



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:17 AM UTC

PDB ID : 1HU0 / pdb_00001hu0
Title : CRYSTAL STRUCTURE OF AN HOGG1-DNA BOROHYDRIDE
TRAPPED INTERMEDIATE COMPLEX
Authors : Fromme, J.C.; Bruner, S.D.; Yang, W.; Karplus, M.; Verdine, G.L.
Deposited on : 2001-01-03
Resolution : 2.35 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

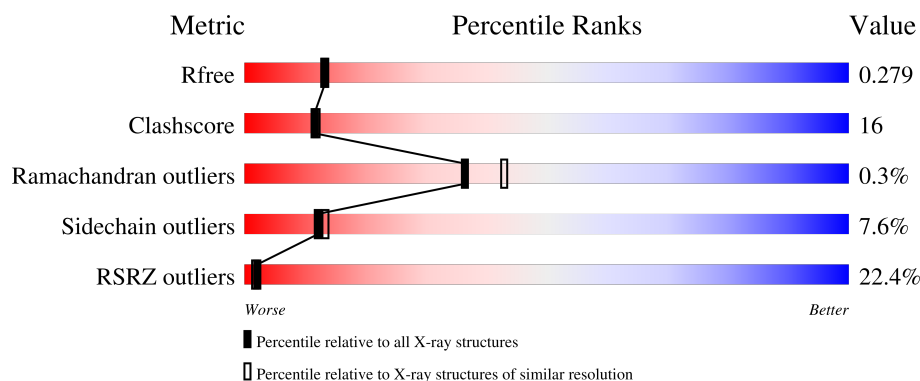
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	D	15	67% 33% 67%
2	E	15	60% 60% 40%
3	A	324	18% 68% 24% 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3200 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*GP*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*C)-3'.

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	15	Total	C	N	O	P	0	0	0
			290	139	49	88	14			

- Molecule 3 is a protein called 8-OXOGUANINE DNA GLYCOSYLASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	314	Total	C	N	O	S	0	0	0
			2491	1579	448	453	11			

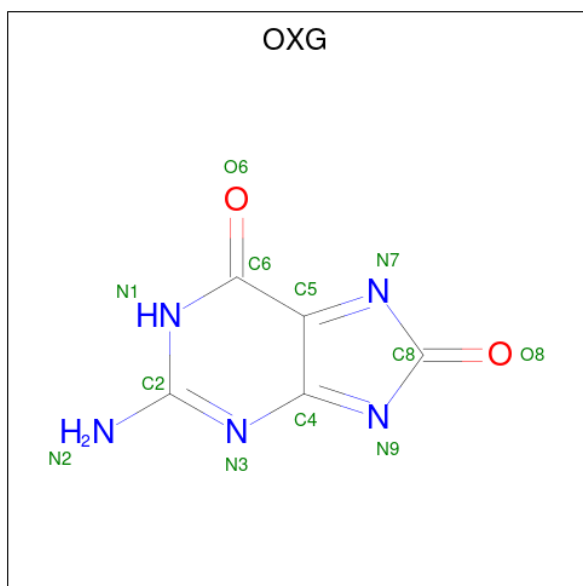
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	ALA	-	cloning artifact	UNP O15527
A	5	MET	-	cloning artifact	UNP O15527
A	6	ALA	-	cloning artifact	UNP O15527
A	7	ASP	-	cloning artifact	UNP O15527
A	8	ILE	-	cloning artifact	UNP O15527
A	9	GLY	-	cloning artifact	UNP O15527
A	10	SER	-	cloning artifact	UNP O15527
A	11	GLU	-	cloning artifact	UNP O15527

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

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

- Molecule 5 is 8-OXOGUANINE (CCD ID: OXG) (formula: $C_5H_3N_5O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			12	5	5	2		

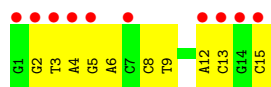
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	2	Total	O	0	0
			2	2		
6	E	5	Total	O	0	0
			5	5		
6	A	91	Total	O	0	0
			91	91		

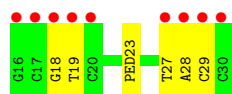
3 Residue-property plots

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.

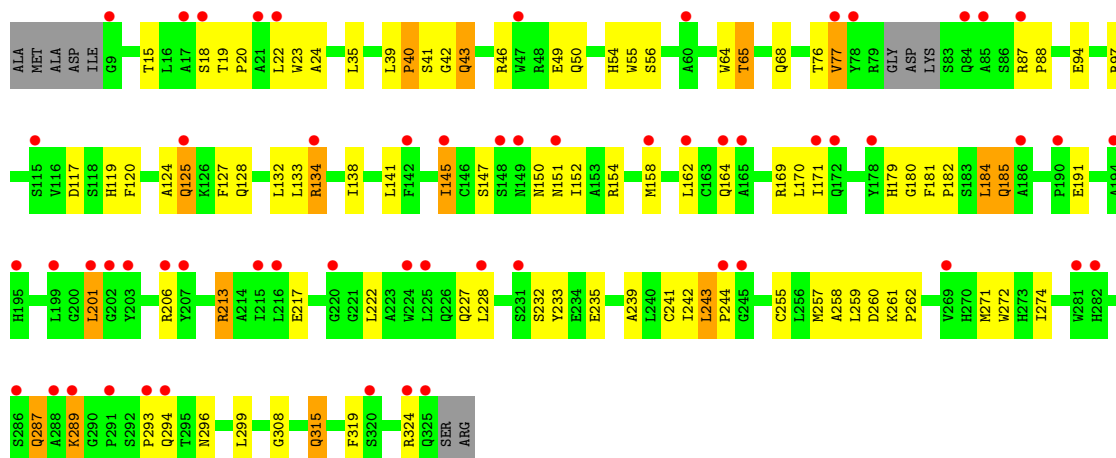
- Molecule 1: 5'-D(*GP*GP*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*C)-3'



- Molecule 2: 5'-D(*GP*CP*GP*TP*CP*CP*AP*(PED)P*GP*TP*CP*TP*AP*CP*C)-3',



- Molecule 3: 8-OXOGUANINE DNA GLYCOSYLASE 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	91.99Å 91.99Å 211.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.14 – 2.35 42.14 – 2.35	Depositor EDS
% Data completeness (in resolution range)	81.5 (42.14-2.35) 91.5 (42.14-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.75 (at 2.34Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.210 , 0.245 0.254 , 0.279	Depositor DCC
R_{free} test set	1860 reflections (8.13%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.2	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	0.90	EDS
Total number of atoms	3200	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% 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: CA, PED, OXG

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	D	0.27	0/346	0.57	0/533
2	E	0.27	0/310	0.58	0/473
3	A	0.43	0/2559	0.97	14/3482 (0.4%)
All	All	0.40	0/3215	0.89	14/4488 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	145	ILE	N-CA-C	-6.62	102.81	111.09
3	A	43	GLN	N-CA-C	-6.59	101.48	111.56
3	A	308	GLY	CA-C-N	6.55	126.48	119.28
3	A	308	GLY	C-N-CA	6.55	126.48	119.28
3	A	243	LEU	CA-C-N	6.23	126.56	119.83
3	A	243	LEU	C-N-CA	6.23	126.56	119.83
3	A	180	GLY	N-CA-C	-5.94	103.28	112.85
3	A	147	SER	N-CA-C	5.72	119.99	112.89
3	A	39	LEU	N-CA-C	5.31	120.89	113.57
3	A	65	THR	N-CA-C	-5.30	99.81	108.76
3	A	41	SER	N-CA-C	5.26	118.82	112.93
3	A	46	ARG	N-CA-C	5.17	119.08	112.26
3	A	127	PHE	N-CA-C	5.14	115.79	108.54
3	A	15	THR	N-CA-C	-5.02	101.75	109.52

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	D	308	0	169	5	1
2	E	290	0	167	4	0
3	A	2491	0	2423	86	1
4	E	1	0	0	0	0
5	A	12	0	3	1	0
6	A	91	0	0	2	0
6	D	2	0	0	0	0
6	E	5	0	0	0	0
All	All	3200	0	2762	94	2

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 16.

All (94) 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:87:ARG:NE	3:A:87:ARG:HA	1.65	1.09
3:A:87:ARG:NH2	3:A:88:PRO:HD3	1.80	0.96
3:A:87:ARG:CZ	3:A:88:PRO:HD3	2.07	0.84
3:A:185:GLN:NE2	3:A:185:GLN:H	1.75	0.82
3:A:76:THR:HA	3:A:87:ARG:HH22	1.46	0.78
3:A:289:LYS:HB3	3:A:289:LYS:NZ	2.04	0.72
3:A:158:MET:HE3	3:A:201:LEU:HA	1.72	0.72
1:D:3:DT:H2''	1:D:4:DA:C8	2.25	0.71
3:A:23:TRP:N	3:A:87:ARG:HD2	2.05	0.70
3:A:35:LEU:H	3:A:68:GLN:HE22	1.40	0.70
3:A:233:TYR:HA	3:A:261:LYS:HD2	1.73	0.69
3:A:87:ARG:NE	3:A:87:ARG:CA	2.52	0.69
3:A:272:TRP:HE1	3:A:296:ASN:ND2	1.92	0.68
3:A:35:LEU:H	3:A:68:GLN:NE2	1.94	0.65
3:A:289:LYS:H	3:A:289:LYS:HD2	1.62	0.65
3:A:87:ARG:HA	3:A:87:ARG:CZ	2.28	0.63
3:A:119:HIS:HE1	6:A:341:HOH:O	1.82	0.62
3:A:134:ARG:NH2	3:A:260:ASP:H	1.98	0.61
3:A:22:LEU:C	3:A:87:ARG:CZ	2.74	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:138:ILE:HD12	3:A:184:LEU:HD13	1.82	0.61
3:A:43:GLN:HE22	3:A:133:LEU:H	1.49	0.61
3:A:141:LEU:O	3:A:145:ILE:HG13	2.01	0.60
3:A:233:TYR:CA	3:A:261:LYS:HD2	2.33	0.59
2:E:28:DA:H2''	2:E:29:DC:O5'	2.04	0.58
3:A:64:TRP:CZ3	3:A:77:VAL:HG22	2.38	0.58
3:A:134:ARG:HH22	3:A:258:ALA:C	2.11	0.58
2:E:23:PED:O1P	3:A:152:ILE:HG13	2.03	0.57
3:A:94:GLU:OE2	3:A:97:ARG:HD3	2.05	0.56
3:A:87:ARG:HA	3:A:87:ARG:HE	1.62	0.56
3:A:185:GLN:H	3:A:185:GLN:HE21	1.52	0.55
3:A:271:MET:HE3	3:A:319:PHE:HD1	1.70	0.55
3:A:287:GLN:H	3:A:287:GLN:CD	2.14	0.55
3:A:43:GLN:NE2	3:A:133:LEU:H	2.03	0.55
3:A:289:LYS:HB3	3:A:289:LYS:HZ3	1.71	0.55
1:D:2:DG:H2''	1:D:3:DT:OP2	2.07	0.55
3:A:22:LEU:HA	3:A:87:ARG:NE	2.22	0.55
3:A:22:LEU:O	3:A:87:ARG:NH2	2.39	0.55
3:A:50:GLN:NE2	3:A:50:GLN:HA	2.22	0.54
2:E:18:DG:H2''	2:E:19:DT:O5'	2.09	0.53
3:A:87:ARG:NH2	3:A:88:PRO:CD	2.65	0.52
3:A:150:ASN:HD22	3:A:151:ASN:H	1.57	0.52
3:A:206:ARG:HH21	3:A:206:ARG:HG3	1.75	0.52
2:E:27:DT:H2''	2:E:28:DA:C8	2.45	0.52
3:A:134:ARG:NE	3:A:260:ASP:OD1	2.42	0.52
3:A:287:GLN:H	3:A:287:GLN:NE2	2.08	0.52
3:A:77:VAL:HG23	3:A:88:PRO:HG3	1.92	0.51
3:A:289:LYS:HD2	3:A:289:LYS:N	2.23	0.51
3:A:255:CYS:HB3	3:A:261:LYS:HG3	1.93	0.51
3:A:22:LEU:O	3:A:87:ARG:CZ	2.59	0.51
3:A:18:SER:C	3:A:20:PRO:HD3	2.35	0.50
3:A:23:TRP:HA	3:A:87:ARG:NH2	2.27	0.50
3:A:87:ARG:HH21	3:A:88:PRO:HD3	1.74	0.49
3:A:42:GLY:O	5:A:328:OXG:N7	2.46	0.48
1:D:12:DA:H2''	1:D:13:DC:O5'	2.13	0.48
3:A:255:CYS:HA	3:A:259:LEU:HB2	1.94	0.48
3:A:23:TRP:HA	3:A:87:ARG:HH21	1.77	0.48
3:A:239:ALA:O	3:A:242:ILE:HB	2.13	0.48
3:A:150:ASN:ND2	3:A:151:ASN:H	2.11	0.48
3:A:232:SER:OG	3:A:235:GLU:HB2	2.13	0.48
3:A:293:PRO:HG2	3:A:294:GLN:NE2	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:171:ILE:C	3:A:171:ILE:HD12	2.40	0.46
3:A:162:LEU:CD2	3:A:182:PRO:HG2	2.46	0.46
3:A:24:ALA:HB2	3:A:88:PRO:HD2	1.98	0.45
3:A:191:GLU:HA	6:A:393:HOH:O	2.15	0.45
1:D:8:DC:H2''	1:D:9:DT:O5'	2.17	0.45
3:A:56:SER:HA	3:A:64:TRP:O	2.15	0.45
3:A:55:TRP:O	3:A:65:THR:HA	2.16	0.45
3:A:19:THR:N	3:A:20:PRO:HD3	2.30	0.45
3:A:158:MET:HE3	3:A:201:LEU:CA	2.43	0.44
3:A:120:PHE:CE2	3:A:124:ALA:HB2	2.53	0.44
3:A:181:PHE:CD2	3:A:182:PRO:HD2	2.53	0.44
3:A:23:TRP:N	3:A:87:ARG:CD	2.80	0.44
3:A:132:LEU:HD12	3:A:257:MET:HG2	2.00	0.43
3:A:324:ARG:N	3:A:324:ARG:HD2	2.33	0.43
3:A:206:ARG:HG3	3:A:206:ARG:NH2	2.34	0.43
3:A:164:GLN:HA	3:A:179:HIS:CD2	2.53	0.43
3:A:289:LYS:HB3	3:A:289:LYS:HZ2	1.83	0.43
3:A:117:ASP:OD1	3:A:117:ASP:C	2.61	0.43
3:A:22:LEU:C	3:A:87:ARG:NH2	2.76	0.43
3:A:185:GLN:NE2	3:A:185:GLN:N	2.57	0.43
3:A:64:TRP:CE2	3:A:77:VAL:HG13	2.54	0.43
3:A:87:ARG:HD3	3:A:88:PRO:HD3	2.01	0.43
3:A:293:PRO:HG2	3:A:294:GLN:HE22	1.82	0.43
3:A:271:MET:CE	3:A:274:ILE:HD12	2.49	0.42
3:A:50:GLN:HA	3:A:50:GLN:HE21	1.84	0.42
3:A:260:ASP:O	3:A:262:PRO:HD3	2.20	0.42
3:A:271:MET:HE3	3:A:319:PHE:CD1	2.53	0.42
1:D:5:DG:H2''	1:D:6:DA:C8	2.55	0.42
3:A:125:GLN:HE21	3:A:125:GLN:HB3	1.58	0.42
3:A:158:MET:HG2	3:A:201:LEU:HD22	2.02	0.42
3:A:243:LEU:HA	3:A:244:PRO:HD3	1.81	0.41
3:A:49:GLU:HA	3:A:54:HIS:O	2.20	0.41
3:A:35:LEU:N	3:A:68:GLN:HE22	2.15	0.41
3:A:315:GLN:HE21	3:A:315:GLN:HB3	1.69	0.40

All (2) 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)
1:D:15:DC:O3'	1:D:15:DC:O3'[11_655]	0.81	1.39
3:A:213:ARG:NH1	3:A:217:GLU:OE2[11_555]	1.88	0.32

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	310/324 (96%)	297 (96%)	12 (4%)	1 (0%)	36	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	40	PRO

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	264/271 (97%)	244 (92%)	20 (8%)	12	13

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	40	PRO
3	A	77	VAL
3	A	125	GLN
3	A	128	GLN
3	A	134	ARG
3	A	154	ARG
3	A	169	ARG
3	A	170	LEU
3	A	184	LEU
3	A	185	GLN

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Mol	Chain	Res	Type
3	A	201	LEU
3	A	213	ARG
3	A	222	LEU
3	A	227	GLN
3	A	228	LEU
3	A	241	CYS
3	A	287	GLN
3	A	289	LYS
3	A	299	LEU
3	A	315	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	43	GLN
3	A	50	GLN
3	A	68	GLN
3	A	108	GLN
3	A	119	HIS
3	A	125	GLN
3	A	150	ASN
3	A	172	GLN
3	A	185	GLN
3	A	219	GLN
3	A	226	GLN
3	A	227	GLN
3	A	273	HIS
3	A	276	GLN
3	A	287	GLN
3	A	294	GLN
3	A	296	ASN

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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	OXG	A	328	-	13,13,13	1.01	1 (7%)	14,19,19	0.94	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	OXG	A	328	-	-	-	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	328	OXG	C8-N7	-2.82	1.30	1.39

There are no bond angle outliers.

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
5	A	328	OXG	1	0

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	D	15/15 (100%)	2.62	10 (66%) 0 0	45, 74, 105, 115	0
2	E	14/15 (93%)	2.66	9 (64%) 0 0	40, 79, 114, 119	0
3	A	314/324 (96%)	1.34	58 (18%) 3 3	21, 34, 52, 73	0
All	All	343/354 (96%)	1.45	77 (22%) 2 2	21, 35, 73, 119	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	15	DC	4.8
2	E	16	DG	4.8
3	A	87	ARG	4.6
1	D	14	DG	4.5
3	A	22	LEU	4.4
1	D	1	DG	4.3
3	A	294	GLN	3.7
2	E	18	DG	3.6
2	E	17	DC	3.6
1	D	2	DG	3.5
2	E	29	DC	3.3
3	A	228	LEU	3.3
3	A	216	LEU	3.2
3	A	325	GLN	3.2
2	E	28	DA	3.2
3	A	158	MET	3.1
2	E	30	DC	3.0
3	A	201	LEU	3.0
3	A	195	HIS	3.0
3	A	288	ALA	3.0
1	D	13	DC	3.0
3	A	9	GLY	2.9
1	D	3	DT	2.9

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Mol	Chain	Res	Type	RSRZ
3	A	21	ALA	2.9
3	A	202	GLY	2.8
3	A	199	LEU	2.8
2	E	19	DT	2.8
3	A	134	ARG	2.8
3	A	85	ALA	2.7
3	A	171	ILE	2.7
3	A	60	ALA	2.6
3	A	190	PRO	2.6
3	A	142	PHE	2.6
3	A	203	TYR	2.6
2	E	27	DT	2.6
3	A	224	TRP	2.5
3	A	293	PRO	2.5
3	A	115	SER	2.5
3	A	286	SER	2.5
3	A	77	VAL	2.5
3	A	151	ASN	2.5
3	A	186	ALA	2.5
3	A	84	GLN	2.5
3	A	164	GLN	2.5
3	A	220	GLY	2.5
1	D	7	DC	2.4
3	A	282	HIS	2.4
3	A	324	ARG	2.4
3	A	194	ALA	2.3
2	E	20	DC	2.3
3	A	207	TYR	2.3
3	A	320	SER	2.3
1	D	5	DG	2.3
3	A	231	SER	2.3
3	A	149	ASN	2.3
3	A	206	ARG	2.3
1	D	12	DA	2.2
3	A	244	PRO	2.2
3	A	162	LEU	2.2
3	A	78	TYR	2.2
3	A	269	VAL	2.2
3	A	178	TYR	2.2
3	A	225	LEU	2.2
3	A	291	PRO	2.2
3	A	172	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
3	A	18	SER	2.2
3	A	245	GLY	2.1
3	A	165	ALA	2.1
3	A	289	LYS	2.1
3	A	47	TRP	2.1
1	D	4	DA	2.1
3	A	148	SER	2.0
3	A	125	GLN	2.0
3	A	145	ILE	2.0
3	A	215	ILE	2.0
3	A	281	TRP	2.0
3	A	17	ALA	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](#)

There are no oligosaccharides in this entry.

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	OXG	A	328	12/12	0.72	0.19	62,66,66,67	0
4	CA	E	100	1/1	0.90	0.24	63,63,63,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.