



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 03:14 PM UTC

PDB ID : 1UA0 / pdb_00001ua0
Title : Aminofluorene DNA adduct at the pre-insertion site of a DNA polymerase
Authors : Hsu, G.W.; Kiefer, J.R.; Becherel, O.J.; Fuchs, R.P.P.; Beese, L.S.
Deposited on : 2004-08-11
Resolution : 2.10 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 ⓘ 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

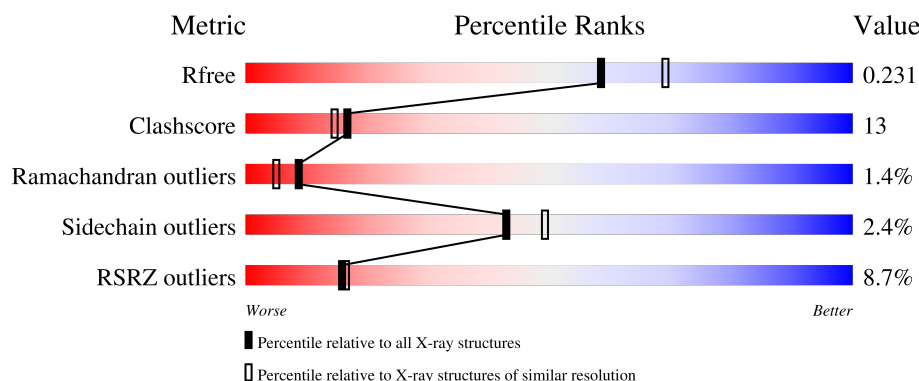
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 $\leq 5\%$. 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	
2	C	14	
3	A	580	
4	D	2	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5438 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 DNA primer strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	9	Total	C	N	O	P	0	0	0
			190	90	39	53	8			

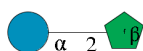
- Molecule 2 is a DNA chain called DNA template strand with aminofluorene adduct.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	P	0	0	0
			235	114	39	71	11			

- Molecule 3 is a protein called DNA polymerase I.

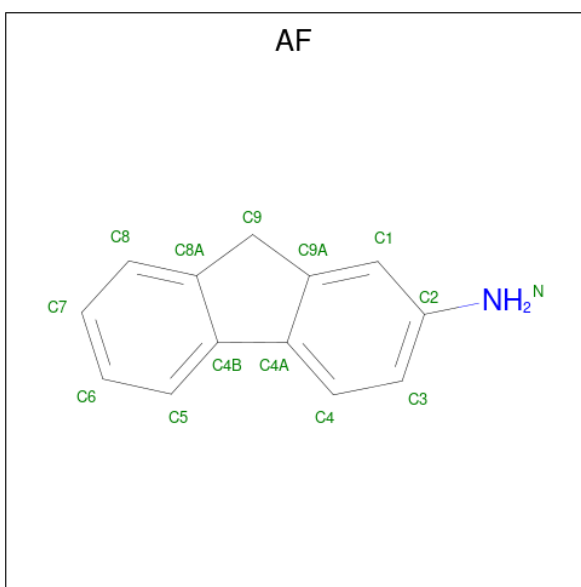
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	580	Total	C	N	O	S	0	0	0
			4650	2956	807	870	17			

- 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 2-AMINOFLUORENE (CCD ID: AF) (formula: C₁₃H₁₁N).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	N	0	0
			14	13	1		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

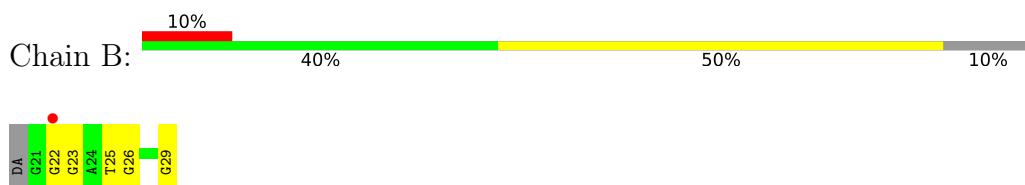
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	15	Total 15	O 15	0	0
7	C	23	Total 23	O 23	0	0
7	A	278	Total 278	O 278	0	0

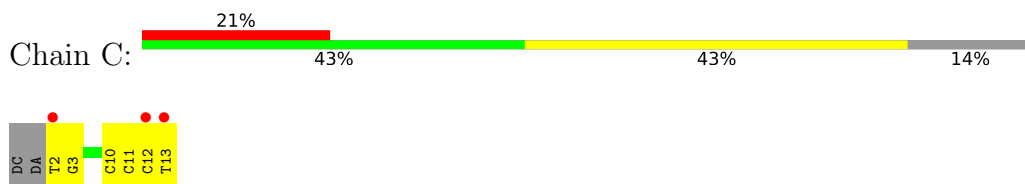
3 Residue-property plots

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.

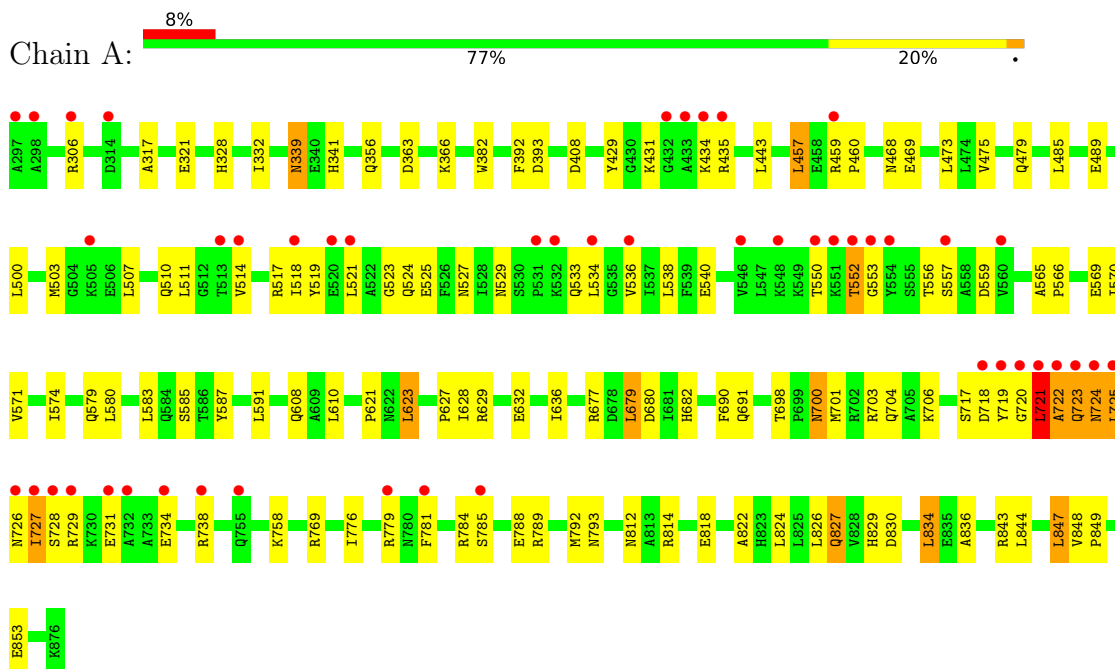
- Molecule 1: DNA primer strand



- Molecule 2: DNA template strand with aminofluorene adduct



- Molecule 3: DNA polymerase I



- Molecule 4: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain D:



GLC1
FR02

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.87Å 93.47Å 104.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.17 – 2.10 37.17 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.5 (37.17-2.10) 93.5 (37.17-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.14 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.210 , 0.242 0.200 , 0.231	Depositor DCC
R_{free} test set	2414 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5438	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% 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: GLC, AF, FRU, SO4

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	0.27	0/214	0.68	0/331
2	C	0.31	0/261	0.75	0/399
3	A	0.38	0/4733	0.87	11/6395 (0.2%)
All	All	0.37	0/5208	0.85	11/7125 (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	-6.62	99.61	109.15
3	A	321	GLU	N-CA-C	6.55	119.22	109.59
3	A	475	VAL	N-CA-C	5.72	116.50	110.72
3	A	317	ALA	N-CA-C	-5.52	100.29	109.46
3	A	585	SER	N-CA-C	5.52	116.98	111.07
3	A	793	ASN	N-CA-C	5.49	118.44	111.69
3	A	769	ARG	N-CA-C	-5.42	102.46	110.48
3	A	332	ILE	N-CA-C	-5.38	99.49	107.51
3	A	457	LEU	N-CA-C	5.21	119.49	113.18
3	A	538	LEU	N-CA-C	5.13	116.72	111.03
3	A	847	LEU	N-CA-C	5.12	116.67	111.14

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	190	0	103	6	0
2	C	235	0	137	7	0
3	A	4650	0	4700	117	0
4	D	23	0	21	1	0
5	C	14	0	10	2	0
6	A	10	0	0	0	0
7	A	278	0	0	2	0
7	B	15	0	0	0	0
7	C	23	0	0	0	0
All	All	5438	0	4971	129	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 13.

All (129) 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:718:ASP:HB3	3:A:719:TYR:HE2	1.35	0.91
3:A:719:TYR:CD1	3:A:785:SER:HB2	2.14	0.81
3:A:734:GLU:O	3:A:738:ARG:HG2	1.84	0.77
3:A:727:ILE:HG12	3:A:731:GLU:CD	2.13	0.73
3:A:328:HIS:HD2	3:A:382:TRP:HE1	1.37	0.72
3:A:724:ASN:C	3:A:725:LEU:HD23	2.16	0.70
3:A:719:TYR:HE1	3:A:781:PHE:O	1.75	0.69
3:A:700:ASN:HD22	3:A:703:ARG:HH11	1.39	0.69
3:A:719:TYR:CE1	3:A:785:SER:HB2	2.28	0.69
3:A:459:ARG:HB3	3:A:460:PRO:HD3	1.76	0.68
3:A:621:PRO:HG2	3:A:623:LEU:HD13	1.76	0.67
3:A:431:LYS:H	3:A:434:LYS:NZ	1.92	0.67
3:A:408:ASP:HB2	4:D:2:FRU:H11	1.76	0.67
3:A:339:ASN:C	3:A:339:ASN:HD22	2.03	0.66
3:A:723:GLN:C	3:A:725:LEU:H	2.02	0.66
3:A:717:SER:HG	3:A:719:TYR:N	1.95	0.65
3:A:629:ARG:HH11	3:A:629:ARG:HG2	1.61	0.65
3:A:722:ALA:O	3:A:724:ASN:N	2.30	0.65
3:A:727:ILE:HG12	3:A:731:GLU:OE2	1.97	0.65
3:A:392:PHE:HA	3:A:479:GLN:HE22	1.62	0.64
1:B:22:DG:H2'	1:B:23:DG:C8	2.34	0.63
3:A:519:TYR:CD1	3:A:525:GLU:HA	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:847:LEU:C	3:A:847:LEU:HD23	2.25	0.62
3:A:721:LEU:O	3:A:722:ALA:HB2	1.99	0.60
3:A:677:ARG:HB2	3:A:679:LEU:HD13	1.83	0.60
3:A:718:ASP:HB3	3:A:719:TYR:CE2	2.25	0.60
3:A:721:LEU:O	3:A:722:ALA:CB	2.49	0.60
3:A:827:GLN:HE21	3:A:827:GLN:C	2.10	0.59
1:B:22:DG:H2''	1:B:23:DG:H5'	1.84	0.59
3:A:356:GLN:NE2	3:A:356:GLN:H	2.00	0.59
3:A:393:ASP:H	3:A:479:GLN:NE2	2.02	0.58
3:A:719:TYR:CE1	3:A:781:PHE:O	2.54	0.58
3:A:700:ASN:ND2	3:A:703:ARG:HH11	2.03	0.57
3:A:727:ILE:HG21	3:A:731:GLU:OE1	2.04	0.57
3:A:827:GLN:NE2	3:A:829:HIS:H	2.03	0.57
3:A:727:ILE:HG21	3:A:731:GLU:CD	2.30	0.56
3:A:677:ARG:CB	3:A:679:LEU:HD13	2.36	0.56
3:A:485:LEU:O	3:A:489:GLU:HG3	2.06	0.55
3:A:550:THR:HG22	3:A:553:GLY:O	2.06	0.55
3:A:534:LEU:HD12	3:A:556:THR:HG21	1.87	0.55
3:A:727:ILE:HG12	3:A:731:GLU:CG	2.36	0.55
3:A:550:THR:CG2	3:A:553:GLY:H	2.20	0.55
3:A:701:MET:HE1	7:A:2954:HOH:O	2.07	0.54
3:A:758:LYS:HA	3:A:776:ILE:HG21	1.90	0.53
3:A:570:ILE:O	3:A:574:ILE:HG12	2.09	0.53
3:A:623:LEU:HD23	3:A:826:LEU:HD21	1.91	0.52
3:A:719:TYR:N	3:A:719:TYR:CD2	2.76	0.52
3:A:720:GLY:C	3:A:721:LEU:HG	2.35	0.52
3:A:822:ALA:CB	3:A:836:ALA:HB2	2.39	0.52
3:A:431:LYS:N	3:A:434:LYS:NZ	2.58	0.52
3:A:510:GLN:O	3:A:514:VAL:HG23	2.09	0.52
3:A:536:VAL:O	3:A:540:GLU:HB2	2.11	0.51
3:A:507:LEU:O	3:A:511:LEU:HB2	2.10	0.51
3:A:363:ASP:HB3	3:A:366:LYS:HG2	1.93	0.51
3:A:550:THR:HG23	3:A:553:GLY:H	1.75	0.50
3:A:698:THR:H	3:A:701:MET:HE3	1.77	0.49
3:A:718:ASP:C	3:A:719:TYR:CD2	2.91	0.49
3:A:518:ILE:HD13	3:A:574:ILE:HD13	1.95	0.49
3:A:726:ASN:O	3:A:728:SER:N	2.40	0.49
3:A:557:SER:HB2	3:A:559:ASP:OD1	2.12	0.49
3:A:565:ALA:HA	3:A:571:VAL:HB	1.94	0.49
2:C:12:DC:H2''	2:C:13:DT:C7	2.43	0.49
3:A:527:ASN:H	3:A:533:GLN:NE2	2.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:723:GLN:C	3:A:725:LEU:N	2.71	0.49
3:A:848:VAL:HB	3:A:849:PRO:HD3	1.95	0.48
1:B:22:DG:H2''	1:B:23:DG:C5'	2.42	0.48
3:A:691:GLN:NE2	3:A:738:ARG:NH2	2.62	0.48
3:A:726:ASN:O	3:A:727:ILE:HG22	2.14	0.48
3:A:431:LYS:H	3:A:434:LYS:HZ2	1.58	0.48
3:A:727:ILE:HG23	3:A:728:SER:N	2.28	0.48
2:C:3:DG:N7	5:C:333:AF:H1	2.29	0.47
3:A:587:TYR:CE1	3:A:627:PRO:HD3	2.49	0.47
3:A:784:ARG:O	3:A:788:GLU:HG3	2.14	0.47
3:A:700:ASN:HD22	3:A:703:ARG:NH1	2.10	0.47
3:A:690:PHE:CD2	3:A:701:MET:HG2	2.49	0.47
1:B:25:DT:H2''	1:B:26:DG:C5'	2.44	0.47
3:A:719:TYR:CD1	3:A:785:SER:CB	2.93	0.47
3:A:565:ALA:HB3	3:A:566:PRO:HD3	1.97	0.47
3:A:723:GLN:O	3:A:725:LEU:N	2.47	0.47
3:A:718:ASP:C	3:A:719:TYR:HD2	2.23	0.47
3:A:610:LEU:HD23	3:A:610:LEU:C	2.40	0.46
3:A:429:TYR:O	3:A:435:ARG:HA	2.15	0.46
3:A:431:LYS:H	3:A:434:LYS:HZ1	1.60	0.46
3:A:511:LEU:HD13	3:A:580:LEU:HB2	1.98	0.46
3:A:523:GLY:O	3:A:524:GLN:HB3	2.16	0.46
3:A:824:LEU:HD23	3:A:834:LEU:HD22	1.96	0.46
3:A:468:ASN:O	3:A:469:GLU:HB2	2.15	0.46
3:A:431:LYS:HG2	3:A:434:LYS:HZ1	1.80	0.46
3:A:629:ARG:HH11	3:A:629:ARG:CG	2.29	0.46
2:C:12:DC:H2''	2:C:13:DT:C5	2.51	0.45
3:A:457:LEU:C	3:A:460:PRO:HD2	2.41	0.45
1:B:25:DT:H2''	1:B:26:DG:H5'	1.98	0.45
3:A:339:ASN:ND2	3:A:341:HIS:H	2.15	0.45
2:C:2:DT:H73	2:C:3:DG:O6	2.17	0.45
3:A:339:ASN:C	3:A:339:ASN:ND2	2.73	0.45
3:A:328:HIS:CD2	3:A:382:TRP:HE1	2.25	0.45
3:A:550:THR:C	3:A:552:THR:H	2.25	0.45
3:A:719:TYR:HE1	3:A:781:PHE:C	2.24	0.45
3:A:500:LEU:HD21	3:A:591:LEU:HD23	2.00	0.44
3:A:503:MET:HE1	3:A:636:ILE:HD13	1.98	0.44
3:A:527:ASN:OD1	3:A:529:ASN:HB2	2.17	0.43
2:C:12:DC:H2''	2:C:13:DT:H71	2.00	0.43
5:C:333:AF:H8	3:A:704:GLN:NE2	2.34	0.43
3:A:431:LYS:HG2	3:A:434:LYS:NZ	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:727:ILE:O	3:A:728:SER:C	2.62	0.42
3:A:534:LEU:C	3:A:534:LEU:HD13	2.44	0.42
3:A:843:ARG:HD2	7:A:2503:HOH:O	2.19	0.42
3:A:608:GLN:O	3:A:608:GLN:HG2	2.19	0.42
3:A:718:ASP:CB	3:A:719:TYR:HE2	2.18	0.42
3:A:849:PRO:O	3:A:853:GLU:HG3	2.20	0.41
3:A:719:TYR:CE1	3:A:781:PHE:C	2.98	0.41
3:A:579:GLN:NE2	3:A:632:GLU:OE1	2.53	0.41
2:C:2:DT:H2'	2:C:3:DG:C4	2.55	0.41
3:A:363:ASP:HB3	3:A:366:LYS:CG	2.51	0.41
3:A:517:ARG:NH2	3:A:521:LEU:HD21	2.36	0.41
3:A:690:PHE:CE2	3:A:701:MET:HG2	2.55	0.41
3:A:789:ARG:HA	3:A:792:MET:HE3	2.01	0.41
3:A:517:ARG:HH21	3:A:569:GLU:CD	2.29	0.41
3:A:779:ARG:N	3:A:779:ARG:HD3	2.36	0.41
3:A:306:ARG:HH11	3:A:306:ARG:HG2	1.86	0.41
3:A:500:LEU:HD12	3:A:500:LEU:HA	1.91	0.41
3:A:720:GLY:O	3:A:721:LEU:O	2.39	0.41
3:A:814:ARG:NH2	3:A:818:GLU:OE2	2.53	0.41
3:A:822:ALA:HB2	3:A:836:ALA:HB2	2.02	0.41
2:C:10:DC:H2''	2:C:11:DC:O5'	2.20	0.40
3:A:728:SER:O	3:A:729:ARG:C	2.64	0.40
1:B:29:DG:O3'	3:A:830:ASP:OD1	2.37	0.40
3:A:682:HIS:CE1	3:A:706:LYS:HG3	2.56	0.40
3:A:721:LEU:HD23	3:A:721:LEU:N	2.35	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	576/580 (99%)	547 (95%)	21 (4%)	8 (1%)	9 5

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	721	LEU
3	A	722	ALA
3	A	723	GLN
3	A	725	LEU
3	A	727	ILE
3	A	724	ASN
3	A	552	THR
3	A	628	ILE

5.3.2 Protein sidechains ⓘ

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	495/495 (100%)	483 (98%)	12 (2%)	43	49

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	339	ASN
3	A	443	LEU
3	A	473	LEU
3	A	583	LEU
3	A	623	LEU
3	A	679	LEU
3	A	700	ASN
3	A	721	LEU
3	A	812	ASN
3	A	827	GLN
3	A	834	LEU
3	A	844	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	328	HIS
3	A	339	ASN
3	A	356	GLN
3	A	405	GLN
3	A	479	GLN
3	A	502	GLN
3	A	533	GLN
3	A	576	HIS
3	A	579	GLN
3	A	624	GLN
3	A	656	GLN
3	A	691	GLN
3	A	700	ASN
3	A	724	ASN
3	A	726	ASN
3	A	759	GLN
3	A	782	ASN
3	A	827	GLN
3	A	854	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	3.24	4 (36%)	15,15,17	1.47	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FRU	D	2	4	11,12,12	1.49	1 (9%)	10,18,18	0.83	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	9.48	1.67	1.52
4	D	2	FRU	O2-C2	4.16	1.47	1.40
4	D	1	GLC	O5-C5	2.90	1.49	1.43
4	D	1	GLC	O5-C1	2.80	1.48	1.43
4	D	1	GLC	C4-C5	2.09	1.57	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	4.30	117.95	112.19
4	D	1	GLC	C1-C2-C3	-2.34	106.24	109.64

There are no chirality outliers.

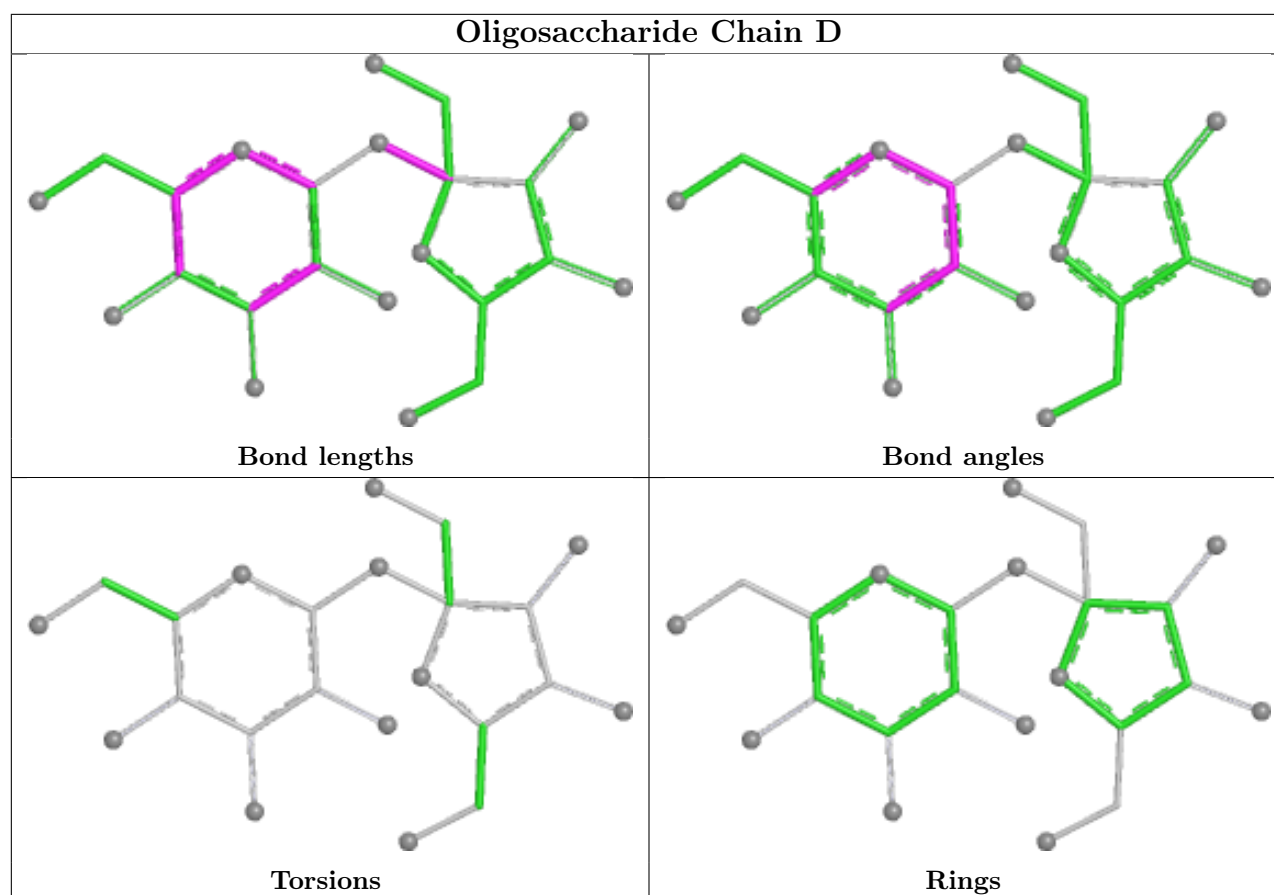
There are no torsion outliers.

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](#)

3 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	$\# Z > 2$
5	AF	C	333	2	16,16,16	2.61	12 (75%)	23,23,23	1.13	3 (13%)
6	SO4	A	903	-	4,4,4	0.38	0	6,6,6	0.09	0
6	SO4	A	902	-	4,4,4	0.36	0	6,6,6	0.10	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
5	AF	C	333	2	-	-	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	333	AF	C1-C9A	4.01	1.45	1.39
5	C	333	AF	C5-C4B	3.74	1.45	1.40
5	C	333	AF	C4B-C4A	-3.16	1.38	1.46
5	C	333	AF	C8-C8A	3.14	1.44	1.39
5	C	333	AF	C1-C2	2.98	1.44	1.39
5	C	333	AF	C2-N	2.97	1.48	1.38
5	C	333	AF	C4A-C9A	2.64	1.46	1.40
5	C	333	AF	C4-C4A	2.42	1.43	1.40
5	C	333	AF	C7-C8	2.35	1.42	1.38
5	C	333	AF	C4B-C8A	2.34	1.45	1.40
5	C	333	AF	C6-C5	2.17	1.42	1.38
5	C	333	AF	C7-C6	2.13	1.42	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	333	AF	C3-C2-N	-2.68	115.96	120.90
5	C	333	AF	C1-C2-N	2.57	124.89	120.56
5	C	333	AF	C2-C1-C9A	-2.09	119.01	121.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	333	AF	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	718:ASP	C	719:TYR	N	3.80

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	9/10 (90%)	0.27	1 (11%)	10 10	34, 40, 73, 78	0
2	C	12/14 (85%)	0.49	3 (25%)	2 2	29, 38, 73, 84	0
3	A	580/580 (100%)	0.49	48 (8%)	17 18	25, 35, 62, 94	0
All	All	601/604 (99%)	0.49	52 (8%)	16 16	25, 35, 62, 94	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	721	LEU	8.9
3	A	719	TYR	8.8
3	A	725	LEU	7.7
3	A	727	ILE	6.9
3	A	718	ASP	5.8
3	A	726	ASN	5.0
3	A	720	GLY	4.9
3	A	724	ASN	4.8
3	A	550	THR	4.5
3	A	554	TYR	4.4
3	A	781	PHE	4.4
3	A	551	LYS	4.3
3	A	297	ALA	4.2
3	A	722	ALA	3.8
2	C	13	DT	3.6
3	A	728	SER	3.4
3	A	531	PRO	3.4
3	A	738	ARG	3.3
3	A	433	ALA	3.3
3	A	723	GLN	3.3
3	A	552	THR	3.3
3	A	729	ARG	3.2
3	A	532	LYS	3.1

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Mol	Chain	Res	Type	RSRZ
3	A	314	ASP	2.9
3	A	732	ALA	2.9
3	A	518	ILE	2.8
3	A	505	LYS	2.8
3	A	534	LEU	2.8
3	A	731	GLU	2.7
2	C	2	DT	2.7
3	A	298	ALA	2.5
3	A	553	GLY	2.5
3	A	557	SER	2.5
3	A	435	ARG	2.5
3	A	521	LEU	2.4
3	A	546	VAL	2.4
3	A	536	VAL	2.4
3	A	520	GLU	2.4
3	A	432	GLY	2.4
3	A	434	LYS	2.3
3	A	459	ARG	2.3
3	A	779	ARG	2.3
3	A	785	SER	2.2
3	A	755	GLN	2.2
3	A	514	VAL	2.2
1	B	22	DG	2.2
2	C	12	DC	2.1
3	A	548	LYS	2.1
3	A	734	GLU	2.0
3	A	306	ARG	2.0
3	A	560	VAL	2.0
3	A	513	THR	2.0

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

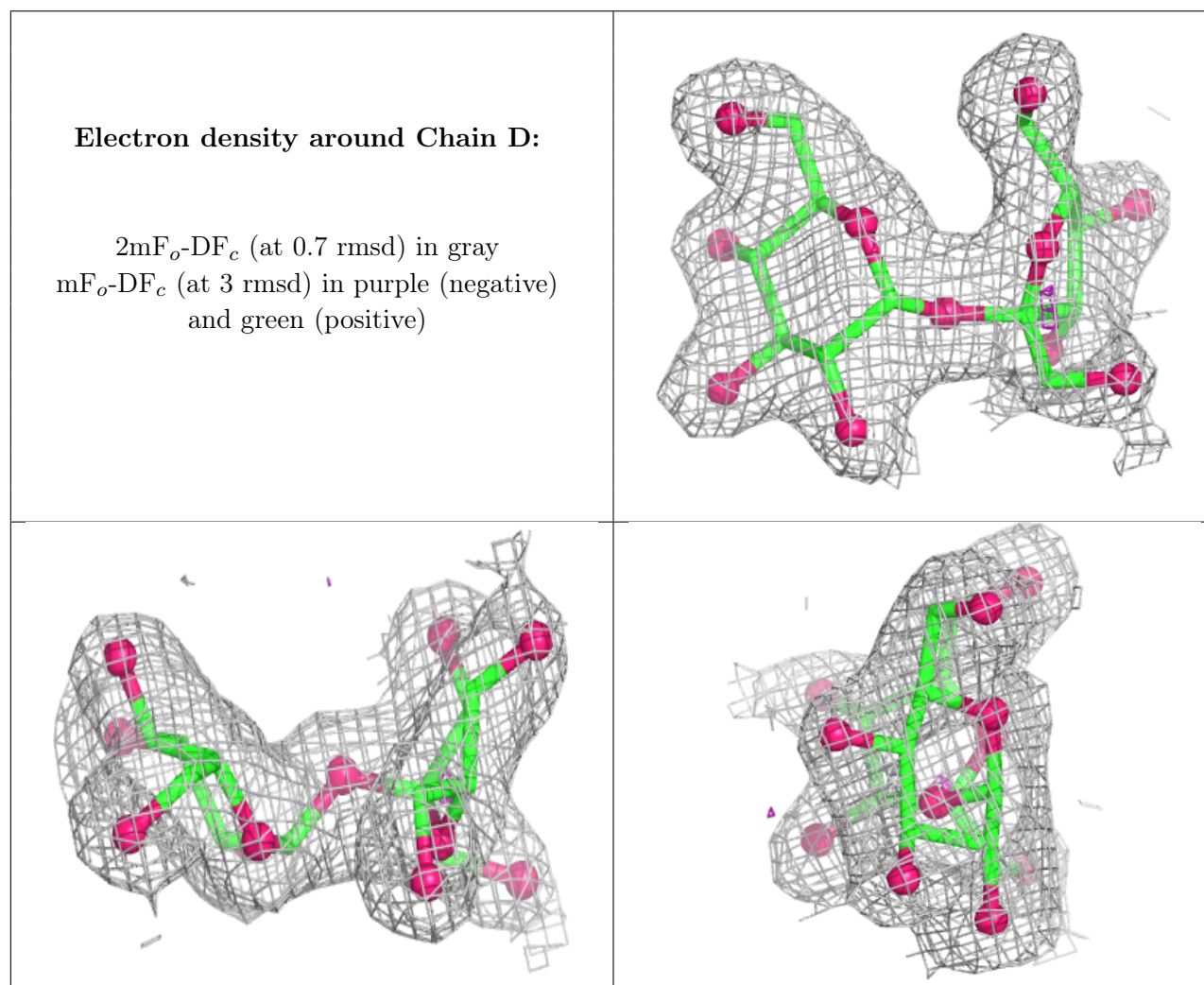
There are no non-standard protein/DNA/RNA residues in this entry.

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($\text{\AA}^2$)	Q<0.9
4	GLC	D	1	11/12	0.92	0.09	43,44,44,46	0
4	FRU	D	2	12/12	0.94	0.09	39,43,43,43	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
6	SO4	A	903	5/5	0.76	0.11	80,80,80,80	0
6	SO4	A	902	5/5	0.88	0.13	64,64,64,65	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	AF	C	333	14/14	0.92	0.11	45,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.