



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:45 AM GMT

PDB ID : 3PX7
Title : Crystal Structure of covalent complex of topoisomerase 1A with substrate
Authors : Zhang, Z.; Tse-dinh, Y.C.; Cheng, B.
Deposited on : 2010-12-09
Resolution : 2.30 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

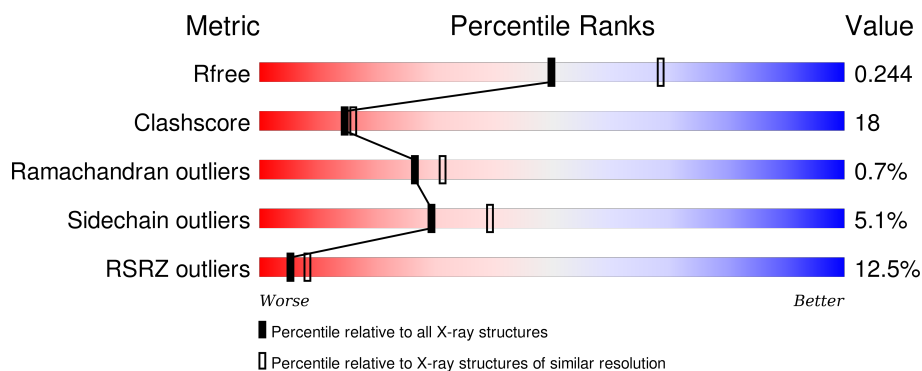
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	595	
2	B	5	
3	C	8	

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
1	PTR	A	319	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4726 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.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	P	S	0	0	0
			4407	2772	786	831	1	17			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	111	ASN	ASP	ENGINEERED MUTATION	UNP C9QXS7

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	2	Total	C	N	O	P	0	0	0
			36	20	4	11	1			

- Molecule 3 is a DNA chain called DNA 5'-D(*A*AP*TP*GP*CP*GP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	7	Total	C	N	O	P	0	0	0
			140	68	25	41	6			

- Molecule 4 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

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

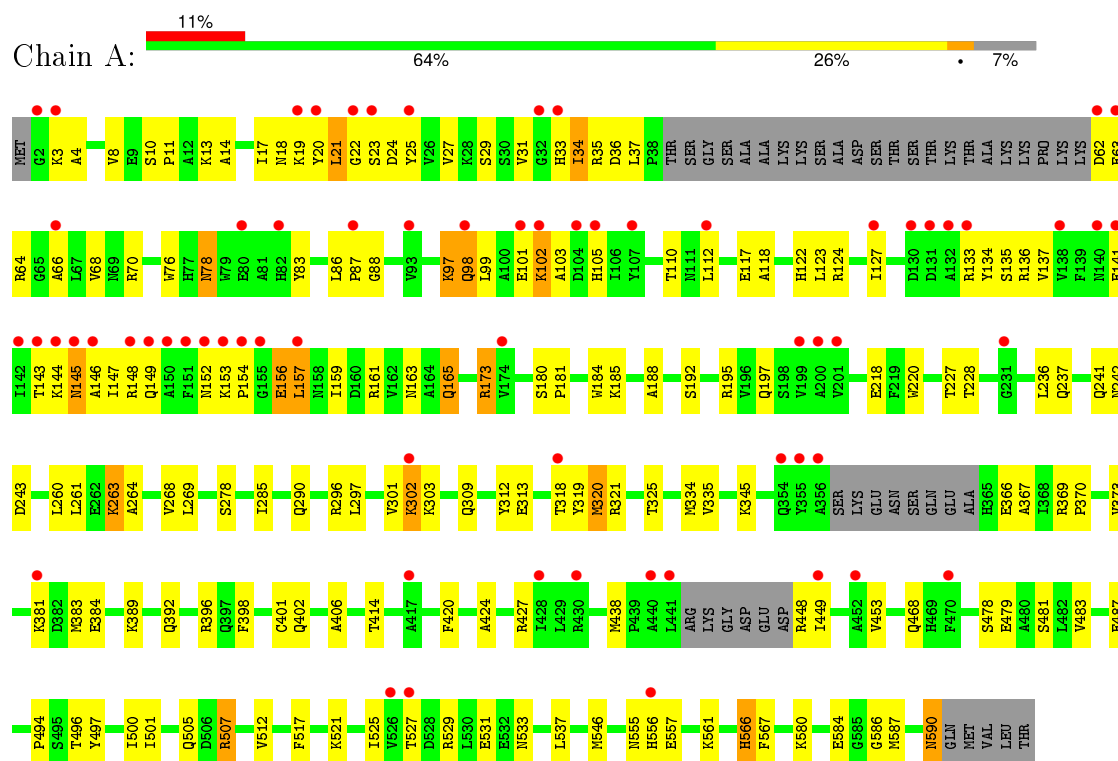
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	133	Total 133	O 133	0	0
5	C	7	Total 7	O 7	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

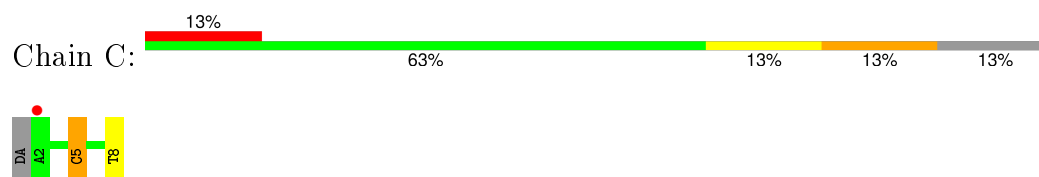
• Molecule 1: DNA topoisomerase



• Molecule 2: DNA 5'-D(*TP*TP*GP*GP*G)-3'



• Molecule 3: DNA 5'-D(*A*AP*TP*GP*CP*GP*CP*T)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.60 Å 91.76 Å 141.98 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	500.00 – 2.30 32.94 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.2 (500.00-2.30) 99.0 (32.94-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.57 (at 2.29 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.230 , 0.269 0.245 , 0.244	Depositor DCC
R_{free} test set	1811 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.2	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	0 of 36231 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4726	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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: HG, PTR

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	0.35	0/4475	0.56	0/6043
2	B	0.64	0/39	1.69	1/59 (1.7%)
3	C	0.37	0/156	0.88	1/239 (0.4%)
All	All	0.36	0/4670	0.59	2/6341 (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	B	1	DT	P-O3'-C3'	8.46	129.86	119.70
3	C	5	DC	O4'-C1'-N1	6.35	112.44	108.00

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	4407	0	4378	160	0
2	B	36	0	23	5	0
3	C	140	0	80	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	133	0	0	2	0
5	C	7	0	0	0	0
All	All	4726	0	4481	161	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 18.

All (161) 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:118:ALA:HB2	1:A:165:GLN:HG2	1.44	1.00
1:A:21:LEU:HD21	1:A:25:TYR:HB2	1.53	0.91
1:A:320:MET:H	1:A:320:MET:CE	1.84	0.90
1:A:302:LYS:HE2	1:A:302:LYS:H	1.35	0.89
1:A:521:LYS:HE3	1:A:525:ILE:HD11	1.57	0.86
1:A:263:LYS:C	1:A:263:LYS:HD2	1.97	0.85
1:A:260:LEU:O	1:A:263:LYS:HE3	1.78	0.83
1:A:33:HIS:O	1:A:34:ILE:HG12	1.79	0.82
1:A:112:LEU:H	1:A:112:LEU:HD23	1.43	0.81
1:A:263:LYS:HD2	1:A:264:ALA:N	1.95	0.81
1:A:320:MET:H	1:A:320:MET:HE3	1.48	0.78
1:A:478:SER:H	1:A:481:SER:HB3	1.48	0.77
1:A:309:GLN:HE22	1:A:321:ARG:NH2	1.85	0.74
1:A:302:LYS:CE	1:A:302:LYS:H	2.01	0.74
1:A:449:ILE:HG13	1:A:449:ILE:O	1.86	0.74
1:A:319:PTR:O1P	1:A:321:ARG:HD2	1.90	0.70
1:A:319:PTR:P	2:B:1:DT:C5'	2.79	0.70
1:A:334:MET:SD	1:A:373:VAL:HG23	2.30	0.70
1:A:118:ALA:CB	1:A:165:GLN:HG2	2.21	0.70
1:A:63:GLU:HB3	1:A:76:TRP:CD2	2.29	0.68
1:A:373:VAL:HG22	1:A:398:PHE:HE2	1.59	0.67
1:A:366:GLU:HG2	1:A:367:ALA:H	1.59	0.67
1:A:147:ILE:HD12	1:A:148:ARG:N	2.10	0.67
1:A:320:MET:N	1:A:320:MET:CE	2.58	0.65
1:A:319:PTR:O3P	3:C:8:DT:H2''	1.97	0.65
1:A:302:LYS:HE3	1:A:479:GLU:OE1	1.97	0.64
1:A:37:LEU:CD1	1:A:173:ARG:HG3	2.27	0.64
1:A:124:ARG:HE	1:A:157:LEU:HD21	1.63	0.64
1:A:35:ARG:NH1	1:A:83:TYR:HB2	2.14	0.63
1:A:309:GLN:HE22	1:A:321:ARG:HH22	1.46	0.63
1:A:319:PTR:HA	1:A:320:MET:CE	2.29	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:LEU:HD23	1:A:424:ALA:HB2	1.81	0.62
1:A:66:ALA:O	1:A:70:ARG:HG2	2.01	0.61
1:A:483:VAL:HG23	1:A:501:ILE:HD11	1.82	0.60
1:A:97:LYS:NZ	1:A:127:ILE:HA	2.16	0.60
1:A:278:SER:HB3	1:A:406:ALA:HB3	1.82	0.60
1:A:319:PTR:HA	1:A:320:MET:HE3	1.84	0.59
1:A:312:TYR:CD2	1:A:320:MET:HE1	2.36	0.59
1:A:23:SER:O	1:A:24:ASP:HB3	2.03	0.58
1:A:21:LEU:HD23	1:A:22:GLY:O	2.04	0.58
1:A:112:LEU:H	1:A:112:LEU:CD2	2.14	0.58
1:A:269:LEU:HD12	1:A:414:THR:HG22	1.85	0.58
1:A:156:GLU:OE1	1:A:156:GLU:N	2.36	0.57
1:A:144:LYS:HA	1:A:147:ILE:HG13	1.86	0.57
1:A:19:LYS:C	1:A:21:LEU:H	2.07	0.57
1:A:496:THR:HG22	1:A:500:ILE:HG12	1.87	0.57
1:A:11:PRO:HD2	2:B:2:DT:OP1	2.05	0.57
1:A:313:GLU:OE1	2:B:2:DT:H5'	2.05	0.56
1:A:147:ILE:HD12	1:A:148:ARG:HG3	1.89	0.55
1:A:18:ASN:HA	1:A:21:LEU:HD22	1.89	0.55
1:A:161:ARG:HG3	1:A:161:ARG:HH11	1.72	0.55
1:A:143:THR:HG22	1:A:145:ASN:H	1.72	0.54
1:A:533:ASN:HD21	1:A:587:MET:HA	1.73	0.54
1:A:102:LYS:O	1:A:102:LYS:HG2	2.08	0.54
1:A:285:ILE:HG23	1:A:325:THR:HG22	1.90	0.54
1:A:34:ILE:HD13	1:A:123:LEU:HD23	1.90	0.53
1:A:555:ASN:O	1:A:557:GLU:HG3	2.09	0.53
1:A:260:LEU:O	1:A:263:LYS:CE	2.54	0.53
1:A:527:THR:O	1:A:531:GLU:HG3	2.09	0.53
1:A:124:ARG:HH11	1:A:124:ARG:HG2	1.74	0.52
1:A:144:LYS:HA	1:A:147:ILE:CG1	2.40	0.52
1:A:37:LEU:HD12	1:A:173:ARG:HG3	1.92	0.52
1:A:149:GLN:HA	1:A:152:ASN:OD1	2.09	0.52
1:A:319:PTR:OH	2:B:1:DT:H3'	2.11	0.51
1:A:3:LYS:HD3	1:A:102:LYS:HE2	1.90	0.51
1:A:302:LYS:H	1:A:302:LYS:CD	2.23	0.51
1:A:97:LYS:HZ3	1:A:127:ILE:HA	1.74	0.51
1:A:35:ARG:HG2	1:A:83:TYR:HB3	1.93	0.51
1:A:366:GLU:HG2	1:A:367:ALA:N	2.26	0.51
1:A:320:MET:HE3	1:A:320:MET:N	2.19	0.50
1:A:13:LYS:O	1:A:17:ILE:HG12	2.11	0.50
1:A:296:ARG:O	1:A:297:LEU:HD23	2.11	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ARG:HG2	1:A:133:ARG:O	2.11	0.50
1:A:36:ASP:CG	1:A:37:LEU:H	2.16	0.49
1:A:184:TRP:HA	1:A:188:ALA:O	2.12	0.49
1:A:78:ASN:ND2	1:A:561:LYS:HD3	2.27	0.49
1:A:180:SER:HB2	1:A:181:PRO:HD3	1.95	0.49
1:A:97:LYS:NZ	1:A:127:ILE:HG12	2.28	0.48
1:A:8:VAL:O	1:A:29:SER:HA	2.13	0.48
1:A:112:LEU:HD23	1:A:112:LEU:N	2.22	0.48
1:A:521:LYS:O	1:A:525:ILE:HD13	2.13	0.48
1:A:242:ASN:O	1:A:243:ASP:HB2	2.12	0.48
1:A:302:LYS:HD2	1:A:303:LYS:H	1.79	0.48
1:A:290:GLN:CD	1:A:301:VAL:HG13	2.33	0.48
1:A:335:VAL:HG11	1:A:402:GLN:HG3	1.96	0.48
1:A:580:LYS:O	1:A:586:GLY:HA3	2.14	0.48
1:A:590:ASN:O	1:A:590:ASN:ND2	2.47	0.48
1:A:525:ILE:O	1:A:529:ARG:HG2	2.13	0.47
1:A:505:GLN:NE2	1:A:512:VAL:HG23	2.29	0.47
1:A:105:HIS:NE2	1:A:135:SER:HB2	2.30	0.47
1:A:147:ILE:CD1	1:A:148:ARG:HG3	2.45	0.47
1:A:3:LYS:CD	1:A:102:LYS:HE2	2.44	0.47
1:A:302:LYS:N	1:A:302:LYS:CD	2.78	0.47
1:A:507:ARG:NH2	3:C:5:DC:OP1	2.41	0.47
1:A:218:GLU:OE2	1:A:468:GLN:NE2	2.44	0.47
1:A:117:GLU:HG2	1:A:161:ARG:HB2	1.97	0.47
1:A:195:ARG:HD3	1:A:546:MET:HE1	1.95	0.47
1:A:62:ASP:OD1	1:A:63:GLU:N	2.48	0.46
1:A:507:ARG:HB2	1:A:507:ARG:NH1	2.31	0.46
1:A:384:GLU:HG3	5:A:659:HOH:O	2.13	0.46
1:A:192:SER:OG	1:A:197:GLN:NE2	2.49	0.46
1:A:64:ARG:O	1:A:68:VAL:HG23	2.16	0.46
1:A:268:VAL:HG21	1:A:453:VAL:CG1	2.46	0.46
1:A:31:VAL:O	1:A:31:VAL:HG12	2.15	0.46
1:A:318:THR:O	1:A:319:PTR:C	2.64	0.45
1:A:105:HIS:CE1	1:A:135:SER:HB2	2.51	0.45
1:A:369:ARG:HB2	1:A:370:PRO:HD2	1.98	0.45
1:A:227:THR:HG22	1:A:228:THR:O	2.16	0.45
1:A:4:ALA:HB3	1:A:25:TYR:CD2	2.51	0.45
1:A:124:ARG:NH1	1:A:124:ARG:HG2	2.32	0.45
1:A:220:TRP:CH2	1:A:468:GLN:HB2	2.52	0.45
1:A:479:GLU:HB3	1:A:501:ILE:HD13	1.97	0.45
1:A:14:ALA:HB1	1:A:27:VAL:HG12	1.97	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LYS:O	1:A:21:LEU:N	2.48	0.45
2:B:1:DT:C7	2:B:2:DT:H73	2.47	0.44
1:A:18:ASN:O	1:A:22:GLY:N	2.49	0.44
1:A:97:LYS:NZ	1:A:97:LYS:HA	2.33	0.44
1:A:345:LYS:HB3	1:A:345:LYS:HZ2	1.81	0.44
1:A:124:ARG:HG3	1:A:134:TYR:CE2	2.53	0.44
1:A:161:ARG:HG3	1:A:161:ARG:NH1	2.32	0.44
1:A:23:SER:C	1:A:25:TYR:H	2.20	0.44
1:A:35:ARG:NE	1:A:122:HIS:HD2	2.15	0.43
1:A:496:THR:O	1:A:500:ILE:HG12	2.18	0.43
1:A:159:ILE:HG12	1:A:163:ASN:ND2	2.32	0.43
1:A:302:LYS:N	1:A:302:LYS:HD2	2.33	0.43
1:A:101:GLU:C	1:A:103:ALA:H	2.21	0.43
1:A:86:LEU:C	1:A:88:GLY:H	2.20	0.43
1:A:392:GLN:CG	1:A:396:ARG:NH1	2.81	0.43
1:A:392:GLN:CD	1:A:396:ARG:NH1	2.72	0.43
1:A:580:LYS:HE3	1:A:584:GLU:OE1	2.18	0.43
1:A:261:LEU:HD21	1:A:420:PHE:CG	2.53	0.43
1:A:185:LYS:HB2	1:A:185:LYS:HE3	1.83	0.43
1:A:533:ASN:ND2	1:A:587:MET:HA	2.34	0.43
1:A:427:ARG:O	1:A:448:ARG:HB3	2.19	0.43
1:A:11:PRO:HG3	1:A:31:VAL:HG22	2.00	0.42
1:A:479:GLU:HA	1:A:517:PHE:HE1	1.84	0.42
1:A:147:ILE:HD12	1:A:147:ILE:C	2.39	0.42
1:A:149:GLN:HA	1:A:152:ASN:CG	2.38	0.42
1:A:590:ASN:N	1:A:590:ASN:HD22	2.17	0.42
1:A:10:SER:HA	1:A:11:PRO:HD3	1.94	0.42
1:A:268:VAL:HG21	1:A:453:VAL:HG11	2.01	0.42
1:A:494:PRO:HA	1:A:497:TYR:CE2	2.55	0.42
1:A:117:GLU:OE1	1:A:136:ARG:NH1	2.53	0.42
1:A:396:ARG:HB3	1:A:438:MET:HG3	2.01	0.42
1:A:566:HIS:ND1	1:A:567:PHE:N	2.67	0.42
1:A:241:GLN:HG2	1:A:242:ASN:ND2	2.35	0.42
1:A:110:THR:O	1:A:136:ARG:NH2	2.53	0.41
1:A:145:ASN:O	1:A:145:ASN:ND2	2.53	0.41
1:A:97:LYS:HZ1	1:A:127:ILE:HG12	1.86	0.41
1:A:86:LEU:O	1:A:88:GLY:N	2.54	0.41
1:A:98:GLN:HG3	1:A:99:LEU:N	2.35	0.41
1:A:137:VAL:HG12	1:A:154:PRO:HB3	2.01	0.41
1:A:319:PTR:HE2	1:A:319:PTR:O2P	2.20	0.41
1:A:260:LEU:C	1:A:263:LYS:HE3	2.37	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:LEU:HD13	1:A:173:ARG:HG3	2.03	0.41
1:A:566:HIS:HE1	5:A:666:HOH:O	2.03	0.41
1:A:345:LYS:NZ	1:A:345:LYS:HB3	2.35	0.40
1:A:141:GLU:H	1:A:146:ALA:HB1	1.86	0.40
1:A:19:LYS:C	1:A:21:LEU:N	2.74	0.40
1:A:78:ASN:HD22	1:A:78:ASN:HA	1.61	0.40
1:A:521:LYS:HE3	1:A:525:ILE:CD1	2.41	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	543/595 (91%)	517 (95%)	22 (4%)	4 (1%)	26	31

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	TYR
1	A	34	ILE
1	A	87	PRO
1	A	102	LYS

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	467/503 (93%)	443 (95%)	24 (5%)	29	39

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LEU
1	A	78	ASN
1	A	97	LYS
1	A	98	GLN
1	A	145	ASN
1	A	153	LYS
1	A	156	GLU
1	A	157	LEU
1	A	165	GLN
1	A	173	ARG
1	A	237	GLN
1	A	263	LYS
1	A	302	LYS
1	A	320	MET
1	A	381	LYS
1	A	383	MET
1	A	389	LYS
1	A	401	CYS
1	A	487	GLU
1	A	507	ARG
1	A	537	LEU
1	A	556	HIS
1	A	566	HIS
1	A	590	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	ASN
1	A	111	ASN
1	A	122	HIS
1	A	140	ASN
1	A	145	ASN
1	A	152	ASN
1	A	163	ASN
1	A	165	GLN
1	A	166	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	197	GLN
1	A	237	GLN
1	A	242	ASN
1	A	253	GLN
1	A	309	GLN
1	A	333	ASN
1	A	353	ASN
1	A	505	GLN
1	A	533	ASN
1	A	539	ASN
1	A	548	ASN
1	A	590	ASN

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	$\# Z > 2$
1	PTR	A	319	1,2	14,16,17	1.94	1 (7%)	18,22,24	0.71	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
1	PTR	A	319	1,2	-	0/9/11/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	319	PTR	OH-CZ	-7.14	1.23	1.40

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 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	319	PTR	8	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are 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 [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	A	551/595 (92%)	0.79	67 (12%) 5 9	23, 45, 93, 106	0
2	B	2/5 (40%)	3.07	2 (100%) 0 0	78, 78, 78, 79	0
3	C	7/8 (87%)	1.08	1 (14%) 4 6	37, 40, 79, 100	0
All	All	560/608 (92%)	0.80	70 (12%) 5 8	23, 45, 93, 106	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	GLY	8.3
1	A	140	ASN	7.8
1	A	145	ASN	7.4
1	A	152	ASN	6.1
1	A	142	ILE	5.8
1	A	153	LYS	5.4
1	A	62	ASP	5.1
1	A	102	LYS	5.1
1	A	87	PRO	5.0
1	A	441	LEU	5.0
1	A	154	PRO	4.8
3	C	2	DA	4.8
1	A	356	ALA	4.7
1	A	23	SER	4.7
1	A	449	ILE	4.7
1	A	151	PHE	4.6
1	A	381	LYS	4.2
1	A	148	ARG	4.1
1	A	144	LYS	4.1
1	A	150	ALA	4.0
1	A	80	GLU	3.8
1	A	143	THR	3.8
1	A	440	ALA	3.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	354	GLN	3.5
1	A	146	ALA	3.5
1	A	149	GLN	3.3
1	A	155	GLY	3.3
1	A	355	TYR	3.2
1	A	3	LYS	3.2
2	B	2	DT	3.1
1	A	22	GLY	3.1
1	A	105	HIS	3.1
1	A	428	ILE	3.1
1	A	19	LYS	3.1
2	B	1	DT	3.0
1	A	63	GLU	3.0
1	A	20	TYR	2.9
1	A	25	TYR	2.9
1	A	32	GLY	2.9
1	A	470	PHE	2.9
1	A	141	GLU	2.9
1	A	430	ARG	2.8
1	A	66	ALA	2.8
1	A	302	LYS	2.8
1	A	131	ASP	2.8
1	A	93	VAL	2.7
1	A	132	ALA	2.7
1	A	33	HIS	2.6
1	A	98	GLN	2.6
1	A	231	GLY	2.5
1	A	112	LEU	2.5
1	A	130	ASP	2.5
1	A	101	GLU	2.4
1	A	133	ARG	2.4
1	A	200	ALA	2.4
1	A	104	ASP	2.4
1	A	526	VAL	2.3
1	A	452	ALA	2.3
1	A	82	HIS	2.3
1	A	527	THR	2.2
1	A	556	HIS	2.2
1	A	138	VAL	2.1
1	A	157	LEU	2.1
1	A	417	ALA	2.1
1	A	127	ILE	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	107	TYR	2.1
1	A	199	VAL	2.1
1	A	174	VAL	2.1
1	A	201	VAL	2.0
1	A	318	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	PTR	A	319	16/17	0.82	0.22	-	42,49,52,52	0

6.3 Carbohydrates [i](#)

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	HG	A	596	1/1	0.99	0.15	0.08	34,34,34,34	1
4	HG	C	9	1/1	0.99	0.22	-	51,51,51,51	1
4	HG	B	6	1/1	0.94	0.11	-	52,52,52,52	1

6.5 Other polymers [i](#)

There are no such residues in this entry.