



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:55 PM UTC

PDB ID : 1U48 / pdb_00001u48
Title : Extension of a cytosine-8-oxoguanine base pair
Authors : Hsu, G.W.; Ober, M.; Carell, T.; Beese, L.S.
Deposited on : 2004-07-23
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

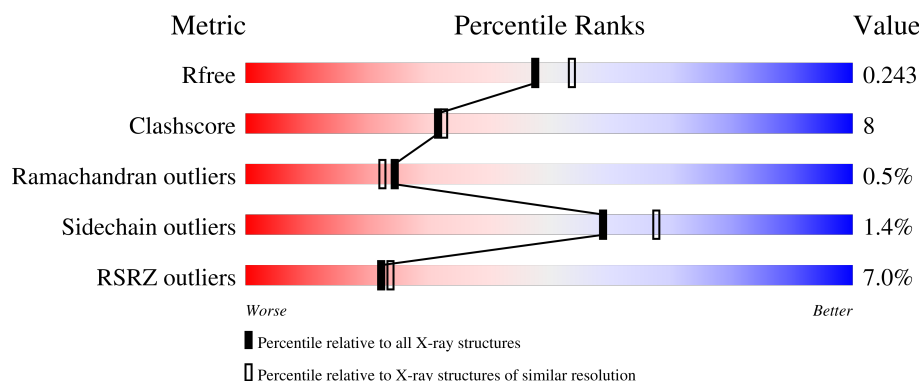
1 Overall quality at a glance

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	12	8% 42% 42% 17%
2	C	15	20% 27% 60% 13%
3	A	580	7% 79% 19% .
4	D	2	50% 50%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5628 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	10	Total	C	N	O	P	0	0	0
			199	96	36	58	9			

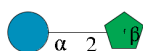
- Molecule 2 is a DNA chain called DNA template strand with 8-oxoguanine.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	P	0	0	0
			268	127	53	76	12			

- Molecule 3 is a protein called DNA polymerase I.

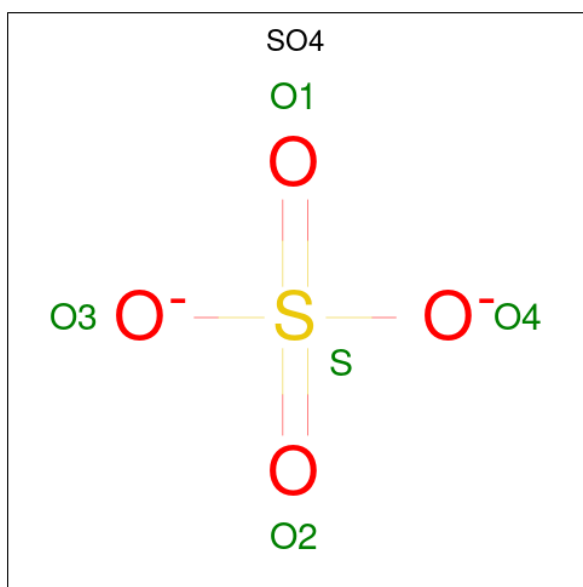
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	580	Total	C	N	O	S	0	0	0
			4650	2956	807	870	17			

- 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		

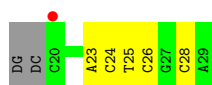
- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

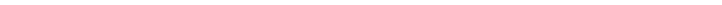
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		

- Molecule 7 is water.

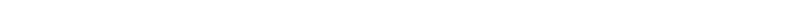
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	18	Total	O	0	0
			18	18		
7	C	31	Total	O	0	0
			31	31		
7	A	423	Total	O	0	0
			423	423		

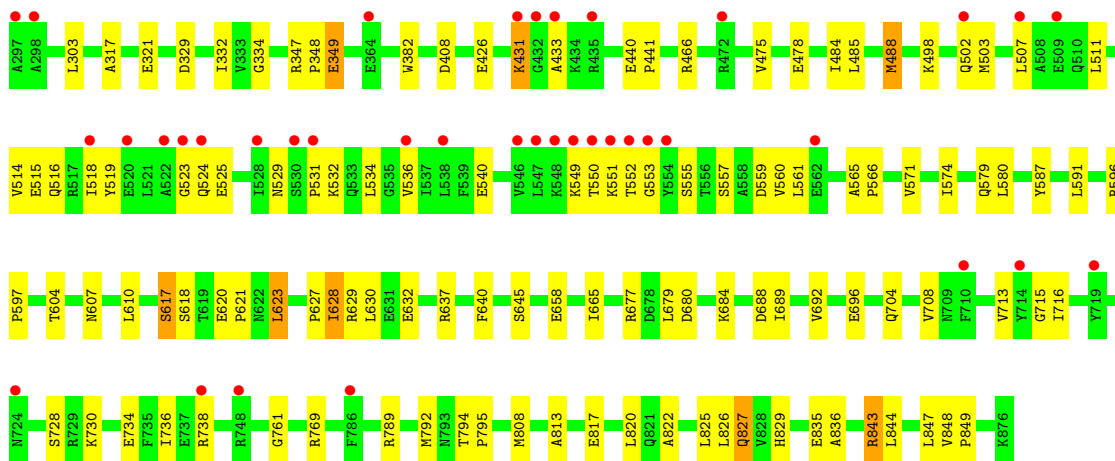
- Molecule 1: DNA primer strand

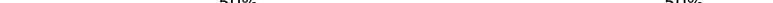


- Chain C:  20% 27% 60% 13%



- Chain A:  7% 79% 19%



- Chain D:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.34Å 93.27Å 106.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.78 – 2.10 40.78 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.5 (40.78-2.10) 93.5 (40.78-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.205 , 0.239 0.207 , 0.243	Depositor DCC
R_{free} test set	2416 reflections (4.44%)	wwPDB-VP
Wilson B-factor (Å ²)	24.1	Xtriage
Anisotropy	0.114	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.3	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	5628	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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: FRU, SOG, SO4, GLC, MG

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	0/222	0.69	0/340
2	C	0.28	0/274	0.66	0/419
3	A	0.40	1/4734 (0.0%)	0.86	11/6398 (0.2%)
All	All	0.39	1/5230 (0.0%)	0.84	11/7157 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	488	MET	SD-CE	-8.18	1.59	1.79

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	321	GLU	N-CA-C	7.13	120.16	109.25
3	A	769	ARG	N-CA-C	-6.68	101.08	110.50
3	A	475	VAL	N-CA-C	6.40	117.19	110.72
3	A	317	ALA	N-CA-C	-6.03	99.45	109.46
3	A	620	GLU	CB-CA-C	-5.97	106.96	111.20
3	A	847	LEU	N-CA-C	5.94	117.55	111.14
3	A	680	ASP	N-CA-C	-5.83	100.75	109.15
3	A	617	SER	N-CA-C	-5.45	100.89	109.50
3	A	665	ILE	N-CA-C	5.17	115.90	110.62
3	A	332	ILE	N-CA-C	-5.04	99.91	107.37
3	A	478	GLU	N-CA-C	5.03	116.45	111.07

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	199	0	114	5	0
2	C	268	0	147	6	0
3	A	4650	0	4701	74	0
4	D	23	0	21	1	0
5	A	15	0	0	0	0
6	A	1	0	0	0	0
7	A	423	0	0	2	0
7	B	18	0	0	2	0
7	C	31	0	0	0	0
All	All	5628	0	4983	82	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 (82) 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:484:ILE:HG22	3:A:488:MET:HE2	1.38	1.06
3:A:843:ARG:HH11	3:A:843:ARG:HB3	1.43	0.83
3:A:485:LEU:HA	3:A:488:MET:HE3	1.63	0.80
3:A:623:LEU:HD23	3:A:826:LEU:HD21	1.67	0.76
3:A:557:SER:O	3:A:560:VAL:HG12	1.86	0.75
2:C:3:DA:H2'	2:C:3:DA:N3	2.03	0.72
3:A:621:PRO:HG2	3:A:623:LEU:HD13	1.74	0.68
2:C:10:DT:H2''	2:C:11:DC:H5'	1.78	0.65
3:A:534:LEU:HD11	3:A:574:ILE:HD13	1.79	0.65
7:B:3354:HOH:O	3:A:629:ARG:HG3	1.96	0.64
3:A:827:GLN:NE2	3:A:829:HIS:H	1.95	0.64
3:A:730:LYS:O	3:A:734:GLU:HG3	1.98	0.64
3:A:813:ALA:O	3:A:817:GLU:HG3	1.98	0.63
3:A:488:MET:HE1	3:A:808:MET:CE	2.32	0.60
3:A:789:ARG:HA	3:A:792:MET:HE3	1.85	0.58
3:A:349:GLU:H	3:A:349:GLU:CD	2.10	0.58
3:A:514:VAL:O	3:A:518:ILE:HG13	2.05	0.57
3:A:761:GLY:HA2	7:A:3529:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:843:ARG:HH11	3:A:843:ARG:CB	2.17	0.56
3:A:848:VAL:HB	3:A:849:PRO:HD3	1.86	0.56
2:C:10:DT:H2''	2:C:11:DC:C5'	2.36	0.55
3:A:519:TYR:CD1	3:A:525:GLU:HA	2.41	0.54
1:B:24:DC:OP1	3:A:552:THR:HG22	2.07	0.54
3:A:692:VAL:HB	3:A:696:GLU:HB2	1.89	0.54
3:A:426:GLU:OE2	3:A:431:LYS:HE2	2.08	0.53
3:A:536:VAL:O	3:A:540:GLU:HB2	2.08	0.53
3:A:591:LEU:HD21	3:A:640:PHE:CZ	2.44	0.53
3:A:728:SER:OG	3:A:730:LYS:HG2	2.09	0.53
3:A:689:ILE:O	3:A:738:ARG:NH1	2.43	0.52
2:C:14:DG:H2''	2:C:15:DC:C5	2.45	0.52
3:A:503:MET:O	3:A:507:LEU:HD23	2.10	0.52
3:A:334:GLY:HA2	3:A:348:PRO:HD3	1.93	0.51
3:A:579:GLN:NE2	3:A:632:GLU:OE1	2.43	0.51
3:A:550:THR:HG23	3:A:555:SER:HA	1.94	0.50
3:A:704:GLN:O	3:A:708:VAL:HG23	2.11	0.50
1:B:28:DC:OP2	3:A:629:ARG:HD2	2.12	0.50
3:A:632:GLU:H	3:A:632:GLU:CD	2.20	0.49
3:A:607:ASN:HB3	3:A:617:SER:OG	2.13	0.49
3:A:621:PRO:HG2	3:A:623:LEU:CD1	2.42	0.49
2:C:12:DA:H2'	2:C:13:DG:C8	2.47	0.49
3:A:523:GLY:O	3:A:524:GLN:HB3	2.13	0.48
3:A:498:LYS:O	3:A:502:GLN:HG3	2.13	0.48
3:A:440:GLU:HB3	3:A:441:PRO:HD3	1.96	0.48
3:A:529:ASN:O	3:A:531:PRO:HD3	2.15	0.47
3:A:549:LYS:HG3	3:A:553:GLY:O	2.14	0.47
3:A:408:ASP:HB2	4:D:2:FRU:H11	1.97	0.45
1:B:25:DT:H2''	1:B:26:DC:H5'	1.98	0.45
3:A:820:LEU:HD21	3:A:843:ARG:NH1	2.31	0.45
3:A:484:ILE:HG22	3:A:488:MET:CE	2.26	0.45
1:B:25:DT:H2''	1:B:26:DC:C5'	2.46	0.45
3:A:511:LEU:HD13	3:A:580:LEU:HB2	1.99	0.45
3:A:645:SER:HA	7:A:2638:HOH:O	2.16	0.45
3:A:715:GLY:O	3:A:789:ARG:HG2	2.17	0.45
3:A:485:LEU:HD23	3:A:488:MET:CE	2.47	0.44
3:A:347:ARG:HB3	3:A:349:GLU:OE2	2.17	0.44
3:A:466:ARG:HH11	3:A:466:ARG:HG3	1.83	0.44
3:A:515:GLU:HG2	3:A:519:TYR:CE2	2.52	0.44
3:A:794:THR:HB	3:A:795:PRO:HD3	2.00	0.43
3:A:488:MET:HE1	3:A:808:MET:HE1	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:658:GLU:OE2	3:A:713:VAL:HG11	2.17	0.43
3:A:587:TYR:CE2	3:A:627:PRO:HD3	2.54	0.43
3:A:827:GLN:HE21	3:A:827:GLN:C	2.27	0.42
3:A:579:GLN:HE21	3:A:630:LEU:HD13	1.84	0.42
3:A:628:ILE:HD12	3:A:637:ARG:NH1	2.33	0.42
3:A:532:LYS:O	3:A:536:VAL:HG23	2.19	0.42
3:A:561:LEU:O	3:A:571:VAL:HG11	2.20	0.42
3:A:623:LEU:HD23	3:A:826:LEU:CD2	2.43	0.42
3:A:604:THR:HB	3:A:623:LEU:HD22	2.02	0.42
3:A:684:LYS:HE3	3:A:688:ASP:OD2	2.20	0.42
1:B:23:DA:H2''	1:B:24:DC:O5'	2.19	0.41
3:A:329:ASP:H	3:A:382:TRP:CD1	2.38	0.41
3:A:825:LEU:HD11	3:A:835:GLU:HB3	2.01	0.41
3:A:716:ILE:HG13	3:A:736:ILE:HD13	2.03	0.41
3:A:591:LEU:HD21	3:A:640:PHE:HZ	1.85	0.41
3:A:677:ARG:HB2	3:A:679:LEU:HG	2.02	0.41
7:B:2804:HOH:O	3:A:433:ALA:HB2	2.20	0.41
3:A:596:ARG:HA	3:A:597:PRO:HD3	1.96	0.40
3:A:610:LEU:HD12	3:A:610:LEU:O	2.21	0.40
3:A:565:ALA:N	3:A:566:PRO:HD2	2.37	0.40
2:C:7:DG:H4'	3:A:618:SER:O	2.21	0.40
3:A:822:ALA:CB	3:A:836:ALA:HB2	2.51	0.40
3:A:516:GLN:NE2	3:A:516:GLN:HA	2.36	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
3	A	578/580 (100%)	556 (96%)	19 (3%)	3 (0%)	24 22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	551	LYS
3	A	431	LYS
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	495/495 (100%)	488 (99%)	7 (1%)	59 67

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	303	LEU
3	A	349	GLU
3	A	559	ASP
3	A	623	LEU
3	A	827	GLN
3	A	843	ARG
3	A	844	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	405	GLN
3	A	502	GLN
3	A	516	GLN
3	A	527	ASN
3	A	579	GLN
3	A	584	GLN
3	A	624	GLN
3	A	656	GLN
3	A	723	GLN
3	A	752	ASN
3	A	759	GLN
3	A	827	GLN

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Mol	Chain	Res	Type
3	A	867	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	8OG	C	5	2,1	22,25,26	2.34	8 (36%)	26,37,40	2.10	8 (30%)

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	8OG	C	5	2,1	-	0/7/21/22	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	5	8OG	C8-N9	-5.28	1.31	1.40
2	C	5	8OG	C5-C4	4.79	1.43	1.37
2	C	5	8OG	C8-N7	-4.59	1.30	1.38
2	C	5	8OG	C2-N3	3.35	1.41	1.33
2	C	5	8OG	C4-N3	3.06	1.41	1.34
2	C	5	8OG	C4-N9	-2.91	1.33	1.39
2	C	5	8OG	C6-N1	2.74	1.44	1.38
2	C	5	8OG	C2-N1	2.31	1.43	1.37

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	5	8OG	N9-C4-N3	5.85	133.44	126.13
2	C	5	8OG	C2-N3-C4	4.42	119.91	112.30
2	C	5	8OG	N7-C8-N9	3.70	110.73	106.61
2	C	5	8OG	O6-C6-C5	-2.97	120.11	127.26
2	C	5	8OG	C5-C6-N1	2.68	119.50	112.13
2	C	5	8OG	C2-N1-C6	-2.52	120.55	125.11
2	C	5	8OG	C1'-N9-C4	-2.31	122.81	126.88
2	C	5	8OG	N1-C2-N3	-2.13	119.42	123.32

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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
4	GLC	D	1	4	11,11,12	3.31	4 (36%)	15,15,17	1.49	2 (13%)
4	FRU	D	2	4	11,12,12	1.37	1 (9%)	10,18,18	0.88	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	9.72	1.67	1.52
4	D	2	FRU	O2-C2	3.80	1.47	1.40
4	D	1	GLC	O5-C5	2.74	1.48	1.43
4	D	1	GLC	C4-C5	2.54	1.58	1.53
4	D	1	GLC	O5-C1	2.53	1.47	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1	GLC	C1-O5-C5	3.93	117.45	112.19
4	D	1	GLC	C1-C2-C3	-2.86	105.48	109.64

There are no chirality outliers.

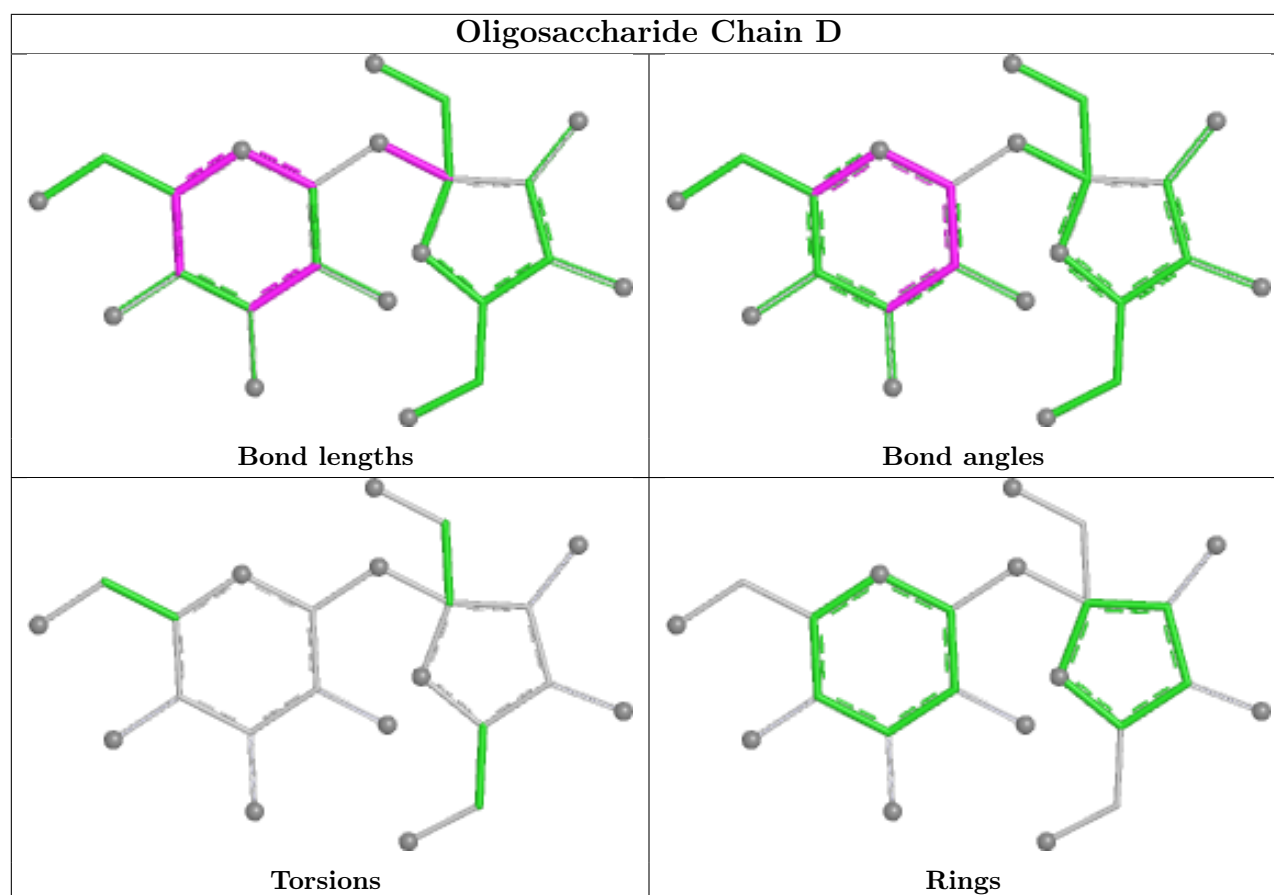
There are no torsion outliers.

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 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	910	-	4,4,4	0.37	0	6,6,6	0.06	0
5	SO4	A	911	-	4,4,4	0.40	0	6,6,6	0.08	0
5	SO4	A	912	-	4,4,4	0.36	0	6,6,6	0.11	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

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, 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	10/12 (83%)	0.44	1 (10%) 12 13	26, 33, 64, 66	0
2	C	12/15 (80%)	1.00	3 (25%) 2 2	24, 39, 82, 91	0
3	A	580/580 (100%)	0.28	38 (6%) 24 26	14, 24, 48, 64	0
All	All	602/607 (99%)	0.29	42 (6%) 22 24	14, 25, 50, 91	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	297	ALA	4.7
3	A	298	ALA	4.2
3	A	433	ALA	3.9
3	A	432	GLY	3.9
3	A	552	THR	3.8
3	A	431	LYS	3.8
3	A	719	TYR	3.6
3	A	520	GLU	3.6
3	A	549	LYS	3.6
2	C	15	DC	3.5
3	A	710	PHE	3.4
2	C	14	DG	3.4
3	A	550	THR	3.2
3	A	553	GLY	3.1
3	A	502	GLN	3.1
3	A	551	LYS	2.8
3	A	554	TYR	2.8
2	C	13	DG	2.7
3	A	522	ALA	2.6
3	A	509	GLU	2.6
3	A	786	PHE	2.5
3	A	435	ARG	2.5
3	A	524	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
3	A	523	GLY	2.4
3	A	364	GLU	2.4
3	A	538	LEU	2.3
3	A	547	LEU	2.2
3	A	530	SER	2.2
3	A	724	ASN	2.2
3	A	528	ILE	2.2
3	A	738	ARG	2.1
3	A	714	TYR	2.1
3	A	531	PRO	2.1
3	A	546	VAL	2.1
3	A	748	ARG	2.1
3	A	562	GLU	2.1
3	A	518	ILE	2.1
3	A	548	LYS	2.1
3	A	536	VAL	2.1
1	B	20	DC	2.0
3	A	507	LEU	2.0
3	A	472	ARG	2.0

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

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.

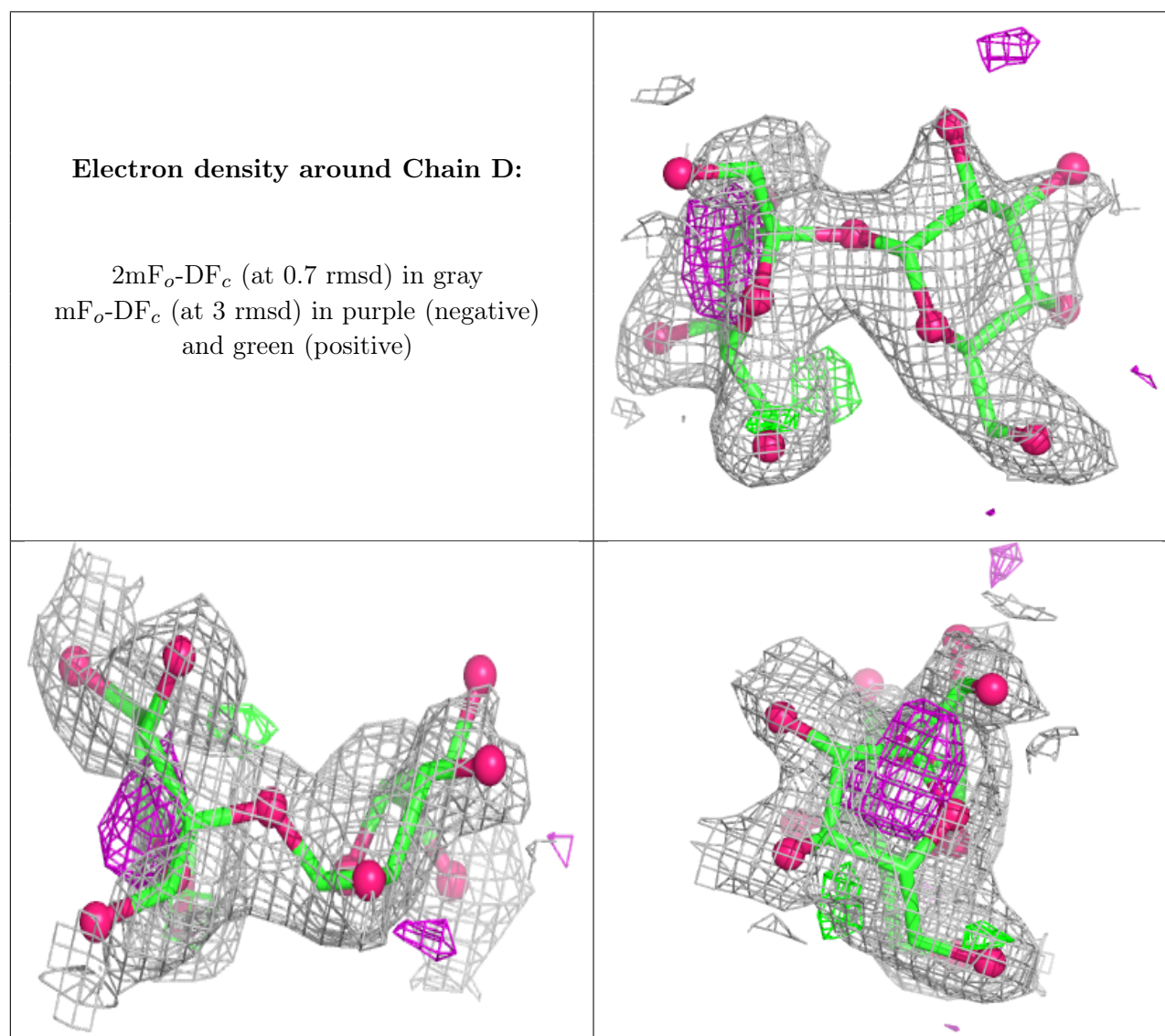
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	8OG	C	5	23/24	0.93	0.09	29,32,33,34	0

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(Å ²)	Q<0.9
4	GLC	D	1	11/12	0.79	0.16	55,57,58,58	0
4	FRU	D	2	12/12	0.79	0.19	53,55,57,57	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
5	SO4	A	911	5/5	0.86	0.12	70,70,70,71	0
5	SO4	A	910	5/5	0.88	0.11	61,61,61,62	0
5	SO4	A	912	5/5	0.89	0.13	48,48,49,49	0
6	MG	A	920	1/1	0.96	0.13	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.