



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:03 PM GMT

PDB ID : 4BDP  
Title : CRYSTAL STRUCTURE OF BACILLUS DNA POLYMERASE I FRAGMENT COMPLEXED TO 11 BASE PAIRS OF DUPLEX DNA AFTER ADDITION OF TWO DATP RESIDUES  
Authors : Kiefer, J.R.; Mao, C.; Beese, L.S.  
Deposited on : 1997-11-17  
Resolution : 1.80 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

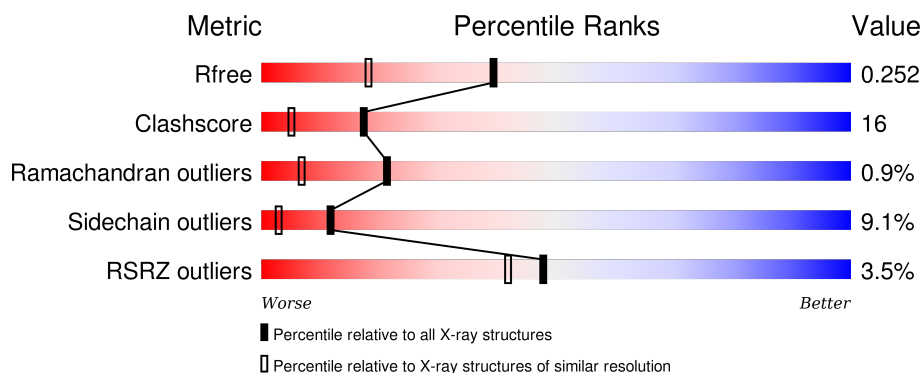
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	11	
2	T	13	
3	A	580	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	A	910	-	X	-	-
4	SO4	A	911	-	X	-	-
4	SO4	A	912	-	X	-	-

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5707 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\*AP\*TP\*CP\*AP\*TP\*GP\*CP\*AP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	P	11	Total	C	N	O	P	0	0	0
			222	107	43	62	10			

- Molecule 2 is a DNA chain called DNA (5'-D(\*TP\*AP\*TP\*TP\*GP\*CP\*AP\*TP\*GP\*AP\*TP\*GP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	13	Total	C	N	O	P	0	0	0
			264	128	46	78	12			

- Molecule 3 is a protein called PROTEIN (DNA POLYMERASE I).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	580	Total	C	N	O	S	0	0	0
			4650	2956	807	870	17			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	298	ALA	LYS	CONFLICT	UNP P52026
A	411	ARG	ALA	CONFLICT	UNP P52026
A	456	GLU	ALA	CONFLICT	UNP P52026
A	505	LYS	GLU	CONFLICT	UNP P52026
A	512	GLY	ARG	CONFLICT	UNP P52026
A	550	THR	SER	CONFLICT	UNP P52026
A	?	-	GLN	DELETION	UNP P52026
A	823	HIS	ARG	CONFLICT	UNP P52026

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

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

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

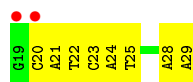
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	496	Total	O	0	0
			496	496		
6	P	28	Total	O	0	0
			28	28		
6	T	30	Total	O	0	0
			30	30		

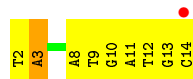
### 3 Residue-property plots

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.

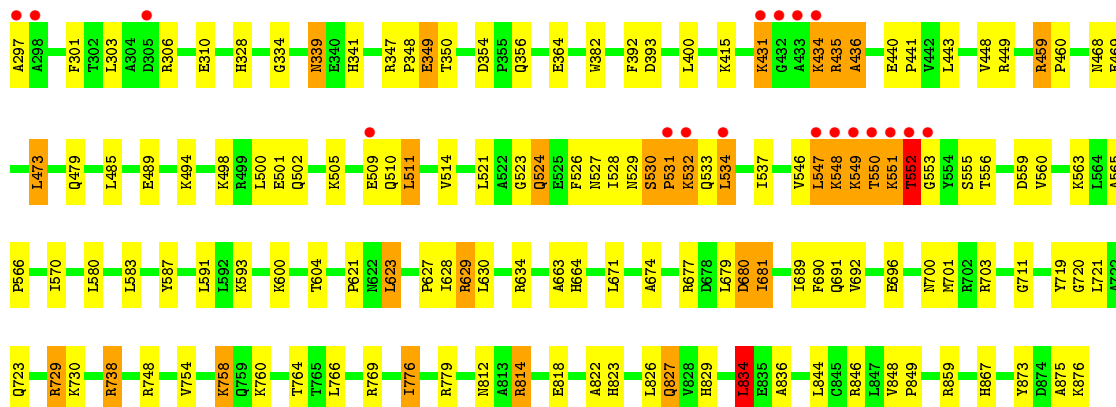
- Molecule 1: DNA (5'-D(\*GP\*CP\*AP\*TP\*CP\*AP\*TP\*GP\*CP\*AP\*A)-3')



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



- Molecule 3: PROTEIN (DNA POLYMERASE I)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.24Å 93.28Å 106.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.80 19.57 – 1.80	Depositor EDS
% Data completeness (in resolution range)	85.9 (20.00-1.80) 90.2 (19.57-1.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.85 (at 1.80Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.248 , 0.296 0.203 , 0.252	Depositor DCC
$R_{free}$ test set	3680 reflections (5.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.8	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 64.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 72389 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SO4

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	0.57	0/249	0.83	1/382 (0.3%)
2	T	0.70	0/295	0.84	0/454
3	A	0.46	0/4734	0.68	1/6398 (0.0%)
All	All	0.48	0/5278	0.70	2/7234 (0.0%)

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
2	T	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	834	LEU	CA-CB-CG	8.45	134.72	115.30
1	P	25	DT	C1'-O4'-C4'	-5.97	104.13	110.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	T	3	DA	Sidechain



## 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	P	222	0	125	11	0
2	T	264	0	150	22	0
3	A	4650	0	4698	143	0
4	A	15	0	0	0	0
5	A	2	0	0	0	0
6	A	496	0	0	11	0
6	P	28	0	0	0	0
6	T	30	0	0	2	0
All	All	5707	0	4973	163	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 16.

All (163) 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:548:LYS:HD3	3:A:560:VAL:HG22	1.35	1.03
2:T:2:DT:H4'	2:T:3:DA:C5'	1.91	1.00
2:T:2:DT:H4'	2:T:3:DA:H5''	1.44	0.98
2:T:2:DT:C4'	2:T:3:DA:H5''	1.93	0.98
3:A:549:LYS:HG2	3:A:553:GLY:O	1.67	0.95
3:A:347:ARG:HB3	3:A:349:GLU:OE1	1.69	0.93
3:A:700:ASN:HD21	3:A:703:ARG:HH21	1.13	0.93
3:A:591:LEU:HD22	6:A:2713:HOH:O	1.71	0.90
3:A:431:LYS:H	3:A:434:LYS:HD2	1.39	0.86
3:A:867:HIS:HB2	3:A:875:ALA:O	1.76	0.86
1:P:22:DT:H2''	1:P:23:DC:H5'	1.58	0.85
3:A:677:ARG:HB2	3:A:679:LEU:HD13	1.58	0.82
3:A:500:LEU:HD21	3:A:591:LEU:HD23	1.61	0.81
3:A:328:HIS:HD2	3:A:382:TRP:HE1	1.29	0.81
3:A:339:ASN:ND2	3:A:341:HIS:H	1.79	0.81
3:A:550:THR:CG2	3:A:552:THR:HG23	2.12	0.79
2:T:9:DT:H2'	2:T:10:DG:C8	2.20	0.77
2:T:2:DT:H4'	2:T:3:DA:H5'	1.66	0.76
3:A:719:TYR:CE1	3:A:729:ARG:HG2	2.24	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:24:DA:OP1	3:A:551:LYS:HG2	1.90	0.71
3:A:677:ARG:CB	3:A:679:LEU:HD13	2.19	0.71
3:A:547:LEU:O	3:A:548:LYS:HG3	1.90	0.70
1:P:28:DA:P	3:A:629:ARG:HG3	2.31	0.70
2:T:2:DT:C5'	2:T:3:DA:H5''	2.22	0.70
3:A:700:ASN:ND2	3:A:703:ARG:HH21	1.88	0.69
3:A:339:ASN:HD22	3:A:339:ASN:C	1.95	0.69
3:A:485:LEU:O	3:A:489:GLU:HG3	1.94	0.67
3:A:532:LYS:NZ	3:A:532:LYS:HB3	2.10	0.66
3:A:621:PRO:HG2	3:A:623:LEU:HD13	1.77	0.66
3:A:431:LYS:N	3:A:434:LYS:HD2	2.10	0.66
2:T:2:DT:H5''	2:T:3:DA:H5''	1.78	0.65
3:A:758:LYS:HD2	3:A:776:ILE:HG12	1.76	0.65
3:A:546:VAL:O	3:A:546:VAL:HG12	1.96	0.65
3:A:551:LYS:C	3:A:552:THR:HG22	2.15	0.65
1:P:22:DT:H2''	1:P:23:DC:C5'	2.25	0.65
6:T:1052:HOH:O	3:A:593:LYS:HE3	1.96	0.64
3:A:400:LEU:HD22	3:A:473:LEU:HD13	1.79	0.64
3:A:498:LYS:O	3:A:502:GLN:HG3	1.97	0.64
3:A:435:ARG:O	3:A:436:ALA:HB2	1.97	0.64
3:A:339:ASN:ND2	3:A:341:HIS:N	2.44	0.64
2:T:3:DA:H5'	3:A:720:GLY:HA3	1.78	0.64
3:A:431:LYS:O	3:A:434:LYS:HG3	1.98	0.64
2:T:12:DT:H5''	3:A:532:LYS:HZ3	1.61	0.63
3:A:530:SER:O	3:A:532:LYS:N	2.31	0.63
3:A:690:PHE:CD2	3:A:701:MET:HE3	2.35	0.62
1:P:24:DA:OP1	3:A:550:THR:HG23	2.00	0.62
3:A:548:LYS:CD	3:A:560:VAL:HG22	2.21	0.61
2:T:3:DA:H5'	3:A:720:GLY:CA	2.30	0.61
3:A:431:LYS:HD3	3:A:434:LYS:HZ2	1.65	0.61
3:A:551:LYS:HG3	3:A:551:LYS:O	2.00	0.60
3:A:738:ARG:HE	3:A:738:ARG:HA	1.66	0.60
3:A:623:LEU:HD23	3:A:826:LEU:HD21	1.83	0.60
3:A:674:ALA:HA	3:A:679:LEU:HD22	1.83	0.60
3:A:339:ASN:HD22	3:A:341:HIS:N	2.00	0.60
3:A:550:THR:C	3:A:552:THR:H	2.05	0.59
3:A:534:LEU:HD13	3:A:556:THR:HG21	1.84	0.59
3:A:550:THR:HG23	3:A:551:LYS:N	2.17	0.59
3:A:827:GLN:NE2	3:A:829:HIS:H	2.00	0.59
2:T:3:DA:C2	3:A:711:GLY:HA3	2.38	0.58
3:A:549:LYS:HA	3:A:555:SER:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:2:DT:C3'	2:T:3:DA:H5"	2.34	0.58
3:A:339:ASN:HD22	3:A:341:HIS:H	1.49	0.58
3:A:431:LYS:HD3	3:A:434:LYS:NZ	2.19	0.56
3:A:526:PHE:HE2	3:A:534:LEU:HD23	1.70	0.56
3:A:534:LEU:CD1	3:A:556:THR:HG21	2.36	0.56
3:A:529:ASN:O	3:A:531:PRO:HD3	2.06	0.55
3:A:339:ASN:HD21	3:A:341:HIS:HB2	1.70	0.55
1:P:28:DA:OP2	3:A:629:ARG:HG3	2.06	0.55
3:A:729:ARG:H	3:A:729:ARG:HD2	1.72	0.55
3:A:532:LYS:CB	3:A:532:LYS:NZ	2.70	0.54
3:A:680:ASP:HB3	6:A:2666:HOH:O	2.07	0.54
3:A:846:ARG:HG3	6:A:2171:HOH:O	2.08	0.54
3:A:621:PRO:HG2	3:A:623:LEU:CD1	2.38	0.53
1:P:28:DA:H2'	1:P:29:DA:C8	2.44	0.53
3:A:494:LYS:HE3	3:A:600:LYS:HB2	1.89	0.53
3:A:530:SER:O	3:A:533:GLN:N	2.42	0.53
3:A:524:GLN:NE2	3:A:537:ILE:HD11	2.24	0.52
3:A:664:HIS:O	3:A:859:ARG:NH1	2.42	0.52
3:A:634:ARG:HD2	6:A:2649:HOH:O	2.09	0.52
3:A:729:ARG:CD	3:A:729:ARG:H	2.20	0.52
3:A:393:ASP:H	3:A:479:GLN:NE2	2.08	0.52
3:A:328:HIS:CD2	3:A:382:TRP:HE1	2.17	0.52
3:A:547:LEU:HB2	3:A:548:LYS:HD2	1.92	0.52
3:A:548:LYS:HD3	3:A:560:VAL:CG2	2.25	0.51
3:A:521:LEU:HD13	3:A:570:ILE:HA	1.91	0.51
1:P:20:DC:H2"	1:P:21:DA:OP2	2.09	0.51
3:A:550:THR:CG2	3:A:551:LYS:N	2.73	0.51
3:A:738:ARG:NE	3:A:738:ARG:HA	2.25	0.51
3:A:729:ARG:HD2	6:A:2568:HOH:O	2.11	0.50
3:A:440:GLU:HB3	3:A:441:PRO:HD3	1.93	0.50
3:A:459:ARG:HB3	3:A:460:PRO:HD3	1.93	0.50
3:A:550:THR:N	3:A:553:GLY:O	2.44	0.50
2:T:13:DG:H2"	2:T:14:DC:C5	2.47	0.50
3:A:550:THR:O	3:A:552:THR:N	2.43	0.50
3:A:551:LYS:CG	3:A:551:LYS:O	2.60	0.50
3:A:692:VAL:HB	3:A:696:GLU:HB2	1.93	0.50
3:A:392:PHE:HA	3:A:479:GLN:HE22	1.76	0.49
3:A:468:ASN:O	3:A:469:GLU:HB2	2.12	0.49
3:A:511:LEU:HD13	3:A:580:LEU:HB2	1.94	0.48
3:A:867:HIS:CG	3:A:876:LYS:HE2	2.48	0.48
1:P:28:DA:OP1	3:A:629:ARG:HG3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:531:PRO:N	6:A:2633:HOH:O	2.47	0.48
3:A:550:THR:HG22	3:A:552:THR:N	2.27	0.48
3:A:510:GLN:O	3:A:514:VAL:HG23	2.12	0.48
3:A:551:LYS:O	3:A:552:THR:HG22	2.13	0.48
3:A:754:VAL:O	3:A:758:LYS:HD3	2.13	0.47
3:A:550:THR:HG22	3:A:552:THR:HG23	1.96	0.47
3:A:551:LYS:C	3:A:552:THR:CG2	2.82	0.47
3:A:339:ASN:ND2	3:A:339:ASN:C	2.67	0.47
3:A:681:ILE:H	3:A:681:ILE:HG13	1.54	0.47
2:T:10:DG:H1'	2:T:11:DA:H5''	1.95	0.47
3:A:681:ILE:HG13	6:A:2123:HOH:O	2.15	0.47
2:T:3:DA:N1	3:A:721:LEU:HD13	2.30	0.47
3:A:530:SER:C	3:A:532:LYS:N	2.69	0.46
3:A:689:ILE:O	3:A:738:ARG:NH1	2.49	0.46
3:A:523:GLY:O	3:A:524:GLN:HB3	2.14	0.46
3:A:334:GLY:HA2	3:A:348:PRO:HD3	1.98	0.46
3:A:814:ARG:NH2	3:A:818:GLU:OE2	2.47	0.46
3:A:301:PHE:CG	3:A:448:VAL:HG21	2.51	0.46
3:A:527:ASN:OD1	3:A:529:ASN:HB2	2.16	0.45
3:A:549:LYS:HA	3:A:555:SER:N	2.32	0.45
3:A:634:ARG:HD3	6:A:2531:HOH:O	2.16	0.45
3:A:822:ALA:CB	3:A:836:ALA:HB2	2.46	0.45
3:A:559:ASP:O	3:A:563:LYS:HG3	2.16	0.45
3:A:634:ARG:HG2	3:A:873:TYR:CE2	2.52	0.45
3:A:459:ARG:HB3	3:A:460:PRO:CD	2.46	0.45
3:A:760:LYS:HE2	6:A:2591:HOH:O	2.16	0.45
3:A:527:ASN:O	3:A:529:ASN:N	2.50	0.44
2:T:10:DG:H2''	2:T:11:DA:H5'	1.98	0.44
3:A:297:ALA:O	3:A:449:ARG:NH2	2.51	0.44
2:T:9:DT:H1'	6:T:2361:HOH:O	2.16	0.44
3:A:364:GLU:O	3:A:364:GLU:HG3	2.16	0.44
3:A:848:VAL:HB	3:A:849:PRO:HD3	1.98	0.44
3:A:310:GLU:HB2	3:A:341:HIS:CD2	2.53	0.44
2:T:8:DA:H4'	6:A:2571:HOH:O	2.17	0.44
3:A:764:THR:HA	3:A:769:ARG:O	2.18	0.44
3:A:551:LYS:HG2	3:A:552:THR:HG22	2.00	0.43
2:T:12:DT:OP1	3:A:530:SER:HB3	2.18	0.43
3:A:550:THR:HG22	3:A:552:THR:H	1.83	0.43
3:A:876:LYS:HG2	6:A:2601:HOH:O	2.17	0.43
3:A:532:LYS:HB3	3:A:532:LYS:HZ1	1.79	0.43
3:A:587:TYR:CE1	3:A:627:PRO:HD3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:10:DG:H2''	2:T:11:DA:C5'	2.49	0.43
3:A:719:TYR:CD1	3:A:729:ARG:HG2	2.54	0.42
1:P:24:DA:OP2	3:A:551:LYS:HE3	2.19	0.42
3:A:527:ASN:C	3:A:529:ASN:N	2.72	0.42
3:A:691:GLN:HE21	3:A:691:GLN:HB2	1.53	0.42
3:A:511:LEU:CD1	3:A:580:LEU:HB2	2.50	0.42
3:A:565:ALA:HB3	3:A:566:PRO:HD3	2.02	0.42
3:A:550:THR:HG21	3:A:552:THR:HG23	1.97	0.41
3:A:738:ARG:NE	3:A:738:ARG:CA	2.82	0.41
3:A:604:THR:HB	3:A:623:LEU:HD22	2.02	0.41
3:A:435:ARG:O	3:A:436:ALA:CB	2.64	0.41
1:P:22:DT:H2'	1:P:23:DC:C6	2.55	0.41
3:A:354:ASP:OD1	3:A:356:GLN:HB2	2.21	0.40
3:A:663:ALA:HB2	3:A:671:LEU:HG	2.03	0.40
2:T:11:DA:H2''	2:T:12:DT:O5'	2.21	0.40
3:A:823:HIS:O	3:A:834:LEU:HB3	2.21	0.40
3:A:549:LYS:HB3	3:A:549:LYS:HE3	1.48	0.40
3:A:677:ARG:CB	3:A:679:LEU:CD1	2.96	0.40
2:T:3:DA:H5'	3:A:720:GLY:HA2	2.03	0.40
3:A:431:LYS:HD3	3:A:434:LYS:CE	2.52	0.40
3:A:528:ILE:H	3:A:528:ILE:HG13	1.71	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	578/580 (100%)	558 (96%)	15 (3%)	5 (1%)	21 7

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	552	THR
3	A	531	PRO
3	A	551	LYS
3	A	436	ALA
3	A	628	ILE

### 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
3	A	495/496 (100%)	450 (91%)	45 (9%)	12 3

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	303	LEU
3	A	306	ARG
3	A	339	ASN
3	A	349	GLU
3	A	350	THR
3	A	415	LYS
3	A	431	LYS
3	A	434	LYS
3	A	435	ARG
3	A	443	LEU
3	A	459	ARG
3	A	473	LEU
3	A	501	GLU
3	A	505	LYS
3	A	509	GLU
3	A	511	LEU
3	A	524	GLN
3	A	530	SER
3	A	532	LYS
3	A	534	LEU
3	A	547	LEU
3	A	548	LYS

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Mol	Chain	Res	Type
3	A	549	LYS
3	A	550	THR
3	A	552	THR
3	A	583	LEU
3	A	623	LEU
3	A	629	ARG
3	A	630	LEU
3	A	680	ASP
3	A	681	ILE
3	A	723	GLN
3	A	729	ARG
3	A	730	LYS
3	A	738	ARG
3	A	748	ARG
3	A	758	LYS
3	A	766	LEU
3	A	776	ILE
3	A	779	ARG
3	A	812	ASN
3	A	814	ARG
3	A	827	GLN
3	A	834	LEU
3	A	844	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	328	HIS
3	A	339	ASN
3	A	341	HIS
3	A	479	GLN
3	A	624	GLN
3	A	656	GLN
3	A	691	GLN
3	A	700	ASN
3	A	704	GLN
3	A	827	GLN
3	A	854	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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	SO4	A	910	-	4,4,4	4.80	4 (100%)	6,6,6	0.18	0
4	SO4	A	911	5	4,4,4	4.63	4 (100%)	6,6,6	0.26	0
4	SO4	A	912	-	4,4,4	4.72	4 (100%)	6,6,6	0.47	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
4	SO4	A	910	-	-	0/0/0/0	0/0/0/0
4	SO4	A	911	5	-	0/0/0/0	0/0/0/0
4	SO4	A	912	-	-	0/0/0/0	0/0/0/0

All (12) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	911	SO4	O2-S	2.85	1.56	1.47
4	A	912	SO4	O2-S	3.02	1.57	1.47
4	A	910	SO4	O2-S	3.43	1.58	1.47
4	A	911	SO4	O1-S	3.47	1.59	1.47
4	A	912	SO4	O1-S	3.50	1.59	1.47
4	A	910	SO4	O1-S	3.65	1.59	1.47
4	A	912	SO4	O4-S	5.55	1.67	1.47
4	A	911	SO4	O3-S	5.65	1.67	1.47
4	A	910	SO4	O3-S	5.76	1.68	1.47
4	A	911	SO4	O4-S	5.81	1.68	1.47
4	A	910	SO4	O4-S	5.81	1.68	1.47
4	A	912	SO4	O3-S	6.08	1.69	1.47

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, 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	11/11 (100%)	0.22	2 (18%) 2 1	21, 26, 67, 71	0
2	T	13/13 (100%)	0.10	1 (7%) 16 13	19, 28, 64, 71	0
3	A	580/580 (100%)	0.02	18 (3%) 52 47	16, 23, 42, 66	0
All	All	604/604 (100%)	0.03	21 (3%) 48 42	16, 24, 45, 71	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	550	THR	7.2
3	A	549	LYS	7.0
3	A	552	THR	6.9
3	A	433	ALA	6.6
3	A	551	LYS	5.9
3	A	297	ALA	5.2
3	A	298	ALA	4.3
3	A	434	LYS	3.5
3	A	548	LYS	3.2
1	P	19	DG	3.2
3	A	547	LEU	3.2
2	T	14	DC	3.0
3	A	531	PRO	2.7
3	A	553	GLY	2.7
3	A	534	LEU	2.7
3	A	431	LYS	2.5
3	A	532	LYS	2.5
3	A	305	ASP	2.1
3	A	432	GLY	2.1
1	P	20	DC	2.0
3	A	509	GLU	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, 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
4	SO4	A	912	5/5	0.92	0.14	0.08	47,48,49,49	0
5	MG	A	950	1/1	0.93	0.09	-	36,36,36,36	0
4	SO4	A	910	5/5	0.94	0.16	-	47,47,48,48	0
5	MG	A	952	1/1	0.89	0.19	-	44,44,44,44	0
4	SO4	A	911	5/5	0.98	0.10	-	34,35,35,36	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.