Full wwPDB X-ray Structure Validation Report (i)

Feb 18, 2024 – 08:15 AM EST

PDB ID	:	4DTX
Title	:	RB69 DNA Polymerase Ternary Complex with dTTP Opposite an Abasic Site and ddC/dG as the Penultimate Base-pair
Authors	:	Xia, S.; Wang, J.; Konigsberg, W.H.
Deposited on	:	2012-02-21
Resolution	:	1.84 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 (i) 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 (i)) were used in the production of this report:

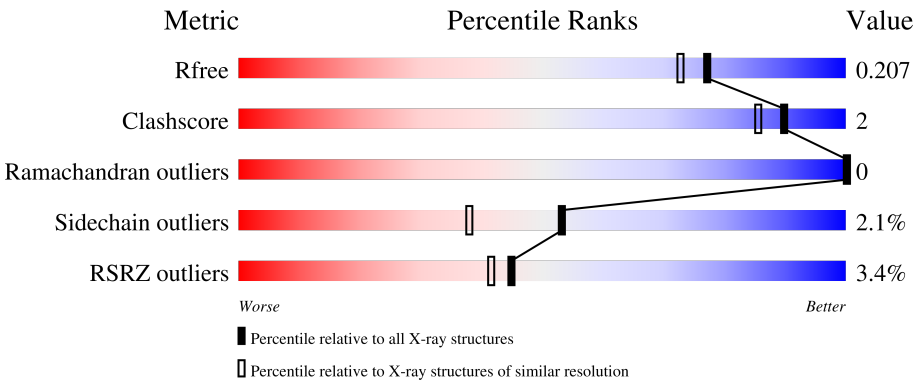
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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}	130704	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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 <=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	903	3%, 94%, 6%
2	T	18	6%, 67%, 22%, 11%
3	P	13	62%, 38%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 8867 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 polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	901	7407	4759	1234	1380	34	0	13	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	ALA	ASP	engineered mutation	UNP Q38087
A	327	ALA	ASP	engineered mutation	UNP Q38087
A	561	ALA	LEU	engineered mutation	UNP Q38087
A	565	GLY	SER	engineered mutation	UNP Q38087
A	567	ALA	TYR	engineered mutation	UNP Q38087

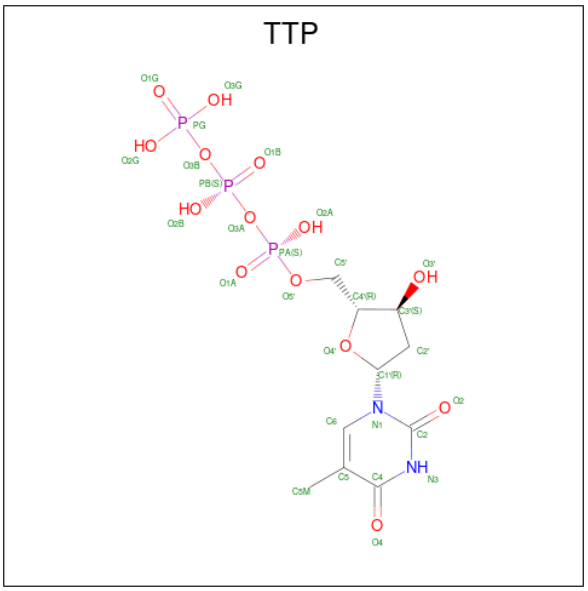
- Molecule 2 is a DNA chain called DNA tempalte.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	T	18	358	170	66	105	17	0	0	0

- Molecule 3 is a DNA chain called DNA primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	P	13	262	126	48	76	12	0	0	0

- Molecule 4 is THYMIDINE-5'-TRIPHOSPHATE (three-letter code: TTP) (formula: $C_{10}H_{17}N_2O_{14}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	29	10	2	14	3	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
5	A	4	4	4	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	711	721	721	0	10
			Total	O		
6	T	58	59	59	0	1
			Total	O		
6	P	26	27	27	0	1

3 Residue-property plots (i)

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: DNA polymerase

Chain A: 3% 94% 6%

M1 D21 R27 K34 E43 S44 Q45 A46 R66 L90 N98 N112 D121 H131 Y156 L170 E177 W216 N217 A222 V252 I253 E254 N255 M256 Y257 K302 L303 K304 I309 R330 V331 L332 Q333 S360 K363 R403

L415 D452 L508 S509 V510 L514 W574 L580 L605 V608 C609 G610 D623 K640 W647 D664 R668 L693 L760 I779 K816 G817 N818 I819 D820 A821 L833 P834 L835 R836 E837 W848 I854 L857 V862 W865 M866

V870 L871 L872 E873 K874 L889 D890 Y891 K894 A895 S896 L897 F898 D899 M900 F901 ASP PHE

• Molecule 2: DNA tempalte

Chain T: 6% 67% 22% 11%

T1 N4 G5 T6 G12 T13 C14 C15 G16 C17 G18

• Molecule 3: DNA primer

Chain P: 62% 38%

G103 C104 T112 T113 A114 C115

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	75.41Å 120.21Å 131.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.38 – 1.84 35.38 – 1.84	Depositor EDS
% Data completeness (in resolution range)	96.0 (35.38-1.84) 96.0 (35.38-1.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.23 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.170 , 0.206 0.170 , 0.207	Depositor DCC
R_{free} test set	4993 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	27.4	Xtriage
Anisotropy	0.148	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 44.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8867	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: TTP, DOC, CA, 3DR

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	0.49	5/7625 (0.1%)	0.54	1/10303 (0.0%)
2	T	0.33	0/388	0.85	2/595 (0.3%)
3	P	0.37	0/273	0.91	1/420 (0.2%)
All	All	0.48	5/8286 (0.1%)	0.58	4/11318 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	216	TRP	CD2-CE2	5.20	1.47	1.41
1	A	647	TRP	CD2-CE2	5.19	1.47	1.41
1	A	848	TRP	CD2-CE2	5.10	1.47	1.41
1	A	574	TRP	CD2-CE2	5.08	1.47	1.41
1	A	865	TRP	CD2-CE2	5.01	1.47	1.41

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	P	112	DT	O3'-P-O5'	-5.93	92.73	104.00
2	T	6	DT	P-O3'-C3'	5.82	126.68	119.70
2	T	15	DC	P-O3'-C3'	5.19	125.93	119.70
1	A	760	LEU	CA-CB-CG	5.12	127.09	115.30

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	7407	0	7345	25	0
2	T	358	0	200	3	0
3	P	262	0	148	2	0
4	A	29	0	13	0	0
5	A	4	0	0	0	0
6	A	721	0	0	5	0
6	P	27	0	0	0	0
6	T	59	0	0	0	0
All	All	8867	0	7706	29	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 2.

All (29) 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:330[B]:ARG:HH11	1:A:333:GLN:HE22	1.25	0.84
1:A:403:ARG:NH2	1:A:889[A]:LEU:HD13	2.08	0.69
1:A:360:SER:OG	1:A:363:LYS:HB3	1.93	0.67
1:A:415:LEU:HD22	1:A:623:ASP:HB3	1.78	0.66
1:A:779:ILE:HB	1:A:871:LEU:HD21	1.81	0.62
1:A:873:GLU:HG2	6:A:1319:HOH:O	2.03	0.59
1:A:821:ALA:HA	6:A:1201:HOH:O	2.04	0.57
1:A:98:ASN:HB3	6:A:1378:HOH:O	2.04	0.57
1:A:664:ASP:O	1:A:668[B]:ARG:HG3	2.07	0.54
1:A:170:LEU:HA	1:A:177:GLU:HG3	1.90	0.53
1:A:605:LEU:HA	1:A:608:VAL:HG22	1.92	0.51
1:A:131:HIS:HD2	1:A:156:TYR:OH	1.94	0.50
1:A:870:VAL:HG13	1:A:874:LYS:HD3	1.96	0.48
1:A:760:LEU:HD13	1:A:891:TYR:HA	1.96	0.46
1:A:403:ARG:HH22	1:A:889[A]:LEU:HD13	1.80	0.46
1:A:112[B]:ASN:HD22	1:A:332:LEU:HG	1.82	0.45
3:P:103:DG:H2"	3:P:104:DC:C6	2.52	0.45
1:A:360:SER:HB2	2:T:4:3DR:OP2	2.17	0.44
1:A:854:ILE:HD13	1:A:862:VAL:HG21	2.00	0.42
1:A:112[A]:ASN:ND2	6:A:1118:HOH:O	2.50	0.42
3:P:113:DT:H2"	3:P:114:DA:H5'	2.02	0.42
1:A:27:ARG:HG3	6:A:1470:HOH:O	2.18	0.42
2:T:15:DC:H2"	2:T:16:DG:C8	2.55	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:LEU:HD11	1:A:363:LYS:HD2	2.02	0.41
1:A:21:ASP:OD2	1:A:27:ARG:NE	2.54	0.41
1:A:862:VAL:O	1:A:866:MET:HG3	2.20	0.41
2:T:12:DG:H2'	2:T:13:DT:H71	2.02	0.41
1:A:833:LEU:HD13	1:A:866:MET:HG2	2.03	0.40
1:A:415:LEU:HD22	1:A:623:ASP:CB	2.48	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	912/903 (101%)	893 (98%)	19 (2%)	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	808/797 (101%)	789 (98%)	19 (2%)	49	32

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ARG
1	A	34	LYS
1	A	217[A]	ASN
1	A	217[B]	ASN
1	A	256	MET
1	A	304	LYS
1	A	309	ILE
1	A	452	ASP
1	A	514	LEU
1	A	580	LEU
1	A	640	LYS
1	A	693	LEU
1	A	760	LEU
1	A	835	LEU
1	A	837	GLU
1	A	889[A]	LEU
1	A	889[B]	LEU
1	A	894	LYS
1	A	900	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	HIS
1	A	203	ASN
1	A	333	GLN
1	A	339	GLN
1	A	546	GLN
1	A	761	GLN
1	A	773	GLN
1	A	823	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
2	3DR	T	4	2	8,11,12	0.51	0	9,14,17	1.37	1 (11%)
3	DOC	P	115	3,2	16,19,20	0.43	0	20,26,29	0.53	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
2	3DR	T	4	2	-	0/3/15/16	0/1/1/1
3	DOC	P	115	3,2	-	0/7/18/19	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	T	4	3DR	O4'-C4'-C3'	2.71	107.71	103.73

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
2	T	4	3DR	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 4 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	$\#\|Z\| > 2$
4	TTP	A	1001	5	26,30,30	1.23	4 (15%)	39,47,47	1.93	9 (23%)

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	TTP	A	1001	5	-	4/22/34/34	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1001	TTP	C6-C5	2.86	1.39	1.34
4	A	1001	TTP	C4-C5	2.59	1.49	1.44
4	A	1001	TTP	C4-N3	-2.48	1.34	1.38
4	A	1001	TTP	C2-N1	2.28	1.42	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	1001	TTP	C4-N3-C2	-4.93	120.97	127.35
4	A	1001	TTP	N3-C2-N1	4.86	121.34	114.89
4	A	1001	TTP	C5-C4-N3	4.51	119.16	115.31
4	A	1001	TTP	O4-C4-C5	-3.38	120.99	124.90
4	A	1001	TTP	C5-C6-N1	-3.31	119.93	123.34
4	A	1001	TTP	C5M-C5-C4	2.97	122.03	118.77
4	A	1001	TTP	O2-C2-N1	-2.65	119.26	122.79
4	A	1001	TTP	C2'-C1'-N1	-2.64	107.70	113.77
4	A	1001	TTP	C5M-C5-C6	-2.36	119.70	122.85

There are no chirality outliers.

All (4) torsion outliers are listed below:

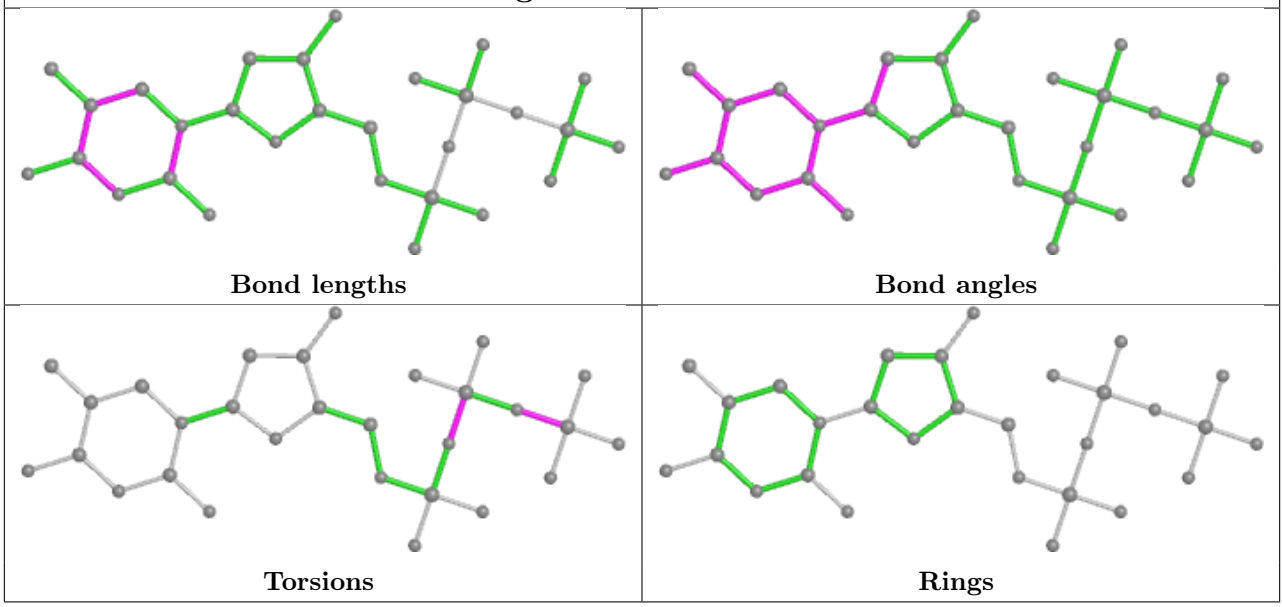
Mol	Chain	Res	Type	Atoms
4	A	1001	TTP	PB-O3B-PG-O1G
4	A	1001	TTP	PA-O3A-PB-O2B
4	A	1001	TTP	PB-O3B-PG-O2G
4	A	1001	TTP	PB-O3B-PG-O3G

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand TTP A 1001



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)

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	901/903 (99%)	-0.16	31 (3%) 45 41	18, 28, 53, 92	1 (0%)
2	T	17/18 (94%)	-0.49	1 (5%) 22 20	22, 31, 47, 58	0
3	P	12/13 (92%)	-0.30	0 100 100	22, 40, 60, 61	0
All	All	930/934 (99%)	-0.17	32 (3%) 45 41	18, 28, 54, 92	1 (0%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	819	ILE	8.7
1	A	44	SER	5.9
1	A	256	MET	5.4
1	A	817	GLY	5.4
1	A	257	TYR	5.2
1	A	897	LEU	4.5
1	A	857	LEU	4.2
1	A	508	LEU	4.1
1	A	895	ALA	3.6
1	A	43	GLU	3.6
1	A	46	ALA	3.3
1	A	816	LYS	3.3
1	A	510	VAL	3.3
1	A	121	ASP	3.0
1	A	255	ASN	2.9
1	A	303	LEU	2.9
1	A	252	VAL	2.9
2	T	1	DT	2.8
1	A	66	ARG	2.7
1	A	898	PHE	2.7
1	A	818[A]	ASN	2.6
1	A	899	ASP	2.5
1	A	901	PHE	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	254	GLU	2.5
1	A	222	ALA	2.4
1	A	302	LYS	2.4
1	A	608	VAL	2.3
1	A	820	ASP	2.2
1	A	610	GLY	2.0
1	A	894	LYS	2.0
1	A	45	GLN	2.0
1	A	896	SER	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. 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(Å^2)	Q<0.9
3	DOC	P	115	18/19	0.97	0.09	21,23,29,30	0
2	3DR	T	4	11/12	0.97	0.09	28,32,37,43	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

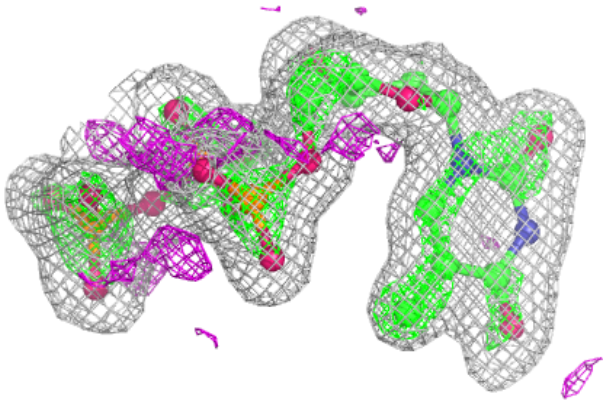
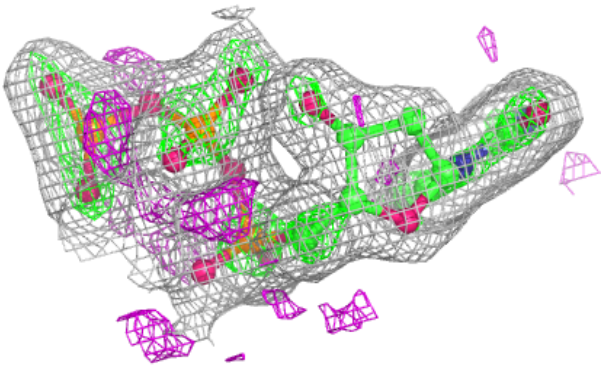
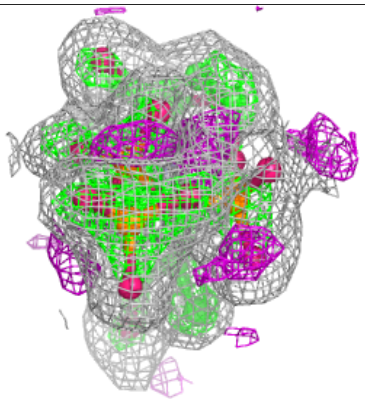
6.4 Ligands (i)

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(Å^2)	Q<0.9
5	CA	A	1003	1/1	0.95	0.06	49,49,49,49	0
5	CA	A	1005	1/1	0.96	0.04	46,46,46,46	0
4	TTP	A	1001	29/29	0.97	0.20	30,34,39,41	0
5	CA	A	1004	1/1	0.98	0.04	38,38,38,38	0
5	CA	A	1002	1/1	0.99	0.08	23,23,23,23	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around TTP A 1001: $2mF_o$-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)	
	

6.5 Other polymers (i)

There are no such residues in this entry.