



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:12 AM UTC

PDB ID : 6DSV / pdb_00006dsv
Title : Bst DNA polymerase I post-chemistry (n+2) structure
Authors : Chim, N.; Jackson, L.N.; Chaput, J.C.
Deposited on : 2018-06-14
Resolution : 1.99 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

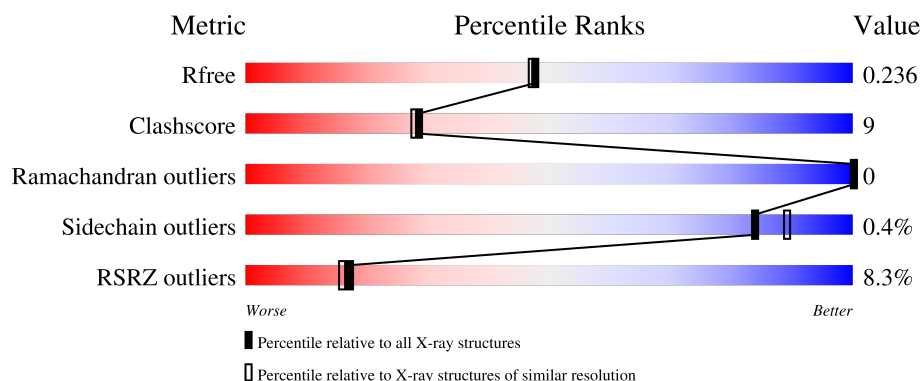
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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 $\leq 5\%$. 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	11	
2	A	580	
3	T	12	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5462 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 DNA (5'-D(*GP*CP*GP*AP*TP*CP*AP*CP*GP*TP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	P	11	Total	C	N	O	P	0	0	0
			223	107	43	63	10			

- Molecule 2 is a protein called DNA polymerase I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	577	Total	C	N	O	S	0	0	0
			4603	2925	801	861	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	550	THR	SER	conflict	UNP E1C9K5

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	12	Total	C	N	O	P	0	0	0
			246	117	45	72	12			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		

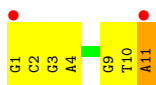
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	P	3	Total	O	0	0
			3	3		
6	A	372	Total	O	0	0
			372	372		
6	T	9	Total	O	0	0
			9	9		

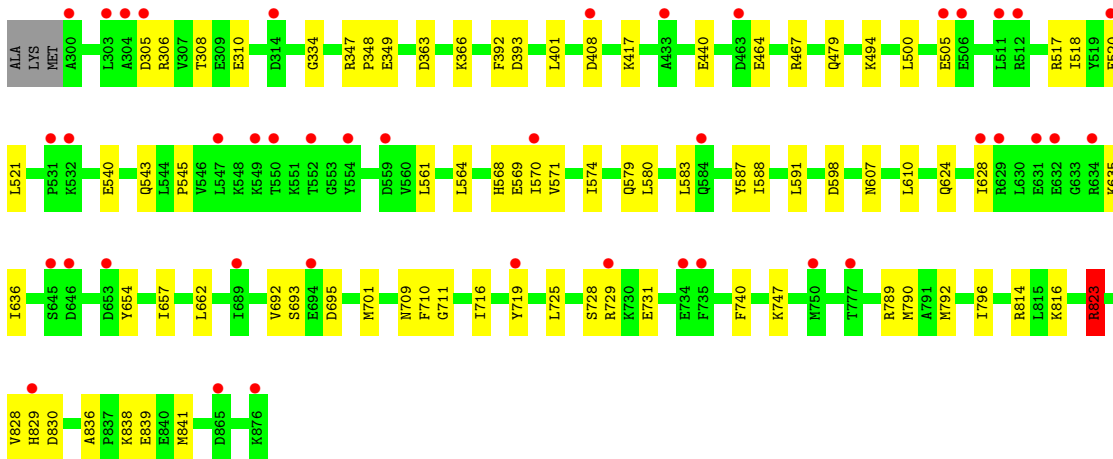
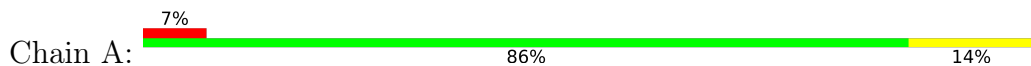
3 Residue-property plots [i](#)

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



- Molecule 2: DNA polymerase I



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.08Å 93.07Å 104.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.54 – 1.99 43.54 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.54-1.99) 94.3 (43.54-1.99)	Depositor EDS
R_{merge}	0.33	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.00Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.193 , 0.228 (Not available) , 0.236	Depositor DCC
R_{free} test set	2003 reflections (3.40%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.506	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5462	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 6.43% 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: SO4, MG

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	0.48	0/250	0.71	1/384 (0.3%)
2	A	0.37	1/4686 (0.0%)	0.53	0/6340
3	T	0.32	0/275	0.57	0/422
All	All	0.38	1/5211 (0.0%)	0.54	1/7146 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	823	ARG	C-O	-5.70	1.17	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	11	DA	C1'-O4'-C4'	-5.86	100.92	109.70

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	P	223	0	124	22	0
2	A	4603	0	4625	64	0
3	T	246	0	136	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	1	0	0	0	0
5	A	5	0	0	0	0
6	A	372	0	0	3	0
6	P	3	0	0	0	0
6	T	9	0	0	0	0
All	All	5462	0	4885	83	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 (83) 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:P:10:DT:H2'	1:P:11:DA:H4'	1.20	1.09
2:A:568:HIS:HD2	2:A:570:ILE:H	1.16	0.92
1:P:11:DA:N7	3:T:4:DT:O4	2.00	0.92
1:P:11:DA:N7	3:T:4:DT:N3	2.20	0.89
1:P:11:DA:N7	3:T:4:DT:C4	2.41	0.88
1:P:10:DT:C2'	1:P:11:DA:H4'	2.06	0.86
1:P:10:DT:H2''	2:A:829:HIS:CE1	2.26	0.70
2:A:823:ARG:NH2	6:A:1002:HOH:O	2.23	0.69
2:A:518:ILE:HG23	2:A:570:ILE:HD11	1.82	0.61
2:A:568:HIS:CD2	2:A:570:ILE:H	2.08	0.60
2:A:540:GLU:O	2:A:543:GLN:NE2	2.28	0.60
3:T:8:DT:H2''	3:T:9:DG:C8	2.37	0.59
2:A:347:ARG:NH2	2:A:440:GLU:OE2	2.34	0.59
1:P:11:DA:O3'	2:A:829:HIS:ND1	2.35	0.59
1:P:2:DC:H42	3:T:13:DG:H1	1.51	0.58
2:A:823:ARG:O	2:A:823:ARG:HG2	2.02	0.57
1:P:9:DG:H1	3:T:6:DC:H42	1.53	0.56
1:P:11:DA:C6	2:A:710:PHE:HZ	2.22	0.56
2:A:517:ARG:NH1	2:A:520:GLU:CD	2.64	0.56
2:A:598:ASP:OD1	2:A:598:ASP:N	2.37	0.55
1:P:10:DT:H2'	1:P:11:DA:C4'	2.14	0.54
2:A:624:GLN:O	2:A:829:HIS:HE1	1.91	0.54
2:A:740:PHE:HB3	2:A:747:LYS:HD2	1.89	0.54
2:A:790:MET:HE2	2:A:790:MET:HA	1.90	0.54
3:T:4:DT:H2'	3:T:5:DA:C8	2.42	0.53
2:A:517:ARG:O	2:A:521:LEU:HD13	2.09	0.53
2:A:693:SER:OG	2:A:695:ASP:OD1	2.21	0.52
2:A:417:LYS:NZ	2:A:464:GLU:OE1	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:624:GLN:O	2:A:829:HIS:CE1	2.63	0.52
2:A:349:GLU:OE2	2:A:349:GLU:N	2.33	0.51
1:P:10:DT:H5''	2:A:628:ILE:HG22	1.93	0.51
2:A:467:ARG:O	6:A:1001:HOH:O	2.20	0.50
3:T:9:DG:H2'	3:T:10:DA:C8	2.47	0.49
2:A:310:GLU:H	2:A:310:GLU:CD	2.20	0.49
2:A:836:ALA:HB3	2:A:841:MET:CE	2.43	0.48
2:A:654:TYR:HB3	2:A:657:ILE:HB	1.94	0.48
2:A:662:LEU:HD13	2:A:796:ILE:HD11	1.95	0.48
3:T:12:DC:H2''	3:T:13:DG:C8	2.49	0.48
2:A:635:LYS:NZ	6:A:1017:HOH:O	2.47	0.48
2:A:517:ARG:HH12	2:A:520:GLU:CD	2.20	0.48
1:P:10:DT:H2''	2:A:829:HIS:NE2	2.30	0.47
1:P:10:DT:H3	3:T:5:DA:H2	1.63	0.47
2:A:393:ASP:H	2:A:479:GLN:NE2	2.12	0.47
2:A:408:ASP:OD2	2:A:408:ASP:N	2.46	0.47
2:A:583:LEU:HD22	2:A:636:ILE:HD11	1.96	0.47
2:A:828:VAL:HG12	2:A:829:HIS:CD2	2.50	0.47
2:A:719:TYR:CE2	2:A:729:ARG:HD3	2.50	0.47
2:A:838:LYS:NZ	2:A:839:GLU:OE1	2.47	0.47
2:A:789:ARG:HA	2:A:792:MET:HE3	1.97	0.46
2:A:829:HIS:HB2	2:A:830:ASP:H	1.54	0.46
1:P:3:DG:H2''	1:P:4:DA:O5'	2.14	0.46
1:P:11:DA:C6	2:A:710:PHE:CZ	3.03	0.46
1:P:11:DA:OP2	2:A:628:ILE:HG21	2.15	0.46
2:A:500:LEU:HD22	2:A:588:ILE:HD12	1.98	0.46
2:A:518:ILE:HD13	2:A:574:ILE:HD13	1.97	0.46
2:A:725:LEU:HD23	2:A:725:LEU:HA	1.78	0.46
2:A:305:ASP:OD1	2:A:305:ASP:N	2.47	0.45
3:T:5:DA:H2''	3:T:6:DC:H5'	1.97	0.45
1:P:11:DA:N6	3:T:4:DT:C7	2.79	0.45
1:P:11:DA:N6	3:T:4:DT:H71	2.31	0.45
2:A:334:GLY:HA2	2:A:348:PRO:HD3	1.99	0.45
2:A:392:PHE:HA	2:A:479:GLN:HE22	1.82	0.45
1:P:11:DA:O3'	2:A:829:HIS:CE1	2.71	0.44
2:A:569:GLU:N	2:A:569:GLU:CD	2.76	0.44
2:A:607:ASN:ND2	2:A:610:LEU:HD12	2.33	0.44
1:P:1:DG:H2''	1:P:2:DC:O5'	2.18	0.44
1:P:11:DA:N6	2:A:710:PHE:HZ	2.16	0.43
2:A:728:SER:OG	2:A:731:GLU:HG3	2.18	0.43
2:A:692:VAL:HG21	2:A:701:MET:HE1	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:711:GLY:HA2	2:A:716:ILE:HG12	2.01	0.42
3:T:9:DG:H2''	3:T:10:DA:H5'	2.01	0.42
2:A:814:ARG:HD2	2:A:814:ARG:HA	1.91	0.42
2:A:363:ASP:OD2	2:A:366:LYS:NZ	2.47	0.41
2:A:545:PRO:HD2	2:A:564:LEU:HD22	2.02	0.41
2:A:579:GLN:HG3	2:A:580:LEU:N	2.36	0.41
2:A:587:TYR:O	2:A:591:LEU:HB2	2.20	0.41
2:A:306:ARG:NH1	2:A:308:THR:HG22	2.35	0.40
2:A:401:LEU:HD23	2:A:401:LEU:HA	1.82	0.40
2:A:494:LYS:HB2	2:A:494:LYS:HE3	1.67	0.40
2:A:709:ASN:HD22	2:A:709:ASN:HA	1.72	0.40
2:A:816:LYS:HA	2:A:816:LYS:HD2	1.88	0.40
2:A:517:ARG:CZ	2:A:521:LEU:HD11	2.51	0.40
2:A:561:LEU:O	2:A:571:VAL:HG11	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	
2	A	575/580 (99%)	560 (97%)	15 (3%)	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
2	A	486/496 (98%)	484 (100%)	2 (0%)	84 89

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	505	GLU
2	A	823	ARG

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

Mol	Chain	Res	Type
2	A	468	ASN
2	A	470	GLN
2	A	479	GLN
2	A	510	GLN
2	A	568	HIS
2	A	579	GLN
2	A	656	GLN
2	A	709	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is 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	SO4	A	902	-	4,4,4	0.26	0	6,6,6	0.15	0

There are no bond length outliers.

There are no bond angle outliers.

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 [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	P	11/11 (100%)	1.32	2 (18%) 3 3	48, 70, 108, 114	0
2	A	577/580 (99%)	0.55	42 (7%) 21 19	13, 30, 62, 97	0
3	T	12/12 (100%)	1.71	6 (50%) 0 1	58, 81, 147, 150	0
All	All	600/603 (99%)	0.59	50 (8%) 17 16	13, 31, 69, 150	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	512	ARG	4.6
2	A	876	LYS	4.4
2	A	505	GLU	4.0
2	A	631	GLU	3.9
2	A	559	ASP	3.7
2	A	632	GLU	3.6
2	A	550	THR	3.4
2	A	634	ARG	3.4
2	A	300	ALA	3.4
2	A	735	PHE	3.3
1	P	11	DA	3.3
3	T	4	DT	3.3
2	A	547	LEU	3.3
2	A	829	HIS	3.0
2	A	532	LYS	3.0
2	A	719	TYR	3.0
2	A	552	THR	2.9
2	A	408	ASP	2.9
2	A	628	ILE	2.9
2	A	314	ASP	2.8
2	A	305	ASP	2.8
2	A	520	GLU	2.7
2	A	777	THR	2.7

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Mol	Chain	Res	Type	RSRZ
2	A	511	LEU	2.7
2	A	734	GLU	2.6
2	A	549	LYS	2.6
3	T	6	DC	2.5
2	A	531	PRO	2.4
2	A	865	ASP	2.3
2	A	694	GLU	2.3
2	A	304	ALA	2.3
2	A	303	LEU	2.3
2	A	463	ASP	2.3
3	T	14	DC	2.3
2	A	750	MET	2.3
2	A	629	ARG	2.2
2	A	645	SER	2.2
2	A	570	ILE	2.2
2	A	689	ILE	2.2
1	P	1	DG	2.2
3	T	13	DG	2.2
2	A	506	GLU	2.2
2	A	653	ASP	2.2
2	A	729	ARG	2.1
2	A	554	TYR	2.1
2	A	584	GLN	2.1
3	T	5	DA	2.1
2	A	433	ALA	2.0
2	A	646	ASP	2.0
3	T	7	DG	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
5	SO4	A	902	5/5	0.87	0.09	70,70,73,75	0
4	MG	A	901	1/1	0.92	0.29	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.