



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:00 PM UTC

PDB ID : 1L3T / pdb_00001l3t
Title : Crystal Structure of Bacillus DNA Polymerase I Fragment product complex with 10 base pairs of duplex DNA following addition of a single dTTP residue
Authors : Johnson, S.J.; Taylor, J.S.; Beese, L.S.
Deposited on : 2002-03-01
Resolution : 1.70 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 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

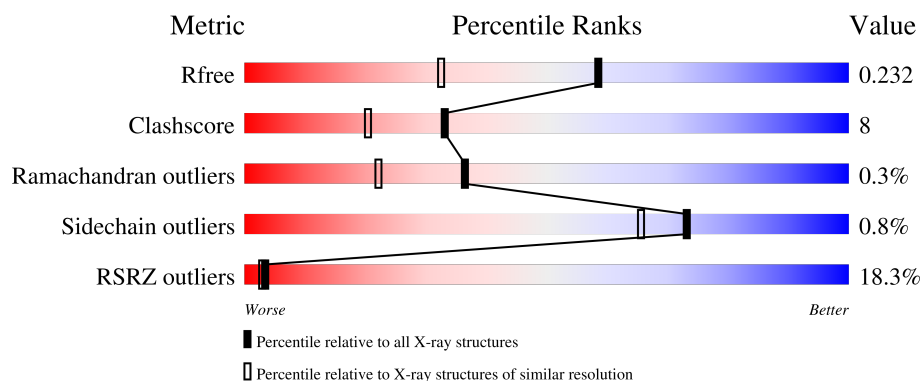
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	B	10	30% 80% 20%
2	C	16	31% 56% 19% 25%
3	A	580	18% 83% 16%
4	D	2	50% 50%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5660 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 DNA chain called 5'-D(*GP*CP*GP*AP*TP*CP*AP*CP*GP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	10	Total	C	N	O	P	0	0	0
			202	97	38	58	9			

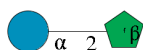
- Molecule 2 is a DNA chain called 5'-D(*GP*AP*CP*G*TP*AP*CP*GP*TP*GP*AP*TP*CP*GP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	P	0	0	0
			243	117	45	70	11			

- Molecule 3 is a protein called DNA Polymerase I.

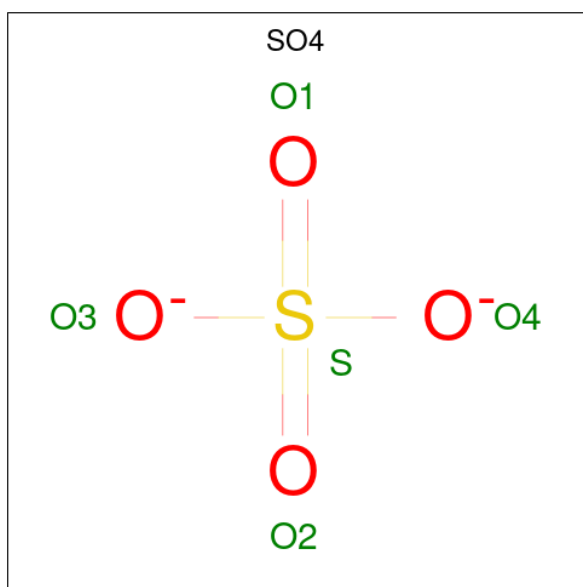
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	580	Total	C	N	O	S	0	0	0
			4650	2956	807	870	17			

- Molecule 4 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

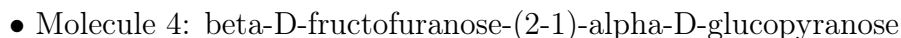
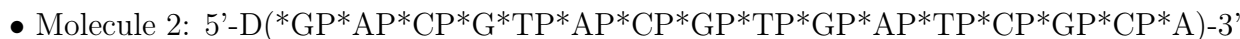
- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	18	Total	O	0	0
			18	18		
7	C	28	Total	O	0	0
			28	28		
7	A	475	Total	O	0	0
			475	475		

- Molecule 1: 5'-D(*GP*CP*GP*AP*TP*CP*AP*CP*GP*T)-3'





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.53Å 93.36Å 106.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.06 – 1.70 35.06 – 1.70	Depositor EDS
% Data completeness (in resolution range)	75.9 (35.06-1.70) 87.3 (35.06-1.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.81 (at 1.70Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.206 , 0.231 0.207 , 0.232	Depositor DCC
R_{free} test set	3636 reflections (4.34%)	wwPDB-VP
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5660	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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: MG, FRU, SO4, GLC

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	B	0.28	0/226	0.62	0/347
2	C	0.28	0/272	0.68	0/418
3	A	0.36	0/4734	0.84	11/6398 (0.2%)
All	All	0.35	0/5232	0.82	11/7163 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	680	ASP	N-CA-C	-7.11	98.91	109.15
3	A	475	VAL	N-CA-C	6.87	117.66	110.72
3	A	317	ALA	N-CA-C	-6.31	98.98	109.46
3	A	321	GLU	N-CA-C	6.28	119.37	109.96
3	A	793	ASN	N-CA-C	5.62	118.60	111.69
3	A	769	ARG	N-CA-C	-5.61	102.59	110.50
3	A	332	ILE	N-CA-C	-5.51	99.94	107.99
3	A	847	LEU	N-CA-C	5.30	116.86	111.14
3	A	457	LEU	N-CA-C	5.25	119.54	113.18
3	A	617	SER	N-CA-C	-5.04	101.72	109.52
3	A	585	SER	N-CA-C	5.01	116.43	110.97

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	B	202	0	114	1	0
2	C	243	0	137	7	0
3	A	4650	0	4698	76	0
4	D	23	0	21	1	0
5	A	20	0	0	0	0
6	A	1	0	0	0	0
7	A	475	0	0	1	0
7	B	18	0	0	0	0
7	C	28	0	0	1	0
All	All	5660	0	4970	79	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 (79) 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)
3:A:727:ILE:HG22	3:A:728:SER:H	1.36	0.91
3:A:729:ARG:H	3:A:729:ARG:HD2	1.43	0.83
3:A:534:LEU:HD11	3:A:574:ILE:HD13	1.62	0.81
3:A:716:ILE:HD11	3:A:721:LEU:HA	1.68	0.75
3:A:408:ASP:HB2	4:D:2:FRU:H11	1.69	0.73
3:A:789:ARG:HA	3:A:792:MET:HE3	1.70	0.71
3:A:813:ALA:O	3:A:817:GLU:HG3	1.91	0.70
3:A:459:ARG:HB3	3:A:460:PRO:HD3	1.76	0.68
3:A:692:VAL:HB	3:A:696:GLU:HB2	1.77	0.66
2:C:3:DT:C6	3:A:716:ILE:HD12	2.32	0.64
3:A:610:LEU:HD23	3:A:610:LEU:C	2.23	0.64
3:A:584:GLN:O	3:A:589:GLU:HG3	1.98	0.63
3:A:727:ILE:HG22	3:A:728:SER:N	2.10	0.61
3:A:716:ILE:HD11	3:A:721:LEU:CA	2.31	0.60
3:A:591:LEU:O	3:A:595:VAL:HG23	2.04	0.58
3:A:722:ALA:HB1	3:A:727:ILE:O	2.04	0.57
3:A:429:TYR:O	3:A:435:ARG:HA	2.07	0.55
3:A:519:TYR:CD1	3:A:525:GLU:HA	2.41	0.55
3:A:708:VAL:O	3:A:712:ILE:HG12	2.07	0.55
2:C:8:DG:H2'	7:C:2742:HOH:O	2.08	0.54
3:A:515:GLU:HG2	3:A:519:TYR:CE2	2.42	0.54
3:A:546:VAL:HG11	3:A:549:LYS:HD3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:722:ALA:O	3:A:726:ASN:N	2.42	0.52
3:A:729:ARG:HD2	3:A:729:ARG:N	2.20	0.52
3:A:728:SER:OG	3:A:729:ARG:HD2	2.10	0.52
3:A:716:ILE:HD13	3:A:721:LEU:HD13	1.92	0.52
3:A:848:VAL:HB	3:A:849:PRO:HD3	1.92	0.51
3:A:558:ALA:O	3:A:562:GLU:HG2	2.10	0.51
3:A:565:ALA:HB3	3:A:566:PRO:HD3	1.93	0.51
3:A:728:SER:HB3	3:A:731:GLU:HG3	1.93	0.51
3:A:629:ARG:NH2	3:A:630:LEU:HD21	2.26	0.50
3:A:692:VAL:HG21	3:A:701:MET:HE1	1.93	0.50
3:A:595:VAL:HG22	3:A:602:VAL:HG13	1.93	0.50
3:A:722:ALA:HB2	3:A:728:SER:O	2.11	0.50
2:C:3:DT:H2'	3:A:714:TYR:HB3	1.93	0.50
3:A:457:LEU:C	3:A:460:PRO:HD2	2.37	0.49
3:A:727:ILE:HG22	3:A:731:GLU:OE1	2.12	0.49
3:A:534:LEU:HD11	3:A:574:ILE:CD1	2.40	0.49
3:A:721:LEU:HB3	3:A:732:ALA:HB1	1.95	0.49
1:B:25:DC:H2''	1:B:26:DA:H5'	1.94	0.49
3:A:355:PRO:HG2	3:A:356:GLN:OE1	2.13	0.48
2:C:3:DT:H6	3:A:716:ILE:HD12	1.77	0.48
3:A:536:VAL:O	3:A:540:GLU:HB2	2.14	0.47
3:A:814:ARG:HD3	3:A:847:LEU:HD11	1.96	0.47
3:A:728:SER:HB3	3:A:731:GLU:CG	2.46	0.46
3:A:354:ASP:OD2	3:A:356:GLN:HB2	2.15	0.46
3:A:583:LEU:HG	3:A:636:ILE:HD11	1.98	0.45
3:A:308:THR:OG1	3:A:311:MET:HE3	2.16	0.45
3:A:349:GLU:H	3:A:349:GLU:CD	2.25	0.45
3:A:729:ARG:HG2	3:A:729:ARG:HH11	1.82	0.45
2:C:3:DT:C2'	3:A:714:TYR:HB3	2.46	0.45
3:A:729:ARG:H	3:A:729:ARG:CD	2.11	0.45
3:A:514:VAL:O	3:A:518:ILE:HG13	2.17	0.44
3:A:595:VAL:HG22	3:A:602:VAL:CG1	2.47	0.44
3:A:727:ILE:CG2	3:A:728:SER:H	2.18	0.44
3:A:593:LYS:HG3	7:A:2690:HOH:O	2.16	0.44
3:A:828:VAL:HB	3:A:831:GLU:CG	2.48	0.43
3:A:524:GLN:HG2	3:A:525:GLU:N	2.33	0.43
3:A:825:LEU:HD11	3:A:835:GLU:HB3	2.00	0.43
3:A:711:GLY:CA	3:A:716:ILE:HB	2.49	0.43
3:A:587:TYR:CE2	3:A:627:PRO:HD3	2.54	0.43
3:A:690:PHE:HB3	3:A:701:MET:HE3	2.01	0.43
3:A:677:ARG:HB2	3:A:679:LEU:HG	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:690:PHE:CB	3:A:701:MET:HE3	2.49	0.42
2:C:3:DT:O2	2:C:3:DT:H3'	2.20	0.42
3:A:549:LYS:HE2	3:A:553:GLY:O	2.18	0.42
3:A:822:ALA:CB	3:A:836:ALA:HB2	2.50	0.42
3:A:579:GLN:NE2	3:A:630:LEU:HD13	2.35	0.42
3:A:728:SER:OG	3:A:729:ARG:N	2.53	0.42
3:A:512:GLY:O	3:A:516:GLN:HG2	2.19	0.41
3:A:561:LEU:HB2	3:A:575:LEU:HD21	2.02	0.41
3:A:728:SER:HB3	3:A:731:GLU:OE1	2.20	0.41
3:A:524:GLN:CG	3:A:525:GLU:N	2.82	0.41
3:A:497:THR:O	3:A:501:GLU:HG3	2.20	0.41
3:A:437:VAL:HA	3:A:438:PRO:HD3	1.90	0.41
2:C:11:DC:H5''	3:A:527:ASN:OD1	2.21	0.41
3:A:329:ASP:H	3:A:382:TRP:CD1	2.39	0.40
3:A:740:PHE:HB3	3:A:747:LYS:HD2	2.04	0.40
3:A:704:GLN:O	3:A:708:VAL:HG23	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	
3	A	578/580 (100%)	558 (96%)	18 (3%)	2 (0%)	36	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	524	GLN
3	A	628	ILE

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	
3	A	495/496 (100%)	491 (99%)	4 (1%)	73	65

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	349	GLU
3	A	583	LEU
3	A	729	ARG
3	A	844	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	502	GLN
3	A	510	GLN
3	A	529	ASN
3	A	638	GLN
3	A	656	GLN
3	A	755	GLN
3	A	768	HIS
3	A	867	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 ⓘ

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
4	GLC	D	1	4	11,11,12	3.21	4 (36%)	15,15,17	1.47	2 (13%)
4	FRU	D	2	4	11,12,12	1.41	1 (9%)	10,18,18	0.78	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
4	GLC	D	1	4	-	0/2/19/22	0/1/1/1
4	FRU	D	2	4	-	0/5/24/24	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	GLC	C2-C3	9.34	1.66	1.52
4	D	2	FRU	O2-C2	3.92	1.47	1.40
4	D	1	GLC	O5-C5	2.76	1.48	1.43
4	D	1	GLC	O5-C1	2.72	1.48	1.43
4	D	1	GLC	C4-C5	2.46	1.58	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1	GLC	C1-O5-C5	4.12	117.71	112.19
4	D	1	GLC	C1-C2-C3	-2.50	106.00	109.64

There are no chirality outliers.

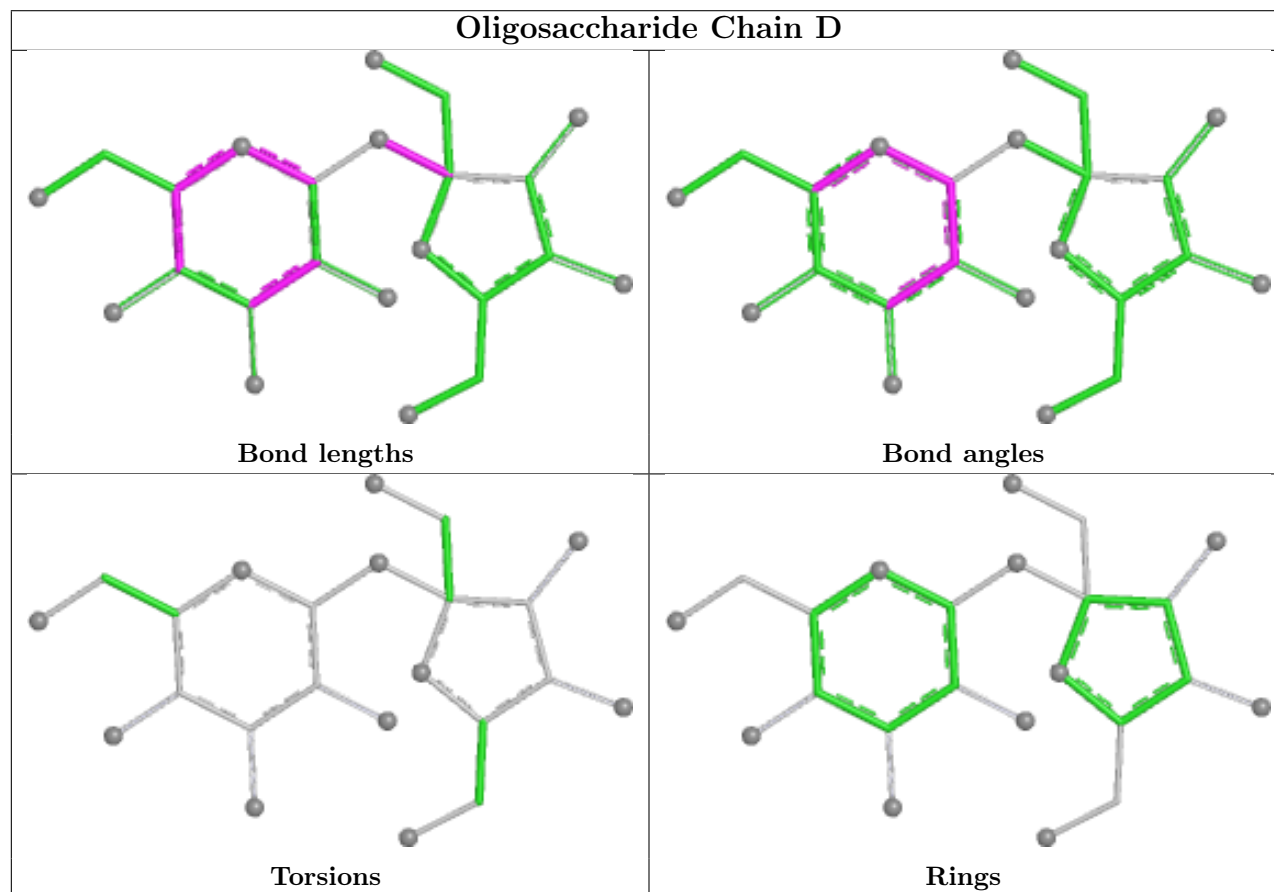
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	2	FRU	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, 1 is monoatomic - leaving 4 for Mogul analysis.

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	$\# Z > 2$
5	SO4	A	913	-	4,4,4	0.40	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	SO4	A	911	-	4,4,4	0.39	0	6,6,6	0.09	0
5	SO4	A	910	-	4,4,4	0.38	0	6,6,6	0.08	0
5	SO4	A	912	-	4,4,4	0.36	0	6,6,6	0.13	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.

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	B	10/10 (100%)	1.15	3 (30%)  	18, 32, 66, 68	0
2	C	12/16 (75%)	1.23	5 (41%)  	16, 33, 64, 70	0
3	A	580/580 (100%)	0.72	102 (17%)  	11, 22, 51, 57	0
All	All	602/606 (99%)	0.73	110 (18%)  	11, 22, 52, 70	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	725	LEU	6.0
3	A	719	TYR	5.9
3	A	721	LEU	5.7
3	A	727	ILE	5.5
3	A	710	PHE	5.4
3	A	297	ALA	5.2
3	A	433	ALA	5.2
3	A	549	LYS	5.1
3	A	716	ILE	5.0
3	A	726	ASN	5.0
3	A	550	THR	5.0
3	A	552	THR	4.9
3	A	729	ARG	4.9
3	A	554	TYR	4.8
3	A	720	GLY	4.8
3	A	551	LYS	4.7
3	A	518	ILE	4.6
3	A	728	SER	4.6
3	A	521	LEU	4.4
3	A	553	GLY	4.4
3	A	722	ALA	4.3
3	A	723	GLN	4.3
3	A	536	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
3	A	434	LYS	4.3
1	B	20	DG	4.2
3	A	531	PRO	4.2
3	A	528	ILE	4.1
3	A	556	THR	4.1
2	C	14	DA	4.0
3	A	511	LEU	4.0
3	A	532	LYS	3.9
3	A	732	ALA	3.9
3	A	539	PHE	3.9
3	A	514	VAL	3.7
3	A	546	VAL	3.7
3	A	431	LYS	3.6
3	A	520	GLU	3.5
3	A	538	LEU	3.5
3	A	735	PHE	3.5
2	C	13	DC	3.4
3	A	559	ASP	3.3
3	A	516	GLN	3.3
3	A	522	ALA	3.3
3	A	523	GLY	3.3
1	B	21	DC	3.2
2	C	3	DT	3.2
3	A	435	ARG	3.2
3	A	537	ILE	3.1
3	A	690	PHE	3.1
3	A	717	SER	3.1
3	A	526	PHE	3.1
1	B	22	DG	3.0
3	A	432	GLY	3.0
3	A	508	ALA	3.0
3	A	509	GLU	3.0
3	A	548	LYS	3.0
2	C	12	DG	3.0
3	A	562	GLU	2.9
3	A	730	LYS	2.9
3	A	534	LEU	2.9
3	A	733	ALA	2.9
3	A	519	TYR	2.9
3	A	734	GLU	2.8
3	A	786	PHE	2.8
3	A	542	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
3	A	781	PHE	2.8
3	A	530	SER	2.8
3	A	558	ALA	2.7
3	A	564	LEU	2.7
3	A	689	ILE	2.7
3	A	691	GLN	2.7
3	A	513	THR	2.7
3	A	571	VAL	2.6
3	A	517	ARG	2.6
3	A	541	LYS	2.6
3	A	718	ASP	2.6
3	A	545	PRO	2.6
3	A	731	GLU	2.5
3	A	561	LEU	2.5
3	A	567	TYR	2.5
3	A	529	ASN	2.5
3	A	724	ASN	2.5
3	A	524	GLN	2.5
3	A	507	LEU	2.5
3	A	715	GLY	2.5
3	A	695	ASP	2.4
3	A	527	ASN	2.4
3	A	543	GLN	2.3
3	A	547	LEU	2.3
3	A	737	GLU	2.3
3	A	779	ARG	2.2
3	A	305	ASP	2.2
3	A	298	LYS	2.2
3	A	692	VAL	2.2
3	A	502	GLN	2.2
3	A	579	GLN	2.2
3	A	505	LYS	2.1
3	A	308	THR	2.1
3	A	544	LEU	2.1
3	A	533	GLN	2.1
3	A	574	ILE	2.1
3	A	515	GLU	2.1
3	A	557	SER	2.1
3	A	306	ARG	2.1
3	A	738	ARG	2.1
3	A	499	ARG	2.0
3	A	540	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
3	A	782	ASN	2.0
2	C	11	DC	2.0
3	A	555	SER	2.0

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

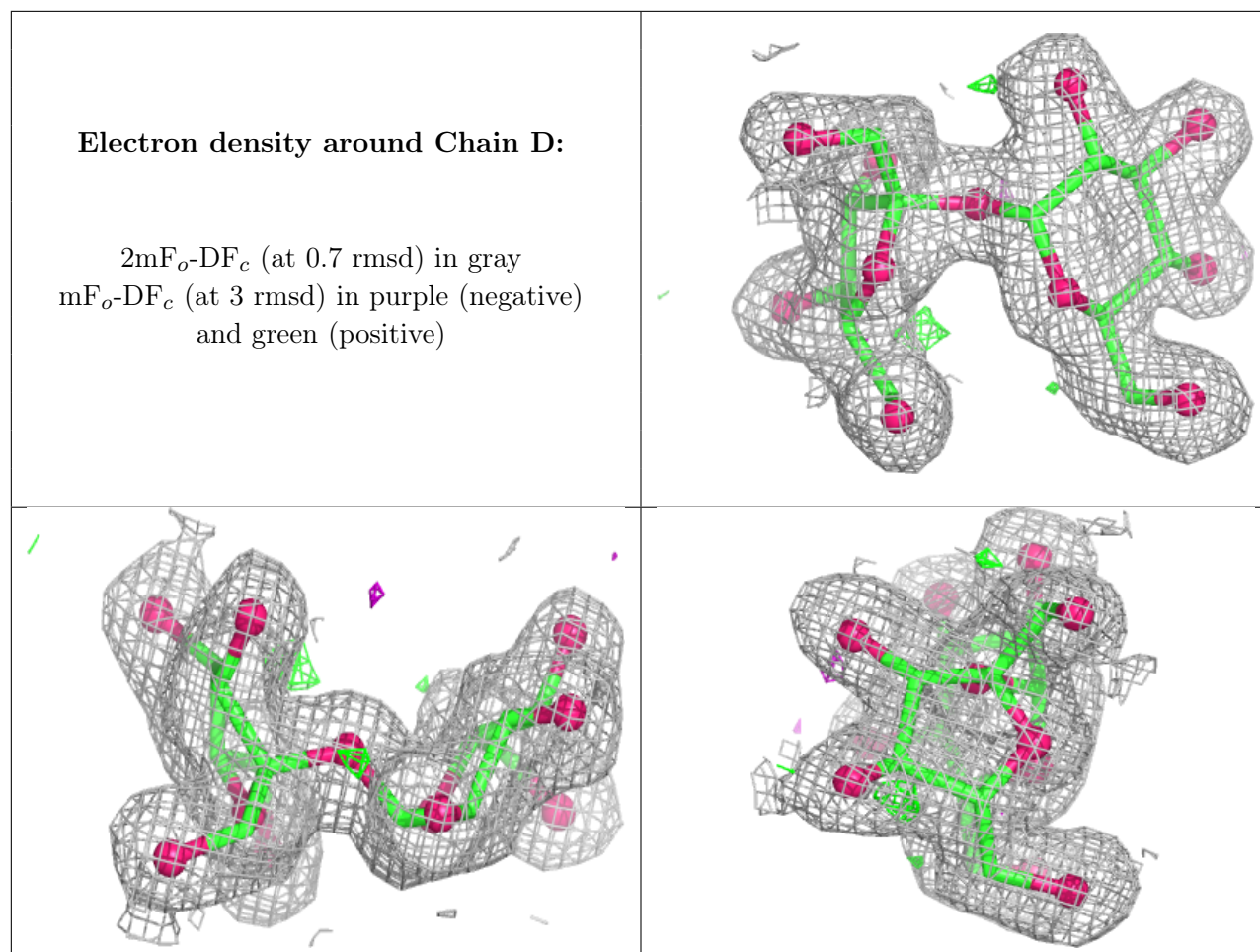
There are no non-standard protein/DNA/RNA residues in this entry.

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
4	FRU	D	2	12/12	0.90	0.09	23,26,26,27	0
4	GLC	D	1	11/12	0.92	0.09	25,27,28,29	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($\text{\AA}^2$)	Q<0.9
5	SO4	A	913	5/5	0.30	0.23	83,83,83,83	0
5	SO4	A	911	5/5	0.77	0.13	66,66,66,66	0
5	SO4	A	912	5/5	0.87	0.14	48,48,48,48	0
5	SO4	A	910	5/5	0.87	0.10	61,61,61,62	0
6	MG	A	920	1/1	0.98	0.13	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.