	0.40
1:B:222:HIS:CE1	1:B:407:SER:C	2.94	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	404/415 (97%)	391 (97%)	12 (3%)	1 (0%)	47	38
1	B	403/415 (97%)	389 (96%)	13 (3%)	1 (0%)	47	38
All	All	807/830 (97%)	780 (97%)	25 (3%)	2 (0%)	41	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	368	ASP
1	A	337	GLY

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	336/337 (100%)	333 (99%)	3 (1%)	78	77
1	B	334/337 (99%)	332 (99%)	2 (1%)	86	85
All	All	670/674 (99%)	665 (99%)	5 (1%)	81	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ASP
1	A	139	MET
1	A	210	HIS
1	B	206	MET
1	B	210	HIS

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	60	HIS
1	A	82	GLN
1	A	149	ASN
1	B	149	ASN
1	B	153	HIS
1	B	185	ASN
1	B	222	HIS

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	$\# Z > 2$
2	PDX	C	1	2	20,20,20	1.70	5 (25%)	23,31,31	1.61	4 (17%)
2	PDX	C	2	2	19,19,20	1.59	5 (26%)	20,29,31	1.26	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	PDX	C	1	2	-	0/12/32/32	0/1/1/1
2	PDX	C	2	2	-	3/12/29/32	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	PDX	O2S-S	4.16	1.63	1.45
2	C	2	PDX	O2-S	-3.23	1.47	1.57
2	C	1	PDX	C1-C2	2.88	1.55	1.52
2	C	2	PDX	O5-C5	2.61	1.48	1.43
2	C	2	PDX	O1S-S	-2.51	1.34	1.45
2	C	1	PDX	O1S-S	-2.44	1.34	1.45
2	C	2	PDX	O3-S'	-2.27	1.50	1.57
2	C	1	PDX	O3-S'	-2.26	1.50	1.57
2	C	1	PDX	O2-S	-2.23	1.50	1.57
2	C	2	PDX	O2X-S'	-2.09	1.36	1.45

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	PDX	C3-O3-S'	5.67	129.83	118.88
2	C	2	PDX	O2-C2-C3	4.21	111.31	106.65
2	C	1	PDX	O4-C4-C5	2.86	116.41	109.30
2	C	1	PDX	O2-C2-C1	2.31	110.68	107.58
2	C	1	PDX	O3S-S-O2S	-2.02	101.47	108.49

There are no chirality outliers.

All (3) torsion outliers are listed below:

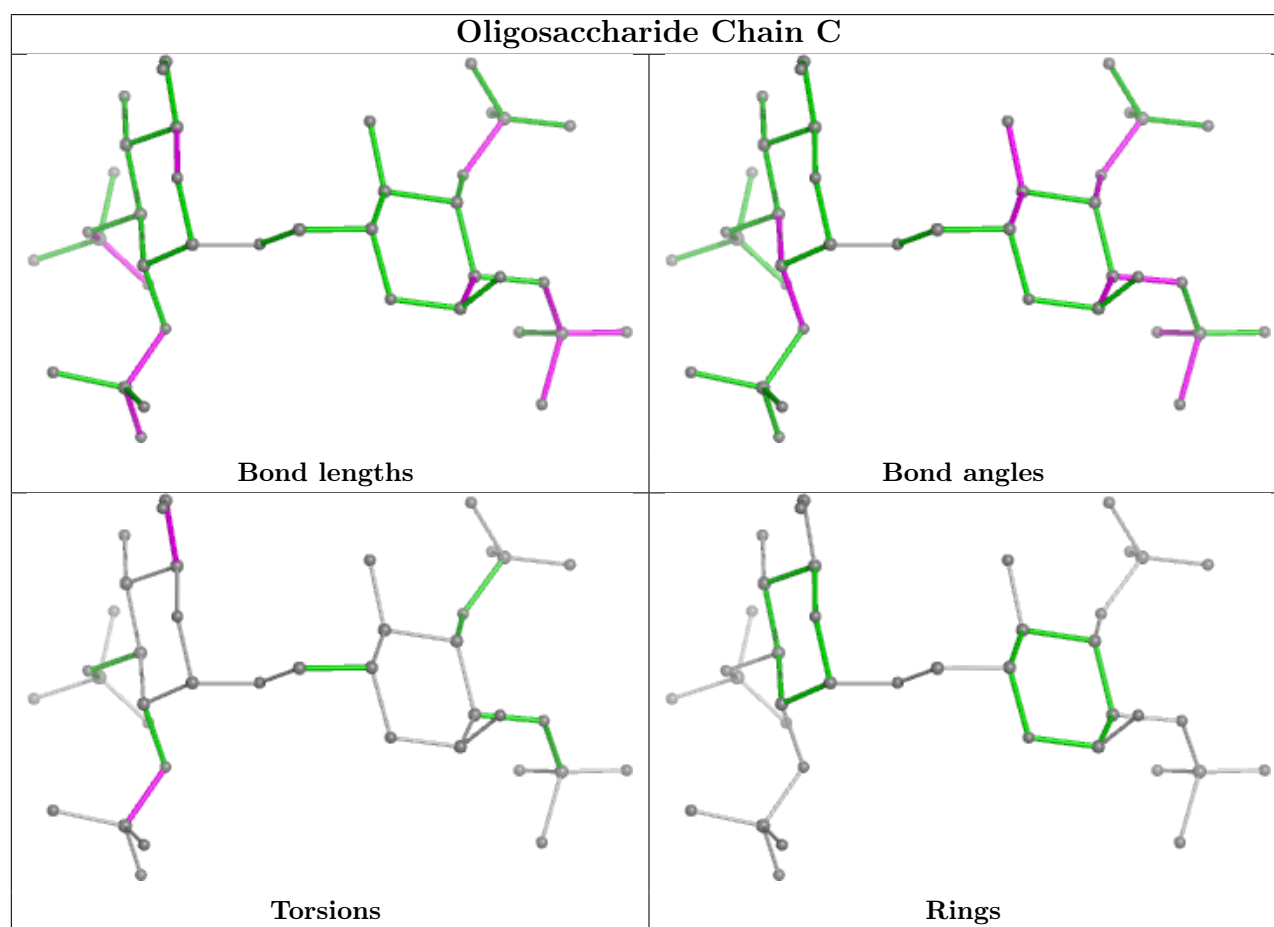
Mol	Chain	Res	Type	Atoms
2	C	2	PDX	O5-C5-C6-O6
2	C	2	PDX	C4-C5-C6-O6
2	C	2	PDX	C2-O2-S-O2S

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	PDX	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 5 ligands modelled in this entry, 5 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	399/415 (96%)	0.49	39 (9%) 7 12	25, 39, 68, 76	0
1	B	402/415 (96%)	0.68	56 (13%) 2 4	24, 38, 76, 89	0
All	All	801/830 (96%)	0.59	95 (11%) 4 7	24, 38, 74, 89	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	381	LEU	7.4
1	B	390	PHE	6.4
1	A	221	TYR	5.6
1	B	391	THR	5.4
1	B	7	GLU	5.3
1	B	382	GLY	5.1
1	B	347	VAL	5.1
1	A	337	GLY	5.1
1	B	385	VAL	5.0
1	B	352	GLY	5.0
1	B	53	VAL	4.8
1	B	343	LEU	4.7
1	B	392	PRO	4.7
1	B	383	GLN	4.7
1	B	412	VAL	4.7
1	A	34	ALA	4.7
1	B	368	ASP	4.7
1	A	255	PHE	4.5
1	A	338	ASN	4.4
1	B	413	VAL	4.3
1	A	58	HIS	4.2
1	A	318	ILE	4.0
1	B	355	ALA	4.0
1	B	359	MET	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	345	GLY	3.9
1	A	258	VAL	3.8
1	A	257	THR	3.8
1	B	255	PHE	3.8
1	B	354	ASP	3.7
1	B	344	LYS	3.7
1	B	357	THR	3.6
1	B	393	GLY	3.5
1	B	93[A]	TRP	3.5
1	A	317	LEU	3.5
1	B	348	LEU	3.4
1	B	358	LEU	3.4
1	B	394	THR	3.4
1	B	386	LEU	3.3
1	A	44	ASP	3.3
1	B	338	ASN	3.3
1	B	258	VAL	3.2
1	B	42	ALA	3.2
1	B	353	ALA	3.2
1	A	252	VAL	3.2
1	A	390	PHE	3.1
1	B	252	VAL	3.1
1	B	318	ILE	3.1
1	A	344	LYS	3.0
1	A	57	THR	3.0
1	B	389	ASP	3.0
1	B	351	GLY	2.9
1	B	69	TYR	2.9
1	B	164	HIS	2.9
1	B	223	SER	2.8
1	A	172	GLU	2.8
1	A	251	TYR	2.7
1	B	366	PHE	2.7
1	B	387	SER	2.7
1	A	82	GLN	2.7
1	B	346	LYS	2.6
1	A	412	VAL	2.6
1	B	396	TYR	2.6
1	A	256	ASN	2.6
1	A	347	VAL	2.6
1	B	350	ILE	2.6
1	B	257	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	339	ILE	2.5
1	A	7	GLU	2.5
1	A	164	HIS	2.5
1	B	8	THR	2.5
1	B	317	LEU	2.5
1	A	294	ILE	2.5
1	B	41	ARG	2.5
1	A	158	VAL	2.4
1	A	336	GLU	2.4
1	A	352	GLY	2.4
1	A	319	ALA	2.4
1	A	335	TYR	2.4
1	A	278	THR	2.4
1	A	316	THR	2.4
1	B	388	ALA	2.3
1	B	126	GLU	2.3
1	A	21	TRP	2.2
1	A	356	GLU	2.2
1	A	334	LYS	2.2
1	A	155	MET	2.1
1	A	315	LEU	2.1
1	A	343	LEU	2.1
1	A	20	TYR	2.0
1	B	371	PRO	2.0
1	A	391	THR	2.0
1	B	256	ASN	2.0
1	B	316	THR	2.0
1	B	222	HIS	2.0
1	B	349	THR	2.0

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

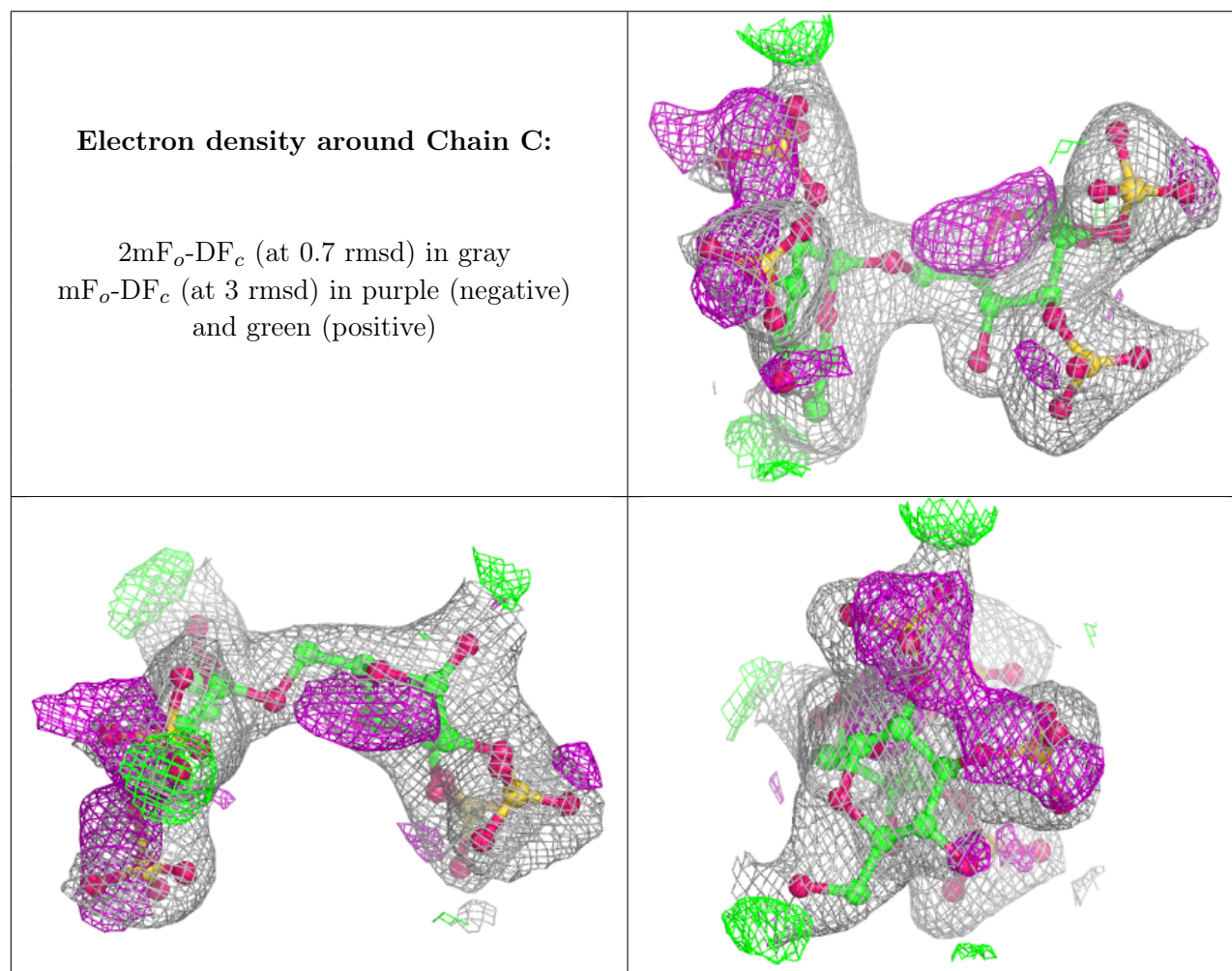
There are no non-standard protein/DNA/RNA residues in this entry.

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(\AA^2)	Q<0.9
2	PDX	C	2	19/20	0.93	0.18	53,56,60,60	0
2	PDX	C	1	20/20	0.94	0.16	52,55,57,58	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 ⓘ

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	XE	A	1416	1/1	0.95	0.06	56,56,56,56	1
3	XE	B	1417	1/1	0.97	0.11	58,58,58,58	1
3	XE	B	1416	1/1	0.98	0.06	45,45,45,45	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	XE	B	1415	1/1	0.98	0.10	34,34,34,34	1
3	XE	A	1415	1/1	1.00	0.08	35,35,35,35	1

6.5 Other polymers [i](#)

There are no such residues in this entry.