



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:01 PM GMT

PDB ID : 1RR8
Title : Structural Mechanisms of Camptothecin Resistance by Mutations in Human Topoisomerase I
Authors : Chrencik, J.E.; Staker, B.L.; Burgin, A.B.; Stewart, L.; Redinbo, M.R.
Deposited on : 2003-12-08
Resolution : 2.60 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

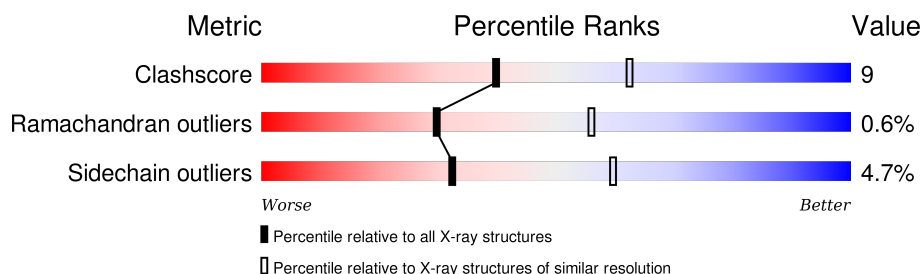
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	22	
2	B	22	
3	C	565	

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
3	PTR	C	723	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5386 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(*AP*AP*AP*AP*AP*GP*AP*CP*TP*T*GP*GP*AP*AP*AP*AP*AP*TP*TP*TP*TP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	22	Total	C	N	O	P	0	0	0
			452	219	87	125	21			

- Molecule 2 is a DNA chain called 5'-D(*AP*AP*AP*AP*AP*TP*TP*TP*TP*TP*CP*CP*AP*AP*GP*TP*CP*TP*TP*TP*TP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	22	Total	C	N	O	P	0	0	0
			443	217	71	134	21			

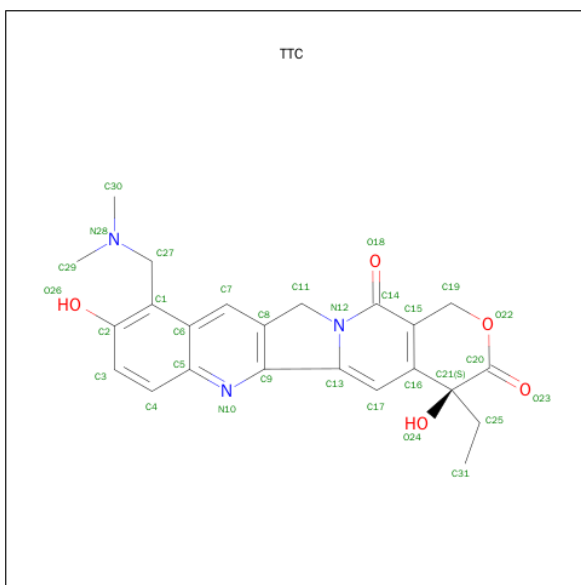
- Molecule 3 is a protein called DNA topoisomerase I.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	494	Total	C	N	O	P	S	0	0	0
			4124	2630	719	752	1	22			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	201	ALA	-	CLONING ARTIFACT	UNP P11387
C	202	ALA	-	CLONING ARTIFACT	UNP P11387
C	361	SER	PHE	ENGINEERED	UNP P11387
C	634	GLN	ARG	ENGINEERED	UNP P11387
C	723	PTR	TYR	MODIFIED RESIDUE	UNP P11387

- Molecule 4 is 2-(1-DIMETHYLAMINOMETHYL-2-HYDROXY-8-HYDROXYMETHYL-9-OXO-9,11-DIHYDRO-INDOLIZINO[1,2-B]QUINOLIN-7-YL)-2-HYDROXY-BUTYRIC ACID (three-letter code: TTC, TTG) (formula: C₂₃H₂₃N₃O₅, C₂₃H₂₅N₃O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			63	46	6	11		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	35	Total	O	0	0
			35	35		
5	B	32	Total	O	0	0
			32	32		
5	C	237	Total	O	0	0
			237	237		

3 Residue-property plots

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.

Note EDS was not executed.

- Molecule 1: 5'-D(*AP*AP*AP*AP*AP*GP*AP*CP*TP*T*GP*GP*AP*AP*AP*AP*TP*TP*TP*TP*T)-3'

Chain A: 



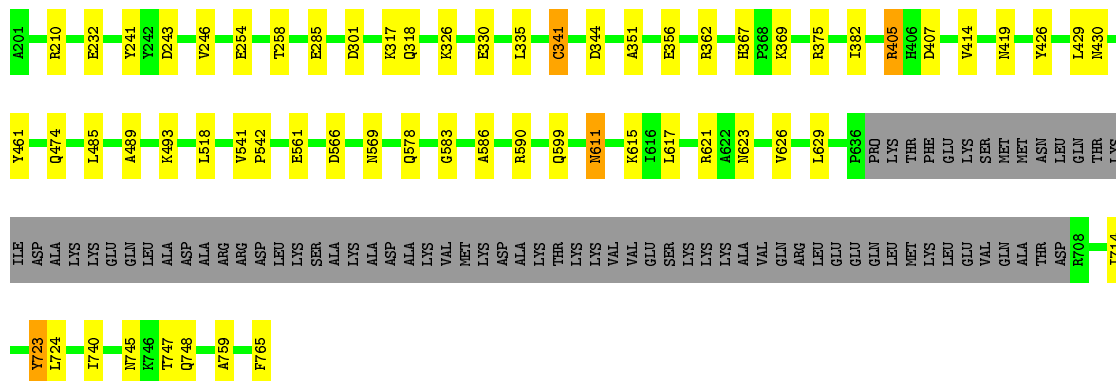
- Molecule 2: 5'-D(*AP*AP*AP*AP*AP*TP*TP*TP*TP*TP*CP*CP*AP*AP*GP*TP*CP*TP*TP*TP*TP*T)-3'

Chain B: 



- Molecule 3: DNA topoisomerase I

Chain C: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.59 Å 115.58 Å 75.01 Å 90.00° 96.29° 90.00°	Depositor
Resolution (Å)	19.70 – 2.60	Depositor
% Data completeness (in resolution range)	10.0 (19.70-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.242 , 0.294	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5386	wwPDB-VP
Average B, all atoms (Å ²)	48.0	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: TTC, TTG, 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	7.07	1/508 (0.2%)	2.15	3/780 (0.4%)
2	B	0.49	0/494	0.79	0/760
3	C	0.66	0/4204	0.72	2/5651 (0.0%)
All	All	2.29	1/5206 (0.0%)	0.99	5/7191 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
3	C	0	2
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	11	DG	P-O5'	159.03	3.18	1.59

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	DG	P-O5'-C5'	-54.29	34.04	120.90
1	A	11	DG	O5'-P-OP2	-14.51	92.64	105.70
3	C	335	LEU	CA-CB-CG	6.03	129.16	115.30
1	A	11	DG	O5'-P-OP1	-5.84	100.44	105.70
3	C	617	LEU	CA-CB-CG	5.79	128.61	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	DA	Sidechain
1	A	9	DT	Sidechain
3	C	723	PTR	Mainchain,Peptide

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	A	452	0	251	36	0
2	B	443	0	255	22	0
3	C	4124	0	4113	30	0
4	A	63	0	40	2	0
5	A	35	0	0	6	0
5	B	32	0	0	3	0
5	C	237	0	0	2	0
All	All	5386	0	4659	85	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 9.

All (85) 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:C:723:PTR:CA	3:C:723:PTR:C	1.76	1.60
1:A:11:DG:OP1	5:A:886:HOH:O	1.67	1.09
1:A:11:DG:P	5:A:886:HOH:O	2.19	0.99
1:A:8:DC:H2''	1:A:9:DT:H5''	1.42	0.99
1:A:17:DA:H2''	1:A:18:DT:H5''	1.50	0.93
3:C:599:GLN:HE22	3:C:765:PHE:H	1.13	0.89
1:A:10:DT:C3'	3:C:723:PTR:O3P	2.24	0.86
1:A:18:DT:H4'	5:A:770:HOH:O	1.83	0.79
1:A:17:DA:C2'	1:A:18:DT:H5''	2.12	0.79
1:A:11:DG:OP1	1:A:11:DG:C8	2.40	0.75
2:B:110:DT:H1'	2:B:111:DC:H5''	1.69	0.72
2:B:120:DT:H6	2:B:120:DT:H5'	1.53	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:DG:H2'	1:A:12:DG:N7	2.06	0.71
1:A:20:DT:C2'	1:A:21:DT:H72	2.24	0.68
1:A:17:DA:H2''	1:A:18:DT:C5'	2.24	0.67
2:B:120:DT:C6	2:B:120:DT:H5'	2.31	0.65
1:A:11:DG:H2'	1:A:12:DG:C8	2.32	0.65
3:C:626:VAL:HG11	3:C:724:LEU:HD21	1.80	0.64
1:A:20:DT:H2''	1:A:21:DT:C7	2.30	0.61
3:C:367:HIS:HD2	3:C:369:LYS:H	1.49	0.60
2:B:112:DC:H1'	5:B:945:HOH:O	2.00	0.60
3:C:375:ARG:H	3:C:419:ASN:HD21	1.48	0.60
1:A:10:DT:H2''	4:A:1991:TTG:O19	2.01	0.60
1:A:8:DC:C2'	1:A:9:DT:H5''	2.24	0.59
2:B:119:DT:H1'	2:B:120:DT:H5''	1.83	0.58
2:B:122:DT:H1'	5:B:846:HOH:O	2.04	0.57
3:C:615:LYS:NZ	5:C:807:HOH:O	2.37	0.57
3:C:382:ILE:HG23	3:C:414:VAL:HG13	1.86	0.57
2:B:121:DT:H3'	5:B:1017:HOH:O	2.05	0.57
1:A:11:DG:C2'	1:A:12:DG:C8	2.90	0.55
3:C:611:ASN:H	3:C:611:ASN:ND2	2.04	0.55
1:A:20:DT:H2''	1:A:21:DT:H72	1.87	0.55
3:C:254:GLU:O	3:C:258:THR:HG23	2.07	0.54
2:B:115:DG:H2'	2:B:116:DT:H72	1.88	0.54
3:C:723:PTR:N	3:C:723:PTR:C	2.61	0.54
1:A:2:DA:H1'	1:A:3:DA:C8	2.42	0.54
2:B:115:DG:H2'	2:B:116:DT:C7	2.38	0.53
3:C:405:ARG:HD2	3:C:407:ASP:HB2	1.90	0.53
1:A:11:DG:H8	1:A:11:DG:OP1	1.88	0.52
3:C:241:TYR:HB2	3:C:301:ASP:HB3	1.91	0.52
3:C:723:PTR:CB	3:C:723:PTR:C	2.79	0.52
1:A:1:DA:H2'	1:A:2:DA:C8	2.46	0.50
1:A:20:DT:H2'	1:A:21:DT:H72	1.95	0.49
2:B:106:DT:H6	2:B:106:DT:H5'	1.77	0.49
1:A:2:DA:H4'	1:A:3:DA:OP1	2.12	0.49
2:B:110:DT:H2''	2:B:111:DC:C5'	2.44	0.48
1:A:11:DG:P	4:A:1991:TTG:H191	2.54	0.48
1:A:17:DA:H1'	1:A:18:DT:H5''	1.96	0.48
3:C:474:GLN:OE1	3:C:566:ASP:O	2.31	0.48
1:A:1:DA:H2''	1:A:2:DA:O5'	2.14	0.47
2:B:105:DA:H1'	2:B:106:DT:H5''	1.96	0.47
1:A:3:DA:C2	1:A:4:DA:C4	3.03	0.47
3:C:341:CYS:SG	3:C:429:LEU:HD21	2.54	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:110:DT:H2''	2:B:111:DC:H5'	1.96	0.47
3:C:326:LYS:O	3:C:330:GLU:HG2	2.15	0.46
3:C:578:GLN:HE22	3:C:583:GLY:H	1.62	0.46
1:A:17:DA:C1'	1:A:18:DT:H5''	2.45	0.46
2:B:104:DA:H2''	2:B:105:DA:OP2	2.16	0.46
2:B:108:DT:H2''	2:B:109:DT:H72	1.97	0.45
3:C:375:ARG:H	3:C:419:ASN:ND2	2.13	0.45
3:C:723:PTR:O	3:C:723:PTR:CA	2.51	0.45
1:A:4:DA:H1'	1:A:5:DA:H5''	1.98	0.45
2:B:108:DT:C2'	2:B:109:DT:H72	2.47	0.45
1:A:21:DT:H5''	5:A:1033:HOH:O	2.17	0.45
2:B:120:DT:H2'	2:B:121:DT:H72	1.99	0.44
1:A:15:DA:C8	1:A:15:DA:H5'	2.52	0.44
3:C:489:ALA:HB2	3:C:586:ALA:HB3	1.99	0.44
3:C:599:GLN:NE2	3:C:765:PHE:H	1.96	0.44
2:B:105:DA:H2''	2:B:106:DT:OP2	2.18	0.43
1:A:1:DA:H2'	1:A:2:DA:H8	1.82	0.43
3:C:246:VAL:HG23	5:C:936:HOH:O	2.17	0.43
2:B:105:DA:OP2	2:B:105:DA:H8	2.01	0.43
1:A:14:DA:H5'	5:A:1029:HOH:O	2.17	0.43
2:B:119:DT:H2''	2:B:120:DT:OP2	2.18	0.43
1:A:10:DT:C3'	3:C:723:PTR:P	3.08	0.42
3:C:541:VAL:HA	3:C:542:PRO:HD3	1.88	0.42
2:B:108:DT:H2''	2:B:109:DT:C7	2.50	0.41
1:A:7:DA:H8	3:C:426:TYR:OH	2.03	0.41
3:C:317:LYS:HG3	3:C:318:GLN:HE21	1.86	0.41
3:C:590:ARG:NH1	3:C:723:PTR:O1P	2.54	0.41
2:B:121:DT:H2'	2:B:122:DT:H72	2.03	0.41
3:C:485:LEU:HD21	3:C:541:VAL:HG21	2.03	0.41
3:C:745:ASN:H	3:C:748:GLN:HE21	1.68	0.40
1:A:15:DA:H8	1:A:15:DA:H5'	1.87	0.40
1:A:3:DA:H1'	5:A:836:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	C	489/565 (86%)	469 (96%)	17 (4%)	3 (1%)	30 56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	344	ASP
3	C	351	ALA
3	C	759	ALA

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	C	444/506 (88%)	423 (95%)	21 (5%)	32 59

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

Mol	Chain	Res	Type
3	C	210	ARG
3	C	232	GLU
3	C	243	ASP
3	C	285	GLU
3	C	341	CYS
3	C	356	GLU
3	C	362	ARG
3	C	405	ARG
3	C	430	ASN
3	C	461	TYR
3	C	493	LYS
3	C	518	LEU
3	C	561	GLU
3	C	569	ASN

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Mol	Chain	Res	Type
3	C	611	ASN
3	C	621	ARG
3	C	623	ASN
3	C	629	LEU
3	C	714	ILE
3	C	740	ILE
3	C	747	THR

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

Mol	Chain	Res	Type
3	C	277	ASN
3	C	318	GLN
3	C	331	ASN
3	C	366	ASN
3	C	367	HIS
3	C	419	ASN
3	C	421	GLN
3	C	430	ASN
3	C	442	GLN
3	C	459	ASN
3	C	569	ASN
3	C	578	GLN
3	C	593	ASN
3	C	599	GLN
3	C	620	ASN
3	C	631	ASN
3	C	748	GLN

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	$\# Z > 2$
3	PTR	C	723	3	14,16,17	1.34	1 (7%)	18,22,24	1.85	2 (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
3	PTR	C	723	3	-	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)
3	C	723	PTR	O-C	2.80	1.32	1.19

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	723	PTR	O-C-CA	-6.71	108.01	125.49
3	C	723	PTR	O2P-P-O1P	2.46	118.49	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	723	PTR	7	0

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

2 ligands 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	TTC	A	100	4	31,35,35	2.39	11 (35%)	26,55,55	2.52	4 (15%)
4	TTG	A	1991	4	27,35,35	2.12	10 (37%)	27,54,54	2.57	6 (22%)

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	TTC	A	100	4	-	0/7/32/32	0/3/5/5
4	TTG	A	1991	4	-	0/15/29/29	0/3/4/4

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1991	TTG	C13-C9	-3.89	1.39	1.47
4	A	100	TTC	C13-C9	-3.24	1.40	1.47
4	A	100	TTC	O22-C20	-3.12	1.30	1.34
4	A	1991	TTG	C4-C5	-2.95	1.36	1.41
4	A	1991	TTG	O24-C21	-2.14	1.41	1.43
4	A	100	TTC	O23-C20	2.01	1.25	1.20
4	A	1991	TTG	C5-N10	2.01	1.41	1.37
4	A	100	TTC	C4-C3	2.29	1.41	1.36
4	A	100	TTC	C7-C6	2.31	1.46	1.42
4	A	1991	TTG	C4-C3	2.35	1.41	1.36
4	A	100	TTC	C21-C20	2.36	1.58	1.53
4	A	1991	TTG	C17-C16	2.49	1.43	1.39
4	A	1991	TTG	C19-C15	2.81	1.55	1.51
4	A	100	TTC	C5-N10	2.82	1.42	1.37
4	A	100	TTC	O22-C19	3.22	1.49	1.45
4	A	100	TTC	C9-C8	3.26	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1991	TTG	C9-C8	3.29	1.47	1.40
4	A	1991	TTG	C27-C1	4.13	1.54	1.51
4	A	1991	TTG	C2-C1	4.79	1.41	1.38
4	A	100	TTC	C27-C1	6.02	1.55	1.51
4	A	100	TTC	C2-C1	6.73	1.43	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1991	TTG	C25-C21-C16	-2.82	107.95	111.33
4	A	100	TTC	C19-C15-C16	-2.74	116.84	120.45
4	A	1991	TTG	C19-C15-C14	-2.67	116.98	121.23
4	A	1991	TTG	C3-C2-C1	-2.53	120.10	121.40
4	A	1991	TTG	C9-C13-N12	-2.38	109.97	113.62
4	A	1991	TTG	C17-C16-C15	-2.33	113.10	118.62
4	A	100	TTC	C3-C2-C1	-2.20	120.27	121.40
4	A	100	TTC	C17-C13-C9	2.05	128.24	124.67
4	A	100	TTC	C13-C9-N10	10.91	130.38	117.79
4	A	1991	TTG	C13-C9-N10	11.18	130.69	117.79

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
4	A	1991	TTG	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.