



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:06 AM GMT

PDB ID : 2PFO
Title : DNA Polymerase lambda in complex with DNA and dUPNPP
Authors : Garcia-Diaz, M.; Bebenek, K.; Krahn, J.M.; Pedersen, L.C.; Kunkel, T.A.
Deposited on : 2007-04-05
Resolution : 2.00 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

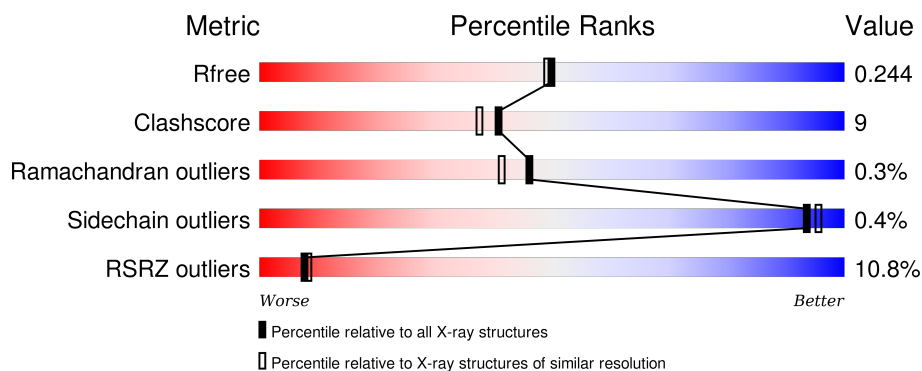
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	T	11	<div> <div>100%</div> </div>
2	P	6	<div> <div>67%</div> <div>33%</div> </div>
3	D	4	<div> <div>75%</div> <div>25%</div> </div>
4	A	335	<div> <div>11%</div> <div>78%</div> <div>18%</div> <div>•</div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 3231 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 Template.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	T	11	Total	C	N	O	P	0	0	0
			224	107	43	64	10			

- Molecule 2 is a DNA chain called Primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	6	Total	C	N	O	P	0	0	0
			119	58	23	33	5			

- Molecule 3 is a DNA chain called Downstream Primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	4	Total	C	N	O	P	0	0	0
			83	38	16	25	4			

- Molecule 4 is a protein called DNA polymerase lambda.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	A	323	Total	C	N	O	S	0	1	0
			2438	1529	438	460	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	MET	-	INITIATING METHIONINE	UNP Q9UGP5
A	543	ALA	CYS	ENGINEERED	UNP Q9UGP5

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

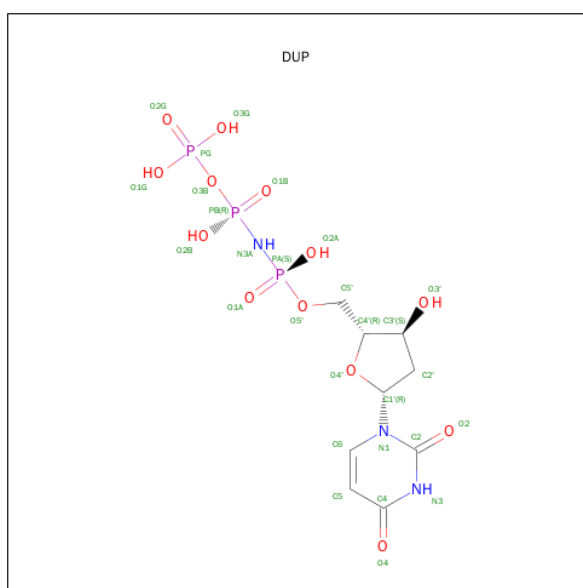
- Molecule 6 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mn	0	0
			1	1		

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

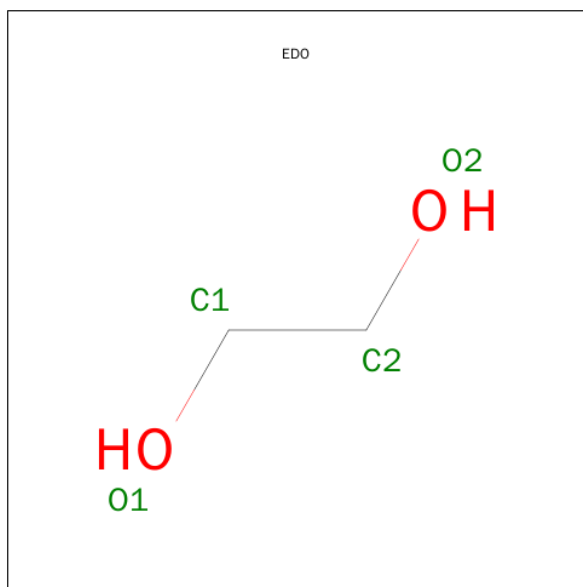
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	4	Total	Na	0	0
			4	4		

- Molecule 8 is 2'-DEOXYURIDINE 5'-ALPHA,BETA-IMIDO-TRIPHOSPHATE (three-letter code: DUP) (formula: C₉H₁₆N₃O₁₃P₃).



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

- Molecule 9 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	245	Total	O	0	0
			245	245		
10	D	11	Total	O	0	0
			11	11		
10	P	29	Total	O	0	0
			29	29		
10	T	44	Total	O	0	0
			44	44		

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: Template

Chain T:  100%


There are no outlier residues recorded for this chain.

- Molecule 2: Primer

Chain P:  67% 33%




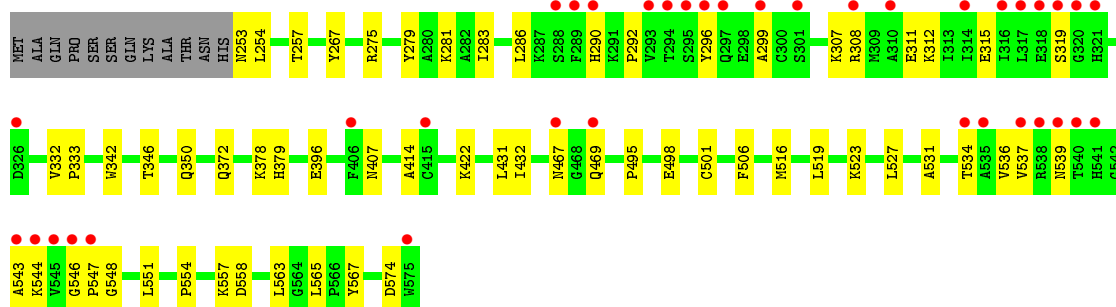
- Molecule 3: Downstream Primer

Chain D:  75% 25%



- Molecule 4: DNA polymerase lambda

Chain A:  11% 78% 18% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.09 Å 61.41 Å 140.31 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 35.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	80.0 (50.00-2.00) 93.1 (35.92-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.00 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.217 , 0.242 0.221 , 0.244	Depositor DCC
R_{free} test set	1556 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 60.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 33156 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3231	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DUP, NA, MG, MN, EDO

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	T	0.42	0/251	0.83	0/386
2	P	0.40	0/133	0.77	0/203
3	D	0.92	1/92 (1.1%)	0.70	0/138
4	A	0.32	0/2487	0.53	0/3372
All	All	0.36	1/2963 (0.0%)	0.58	0/4099

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1	DG	OP3-P	-6.94	1.52	1.61

There are no bond angle outliers.

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	T	224	0	125	0	0
2	P	119	0	68	2	0
3	D	83	0	45	3	0
4	A	2438	0	2324	43	0
5	A	1	0	0	0	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	4	0	0	0	0
8	A	28	0	12	0	0
9	A	4	0	6	0	0
10	A	245	0	0	2	0
10	D	11	0	0	0	0
10	P	29	0	0	1	0
10	T	44	0	0	0	0
All	All	3231	0	2580	47	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 (47) 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)
4:A:467:ASN:HD21	4:A:469:GLN:HB2	1.53	0.74
4:A:537:VAL:HB	4:A:546:GLY:HA3	1.70	0.71
4:A:290:HIS:O	4:A:292:PRO:HD3	1.92	0.69
4:A:534:THR:HB	4:A:551:LEU:HD11	1.75	0.68
4:A:495:PRO:HG2	4:A:498:GLU:HG3	1.76	0.66
4:A:539:ASN:HD21	4:A:543:ALA:HB3	1.60	0.66
4:A:396:GLU:HG3	4:A:414:ALA:HB2	1.80	0.64
4:A:536:VAL:HG11	4:A:544:LYS:HE2	1.84	0.59
4:A:467:ASN:ND2	4:A:469:GLN:HB2	2.21	0.56
4:A:308:ARG:HH11	4:A:308:ARG:HG3	1.71	0.55
4:A:519:LEU:HD23	4:A:565:LEU:HD11	1.88	0.55
4:A:253:ASN:C	4:A:254:LEU:HD12	2.28	0.54
4:A:279:TYR:O	4:A:283:ILE:HG13	2.10	0.51
4:A:254:LEU:HD21	4:A:292:PRO:HA	1.92	0.51
4:A:523:LYS:HE3	4:A:563:LEU:O	2.11	0.50
4:A:534:THR:HG21	4:A:551:LEU:HD21	1.93	0.49
4:A:539:ASN:ND2	4:A:543:ALA:HB3	2.27	0.49
4:A:519:LEU:CD2	4:A:565:LEU:HD11	2.43	0.48
4:A:431:LEU:HD23	4:A:431:LEU:C	2.35	0.47
4:A:431:LEU:HD23	4:A:432:ILE:N	2.29	0.47
4:A:536:VAL:CG1	4:A:544:LYS:HE2	2.43	0.47
4:A:534:THR:CG2	4:A:551:LEU:HD21	2.44	0.47
4:A:308:ARG:O	4:A:312:LYS:HG3	2.13	0.47
4:A:547:PRO:HG2	4:A:548:GLY:H	1.80	0.47
4:A:346[A]:THR:O	4:A:350:GLN:HG3	2.15	0.46
4:A:307:LYS:O	4:A:311:GLU:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:501:CYS:SG	4:A:531:ALA:HA	2.56	0.46
4:A:516:MET:HG2	4:A:527:LEU:CD2	2.46	0.46
4:A:332:VAL:HB	4:A:333:PRO:HD3	1.97	0.45
4:A:267:TYR:CD1	4:A:275:ARG:HD2	2.52	0.45
4:A:554:PRO:HG2	4:A:558:ASP:OD2	2.15	0.45
4:A:267:TYR:CE1	4:A:275:ARG:HD2	2.52	0.45
4:A:296:TYR:O	4:A:299:ALA:HB3	2.18	0.44
3:D:1:DG:H2''	3:D:2:DC:H5'	1.99	0.44
2:P:5:DA:H4'	4:A:342:TRP:CZ2	2.53	0.43
4:A:281:LYS:HB2	10:A:2907:HOH:O	2.17	0.43
2:P:6:DC:H4'	10:P:18:HOH:O	2.17	0.43
3:D:3:DC:H2''	3:D:4:DG:C8	2.53	0.43
4:A:315:GLU:O	4:A:319:SER:HB2	2.19	0.43
4:A:257:THR:HG22	4:A:286:LEU:HD12	2.00	0.43
4:A:422:LYS:HD3	4:A:574:ASP:OD1	2.18	0.42
4:A:346[B]:THR:O	4:A:350:GLN:HG3	2.18	0.42
4:A:308:ARG:NH1	4:A:308:ARG:HG3	2.34	0.41
3:D:1:DG:C2'	3:D:2:DC:H5'	2.51	0.41
4:A:372:GLN:HG3	10:A:2834:HOH:O	2.20	0.40
4:A:557:LYS:HG2	4:A:567:TYR:CD2	2.56	0.40
4:A:378:LYS:HE3	4:A:379:HIS:NE2	2.37	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	
4	A	322/335 (96%)	303 (94%)	18 (6%)	1 (0%)	46	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	A	407	ASN

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	
4	A	242/280 (86%)	241 (100%)	1 (0%)	93	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	A	506	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
4	A	350	GLN
4	A	354	GLN
4	A	464	GLN
4	A	467	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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
9	EDO	A	2801	-	3,3,3	0.60	0	2,2,2	0.54	0
8	DUP	A	952	5,6	23,29,29	1.89	6 (26%)	32,45,45	2.74	6 (18%)

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
9	EDO	A	2801	-	-	0/1/1/1	0/0/0/0
8	DUP	A	952	5,6	-	0/13/34/34	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	952	DUP	PB-O3B	-5.86	1.51	1.59
8	A	952	DUP	PB-O2B	-2.83	1.48	1.56
8	A	952	DUP	PA-O2A	-2.48	1.49	1.56
8	A	952	DUP	PB-O1B	2.37	1.48	1.46
8	A	952	DUP	C4-N3	2.63	1.38	1.33
8	A	952	DUP	PA-O1A	2.69	1.49	1.46

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	952	DUP	C5-C4-N3	-3.24	114.81	123.12
8	A	952	DUP	O5'-PA-O1A	-2.08	104.69	113.31
8	A	952	DUP	O1G-PG-O2G	2.53	118.72	110.58
8	A	952	DUP	O2B-PB-O1B	4.18	118.73	110.00
8	A	952	DUP	O2A-PA-O1A	4.19	118.74	110.00
8	A	952	DUP	C4-N3-C2	13.06	127.07	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	T	11/11 (100%)	-0.36	0 100 100	21, 25, 39, 42	0
2	P	6/6 (100%)	-0.27	0 100 100	18, 20, 21, 22	0
3	D	4/4 (100%)	-0.52	0 100 100	37, 37, 39, 43	0
4	A	323/335 (96%)	0.56	37 (11%) 6 7	16, 36, 78, 86	0
All	All	344/356 (96%)	0.50	37 (10%) 8 8	16, 35, 78, 86	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	A	540	THR	6.3
4	A	543	ALA	6.2
4	A	535	ALA	6.0
4	A	541	HIS	5.2
4	A	319	SER	5.1
4	A	544	LYS	4.9
4	A	545	VAL	4.9
4	A	321	HIS	4.7
4	A	546	GLY	4.2
4	A	293	VAL	4.1
4	A	289	PHE	4.1
4	A	301	SER	3.7
4	A	290	HIS	3.5
4	A	320	GLY	3.5
4	A	406	PHE	3.4
4	A	296	TYR	3.4
4	A	538	ARG	3.4
4	A	299	ALA	3.1
4	A	539	ASN	3.1
4	A	575	TRP	3.1
4	A	547	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
4	A	295	SER	3.0
4	A	318	GLU	2.9
4	A	316	ILE	2.9
4	A	317	LEU	2.9
4	A	310	ALA	2.8
4	A	288	SER	2.8
4	A	297	GLN	2.8
4	A	294	THR	2.7
4	A	534	THR	2.6
4	A	467	ASN	2.5
4	A	308	ARG	2.4
4	A	415	CYS	2.4
4	A	537	VAL	2.3
4	A	314	ILE	2.2
4	A	326	ASP	2.1
4	A	469	GLN	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

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
8	DUP	A	952	28/28	0.98	0.16	0.66	16,23,25,25	0
9	EDO	A	2801	4/4	0.91	0.11	0.17	51,51,52,52	0
7	NA	A	603	1/1	0.88	0.12	-1.10	39,39,39,39	1
7	NA	A	601	1/1	0.98	0.08	-2.16	21,21,21,21	0
7	NA	A	600	1/1	0.97	0.06	-	41,41,41,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	MN	A	951	1/1	1.00	0.11	-	22,22,22,22	0
7	NA	A	602	1/1	0.97	0.17	-	46,46,46,46	0
5	MG	A	950	1/1	0.99	0.17	-	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.