



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2017 – 10:45 AM EST

PDB ID : 4KLQ
Title : Observing a DNA polymerase choose right from wrong.
Authors : Freudenthal, B.D.; Beard, W.A.; Shock, D.D.; Wilson, S.H.
Deposited on : unknown
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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030345

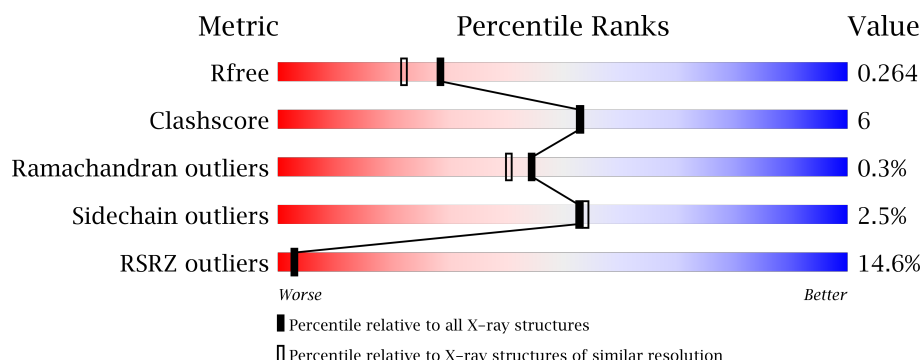
1 Overall quality at a glance

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}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (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	A	335	<div> <div>15%</div> <div> <div></div> <div>85%</div> <div>11%</div> <div>.</div> </div> </div>
2	T	16	<div> <div>6%</div> <div> <div></div> <div>63%</div> <div>38%</div> </div> </div>
3	P	11	<div> <div></div> <div> <div>55%</div> <div>36%</div> <div>9%</div> </div> </div>
4	D	5	<div> <div></div> <div> <div>20%</div> <div>80%</div> </div> </div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 3519 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 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	4	0
			2616	1650	458	499	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	16	Total	C	N	O	P	0	0	0
			323	153	63	92	15			

- Molecule 3 is a DNA chain called 5'-D(*GP*CP*TP*GP*AP*TP*GP*CP*GP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	11	Total	C	N	O	P	0	2	0
			243	116	46	70	11			

- Molecule 4 is a DNA chain called 5'-D(P*GP*TP*CP*GP*G)-3'.

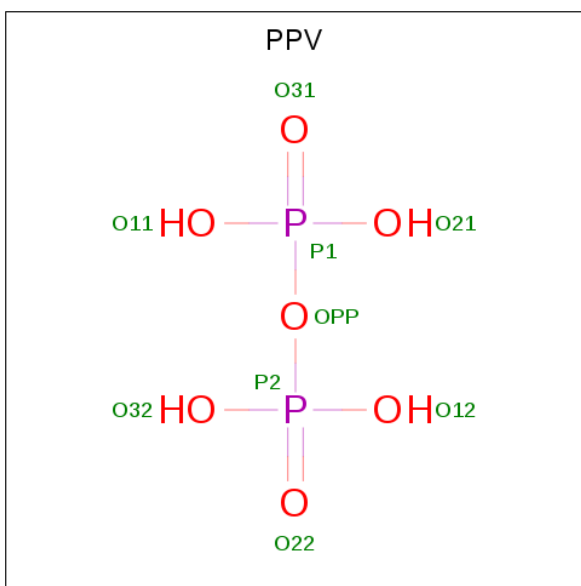
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	5	Total	C	N	O	P	0	0	0
			106	49	20	32	5			

- Molecule 5 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (three-letter code: DTP) (formula: C₁₀H₁₆N₅O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	1
			30	10	5	12	3		

- Molecule 6 is PYROPHOSPHATE (three-letter code: PPV) (formula: $\text{H}_4\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	1
			9	7	2		

- Molecule 7 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	3	Total 3	Mn 3	0	0

- Molecule 8 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total 3	Na 3	0	0

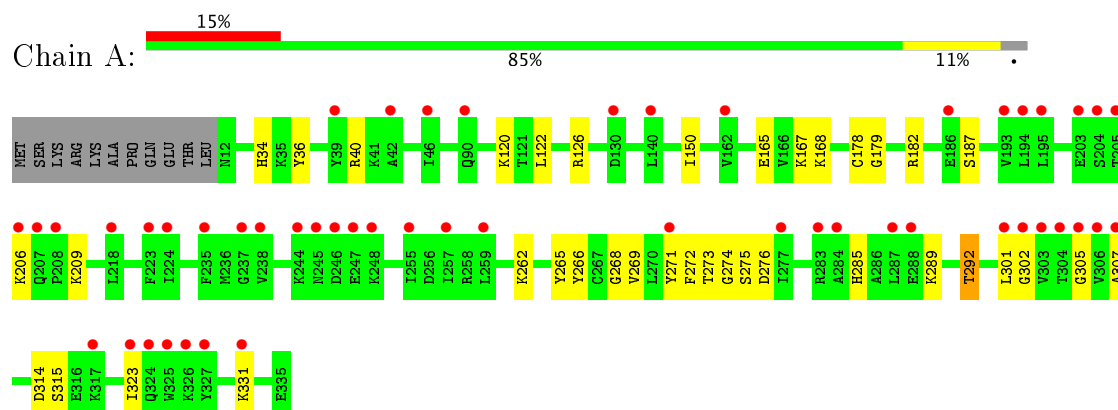
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	134	Total 134	O 134	0	0
9	T	26	Total 26	O 26	0	0
9	P	20	Total 20	O 20	0	1
9	D	6	Total 6	O 6	0	0

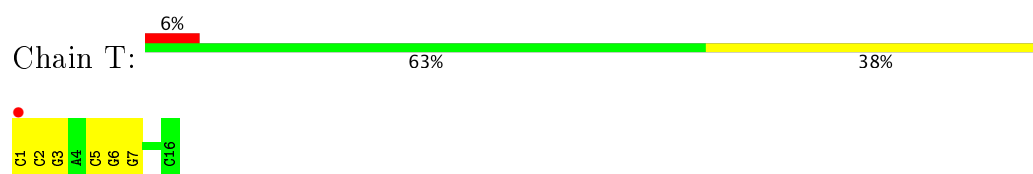
3 Residue-property plots [i](#)

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



- Molecule 2: 5'-D(*CP*CP*GP*AP*CP*GP*GP*CP*GP*CP*AP*TP*CP*AP*GP*C)-3'



- Molecule 3: 5'-D(*GP*CP*TP*GP*AP*TP*GP*CP*GP*CP*A)-3'



- Molecule 4: 5'-D(P*GP*TP*CP*GP*G)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.82Å 78.41Å 54.96Å 90.00° 113.24° 90.00°	Depositor
Resolution (Å)	24.04 – 2.00 24.04 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.0 (24.04-2.00) 89.3 (24.04-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 1.99Å)	Xtriage
Refinement program	PHENIX 1.7 _650	Depositor
R, R_{free}	0.214 , 0.267 0.218 , 0.264	Depositor DCC
R_{free} test set	1910 reflections (7.14%)	DCC
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3519	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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: NA, DTP, PPV, MN

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	0.40	0/2666	0.55	0/3583
2	T	0.91	0/362	1.20	0/556
3	P	1.64	3/272 (1.1%)	1.87	7/418 (1.7%)
4	D	1.25	1/118 (0.8%)	1.57	3/179 (1.7%)
All	All	0.70	4/3418 (0.1%)	0.90	10/4736 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	10[A]	DC	O3'-P	-11.25	1.47	1.61
3	P	10[B]	DC	O3'-P	-11.25	1.47	1.61
3	P	5	DA	C4'-O4'	-5.33	1.39	1.45
4	D	1	DG	C4'-C3'	-5.04	1.47	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	10[A]	DC	P-O3'-C3'	16.63	139.66	119.70
3	P	10[B]	DC	P-O3'-C3'	16.63	139.66	119.70
3	P	10[A]	DC	O3'-P-O5'	-7.01	90.67	104.00
3	P	10[B]	DC	O3'-P-O5'	-7.01	90.67	104.00
3	P	10[A]	DC	O4'-C1'-N1	6.89	112.83	108.00
3	P	10[B]	DC	O4'-C1'-N1	6.89	112.83	108.00
4	D	1	DG	O4'-C1'-N9	-6.14	103.70	108.00
3	P	5	DA	O4'-C1'-N9	-5.65	104.04	108.00
4	D	2	DT	O4'-C1'-N1	-5.28	104.30	108.00
4	D	2	DT	N3-C4-O4	5.13	122.98	119.90

There are no chirality outliers.

There are no planarity outliers.

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	2616	0	2607	25	0
2	T	323	0	179	8	0
3	P	243	0	135	9	0
4	D	106	0	57	3	0
5	A	30	0	9	2	0
6	A	9	0	0	1	0
7	A	3	0	0	0	0
8	A	3	0	0	0	0
9	A	134	0	0	0	0
9	D	6	0	0	0	0
9	P	20	0	0	0	0
9	T	26	0	0	0	0
All	All	3519	0	2987	39	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 6.

All (39) 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:P:11[B]:DA:H8	3:P:11[B]:DA:O5'	1.44	1.00
3:P:11[B]:DA:C8	3:P:11[B]:DA:O5'	2.30	0.82
6:A:402[B]:PPV:O12	3:P:11[B]:DA:H5''	1.81	0.81
1:A:271:TYR:OH	3:P:11[B]:DA:C2	2.35	0.78
2:T:1:DC:H1'	2:T:2:DC:H5'	1.69	0.74
2:T:1:DC:H42	4:D:5:DG:H1	1.39	0.70
4:D:4:DG:H2''	4:D:5:DG:O5'	1.95	0.66
1:A:276:ASP:HB3	5:A:401[A]:DTP:H8	1.78	0.66
1:A:276:ASP:HB3	5:A:401[A]:DTP:C8	2.31	0.60
1:A:179:GLY:H	1:A:273:THR:HA	1.65	0.60
1:A:36:TYR:CZ	1:A:40:ARG:HD2	2.38	0.59
3:P:10[B]:DC:H2''	3:P:11[B]:DA:N7	2.18	0.58
2:T:1:DC:N4	4:D:5:DG:H1	2.02	0.56
1:A:36:TYR:CZ	1:A:40:ARG:CD	2.90	0.54
1:A:292:THR:HG21	1:A:307:ALA:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:11[B]:DA:OP1	3:P:11[B]:DA:O4'	2.30	0.50
2:T:2:DC:H2''	2:T:3:DG:C8	2.47	0.49
1:A:36:TYR:CE1	1:A:40:ARG:HD2	2.47	0.49
1:A:301:LEU:C	1:A:301:LEU:HD12	2.33	0.49
1:A:301:LEU:HB2	1:A:305:GLY:HA2	1.95	0.48
1:A:301:LEU:HD12	1:A:302:GLY:O	2.13	0.47
2:T:1:DC:C1'	2:T:2:DC:H5'	2.41	0.47
1:A:126:ARG:HH11	1:A:126:ARG:HG3	1.79	0.47
1:A:265:TYR:CZ	1:A:269:VAL:HG21	2.51	0.46
3:P:2:DC:H2''	3:P:3:DT:C6	2.51	0.46
1:A:178:CYS:HA	1:A:182:ARG:HB2	1.98	0.45
1:A:285:HIS:HD2	1:A:323:ILE:HD12	1.82	0.45
1:A:150:ILE:O	1:A:187:SER:HA	2.16	0.44
3:P:11[B]:DA:H8	3:P:11[B]:DA:C5'	2.28	0.44
1:A:34:HIS:CD2	2:T:5:DC:C5	3.05	0.43
1:A:268:GLY:O	1:A:272:PHE:HD2	2.01	0.43
1:A:167:LYS:HB3	1:A:167:LYS:HE2	1.79	0.43
1:A:289:LYS:HD3	1:A:289:LYS:HA	1.42	0.42
1:A:266:TYR:CD1	1:A:315:SER:HA	2.55	0.42
1:A:34:HIS:CE1	2:T:5:DC:C6	3.07	0.41
1:A:271:TYR:CZ	3:P:11[B]:DA:C2	3.08	0.41
1:A:274:GLY:HA3	1:A:275:SER:HA	1.81	0.41
2:T:6:DG:H2''	2:T:7:DG:OP2	2.21	0.41
1:A:165:GLU:OE1	1:A:168:LYS:HD2	2.22	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	326/335 (97%)	308 (94%)	17 (5%)	1 (0%)	44 40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	LYS

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	
1	A	284/295 (96%)	277 (98%)	7 (2%)	53	54

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	LYS
1	A	122	LEU
1	A	209	LYS
1	A	262	LYS
1	A	292	THR
1	A	314	ASP
1	A	331	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	264	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

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	$\# Z > 2$
5	DTP	A	401[A]	8,7	26,32,32	1.20	3 (11%)	26,50,50	1.75	2 (7%)
6	PPV	A	402[B]	8,7	8,8,8	1.11	1 (12%)	8,13,13	0.37	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
5	DTP	A	401[A]	8,7	-	0/18/34/34	0/3/3/3
6	PPV	A	402[B]	8,7	-	0/6/6/6	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	401[A]	DTP	PB-O2B	-3.11	1.39	1.55
5	A	401[A]	DTP	C8-N7	-2.91	1.29	1.34
6	A	402[B]	PPV	P2-OPP	-2.04	1.56	1.60
5	A	401[A]	DTP	PA-O1A	2.03	1.58	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	401[A]	DTP	N3-C2-N1	-7.57	122.26	128.86
5	A	401[A]	DTP	C4-C5-N7	-3.36	106.17	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	401[A]	DTP	2	0
6	A	402[B]	PPV	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	324/335 (96%)	0.97	51 (15%) 2 2	21, 38, 75, 99	0
2	T	16/16 (100%)	0.23	1 (6%) 21 21	20, 43, 69, 76	0
3	P	11/11 (100%)	0.04	0 100 100	20, 34, 42, 45	1 (9%)
4	D	5/5 (100%)	0.12	0 100 100	26, 29, 49, 57	0
All	All	356/367 (97%)	0.90	52 (14%) 3 3	20, 38, 75, 99	1 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	LEU	11.1
1	A	306	VAL	8.5
1	A	305	GLY	7.0
1	A	246	ASP	6.7
1	A	204	SER	6.2
1	A	244	LYS	6.1
1	A	304	THR	5.8
1	A	245	ASN	5.3
1	A	323	ILE	5.0
1	A	208	PRO	5.0
1	A	193	VAL	4.5
1	A	303	VAL	4.4
1	A	257	ILE	4.3
1	A	288	GLU	4.3
1	A	302	GLY	4.3
1	A	207	GLN	4.3
1	A	255	ILE	4.0
1	A	324	GLN	3.8
1	A	271	TYR	3.7
1	A	307	ALA	3.7
1	A	325	TRP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	205	THR	3.4
1	A	331	LYS	3.2
1	A	42	ALA	3.2
1	A	327	TYR	3.0
1	A	162	VAL	3.0
1	A	206	LYS	3.0
1	A	287	LEU	2.9
1	A	326	LYS	2.9
1	A	283	ARG	2.9
1	A	223	PHE	2.8
1	A	235	PHE	2.8
1	A	195	LEU	2.8
1	A	90	GLN	2.8
1	A	203	GLU	2.8
2	T	1	DC	2.7
1	A	237	GLY	2.5
1	A	140	LEU	2.4
1	A	317	LYS	2.4
1	A	238	VAL	2.3
1	A	218	LEU	2.2
1	A	259	LEU	2.2
1	A	46	ILE	2.1
1	A	277	ILE	2.1
1	A	248	LYS	2.1
1	A	194	LEU	2.1
1	A	130	ASP	2.1
1	A	284	ALA	2.1
1	A	224	ILE	2.1
1	A	186	GLU	2.0
1	A	39	TYR	2.0
1	A	247	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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
5	DTP	A	401[A]	30/30	0.90	0.19	0.69	14,40,51,53	30
7	MN	A	403	1/1	0.97	0.17	-0.03	25,25,25,25	1
6	PPV	A	402[B]	9/9	0.93	0.12	-0.46	32,35,37,40	9
8	NA	A	407	1/1	0.97	0.13	-0.94	22,22,22,22	0
7	MN	A	404	1/1	0.91	0.08	-1.56	36,36,36,36	0
8	NA	A	408	1/1	0.98	0.07	-4.70	30,30,30,30	0
8	NA	A	406	1/1	0.96	0.12	-	51,51,51,51	0
7	MN	A	405	1/1	0.97	0.32	-	30,30,30,30	0

6.5 Other polymers

There are no such residues in this entry.