



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:34 AM GMT

PDB ID : 2AQ4  
Title : Ternary complex of the catalytic core of REV1 with DNA and dCTP.  
Authors : Nair, D.T.; Johnson, R.E.; Prakash, L.; Prakash, S.; Aggarwal, A.K.  
Deposited on : 2005-08-17  
Resolution : 2.32 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](mailto: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.

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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

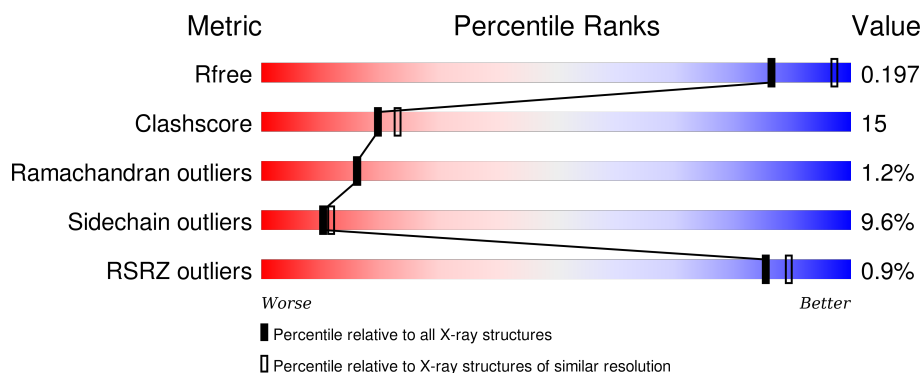
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.32 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	4425 (2.34-2.30)
Clashscore	102246	5057 (2.34-2.30)
Ramachandran outliers	100387	5008 (2.34-2.30)
Sidechain outliers	100360	5007 (2.34-2.30)
RSRZ outliers	91569	4432 (2.34-2.30)

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  $\geq 3$ , 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	P	12	
2	T	16	
3	A	434	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4117 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\*TP\*CP\*CP\*TP\*CP\*CP\*CP\*CP\*TP\*AP\*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	P	12	Total	C	N	O	P	0	0	0
			231	113	37	70	11			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	16	Total	C	N	O	P	0	0	0
			338	160	71	92	15			

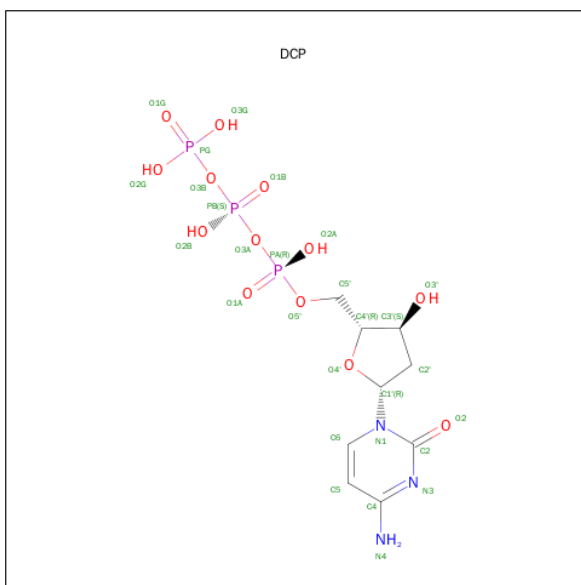
- Molecule 3 is a protein called DNA repair protein REV1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	428	Total	C	N	O	S	0	0	0
			3378	2151	577	627	23			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

- Molecule 5 is 2'-DEOXYCYTIDINE-5'-TRIPHOSPHATE (three-letter code: DCP) (formula: C<sub>9</sub>H<sub>16</sub>N<sub>3</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			28	9	3	13	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	120	Total O 120 120	0	0
6	P	6	Total O 6 6	0	0
6	T	14	Total O 14 14	0	0

### 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.

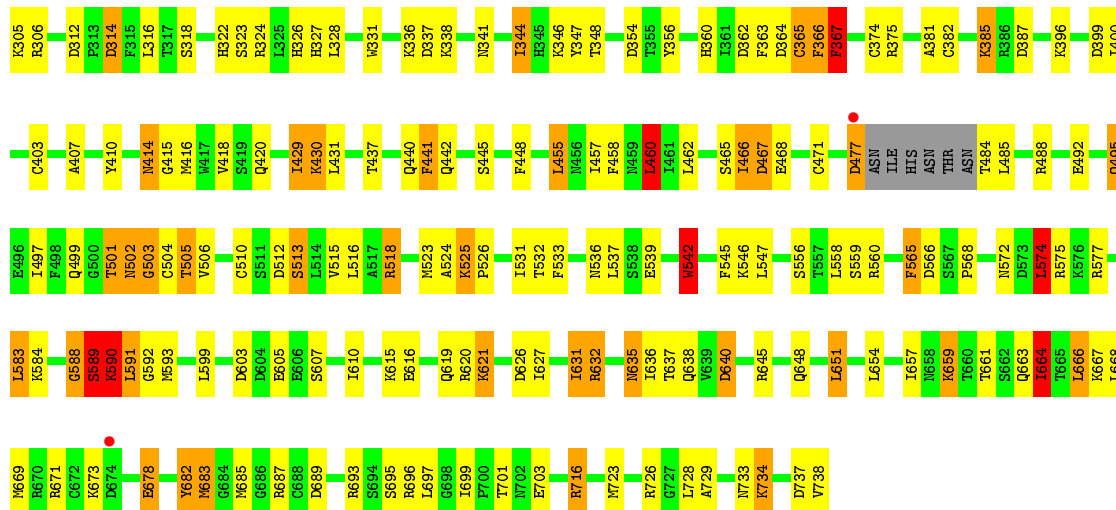
- Molecule 1: 5'-D(\*AP\*TP\*CP\*CP\*TP\*CP\*CP\*CP\*CP\*TP\*AP\*C)-3'



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



- Molecule 3: DNA repair protein REV1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	177.32Å 200.91Å 55.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.32 41.58 – 2.32	Depositor EDS
% Data completeness (in resolution range)	98.5 (50.00-2.32) 98.5 (41.58-2.32)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.13 (at 2.32Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.197 , 0.238 0.200 , 0.197	Depositor DCC
$R_{free}$ test set	2157 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.5	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 35.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 42757 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4117	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: DOC, MG, DCP

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	P	2.56	17/236 (7.2%)	3.44	52/360 (14.4%)
2	T	2.79	27/382 (7.1%)	3.85	80/591 (13.5%)
3	A	1.59	40/3445 (1.2%)	1.50	63/4652 (1.4%)
All	All	1.81	84/4063 (2.1%)	2.05	195/5603 (3.5%)

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
3	A	0	3

All (84) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	T	2	DA	N3-C4	-11.96	1.27	1.34
2	T	2	DA	C8-N7	10.30	1.38	1.31
2	T	2	DA	C6-N6	-10.10	1.25	1.33
2	T	6	DT	C5-C7	10.00	1.56	1.50
3	A	492	GLU	CD-OE1	9.96	1.36	1.25
2	T	2	DA	N9-C8	9.71	1.45	1.37
2	T	2	DA	C3'-O3'	-9.33	1.31	1.44
2	T	4	DG	O5'-C5'	-8.99	1.19	1.42
3	A	546	LYS	CD-CE	8.91	1.73	1.51
1	P	10	DT	C5-C7	8.67	1.55	1.50
2	T	2	DA	N9-C4	-8.27	1.32	1.37
3	A	385	LYS	CD-CE	8.19	1.71	1.51
2	T	7	DA	N7-C5	-8.05	1.34	1.39
2	T	2	DA	C5-C4	8.04	1.44	1.38
1	P	11	DA	N9-C4	-7.88	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	415	GLY	N-CA	7.68	1.57	1.46
3	A	492	GLU	CD-OE2	7.55	1.33	1.25
1	P	5	DT	C5-C6	7.52	1.39	1.34
2	T	2	DA	C5-C6	-7.20	1.34	1.41
2	T	10	DG	N7-C5	7.14	1.43	1.39
3	A	683	MET	SD-CE	7.14	2.17	1.77
2	T	6	DT	N1-C6	-7.14	1.33	1.38
2	T	13	DG	N1-C2	7.12	1.43	1.37
1	P	5	DT	C3'-O3'	-7.12	1.34	1.44
1	P	7	DC	C5-C6	6.93	1.39	1.34
3	A	338	LYS	CE-NZ	6.83	1.66	1.49
2	T	6	DT	O5'-C5'	-6.82	1.25	1.42
2	T	6	DT	C1'-N1	6.74	1.58	1.49
3	A	367	PHE	CB-CG	6.68	1.62	1.51
3	A	488	ARG	CG-CD	6.63	1.68	1.51
3	A	542	TRP	CG-CD1	6.61	1.46	1.36
3	A	418	VAL	CB-CG2	-6.44	1.39	1.52
3	A	546	LYS	CE-NZ	6.41	1.65	1.49
3	A	367	PHE	N-CA	6.40	1.59	1.46
3	A	678	GLU	CB-CG	6.38	1.64	1.52
1	P	6	DC	C5-C6	6.38	1.39	1.34
3	A	682	TYR	C-O	6.36	1.35	1.23
3	A	305	LYS	CA-CB	6.36	1.68	1.53
3	A	532	THR	N-CA	6.33	1.59	1.46
3	A	465	SER	CB-OG	-6.29	1.34	1.42
2	T	7	DA	C6-N6	-6.29	1.28	1.33
3	A	420	GLN	CG-CD	6.29	1.65	1.51
1	P	9	DC	O5'-C5'	-6.21	1.26	1.42
1	P	11	DA	C3'-O3'	-6.09	1.36	1.44
3	A	515	VAL	CB-CG2	-6.02	1.40	1.52
2	T	1	DT	C3'-O3'	5.99	1.51	1.44
3	A	542	TRP	CB-CG	5.93	1.60	1.50
3	A	695	SER	CB-OG	-5.93	1.34	1.42
3	A	678	GLU	CG-CD	5.89	1.60	1.51
1	P	9	DC	P-O5'	5.87	1.65	1.59
1	P	2	DT	C5-C7	5.83	1.53	1.50
2	T	10	DG	C2-N2	-5.80	1.28	1.34
2	T	10	DG	N1-C2	-5.79	1.33	1.37
2	T	1	DT	C5-C7	5.78	1.53	1.50
1	P	10	DT	P-O5'	5.77	1.65	1.59
3	A	441	PHE	CD1-CE1	5.63	1.50	1.39
3	A	659	LYS	CE-NZ	5.61	1.63	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	682	TYR	CG-CD1	5.60	1.46	1.39
3	A	545	PHE	CE2-CZ	5.57	1.48	1.37
3	A	366	PHE	CG-CD1	5.53	1.47	1.38
2	T	4	DG	C4'-C3'	-5.53	1.47	1.52
1	P	9	DC	N3-C4	5.47	1.37	1.33
2	T	5	DG	C5-C4	-5.47	1.34	1.38
3	A	565	PHE	CE2-CZ	5.39	1.47	1.37
3	A	590	LYS	N-CA	5.39	1.57	1.46
3	A	620	ARG	CZ-NH1	5.37	1.40	1.33
1	P	5	DT	N1-C6	5.28	1.42	1.38
2	T	9	DG	C2-N3	5.26	1.36	1.32
1	P	7	DC	O3'-P	5.25	1.67	1.61
3	A	589	SER	N-CA	5.24	1.56	1.46
3	A	430	LYS	CB-CG	5.23	1.66	1.52
2	T	4	DG	O4'-C1'	-5.21	1.35	1.42
3	A	659	LYS	CD-CE	5.21	1.64	1.51
1	P	5	DT	N1-C2	5.20	1.42	1.38
3	A	575	ARG	CB-CG	-5.14	1.38	1.52
2	T	2	DA	C3'-C2'	5.11	1.58	1.52
3	A	593	MET	CG-SD	5.08	1.94	1.81
1	P	11	DA	C5-C4	5.08	1.42	1.38
2	T	8	DG	N7-C5	-5.05	1.36	1.39
3	A	616	GLU	CD-OE1	5.05	1.31	1.25
3	A	687	ARG	CG-CD	5.05	1.64	1.51
1	P	8	DC	C5-C6	5.04	1.38	1.34
3	A	420	GLN	CD-NE2	5.04	1.45	1.32
3	A	356	TYR	CD2-CE2	5.03	1.47	1.39

All (195) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	2	DA	C5-N7-C8	-27.46	90.17	103.90
2	T	5	DG	O4'-C1'-N9	-21.02	93.29	108.00
2	T	2	DA	N7-C8-N9	19.70	123.65	113.80
2	T	2	DA	C4-C5-N7	18.79	120.09	110.70
1	P	10	DT	O4'-C1'-N1	-15.38	97.23	108.00
2	T	15	DA	O4'-C1'-N9	13.25	117.28	108.00
3	A	632	ARG	NE-CZ-NH1	-13.08	113.76	120.30
2	T	2	DA	C2-N3-C4	-13.05	104.08	110.60
2	T	2	DA	C8-N9-C4	-13.01	100.59	105.80
3	A	632	ARG	NE-CZ-NH2	11.82	126.21	120.30
2	T	2	DA	C6-C5-N7	-11.77	124.06	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	414	ASN	C-N-CA	-10.77	99.69	122.30
1	P	11	DA	C6-N1-C2	10.56	124.94	118.60
1	P	9	DC	C5'-C4'-O4'	-10.17	89.98	109.30
3	A	518	ARG	NE-CZ-NH1	10.05	125.33	120.30
2	T	4	DG	O5'-P-OP1	10.01	122.71	110.70
2	T	4	DG	O4'-C1'-N9	-9.92	101.06	108.00
2	T	6	DT	C5'-C4'-C3'	-9.81	96.44	114.10
1	P	7	DC	O5'-P-OP2	-9.80	96.88	105.70
3	A	512	ASP	CB-CG-OD2	9.78	127.10	118.30
1	P	11	DA	O4'-C1'-N9	-9.70	101.21	108.00
2	T	12	DA	OP1-P-OP2	9.62	134.03	119.60
1	P	11	DA	C5-C6-N1	-9.49	112.95	117.70
1	P	11	DA	C2-N3-C4	-9.43	105.88	110.60
2	T	5	DG	C5-C6-N1	9.36	116.18	111.50
2	T	13	DG	N3-C2-N2	-9.27	113.41	119.90
1	P	7	DC	N3-C4-N4	-9.23	111.54	118.00
3	A	502	ASN	C-N-CA	-9.22	102.93	122.30
1	P	8	DC	C6-N1-C2	-9.18	116.63	120.30
2	T	5	DG	C6-N1-C2	-9.00	119.70	125.10
2	T	9	DG	C6-N1-C2	-8.71	119.87	125.10
2	T	1	DT	O4'-C1'-C2'	-8.69	98.95	105.90
2	T	13	DG	N1-C2-N2	8.65	123.98	116.20
3	A	518	ARG	NE-CZ-NH2	-8.60	116.00	120.30
1	P	7	DC	C5-C4-N4	8.48	126.13	120.20
1	P	5	DT	N3-C2-O2	-8.47	117.22	122.30
3	A	664	ILE	CA-CB-CG2	8.47	127.83	110.90
1	P	2	DT	O4'-C1'-N1	-8.44	102.09	108.00
3	A	620	ARG	NE-CZ-NH2	-8.35	116.12	120.30
2	T	4	DG	C4'-C3'-C2'	-8.33	95.60	103.10
3	A	526	PRO	N-CD-CG	-8.27	90.79	103.20
3	A	728	LEU	CA-CB-CG	8.24	134.26	115.30
2	T	2	DA	O4'-C1'-N9	-8.22	102.25	108.00
3	A	367	PHE	N-CA-CB	8.07	125.13	110.60
2	T	2	DA	N3-C4-C5	8.00	132.40	126.80
3	A	414	ASN	O-C-N	-7.95	109.69	123.20
3	A	626	ASP	CB-CG-OD1	7.94	125.44	118.30
2	T	2	DA	N1-C6-N6	7.92	123.35	118.60
2	T	8	DG	P-O3'-C3'	7.91	129.19	119.70
3	A	455	LEU	C-N-CA	-7.88	102.00	121.70
1	P	8	DC	O5'-P-OP2	-7.86	98.63	105.70
2	T	2	DA	N9-C1'-C2'	7.83	127.47	112.60
1	P	11	DA	N3-C4-C5	7.78	132.25	126.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	2	DA	N1-C2-N3	7.70	133.15	129.30
2	T	7	DA	C8-N9-C4	-7.67	102.73	105.80
1	P	10	DT	C2-N3-C4	-7.67	122.60	127.20
2	T	9	DG	N1-C6-O6	-7.67	115.30	119.90
2	T	15	DA	C8-N9-C4	-7.64	102.74	105.80
1	P	2	DT	P-O3'-C3'	7.61	128.83	119.70
3	A	603	ASP	CB-CG-OD2	7.58	125.12	118.30
2	T	13	DG	O4'-C4'-C3'	-7.55	101.47	106.00
3	A	337	ASP	CB-CG-OD2	7.50	125.05	118.30
2	T	15	DA	N7-C8-N9	7.49	117.55	113.80
3	A	515	VAL	CB-CA-C	-7.44	97.26	111.40
2	T	6	DT	C4-C5-C7	7.43	123.46	119.00
3	A	337	ASP	CB-CG-OD1	-7.37	111.66	118.30
2	T	4	DG	C1'-O4'-C4'	-7.19	102.91	110.10
1	P	11	DA	C5-N7-C8	-7.18	100.31	103.90
3	A	460	LEU	CB-CG-CD1	-7.15	98.84	111.00
1	P	6	DC	O4'-C4'-C3'	-7.13	101.65	104.50
3	A	336	LYS	CD-CE-NZ	-7.09	95.39	111.70
3	A	362	ASP	CB-CG-OD2	7.01	124.61	118.30
2	T	7	DA	C5-C6-N1	7.00	121.20	117.70
2	T	5	DG	O4'-C4'-C3'	-6.98	101.71	104.50
2	T	12	DA	O4'-C1'-N9	-6.98	103.11	108.00
2	T	9	DG	C5-C6-N1	6.97	114.98	111.50
1	P	10	DT	N3-C4-C5	6.96	119.38	115.20
2	T	3	DA	C5-C6-N6	6.94	129.25	123.70
2	T	8	DG	N1-C6-O6	6.89	124.04	119.90
3	A	312	ASP	CB-CG-OD2	6.88	124.50	118.30
3	A	306	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	P	8	DC	N1-C2-O2	-6.88	114.77	118.90
3	A	566	ASP	CB-CG-OD2	6.86	124.47	118.30
2	T	16	DT	N3-C4-O4	-6.83	115.80	119.90
3	A	560	ARG	NE-CZ-NH1	6.82	123.71	120.30
3	A	689	ASP	CB-CG-OD2	6.82	124.44	118.30
3	A	666	LEU	CA-CB-CG	6.81	130.96	115.30
3	A	631	ILE	CG1-CB-CG2	-6.81	96.42	111.40
1	P	5	DT	N1-C2-N3	6.75	118.65	114.60
2	T	10	DG	C2-N3-C4	6.72	115.26	111.90
2	T	2	DA	N3-C4-N9	-6.70	122.04	127.40
3	A	728	LEU	CB-CG-CD1	-6.68	99.64	111.00
3	A	366	PHE	O-C-N	-6.64	112.07	122.70
1	P	5	DT	C4-C5-C7	-6.64	115.02	119.00
2	T	6	DT	C5-C6-N1	6.62	127.67	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	4	DG	P-O5'-C5'	6.61	131.48	120.90
2	T	1	DT	C5-C4-O4	-6.54	120.32	124.90
3	A	523	MET	CG-SD-CE	-6.47	89.84	100.20
1	P	5	DT	C4-C5-C6	6.46	121.88	118.00
1	P	11	DA	C4'-C3'-C2'	6.44	108.90	103.10
3	A	591	LEU	CB-CG-CD2	6.43	121.93	111.00
2	T	16	DT	C4-C5-C7	6.41	122.85	119.00
2	T	1	DT	C6-C5-C7	6.40	126.74	122.90
1	P	2	DT	C5-C6-N1	-6.39	119.86	123.70
1	P	5	DT	C4'-C3'-C2'	6.38	108.84	103.10
3	A	399	ASP	CB-CG-OD2	-6.38	112.56	118.30
2	T	16	DT	C6-C5-C7	-6.38	119.08	122.90
1	P	2	DT	O4'-C1'-C2'	6.37	111.00	105.90
1	P	11	DA	N3-C4-N9	-6.35	122.32	127.40
1	P	1	DA	N1-C6-N6	6.32	122.39	118.60
1	P	11	DA	O4'-C1'-C2'	6.32	110.95	105.90
2	T	15	DA	P-O5'-C5'	6.29	130.96	120.90
1	P	7	DC	C2-N1-C1'	-6.28	111.89	118.80
2	T	4	DG	N9-C4-C5	6.24	107.90	105.40
3	A	532	THR	N-CA-C	6.24	127.84	111.00
2	T	13	DG	OP1-P-OP2	6.21	128.92	119.60
3	A	678	GLU	OE1-CD-OE2	-6.18	115.88	123.30
3	A	316	LEU	CB-CG-CD1	-6.17	100.51	111.00
3	A	467	ASP	CB-CG-OD1	6.14	123.82	118.30
1	P	10	DT	C4'-C3'-C2'	6.10	108.59	103.10
2	T	7	DA	C6-N1-C2	-6.09	114.95	118.60
2	T	3	DA	C5'-C4'-O4'	-6.08	97.75	109.30
2	T	10	DG	C5-C6-N1	6.08	114.54	111.50
2	T	9	DG	N3-C4-C5	-6.06	125.57	128.60
1	P	5	DT	P-O3'-C3'	6.05	126.96	119.70
3	A	546	LYS	CD-CE-NZ	6.04	125.59	111.70
2	T	10	DG	N7-C8-N9	-6.04	110.08	113.10
3	A	492	GLU	OE1-CD-OE2	6.03	130.54	123.30
3	A	526	PRO	CA-N-CD	6.00	120.11	111.70
3	A	590	LYS	N-CA-CB	5.99	121.38	110.60
2	T	16	DT	O4'-C1'-C2'	5.98	110.68	105.90
3	A	416	MET	CG-SD-CE	-5.98	90.64	100.20
3	A	460	LEU	CA-CB-CG	5.95	128.99	115.30
1	P	9	DC	O4'-C1'-C2'	-5.95	101.14	105.90
1	P	5	DT	C5-C6-N1	-5.93	120.14	123.70
2	T	4	DG	C2'-C3'-O3'	-5.92	93.07	112.60
2	T	8	DG	C5-C6-O6	-5.90	125.06	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	4	DG	C5'-C4'-O4'	-5.90	98.10	109.30
1	P	10	DT	O4'-C4'-C3'	-5.86	102.16	104.50
2	T	11	DG	OP1-P-O3'	-5.85	92.33	105.20
2	T	13	DG	N9-C4-C5	5.82	107.73	105.40
3	A	344	ILE	CG1-CB-CG2	-5.81	98.62	111.40
1	P	2	DT	N3-C2-O2	-5.80	118.82	122.30
2	T	5	DG	P-O5'-C5'	-5.80	111.62	120.90
3	A	513	SER	CB-CA-C	-5.77	99.14	110.10
1	P	9	DC	O4'-C4'-C3'	-5.75	102.20	104.50
1	P	6	DC	C6-N1-C2	-5.74	118.00	120.30
3	A	314	ASP	CB-CG-OD2	5.74	123.47	118.30
2	T	2	DA	C5-C6-N6	-5.73	119.11	123.70
3	A	306	ARG	NE-CZ-NH2	-5.72	117.44	120.30
3	A	354	ASP	CB-CG-OD2	5.68	123.41	118.30
3	A	525	LYS	N-CA-C	5.65	126.25	111.00
3	A	651	LEU	CB-CG-CD2	-5.62	101.44	111.00
1	P	7	DC	OP1-P-OP2	5.62	128.02	119.60
2	T	11	DG	C5-N7-C8	-5.57	101.51	104.30
1	P	9	DC	N3-C2-O2	-5.56	118.00	121.90
2	T	5	DG	O3'-P-O5'	-5.56	93.43	104.00
2	T	14	DG	O4'-C1'-C2'	-5.55	101.46	105.90
3	A	671	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	P	9	DC	O4'-C1'-N1	-5.52	104.13	108.00
3	A	328	LEU	CB-CG-CD1	5.52	120.38	111.00
2	T	6	DT	O5'-P-OP2	5.51	117.31	110.70
2	T	15	DA	C5-N7-C8	-5.51	101.15	103.90
1	P	5	DT	O4'-C4'-C3'	-5.47	102.31	104.50
2	T	10	DG	C4'-C3'-C2'	5.44	108.00	103.10
1	P	7	DC	C4'-C3'-C2'	5.38	107.94	103.10
2	T	3	DA	C5-C6-N1	-5.37	115.02	117.70
1	P	8	DC	O4'-C1'-N1	-5.35	104.25	108.00
1	P	11	DA	N7-C8-N9	5.34	116.47	113.80
2	T	5	DG	N3-C2-N2	-5.34	116.16	119.90
2	T	4	DG	C4-C5-N7	-5.34	108.67	110.80
3	A	588	GLY	C-N-CA	-5.29	108.49	121.70
3	A	577	ARG	NE-CZ-NH1	-5.26	117.67	120.30
3	A	640	ASP	CB-CG-OD2	5.25	123.02	118.30
3	A	362	ASP	OD1-CG-OD2	-5.23	113.36	123.30
3	A	477	ASP	CB-CG-OD2	5.23	123.00	118.30
3	A	574	LEU	CB-CG-CD2	5.22	119.87	111.00
2	T	3	DA	N9-C4-C5	5.21	107.89	105.80
3	A	503	GLY	N-CA-C	5.21	126.13	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	6	DT	OP2-P-O3'	5.20	116.65	105.20
3	A	429	ILE	CB-CA-C	5.19	121.97	111.60
1	P	6	DC	N3-C4-C5	-5.17	119.83	121.90
2	T	13	DG	OP2-P-O3'	5.13	116.49	105.20
1	P	8	DC	OP1-P-OP2	5.12	127.29	119.60
1	P	8	DC	N1-C2-N3	5.10	122.77	119.20
1	P	6	DC	O3'-P-O5'	5.10	113.69	104.00
3	A	726	ARG	NE-CZ-NH1	-5.09	117.75	120.30
2	T	8	DG	O3'-P-O5'	-5.09	94.33	104.00
2	T	8	DG	C3'-C2'-C1'	-5.09	96.39	102.50
1	P	6	DC	C5-C4-N4	5.07	123.75	120.20
1	P	8	DC	C5-C4-N4	5.06	123.74	120.20
2	T	6	DT	C6-N1-C2	-5.05	118.78	121.30
2	T	9	DG	OP1-P-OP2	5.04	127.16	119.60
3	A	455	LEU	O-C-N	-5.04	114.64	122.70
2	T	5	DG	N9-C1'-C2'	5.02	122.14	112.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	347	TYR	Peptide
3	A	381	ALA	Peptide
3	A	524	ALA	Peptide

## 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	P	231	0	137	6	1
2	T	338	0	175	5	0
3	A	3378	0	3374	105	0
4	A	2	0	0	0	0
5	A	28	0	12	1	0
6	A	120	0	0	3	0
6	P	6	0	0	0	0
6	T	14	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4117	0	3698	114	1

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 15.

All (114) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:723:MET:SD	3:A:723:MET:CE	2.05	1.45
3:A:683:MET:SD	3:A:683:MET:CE	2.17	1.31
3:A:588:GLY:O	3:A:589:SER:CB	1.94	1.12
3:A:366:PHE:O	3:A:367:PHE:HB2	1.55	1.02
3:A:589:SER:O	3:A:590:LYS:HB2	1.62	0.99
3:A:460:LEU:HD12	3:A:460:LEU:C	1.85	0.95
2:T:6:DT:OP1	3:A:326:HIS:HD2	1.48	0.94
3:A:364:ASP:OD2	3:A:505:THR:HG22	1.68	0.94
3:A:588:GLY:O	3:A:589:SER:HB3	1.13	0.93
3:A:501:THR:HG21	3:A:504:CYS:HB2	1.50	0.91
3:A:348:THR:HG21	3:A:615:LYS:HE3	1.53	0.91
1:P:11:DA:H2''	1:P:12:DOC:H5'	1.51	0.90
3:A:533:PHE:H	3:A:536:ASN:HD22	1.16	0.88
1:P:1:DA:H2''	1:P:2:DT:H5'	1.59	0.83
3:A:364:ASP:OD2	3:A:505:THR:CG2	2.29	0.80
3:A:589:SER:O	3:A:590:LYS:CB	2.30	0.80
3:A:533:PHE:H	3:A:536:ASN:ND2	1.80	0.80
3:A:537:LEU:HD13	3:A:542:TRP:CH2	2.16	0.79
3:A:497:ILE:O	3:A:501:THR:HB	1.83	0.79
3:A:348:THR:CG2	3:A:615:LYS:HE3	2.13	0.78
3:A:341:ASN:O	3:A:344:ILE:HG13	1.83	0.78
3:A:664:ILE:HD13	3:A:697:LEU:HD11	1.64	0.77
3:A:318:SER:O	3:A:322:HIS:HD2	1.69	0.76
3:A:537:LEU:HD13	3:A:542:TRP:HH2	1.55	0.72
2:T:6:DT:OP1	3:A:326:HIS:CD2	2.37	0.72
1:P:8:DC:H2''	1:P:9:DC:H5'	1.71	0.70
2:T:15:DA:HI'	2:T:16:DT:H5'	1.72	0.70
3:A:484:THR:HG23	3:A:485:LEU:H	1.56	0.69
3:A:366:PHE:O	3:A:367:PHE:CB	2.19	0.69
3:A:460:LEU:CD1	3:A:460:LEU:C	2.62	0.67
3:A:619:GLN:O	3:A:621:LYS:HE2	1.93	0.67
3:A:696:ARG:HG3	3:A:696:ARG:HH11	1.62	0.64
3:A:501:THR:HG21	3:A:504:CYS:CB	2.25	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:360:HIS:HE1	3:A:468:GLU:OE1	1.80	0.63
3:A:539:GLU:HG2	3:A:572:ASN:HD21	1.63	0.62
3:A:733:ASN:O	3:A:734:LYS:O	2.17	0.61
1:P:6:DC:H2"	1:P:7:DC:H5"	1.82	0.61
3:A:533:PHE:N	3:A:536:ASN:HD22	1.95	0.59
3:A:510:CYS:O	3:A:531:ILE:O	2.19	0.59
3:A:501:THR:HG22	3:A:504:CYS:H	1.66	0.59
3:A:701:THR:HB	3:A:703:GLU:H	1.69	0.58
3:A:374:CYS:SG	3:A:440:GLN:HG3	2.43	0.58
3:A:455:LEU:HB3	3:A:457:ILE:HD12	1.86	0.57
3:A:460:LEU:HD11	3:A:462:LEU:HG	1.85	0.57
2:T:14:DG:H2"	2:T:15:DA:OP2	2.04	0.57
3:A:635:ASN:ND2	3:A:638:GLN:H	2.03	0.57
3:A:640:ASP:OD2	3:A:716:ARG:NH2	2.38	0.57
3:A:348:THR:HG21	3:A:615:LYS:CE	2.30	0.56
3:A:367:PHE:HA	3:A:441:PHE:CE2	2.41	0.56
3:A:460:LEU:O	3:A:460:LEU:HD12	2.04	0.56
3:A:331:TRP:HE1	3:A:442:GLN:HE21	1.54	0.55
3:A:525:LYS:HG3	3:A:525:LYS:O	2.06	0.55
3:A:501:THR:O	3:A:502:ASN:HB2	2.07	0.55
3:A:733:ASN:O	3:A:734:LYS:C	2.43	0.55
3:A:632:ARG:NH1	3:A:678:GLU:OE1	2.40	0.55
3:A:363:PHE:CD1	3:A:467:ASP:HB2	2.43	0.54
3:A:501:THR:HG23	3:A:504:CYS:SG	2.49	0.53
3:A:584:LYS:HG2	3:A:589:SER:HA	1.89	0.53
3:A:495:GLN:OE1	3:A:499:GLN:NE2	2.36	0.53
3:A:645:ARG:NH1	6:A:16:HOH:O	2.42	0.52
3:A:445:SER:HB2	3:A:466:ILE:HD13	1.91	0.52
3:A:502:ASN:O	3:A:504:CYS:N	2.43	0.52
3:A:723:MET:CE	3:A:723:MET:CG	2.85	0.51
3:A:539:GLU:HG2	3:A:572:ASN:ND2	2.26	0.51
3:A:657:ILE:HG13	3:A:659:LYS:HG3	1.91	0.51
3:A:323:SER:O	3:A:327:HIS:HD2	1.94	0.51
3:A:501:THR:CG2	3:A:504:CYS:HB2	2.31	0.51
3:A:360:HIS:CE1	3:A:468:GLU:OE1	2.62	0.51
3:A:458:PHE:CG	3:A:471:CYS:HB3	2.46	0.50
3:A:445:SER:O	3:A:448:PHE:HB3	2.11	0.50
3:A:445:SER:HB2	3:A:466:ILE:CD1	2.42	0.50
3:A:696:ARG:CG	3:A:696:ARG:HH11	2.25	0.49
3:A:460:LEU:CD1	3:A:462:LEU:HG	2.43	0.49
3:A:669:MET:HE2	3:A:685:MET:HE1	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:513:SER:HB2	3:A:516:LEU:H	1.78	0.49
3:A:574:LEU:HD13	3:A:599:LEU:HD11	1.95	0.49
3:A:501:THR:CG2	3:A:504:CYS:CB	2.89	0.49
3:A:635:ASN:ND2	3:A:638:GLN:HG3	2.27	0.49
3:A:437:THR:HB	6:A:95:HOH:O	2.12	0.48
3:A:410:TYR:OH	3:A:429:ILE:HD13	2.15	0.47
3:A:400:ILE:O	3:A:414:ASN:O	2.33	0.47
3:A:669:MET:HE2	3:A:685:MET:CE	2.44	0.47
3:A:669:MET:HE3	3:A:685:MET:HE3	1.96	0.47
3:A:627:ILE:HG22	3:A:631:ILE:HD13	1.96	0.47
3:A:366:PHE:O	5:A:201:DCP:O3'	2.33	0.46
3:A:583:LEU:HD13	3:A:592:GLY:O	2.16	0.45
3:A:669:MET:CE	3:A:685:MET:HE3	2.46	0.45
3:A:318:SER:O	3:A:322:HIS:CD2	2.60	0.45
3:A:645:ARG:HA	3:A:648:GLN:HE21	1.82	0.45
3:A:668:LEU:C	3:A:668:LEU:HD12	2.37	0.44
3:A:667:LYS:HB3	3:A:729:ALA:HB3	2.00	0.44
3:A:547:LEU:HD11	3:A:568:PRO:HG2	1.99	0.44
3:A:696:ARG:NH1	3:A:696:ARG:HG3	2.32	0.44
3:A:607:SER:O	3:A:610:ILE:HG22	2.17	0.44
3:A:502:ASN:O	3:A:503:GLY:C	2.49	0.44
2:T:10:DG:H2''	2:T:11:DG:C8	2.53	0.44
3:A:501:THR:CG2	3:A:504:CYS:H	2.31	0.43
3:A:403:CYS:HB2	3:A:407:ALA:HB3	2.00	0.43
1:P:7:DC:OP1	1:P:7:DC:H4'	2.18	0.43
3:A:382:CYS:O	3:A:387:ASP:OD2	2.37	0.43
3:A:348:THR:CB	3:A:615:LYS:HE3	2.48	0.43
3:A:635:ASN:ND2	3:A:637:THR:HB	2.34	0.43
3:A:734:LYS:HA	6:A:121:HOH:O	2.18	0.42
3:A:632:ARG:HD2	3:A:632:ARG:HH11	1.47	0.42
3:A:466:ILE:HD13	3:A:466:ILE:O	2.20	0.42
1:P:7:DC:H5''	1:P:7:DC:H6	1.85	0.42
3:A:477:ASP:O	3:A:477:ASP:OD1	2.38	0.42
3:A:669:MET:CE	3:A:685:MET:CE	2.98	0.41
3:A:661:THR:HG22	3:A:663:GLN:H	1.85	0.41
3:A:651:LEU:HD23	3:A:651:LEU:HA	1.85	0.41
3:A:696:ARG:CG	3:A:696:ARG:NH1	2.81	0.41
3:A:565:PHE:O	3:A:568:PRO:HD3	2.21	0.40
3:A:682:TYR:O	3:A:683:MET:HB2	2.20	0.40
3:A:364:ASP:O	3:A:365:CYS:C	2.59	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:1:DA:C2	1:P:1:DA:C2[4_566]	2.18	0.02

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	A	424/434 (98%)	402 (95%)	17 (4%)	5 (1%)	16 16

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	367	PHE
3	A	590	LYS
3	A	734	LYS
3	A	365	CYS
3	A	589	SER

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	A	376/391 (96%)	340 (90%)	36 (10%)	10 12

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	314	ASP
3	A	324	ARG
3	A	346	LYS
3	A	375	ARG
3	A	385	LYS
3	A	396	LYS
3	A	430	LYS
3	A	431	LEU
3	A	460	LEU
3	A	466	ILE
3	A	495	GLN
3	A	501	THR
3	A	505	THR
3	A	506	VAL
3	A	518	ARG
3	A	542	TRP
3	A	556	SER
3	A	558	LEU
3	A	559	SER
3	A	574	LEU
3	A	583	LEU
3	A	590	LYS
3	A	591	LEU
3	A	605	GLU
3	A	621	LYS
3	A	635	ASN
3	A	636	ILE
3	A	654	LEU
3	A	664	ILE
3	A	666	LEU
3	A	673	LYS
3	A	693	ARG
3	A	699	ILE
3	A	716	ARG
3	A	737	ASP
3	A	738	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	322	HIS
3	A	326	HIS
3	A	327	HIS

*Continued on next page...*

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Mol	Chain	Res	Type
3	A	360	HIS
3	A	442	GLN
3	A	459	ASN
3	A	536	ASN
3	A	572	ASN
3	A	619	GLN
3	A	635	ASN
3	A	648	GLN
3	A	663	GLN
3	A	733	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	DOC	P	12	1,2	11,19,20	1.20	1 (9%)	14,26,29	2.77	8 (57%)

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	DOC	P	12	1,2	-	0/3/18/19	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	P	12	DOC	C6-C5	-2.42	1.32	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	12	DOC	C3'-C2'-C1'	-5.82	96.20	102.71
1	P	12	DOC	C3'-C4'-C5'	-3.65	101.19	116.05
1	P	12	DOC	C5-C4-N4	-2.70	117.17	121.31
1	P	12	DOC	C2'-C3'-C4'	-2.10	98.49	102.59
1	P	12	DOC	O4'-C4'-C3'	2.10	108.21	104.69
1	P	12	DOC	O4'-C1'-C2'	2.29	109.15	106.67
1	P	12	DOC	N4-C4-N3	2.65	121.34	116.50
1	P	12	DOC	C2-N3-C4	5.21	122.96	115.61

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
1	P	12	DOC	1	0

## 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	DCP	A	201	4	21,29,29	1.27	2 (9%)	33,45,45	1.52	5 (15%)

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	DCP	A	201	4	-	0/18/34/34	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	201	DCP	C4-N3	2.04	1.39	1.35
5	A	201	DCP	C6-N1	3.25	1.40	1.35

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	201	DCP	C2'-C3'-C4'	-2.91	96.75	102.77
5	A	201	DCP	C6-N1-C2	-2.25	117.63	121.28
5	A	201	DCP	C5-C4-N3	-2.23	118.98	121.80
5	A	201	DCP	O3A-PA-O5'	2.65	109.97	102.94
5	A	201	DCP	C2-N3-C4	4.43	121.86	115.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	201	DCP	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	P	11/12 (91%)	0.05	1 (9%) 11 17	27, 35, 65, 67	0
2	T	16/16 (100%)	-0.13	1 (6%) 23 31	21, 28, 58, 67	0
3	A	428/434 (98%)	-0.24	2 (0%) 91 94	16, 31, 49, 69	0
All	All	455/462 (98%)	-0.23	4 (0%) 85 89	16, 31, 50, 69	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	T	16	DT	2.8
3	A	477	ASP	2.6
3	A	674	ASP	2.5
1	P	1	DA	2.2

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	DOC	P	12	18/19	0.86	0.22	-	24,50,68,70	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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
5	DCP	A	201	28/28	0.99	0.15	0.08	16,22,26,28	0
4	MG	A	302	1/1	0.96	0.14	-0.05	16,16,16,16	0
4	MG	A	301	1/1	0.96	0.10	-	41,41,41,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.