



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:39 AM UTC

PDB ID : 1N3C / pdb_00001n3c
Title : Structural and biochemical exploration of a critical amino acid in human 8-oxoguanine glycosylase
Authors : Norman, D.P.; Chung, S.J.; Verdine, G.L.
Deposited on : 2002-10-25
Resolution : 2.70 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

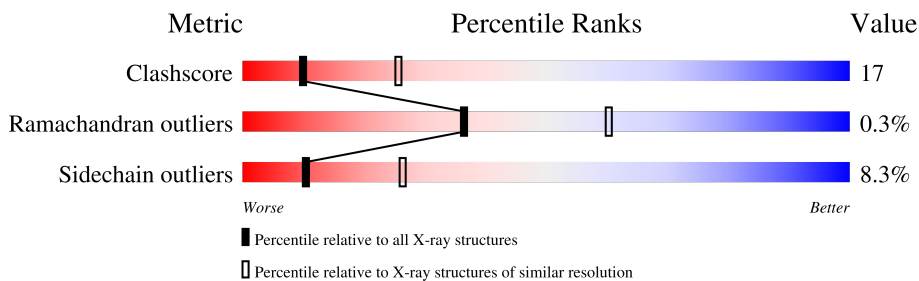
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	B	15	40% 60%
2	C	15	33% 53% 13%
3	A	317	66% 27% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3221 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 complement strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	15	Total	C	N	O	P	0	0	0
			308	146	61	87	14			

- Molecule 2 is a DNA chain called 8-oxoG-containing DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	P	0	0	0
			302	144	54	90	14			

- Molecule 3 is a protein called N-glycosylase/DNA lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	314	Total	C	N	O	S	0	0	0
			2445	1555	438	441	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLY	-	cloning artifact	UNP O15527
A	10	SER	-	cloning artifact	UNP O15527
A	11	GLU	-	cloning artifact	UNP O15527
A	268	ASN	ASP	engineered mutation	UNP O15527

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Ca	0	0
			1	1		

- Molecule 5 is water.

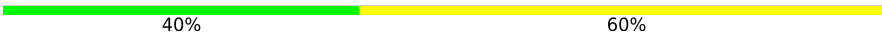
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	9	Total 9	O 9	0	0
5	C	7	Total 7	O 7	0	0
5	A	149	Total 149	O 149	0	0

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. 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. 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.

Note EDS was not executed.

- Molecule 1: DNA complement strand

Chain B: 



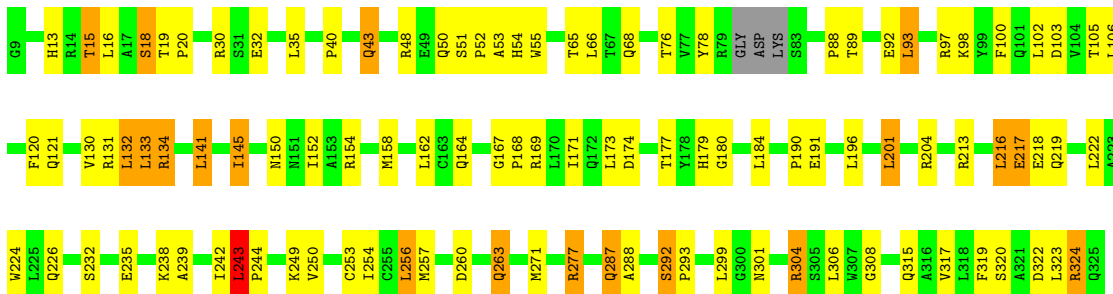
- Molecule 2: 8-oxoG-containing DNA

Chain C: 



- Molecule 3: N-glycosylase/DNA lyase

Chain A: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	92.25Å 92.25Å 210.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.03 – 2.70	Depositor
% Data completeness (in resolution range)	92.6 (29.03-2.70)	Depositor
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.225 , 0.273	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3221	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 8OG

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	0/346	0.51	0/533
2	C	2.18	2/310 (0.6%)	1.57	5/473 (1.1%)
3	A	0.44	0/2513	1.10	21/3425 (0.6%)
All	All	0.79	2/3169 (0.1%)	1.11	26/4431 (0.6%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	30	DC	C4'-O4'	29.86	2.05	1.45
2	C	30	DC	O4'-C1'	23.71	1.89	1.41

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	30	DC	C1'-O4'-C4'	-24.34	73.19	109.70
3	A	134	ARG	NE-CZ-NH2	-17.60	103.36	119.20
3	A	134	ARG	NE-CZ-NH1	14.09	135.59	121.50
2	C	30	DC	O4'-C1'-C2'	12.69	125.43	106.40
2	C	30	DC	O4'-C1'-N1	-11.41	91.29	108.40
2	C	30	DC	O4'-C4'-C3'	10.59	121.29	105.40
3	A	43	GLN	N-CA-C	-7.46	100.14	111.56
3	A	243	LEU	N-CA-C	6.96	118.48	109.64
3	A	308	GLY	CA-C-N	6.95	127.53	119.47
3	A	308	GLY	C-N-CA	6.95	127.53	119.47
3	A	18	SER	N-CA-C	6.67	119.56	111.82
3	A	145	ILE	N-CA-C	-6.59	104.07	110.72
3	A	133	LEU	CA-C-N	-5.96	114.32	123.14
3	A	133	LEU	C-N-CA	-5.96	114.32	123.14
3	A	180	GLY	N-CA-C	-5.46	104.06	112.85
3	A	323	LEU	N-CA-C	-5.42	106.17	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	217	GLU	N-CA-C	5.39	118.96	112.93
3	A	277	ARG	N-CA-C	5.33	117.17	111.36
3	A	243	LEU	CA-C-N	5.33	125.49	119.90
3	A	243	LEU	C-N-CA	5.33	125.49	119.90
3	A	292	SER	CA-C-N	-5.31	113.27	119.32
3	A	292	SER	C-N-CA	-5.31	113.27	119.32
3	A	134	ARG	CD-NE-CZ	5.29	131.81	124.40
3	A	256	LEU	N-CA-C	5.26	116.83	111.14
2	C	30	DC	C5'-C4'-O4'	-5.25	101.52	109.40
3	A	174	ASP	CB-CA-C	-5.08	110.69	116.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	308	0	169	7	0
2	C	302	0	170	13	0
3	A	2445	0	2351	85	1
4	C	1	0	0	0	0
5	A	149	0	0	13	0
5	B	9	0	0	1	0
5	C	7	0	0	0	0
All	All	3221	0	2690	100	1

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 17.

All (100) 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)
2:C:30:DC:C1'	2:C:30:DC:O4'	1.89	1.19
2:C:30:DC:O4'	2:C:30:DC:C4'	2.05	1.05
1:B:14:DG:H1	2:C:17:DC:H42	1.04	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:287:GLN:HE21	3:A:288:ALA:H	1.25	0.84
3:A:169:ARG:HD2	5:A:366:HOH:O	1.76	0.83
3:A:35:LEU:H	3:A:68:GLN:HE22	1.30	0.80
3:A:320:SER:HB2	5:A:357:HOH:O	1.86	0.75
3:A:15:THR:HG22	3:A:18:SER:H	1.49	0.75
3:A:249:LYS:NZ	3:A:253:CYS:SG	2.59	0.75
3:A:132:LEU:HD21	3:A:256:LEU:HG	1.69	0.74
3:A:50:GLN:HE21	3:A:50:GLN:HA	1.55	0.70
3:A:15:THR:HG23	5:A:355:HOH:O	1.95	0.67
3:A:65:THR:HB	3:A:76:THR:CG2	2.26	0.66
1:B:14:DG:H1	2:C:17:DC:N4	1.87	0.65
3:A:50:GLN:HA	3:A:50:GLN:NE2	2.11	0.65
3:A:263:GLN:NE2	3:A:263:GLN:H	1.95	0.65
3:A:30:ARG:HD3	5:A:368:HOH:O	1.97	0.65
3:A:103:ASP:HB2	5:A:389:HOH:O	1.97	0.65
3:A:43:GLN:HE22	3:A:133:LEU:H	1.44	0.64
3:A:158:MET:HE3	3:A:201:LEU:HA	1.79	0.63
3:A:32:GLU:HG3	3:A:131:ARG:HH21	1.62	0.63
3:A:184:LEU:HB3	3:A:216:LEU:HD22	1.81	0.62
2:C:17:DC:H2''	2:C:18:DG:C8	2.35	0.62
3:A:271:MET:HE3	3:A:319:PHE:HD2	1.65	0.62
3:A:32:GLU:OE1	3:A:105:THR:HA	2.00	0.61
3:A:141:LEU:C	3:A:141:LEU:HD23	2.26	0.60
3:A:171:ILE:HD13	3:A:173:LEU:HG	1.83	0.59
3:A:301:ASN:ND2	3:A:304:ARG:HH21	2.00	0.59
3:A:222:LEU:O	3:A:226:GLN:HG3	2.04	0.58
3:A:287:GLN:H	3:A:287:GLN:CD	2.11	0.58
2:C:23:8OG:H5'	3:A:150:ASN:O	2.04	0.58
3:A:249:LYS:C	3:A:249:LYS:HD3	2.28	0.58
3:A:164:GLN:HA	3:A:179:HIS:CD2	2.41	0.56
3:A:169:ARG:HD3	3:A:177:THR:HG21	1.87	0.56
3:A:235:GLU:O	3:A:238:LYS:HB3	2.05	0.56
3:A:287:GLN:HE21	3:A:288:ALA:N	2.00	0.56
1:B:4:DA:H2''	1:B:5:DG:O5'	2.06	0.55
3:A:130:VAL:HB	3:A:317:VAL:HG22	1.90	0.54
3:A:287:GLN:NE2	3:A:288:ALA:H	2.01	0.54
3:A:171:ILE:CD1	3:A:173:LEU:HG	2.37	0.54
3:A:158:MET:CE	3:A:201:LEU:HD13	2.38	0.54
2:C:22:DA:H2''	3:A:150:ASN:O	2.07	0.53
3:A:88:PRO:HB2	3:A:93:LEU:HD13	1.91	0.52
3:A:158:MET:HE3	3:A:201:LEU:HD13	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:DC:H2''	1:B:9:DT:O5'	2.09	0.52
3:A:13:HIS:CD2	3:A:50:GLN:HG3	2.45	0.52
3:A:35:LEU:H	3:A:68:GLN:NE2	2.02	0.52
3:A:43:GLN:NE2	3:A:133:LEU:H	2.07	0.52
2:C:30:DC:O4'	2:C:30:DC:N1	2.42	0.51
3:A:171:ILE:C	3:A:171:ILE:HD12	2.35	0.51
3:A:263:GLN:HE21	3:A:263:GLN:N	2.09	0.51
3:A:32:GLU:OE1	3:A:106:LEU:N	2.41	0.50
3:A:43:GLN:NE2	3:A:133:LEU:HG	2.26	0.50
1:B:12:DA:H2''	1:B:13:DC:O5'	2.11	0.50
3:A:320:SER:O	3:A:324:ARG:HG2	2.11	0.49
3:A:51:SER:HB2	3:A:54:HIS:HB2	1.94	0.49
3:A:97:ARG:HG2	3:A:102:LEU:HD12	1.95	0.48
3:A:219:GLN:O	3:A:224:TRP:HB2	2.14	0.47
3:A:288:ALA:HB1	5:A:463:HOH:O	2.13	0.47
3:A:190:PRO:O	3:A:191:GLU:HB2	2.14	0.47
1:B:5:DG:H2''	1:B:6:DA:C8	2.49	0.47
2:C:23:8OG:OP1	3:A:152:ILE:HG13	2.14	0.47
3:A:19:THR:N	3:A:20:PRO:HD3	2.30	0.47
3:A:263:GLN:NE2	3:A:263:GLN:N	2.62	0.47
3:A:239:ALA:O	3:A:242:ILE:HG12	2.15	0.46
3:A:141:LEU:C	3:A:141:LEU:CD2	2.89	0.46
3:A:257:MET:HA	5:A:346:HOH:O	2.15	0.46
2:C:18:DG:H2''	2:C:19:DT:O5'	2.16	0.46
3:A:89:THR:OG1	3:A:92:GLU:HG3	2.15	0.46
3:A:32:GLU:HG2	5:A:413:HOH:O	2.16	0.46
2:C:27:DT:H2''	2:C:28:DA:C8	2.51	0.45
3:A:100:PHE:O	3:A:131:ARG:HD3	2.17	0.45
3:A:18:SER:C	3:A:20:PRO:HD3	2.42	0.45
3:A:13:HIS:NE2	3:A:50:GLN:HG3	2.32	0.44
3:A:52:PRO:O	3:A:53:ALA:HB3	2.16	0.44
3:A:141:LEU:HD23	3:A:141:LEU:O	2.18	0.44
2:C:19:DT:H2''	2:C:20:DC:O5'	2.18	0.44
3:A:322:ASP:HA	5:A:424:HOH:O	2.16	0.44
3:A:204:ARG:HG3	3:A:204:ARG:NH1	2.33	0.43
3:A:232:SER:OG	3:A:235:GLU:HG2	2.18	0.43
2:C:29:DC:H2''	2:C:30:DC:C5	2.53	0.43
3:A:169:ARG:HD3	3:A:177:THR:CG2	2.48	0.43
1:B:1:DG:H5'	5:B:206:HOH:O	2.19	0.43
3:A:162:LEU:HD13	3:A:196:LEU:HD21	2.01	0.42
3:A:15:THR:HA	3:A:54:HIS:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:134:ARG:NH1	3:A:260:ASP:OD2	2.52	0.42
3:A:78:TYR:N	3:A:78:TYR:CD1	2.87	0.42
3:A:55:TRP:O	3:A:65:THR:HA	2.20	0.42
3:A:292:SER:O	3:A:293:PRO:C	2.62	0.42
3:A:50:GLN:HB2	5:A:358:HOH:O	2.20	0.42
3:A:293:PRO:HG2	5:A:435:HOH:O	2.18	0.42
3:A:243:LEU:HA	3:A:244:PRO:HD3	1.77	0.42
3:A:120:PHE:O	3:A:121:GLN:C	2.63	0.42
3:A:50:GLN:NE2	3:A:50:GLN:CA	2.78	0.41
3:A:93:LEU:HD12	3:A:93:LEU:HA	1.80	0.41
3:A:257:MET:HB3	5:A:354:HOH:O	2.21	0.41
3:A:43:GLN:HE22	3:A:133:LEU:N	2.13	0.41
3:A:250:VAL:O	3:A:254:ILE:HG13	2.21	0.41
3:A:277:ARG:NH2	5:A:454:HOH:O	2.54	0.41
3:A:167:GLY:HA2	3:A:168:PRO:HD3	1.94	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:213:ARG:NH1	3:A:217:GLU:OE2[11_555]	2.08	0.12

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	310/317 (98%)	298 (96%)	11 (4%)	1 (0%)	36 60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	218	GLU

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	252/266 (95%)	231 (92%)	21 (8%)	10	26

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	15	THR
3	A	16	LEU
3	A	40	PRO
3	A	48	ARG
3	A	66	LEU
3	A	93	LEU
3	A	98	LYS
3	A	132	LEU
3	A	141	LEU
3	A	145	ILE
3	A	154	ARG
3	A	201	LEU
3	A	216	LEU
3	A	243	LEU
3	A	263	GLN
3	A	287	GLN
3	A	299	LEU
3	A	304	ARG
3	A	306	LEU
3	A	315	GLN
3	A	324	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	43	GLN
3	A	50	GLN
3	A	68	GLN
3	A	72	GLN

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Mol	Chain	Res	Type
3	A	74	HIS
3	A	108	GLN
3	A	135	GLN
3	A	149	ASN
3	A	151	ASN
3	A	172	GLN
3	A	263	GLN
3	A	276	GLN
3	A	282	HIS
3	A	287	GLN
3	A	296	ASN
3	A	301	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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	23	2	22,25,26	1.01	1 (4%)	26,37,40	1.58	3 (11%)

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	23	2	-	0/7/21/22	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	23	8OG	C8-N7	-3.88	1.31	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	23	8OG	N7-C8-N9	5.38	112.60	106.61
2	C	23	8OG	C5-N7-C8	-3.85	104.17	109.47
2	C	23	8OG	C4-C5-N7	2.57	110.76	106.06

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
2	C	23	8OG	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.