



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:38 PM GMT

PDB ID : 4FM9
Title : Human topoisomerase II alpha bound to DNA
Authors : Wendorff, T.J.; Schmidt, B.H.; Heslop, P.; Austin, C.A.; Berger, J.M.
Deposited on : 2012-06-15
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

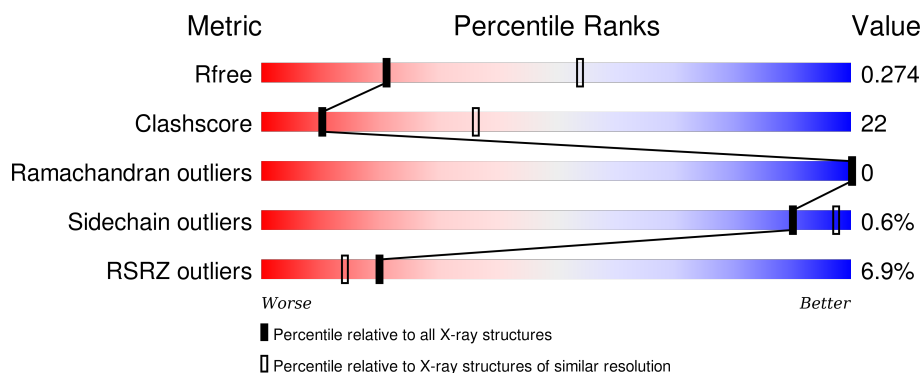
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	A	763	<div> <div>7%</div> <div>55%</div> <div>39%</div> <div>5%</div> </div>
2	C	13	<div> <div>38%</div> <div>54%</div> <div>8%</div> </div>
3	D	17	<div> <div>59%</div> <div>41%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MG	A	1201	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6582 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 protein called DNA topoisomerase 2-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	727	Total	C	N	O	S	0	0	0
			5918	3778	1006	1108	26			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	P	0	0	0
			275	129	57	76	13			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	0	0
			337	162	57	102	16			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		

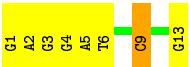
- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	32	Total	O	0	0
			32	32		
6	C	8	Total	O	0	0
			8	8		
6	D	4	Total	O	0	0
			4	4		



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



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	85.92Å 215.08Å 128.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.61 – 2.90 49.61 – 2.90	Depositor EDS
% Data completeness (in resolution range)	94.7 (49.61-2.90) 94.4 (49.61-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.227 , 0.270 0.224 , 0.274	Depositor DCC
R_{free} test set	1292 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	86.4	Xtriage
Anisotropy	0.708	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 53.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 25394 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6582	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 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	A	0.34	0/6039	0.47	0/8139
2	C	0.48	0/310	1.00	2/478 (0.4%)
3	D	0.51	0/375	0.97	0/575
All	All	0.36	0/6724	0.55	2/9192 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	9	DC	O4'-C4'-C3'	-9.10	100.54	106.00
2	C	9	DC	C4'-C3'-C2'	-5.02	98.58	103.10

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	A	5918	0	5969	265	0
2	C	275	0	146	10	0
3	D	337	0	193	13	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	32	0	0	8	0
6	C	8	0	0	1	0
6	D	4	0	0	0	0
All	All	6582	0	6316	277	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 22.

All (277) 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)
1:A:693:LEU:HD11	1:A:697:ASP:HB3	1.63	0.81
1:A:605:HIS:O	1:A:606:LYS:HB2	1.82	0.80
1:A:542:GLN:HG2	1:A:575:THR:O	1.86	0.75
3:D:12:DA:H2"	3:D:13:DT:H5"	1.69	0.74
1:A:903:ALA:HB1	1:A:904:PRO:HD2	1.71	0.73
1:A:602:THR:O	1:A:605:HIS:NE2	2.21	0.72
1:A:1127:TYR:CE2	1:A:1128:LEU:HG	2.26	0.70
1:A:1001:VAL:CG1	1:A:1009:LEU:HD22	2.21	0.70
1:A:542:GLN:CG	1:A:575:THR:O	2.39	0.70
1:A:1053:ARG:HH21	1:A:1082:TYR:HB2	1.56	0.69
1:A:568:ARG:HB3	1:A:637:GLN:NE2	2.07	0.69
1:A:1001:VAL:HG12	1:A:1009:LEU:HD22	1.74	0.69
1:A:1131:MET:HB3	1:A:1132:PRO:HD2	1.73	0.68
1:A:858:THR:HG21	6:A:1311:HOH:O	1.94	0.68
1:A:438:ILE:HD12	1:A:439:PRO:HD2	1.74	0.68
1:A:732:THR:HG21	1:A:751:VAL:HG22	1.76	0.67
1:A:460:THR:HG22	1:A:539:MET:HG2	1.77	0.67
1:A:1138:LYS:O	1:A:1141:LYS:HB3	1.94	0.66
1:A:1047:LYS:O	1:A:1051:GLN:HG3	1.95	0.66
1:A:561:TRP:HB2	1:A:564:LEU:HD12	1.77	0.66
1:A:815:ARG:NH1	1:A:818:PHE:O	2.29	0.65
1:A:1053:ARG:O	1:A:1057:GLU:HG2	1.96	0.65
1:A:522:TYR:CE2	1:A:531:LEU:HD21	2.31	0.65
1:A:1002:LEU:O	1:A:1009:LEU:HD23	1.97	0.65
1:A:1073:LEU:O	1:A:1077:LEU:HG	1.97	0.64
1:A:457:LEU:HB3	1:A:536:ILE:CD1	2.28	0.64
1:A:1056:LEU:O	1:A:1060:ASP:HB3	1.97	0.64
1:A:645:ASP:HB3	1:A:695:TYR:HB2	1.78	0.64
1:A:956:ASP:OD1	1:A:957:TYR:N	2.31	0.63
1:A:1056:LEU:O	1:A:1060:ASP:CB	2.47	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1050:ASN:ND2	1:A:1083:ASP:OD1	2.32	0.62
1:A:1001:VAL:HG22	1:A:1011:LYS:HB2	1.81	0.62
1:A:537:MET:HA	1:A:571:GLU:O	1.99	0.62
1:A:1030:TYR:OH	1:A:1167:TRP:HA	2.00	0.61
1:A:1084:SER:HA	1:A:1127:TYR:OH	1.99	0.61
1:A:1037:LEU:HB3	1:A:1158:LEU:HD21	1.82	0.60
1:A:718:MET:O	1:A:848:VAL:CG2	2.49	0.60
1:A:571:GLU:OE1	1:A:635:ARG:NH1	2.34	0.60
1:A:1053:ARG:NH2	1:A:1082:TYR:HB2	2.17	0.59
1:A:457:LEU:HB3	1:A:536:ILE:HD12	1.82	0.59
1:A:1133:LEU:O	1:A:1135:TYR:CD2	2.55	0.59
1:A:1065:ILE:HA	1:A:1068:LYS:HG3	1.85	0.59
1:A:1004:ASP:OD2	1:A:1010:LYS:HG3	2.01	0.59
2:C:5:DA:H2"	2:C:6:DT:H5'	1.85	0.59
2:C:4:DG:H2"	2:C:5:DA:OP2	2.03	0.59
3:D:12:DA:C2'	3:D:13:DT:H5"	2.33	0.58
1:A:780:ASN:ND2	1:A:866:ASN:HB3	2.17	0.58
1:A:541:ASP:OD1	1:A:616:LEU:HG	2.03	0.58
1:A:1138:LYS:O	1:A:1141:LYS:CB	2.51	0.58
1:A:817:LEU:O	1:A:843:PRO:HA	2.03	0.58
1:A:461:GLU:OE1	2:C:13:DG:O3'	2.22	0.57
1:A:847:MET:HA	1:A:850:ILE:HG22	1.87	0.57
1:A:538:ILE:HD13	1:A:550:LYS:HA	1.86	0.57
1:A:507:ILE:O	1:A:511:ILE:HD12	2.04	0.57
1:A:476:VAL:HG11	1:A:481:TYR:CE2	2.40	0.57
1:A:817:LEU:HD21	1:A:1174:PHE:CE1	2.39	0.57
1:A:524:ASP:O	1:A:528:LEU:HD13	2.04	0.57
1:A:719:VAL:HG11	1:A:840:TRP:CD1	2.41	0.56
1:A:496:ALA:HB3	1:A:501:ILE:HD11	1.87	0.56
1:A:693:LEU:HD11	1:A:697:ASP:CB	2.34	0.56
1:A:1026:ARG:CZ	1:A:1030:TYR:CE2	2.89	0.56
1:A:1030:TYR:CZ	1:A:1167:TRP:HA	2.41	0.56
1:A:736:ARG:HG3	1:A:736:ARG:O	2.05	0.56
1:A:846:PRO:HD3	1:A:876:ILE:HD11	1.86	0.55
1:A:859:GLY:N	3:D:9:DG:H5"	2.21	0.55
1:A:463:ASP:O	1:A:466:LYS:HB3	2.06	0.55
1:A:1048:LEU:HD13	1:A:1147:LEU:HG	1.88	0.55
1:A:602:THR:O	1:A:605:HIS:CE1	2.59	0.55
1:A:507:ILE:HG23	1:A:556:PHE:CE1	2.41	0.55
1:A:1057:GLU:HB3	1:A:1063:ILE:HB	1.89	0.55
1:A:780:ASN:HB2	1:A:850:ILE:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:694:THR:HG1	1:A:697:ASP:CG	2.08	0.54
1:A:996:THR:HG23	3:D:12:DA:OP1	2.06	0.54
1:A:476:VAL:CG1	1:A:481:TYR:CE2	2.91	0.54
1:A:689:THR:O	1:A:689:THR:HG23	2.07	0.54
1:A:599:LYS:HA	1:A:605:HIS:CE1	2.42	0.54
1:A:986:HIS:HB3	1:A:992:GLN:HG3	1.89	0.54
1:A:818:PHE:CD2	1:A:841:TYR:HB3	2.43	0.53
1:A:874:ASN:O	1:A:877:ARG:HG2	2.09	0.53
1:A:802:SER:HB3	1:A:805:TYR:CD2	2.43	0.53
1:A:458:ILE:HA	1:A:537:MET:O	2.08	0.53
2:C:1:DG:H2"	2:C:2:DA:C8	2.44	0.53
1:A:879:LEU:HD23	1:A:1023:PHE:CD2	2.43	0.53
1:A:1053:ARG:HG3	1:A:1054:PHE:N	2.24	0.52
1:A:879:LEU:HD23	1:A:1023:PHE:CG	2.44	0.52
1:A:728:LYS:HB3	1:A:755:SER:HB3	1.90	0.52
1:A:829:LEU:O	1:A:836:VAL:HG22	2.09	0.52
1:A:460:THR:O	1:A:486:LEU:N	2.35	0.52
1:A:1144:LEU:HD23	1:A:1144:LEU:C	2.30	0.52
1:A:892:TYR:CD2	1:A:928:VAL:HG11	2.44	0.52
1:A:541:ASP:O	1:A:543:ASP:N	2.39	0.52
1:A:958:ARG:HD3	1:A:960:TYR:OH	2.10	0.52
1:A:921:ILE:HG22	1:A:970:VAL:HB	1.92	0.52
1:A:718:MET:O	1:A:848:VAL:HG21	2.09	0.52
3:D:13:DT:H2"	3:D:14:DC:C6	2.45	0.51
1:A:460:THR:HB	1:A:465:ALA:HB1	1.92	0.51
1:A:718:MET:HE3	1:A:1021:ASP:HB2	1.92	0.51
1:A:561:TRP:CB	1:A:564:LEU:HD12	2.40	0.51
1:A:757:TYR:CZ	1:A:759:HIS:HB2	2.45	0.51
1:A:868:ASP:O	1:A:872:ILE:HG12	2.10	0.51
1:A:605:HIS:O	1:A:606:LYS:CB	2.58	0.51
1:A:466:LYS:HD2	1:A:485:PRO:HD3	1.92	0.51
2:C:2:DA:H2"	2:C:3:DG:C8	2.46	0.50
1:A:858:THR:CG2	6:A:1311:HOH:O	2.55	0.50
1:A:747:LEU:O	1:A:751:VAL:HG23	2.11	0.50
1:A:849:LEU:HD22	1:A:1015:VAL:HG13	1.94	0.50
2:C:4:DG:P	6:C:205:HOH:O	2.70	0.50
1:A:1044:GLU:O	1:A:1048:LEU:HG	2.11	0.50
1:A:1004:ASP:O	1:A:1007:GLY:N	2.42	0.50
1:A:744:VAL:CG2	1:A:806:ILE:HG13	2.42	0.50
1:A:558:HIS:HA	1:A:565:LEU:HD11	1.93	0.50
1:A:846:PRO:CD	1:A:876:ILE:HD11	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:542:GLN:HG3	1:A:575:THR:O	2.11	0.50
1:A:818:PHE:HD2	1:A:841:TYR:HB3	1.77	0.50
1:A:1026:ARG:NH2	1:A:1030:TYR:CZ	2.80	0.50
1:A:520:LYS:NZ	1:A:527:SER:HA	2.27	0.50
1:A:1138:LYS:O	1:A:1141:LYS:N	2.45	0.49
1:A:673:ARG:NH2	1:A:1006:VAL:O	2.45	0.49
1:A:1026:ARG:NE	1:A:1030:TYR:CE2	2.81	0.49
1:A:860:TRP:C	3:D:9:DG:H4'	2.33	0.49
1:A:920:THR:HG22	1:A:971:LYS:HG2	1.93	0.49
1:A:718:MET:HA	1:A:848:VAL:HG21	1.94	0.49
1:A:990:LYS:NZ	3:D:14:DC:OP1	2.30	0.49
1:A:578:VAL:HG12	1:A:589:PHE:HB2	1.95	0.49
1:A:509:ASN:O	1:A:513:ILE:HG13	2.12	0.49
1:A:906:GLN:HG2	1:A:996:THR:HG22	1.95	0.48
1:A:1171:LEU:O	1:A:1175:ILE:HG12	2.12	0.48
1:A:1070:LYS:HD3	1:A:1073:LEU:HD12	1.95	0.48
1:A:440:LYS:O	1:A:466:LYS:NZ	2.32	0.48
1:A:592:LEU:N	1:A:593:PRO:CD	2.76	0.48
1:A:572:GLU:HB2	1:A:638:PHE:CE2	2.49	0.48
1:A:869:VAL:HG13	1:A:870:ARG:N	2.28	0.48
1:A:1157:THR:HA	1:A:1160:ARG:HD3	1.94	0.48
1:A:736:ARG:HH21	1:A:750:SER:CB	2.27	0.47
1:A:861:SER:OG	3:D:9:DG:C1'	2.63	0.47
1:A:614:LYS:O	6:A:1327:HOH:O	2.20	0.47
1:A:840:TRP:CZ3	1:A:1005:HIS:O	2.67	0.47
1:A:861:SER:OG	3:D:9:DG:O4'	2.32	0.47
1:A:535:LYS:HE3	6:A:1332:HOH:O	2.14	0.47
1:A:456:THR:HA	1:A:535:LYS:O	2.14	0.47
1:A:499:LYS:O	1:A:503:GLU:HG2	2.14	0.47
1:A:1182:GLU:O	1:A:1186:LYS:HG2	2.15	0.47
1:A:741:GLU:OE1	1:A:787:ILE:HG21	2.14	0.47
1:A:568:ARG:HB3	1:A:637:GLN:HE21	1.76	0.47
1:A:776:VAL:HG23	6:A:1312:HOH:O	2.15	0.47
1:A:1133:LEU:O	1:A:1135:TYR:HD2	1.95	0.47
1:A:830:TYR:CE1	1:A:835:ARG:HB2	2.51	0.46
1:A:820:PRO:O	1:A:823:ASP:OD1	2.34	0.46
1:A:695:TYR:O	1:A:699:ILE:HG12	2.15	0.46
1:A:636:ILE:CD1	1:A:686:TYR:CE1	2.98	0.46
1:A:1070:LYS:HA	1:A:1073:LEU:HD12	1.98	0.46
1:A:578:VAL:CG1	1:A:589:PHE:HB2	2.45	0.46
1:A:1055:ILE:O	1:A:1059:ILE:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:VAL:HG23	1:A:587:MET:HB2	1.96	0.46
1:A:452:SER:HB2	1:A:532:ARG:O	2.15	0.46
1:A:739:LYS:O	1:A:809:MET:SD	2.74	0.46
1:A:1085:ASP:HB2	1:A:1127:TYR:CD1	2.51	0.46
2:C:5:DA:H1'	2:C:6:DT:H5'	1.97	0.46
1:A:859:GLY:CA	3:D:9:DG:H5''	2.45	0.46
1:A:997:CYS:HB2	1:A:1000:MET:HE3	1.97	0.46
1:A:1057:GLU:CB	1:A:1063:ILE:HB	2.46	0.46
1:A:1072:GLU:O	1:A:1076:VAL:HG23	2.16	0.46
1:A:867:PHE:CZ	1:A:1015:VAL:HG21	2.51	0.45
1:A:490:ILE:HG22	1:A:491:LEU:N	2.31	0.45
1:A:1085:ASP:H	1:A:1127:TYR:HE1	1.63	0.45
1:A:460:THR:HB	1:A:465:ALA:CB	2.47	0.45
1:A:837:GLU:HB2	1:A:838:PRO:HD2	1.99	0.45
1:A:940:LEU:HD11	1:A:968:PHE:CD1	2.51	0.45
2:C:5:DA:C2'	2:C:6:DT:H5'	2.46	0.45
1:A:471:SER:HB3	1:A:621:SER:HB3	1.97	0.45
1:A:651:LEU:HA	1:A:657:GLN:NE2	2.32	0.45
1:A:870:ARG:NH1	1:A:1179:GLU:OE2	2.49	0.45
1:A:538:ILE:CD1	1:A:550:LYS:HA	2.47	0.45
1:A:626:GLU:O	1:A:629:ALA:HB3	2.17	0.45
1:A:694:THR:OG1	1:A:697:ASP:CG	2.54	0.45
1:A:787:ILE:HB	1:A:807:PHE:O	2.17	0.45
1:A:830:TYR:HE1	1:A:835:ARG:HB2	1.81	0.45
1:A:445:ASN:HB2	1:A:479:ASP:OD1	2.17	0.45
1:A:732:THR:HG23	1:A:754:MET:CE	2.46	0.45
1:A:457:LEU:HB3	1:A:536:ILE:HD13	1.99	0.45
1:A:928:VAL:HG23	1:A:929:ARG:HG2	1.99	0.45
1:A:673:ARG:NH1	1:A:676:LYS:HD3	2.32	0.45
1:A:772:ALA:O	1:A:786:PRO:HG3	2.16	0.45
1:A:941:GLU:N	1:A:942:PRO:CD	2.80	0.44
1:A:934:THR:O	1:A:938:GLN:HG2	2.16	0.44
1:A:921:ILE:CG2	1:A:970:VAL:HB	2.47	0.44
1:A:1085:ASP:N	1:A:1127:TYR:HE1	2.15	0.44
1:A:1055:ILE:HG21	1:A:1141:LYS:HB2	1.98	0.44
1:A:1026:ARG:NH2	1:A:1030:TYR:CE2	2.86	0.44
1:A:847:MET:HA	1:A:850:ILE:CG2	2.48	0.44
1:A:1076:VAL:O	1:A:1080:ARG:HG2	2.18	0.44
1:A:932:THR:HG21	1:A:961:HIS:CD2	2.53	0.44
1:A:785:GLN:O	1:A:787:ILE:HD12	2.17	0.44
1:A:775:PHE:O	1:A:778:SER:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1141:LYS:HD2	1:A:1141:LYS:HA	1.67	0.44
1:A:636:ILE:CD1	1:A:686:TYR:CD1	3.01	0.44
1:A:1002:LEU:C	1:A:1009:LEU:HD23	2.37	0.44
1:A:688:GLN:N	1:A:688:GLN:OE1	2.50	0.44
1:A:450:ARG:O	1:A:451:ASN:HB2	2.17	0.44
1:A:441:LEU:HD11	1:A:443:ASP:HA	2.00	0.43
1:A:494:ARG:HG2	1:A:495:GLU:OE1	2.18	0.43
1:A:912:GLU:H	1:A:925:GLU:HB3	1.83	0.43
1:A:1053:ARG:CZ	1:A:1082:TYR:CD1	3.01	0.43
1:A:874:ASN:HA	1:A:877:ARG:HG2	2.00	0.43
1:A:684:TYR:HE1	1:A:687:GLY:HA2	1.83	0.43
1:A:925:GLU:HG3	1:A:926:LEU:N	2.33	0.43
1:A:689:THR:HG23	6:A:1318:HOH:O	2.17	0.43
1:A:684:TYR:HE1	1:A:687:GLY:CA	2.31	0.43
1:A:837:GLU:HG2	1:A:838:PRO:O	2.18	0.43
1:A:511:ILE:HG23	1:A:516:LEU:O	2.19	0.43
1:A:1147:LEU:O	1:A:1147:LEU:HD12	2.19	0.43
1:A:823:ASP:HA	1:A:826:LEU:HD13	2.00	0.43
1:A:693:LEU:CD1	1:A:697:ASP:CB	2.97	0.43
1:A:1085:ASP:N	1:A:1127:TYR:CE1	2.86	0.43
1:A:732:THR:HG21	1:A:751:VAL:CG2	2.46	0.43
1:A:494:ARG:HD2	1:A:654:SER:HA	2.00	0.42
1:A:568:ARG:HA	6:A:1310:HOH:O	2.18	0.42
1:A:688:GLN:NE2	1:A:690:THR:HG22	2.34	0.42
1:A:949:LYS:HG3	1:A:950:THR:HG23	2.01	0.42
1:A:843:PRO:HB2	1:A:845:ILE:O	2.20	0.42
1:A:553:LEU:O	1:A:556:PHE:HB3	2.18	0.42
1:A:441:LEU:HD22	1:A:513:ILE:HG12	2.01	0.42
1:A:1041:LEU:HD23	1:A:1044:GLU:OE1	2.19	0.42
1:A:1139:GLU:O	1:A:1142:ASP:HB3	2.20	0.42
1:A:593:PRO:HB2	1:A:684:TYR:HB3	2.01	0.42
1:A:923:ILE:HB	1:A:968:PHE:HB2	2.02	0.42
1:A:842:ILE:CG2	1:A:1026:ARG:HD3	2.50	0.42
1:A:543:ASP:OD1	1:A:543:ASP:O	2.37	0.42
1:A:441:LEU:HD12	1:A:442:ASP:N	2.35	0.42
1:A:959:GLU:HB2	1:A:961:HIS:CE1	2.54	0.42
1:A:664:TRP:CE3	1:A:665:LEU:CD1	3.03	0.42
1:A:438:ILE:O	1:A:438:ILE:HG23	2.20	0.42
1:A:985:LEU:H	1:A:985:LEU:HD22	1.85	0.42
1:A:791:GLY:HA3	1:A:797:GLY:HA2	2.02	0.42
1:A:1036:TRP:CZ2	1:A:1090:TRP:CZ2	3.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:903:ALA:HB3	1:A:906:GLN:HB2	2.01	0.41
1:A:1070:LYS:HA	1:A:1073:LEU:CD1	2.50	0.41
1:A:868:ASP:N	1:A:891:SER:HB3	2.34	0.41
1:A:733:CYS:O	1:A:736:ARG:O	2.38	0.41
1:A:1185:GLU:O	1:A:1189:GLU:HG3	2.20	0.41
1:A:1167:TRP:O	1:A:1171:LEU:HG	2.20	0.41
1:A:846:PRO:HD2	1:A:872:ILE:CG2	2.50	0.41
1:A:826:LEU:HD23	1:A:838:PRO:HG3	2.02	0.41
1:A:581:SER:OG	1:A:609:LYS:HE3	2.20	0.41
1:A:730:LEU:HG	1:A:734:PHE:CE2	2.56	0.41
1:A:959:GLU:HB2	1:A:961:HIS:NE2	2.36	0.41
1:A:856:ILE:HD11	2:C:9:DC:C2	2.56	0.41
1:A:764:LEU:O	1:A:767:THR:HB	2.20	0.41
1:A:523:GLU:O	1:A:523:GLU:HG3	2.21	0.41
1:A:824:HIS:NE2	6:A:1324:HOH:O	2.27	0.41
1:A:604:ASN:O	1:A:604:ASN:OD1	2.38	0.41
1:A:813:LEU:HA	1:A:816:LEU:HD12	2.02	0.41
1:A:718:MET:O	1:A:848:VAL:HG23	2.19	0.41
3:D:6:DA:H2'	3:D:7:DT:C6	2.56	0.41
1:A:461:GLU:OE2	1:A:540:THR:HB	2.21	0.41
1:A:1050:ASN:OD1	1:A:1082:TYR:HB3	2.21	0.40
1:A:732:THR:CG2	1:A:751:VAL:HG22	2.49	0.40
1:A:556:PHE:CE2	1:A:557:ILE:HD13	2.56	0.40
1:A:547:SER:HB3	1:A:702:GLU:HG2	2.03	0.40
1:A:668:PHE:HB2	1:A:704:ILE:HD11	2.03	0.40
1:A:1127:TYR:CD2	1:A:1128:LEU:HG	2.55	0.40
1:A:571:GLU:CD	1:A:635:ARG:NH1	2.75	0.40
1:A:919:THR:HG22	1:A:919:THR:O	2.21	0.40
1:A:736:ARG:NH2	1:A:750:SER:CB	2.83	0.40
1:A:859:GLY:H	3:D:9:DG:H5''	1.85	0.40
1:A:577:ILE:HD13	1:A:592:LEU:HD13	2.03	0.40
1:A:856:ILE:HD11	2:C:9:DC:N3	2.37	0.40
1:A:555:ASN:HB2	1:A:653:PHE:CD2	2.57	0.40
1:A:557:ILE:HG13	1:A:570:LEU:HD21	2.03	0.40
1:A:744:VAL:HG21	1:A:806:ILE:HG13	2.02	0.40
1:A:591:SER:OG	1:A:685:LEU:O	2.38	0.40
3:D:11:DC:H2''	3:D:12:DA:OP2	2.22	0.40
1:A:718:MET:SD	1:A:1012:TYR:OH	2.68	0.40
1:A:662:LYS:HE2	1:A:859:GLY:O	2.21	0.40
1:A:648:ALA:O	1:A:651:LEU:HB3	2.22	0.40
1:A:604:ASN:O	1:A:607:LYS:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:716:PRO:HG3	1:A:853:ALA:HB1	2.03	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	
1	A	723/763 (95%)	679 (94%)	44 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	
1	A	655/688 (95%)	651 (99%)	4 (1%)	90	97

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

Mol	Chain	Res	Type
1	A	570	LEU
1	A	945	ASN
1	A	1053	ARG
1	A	1055	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	758	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are 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	GOL	A	1202	-	5,5,5	0.22	0	5,5,5	0.23	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	GOL	A	1202	-	-	0/4/4/4	0/0/0/0

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	A	727/763 (95%)	0.41	52 (7%) 18 12	29, 70, 157, 201	1 (0%)
2	C	13/13 (100%)	-0.02	0 100 100	37, 73, 118, 125	0
3	D	17/17 (100%)	-0.22	0 100 100	39, 57, 110, 113	0
All	All	757/793 (95%)	0.39	52 (6%) 20 14	29, 70, 154, 201	1 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1135	TYR	7.2
1	A	1134	TRP	6.9
1	A	1054	PHE	5.9
1	A	1074	ILE	5.7
1	A	1147	LEU	5.6
1	A	1065	ILE	5.1
1	A	1140	LYS	4.7
1	A	607	LYS	4.4
1	A	1124	THR	4.4
1	A	1139	GLU	4.3
1	A	1072	GLU	4.1
1	A	1051	GLN	3.6
1	A	1125	PHE	3.4
1	A	1071	LYS	3.4
1	A	1142	ASP	3.4
1	A	1145	CYS	3.3
1	A	1070	LYS	3.3
1	A	1126	ASN	3.2
1	A	1136	LEU	3.1
1	A	1138	LYS	3.1
1	A	580	VAL	3.0
1	A	1064	ILE	2.9
1	A	1075	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	1090	TRP	2.9
1	A	1069	PRO	2.8
1	A	1058	LYS	2.8
1	A	1077	LEU	2.8
1	A	1056	LEU	2.8
1	A	1060	ASP	2.8
1	A	1137	THR	2.7
1	A	1063	ILE	2.6
1	A	1059	ILE	2.6
1	A	1036	TRP	2.5
1	A	1084	SER	2.5
1	A	1158	LEU	2.5
1	A	1143	GLU	2.5
1	A	1128	LEU	2.5
1	A	1066	GLU	2.5
1	A	608	TRP	2.4
1	A	1012	TYR	2.3
1	A	1133	LEU	2.2
1	A	1131	MET	2.2
1	A	808	THR	2.2
1	A	1161	LYS	2.2
1	A	692	TYR	2.2
1	A	1001	VAL	2.1
1	A	935	TYR	2.1
1	A	1055	ILE	2.1
1	A	835	ARG	2.1
1	A	1062	LYS	2.0
1	A	589	PHE	2.0
1	A	1129	LEU	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 ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	MG	A	1201	1/1	0.94	0.28	2.08	35,35,35,35	0
5	GOL	A	1202	6/6	0.92	0.22	-0.43	52,62,74,74	0
4	MG	C	101	1/1	0.76	0.69	-	51,51,51,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.