Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:51 AM GMT

PDB ID	:	3CFP
Title	:	Structure of the replicating complex of a POL Alpha family DNA Polymerase, ternary complex 1
Authors	:	Wang, J.; Klimenko, D.; Wang, M.; Steitz, T.A.; Konigsberg, W.H.
Deposited on	:	2008-03-04
Resolution	:	2.50 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

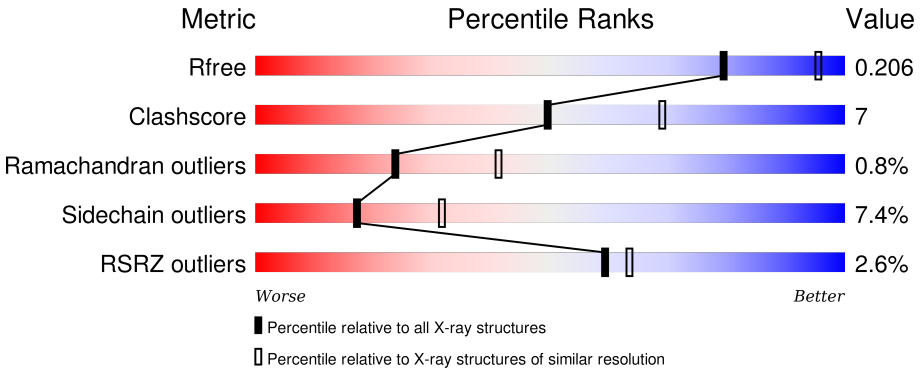
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 <=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	18	44% 44% 11%
2	P	14	43% 36% 21%
3	A	909	3% 79% 17%

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
5	CL	A	1005	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8326 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(*DAP*DCP*DAP*DGP*DGP*DTP*DAP*DAP*DGP*DCP*DAP*DGP*DTP*DCP*DCP*DGP*DCP*DG)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	T	18	369	175	74	103	17	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(*DGP*DCP*DGP*DGP*DAP*DCP*DTP*DGP*DCP*DTP*DTP*DAP*DCP*(DOC))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	P	14	281	135	51	82	13	0	0	0

- Molecule 3 is a protein called DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	901	7343	4714	1224	1372	33	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	561	ALA	LEU	ENGINEERED	UNP Q38087
A	565	GLY	SER	ENGINEERED	UNP Q38087
A	567	ALA	TYR	ENGINEERED	UNP Q38087
A	904	HIS	-	EXPRESSION TAG	UNP Q38087
A	905	HIS	-	EXPRESSION TAG	UNP Q38087
A	906	HIS	-	EXPRESSION TAG	UNP Q38087
A	907	HIS	-	EXPRESSION TAG	UNP Q38087
A	908	HIS	-	EXPRESSION TAG	UNP Q38087
A	909	HIS	-	EXPRESSION TAG	UNP Q38087

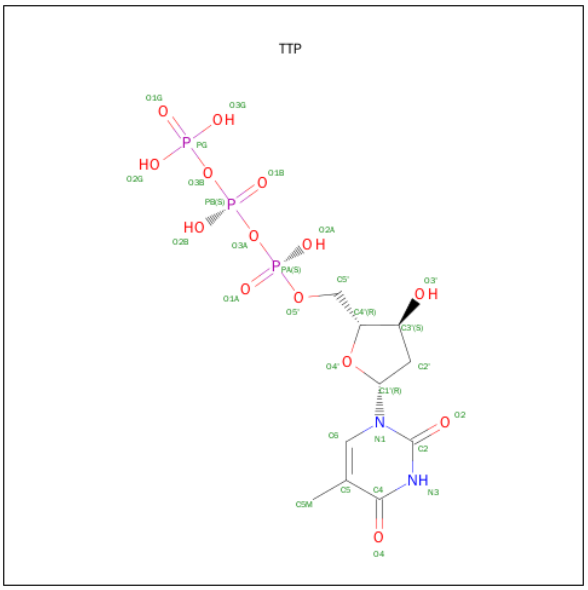
- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total Ca 4 4	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 6 is THYMIDINE-5'-TRIPHOSPHATE (three-letter code: TTP) (formula: $C_{10}H_{17}N_2O_{14}P_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C N O P 29 10 2 14 3	0	0

- Molecule 7 is water.

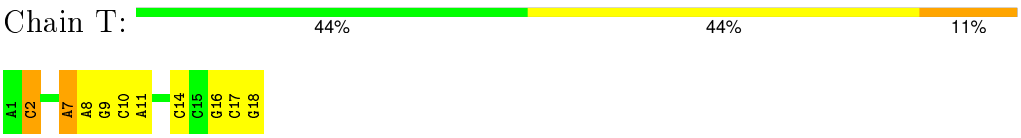
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	260	Total O 260 260	0	0
7	P	17	Total O 17 17	0	0
7	T	22	Total O 22 22	0	0

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 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.

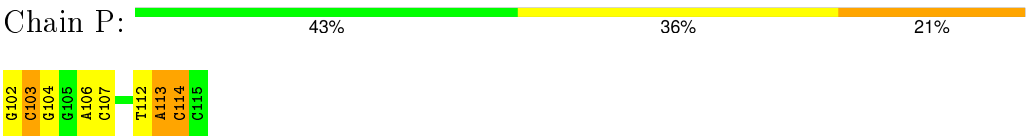
• Molecule 1: DNA (5'-D(*DAP*DCP*DAP*DGP*DGP*DTP*DAP*DAP*DGP*DCP*DAP*DGP*DTP*DCP*DCP*DGP*DCP*DG)-3')

Chain T:



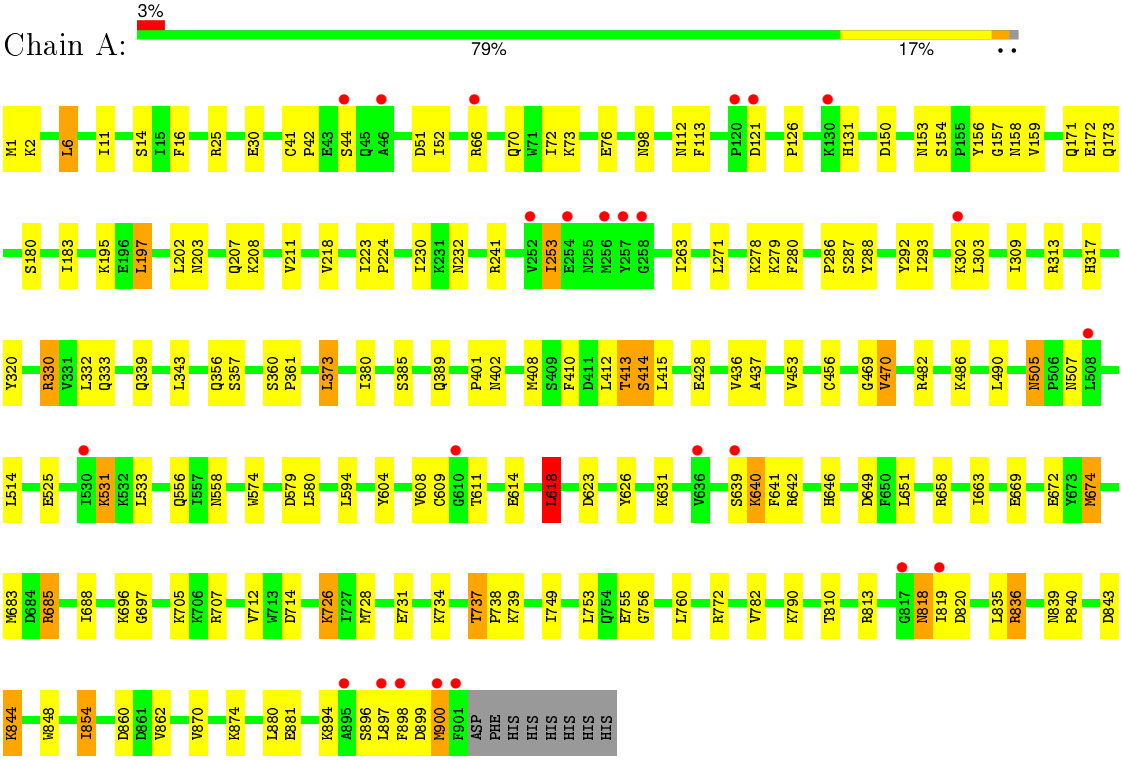
• Molecule 2: DNA (5'-D(*DGP*DCP*DGP*DGP*DAP*DCP*DTP*DGP*DCP*DTP*DTP*DAP*DCP*(DOC))-3')

Chain P:



• Molecule 3: DNA polymerase

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	80.73Å 117.28Å 126.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 38.46 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (50.00-2.50) 99.2 (38.46-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.29 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.197 , 0.267 0.199 , 0.206	Depositor DCC
R_{free} test set	2146 reflections (5.38%)	DCC
Wilson B-factor (Å^2)	52.0	Xtriage
Anisotropy	0.417	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 41.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 42069 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8326	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, CA, TTP, CL

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	1.80	8/415 (1.9%)	1.63	5/639 (0.8%)
2	P	0.91	0/294	1.59	6/452 (1.3%)
3	A	0.48	0/7522	0.60	3/10164 (0.0%)
All	All	0.63	8/8231 (0.1%)	0.76	14/11255 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	T	18	DG	N9-C8	15.97	1.49	1.37
1	T	18	DG	C5-C4	14.89	1.48	1.38
1	T	18	DG	N3-C4	13.77	1.45	1.35
1	T	18	DG	N1-C2	12.23	1.47	1.37
1	T	18	DG	C6-N1	7.82	1.45	1.39
1	T	18	DG	C6-O6	6.60	1.30	1.24
1	T	18	DG	C3'-C2'	6.32	1.59	1.52
1	T	18	DG	C2-N3	5.29	1.36	1.32

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	T	18	DG	C2-N3-C4	-11.46	106.17	111.90
1	T	2	DC	P-O3'-C3'	6.55	127.56	119.70
2	P	104	DG	P-O3'-C3'	6.45	127.44	119.70
2	P	103	DC	P-O3'-C3'	6.32	127.29	119.70
3	A	197	LEU	CA-CB-CG	6.23	129.64	115.30
3	A	618	LEU	CA-CB-CG	6.02	129.15	115.30
1	T	14	DC	O4'-C1'-N1	5.79	112.05	108.00
1	T	7	DA	P-O3'-C3'	5.78	126.64	119.70
2	P	107	DC	C1'-O4'-C4'	-5.66	104.44	110.10
2	P	114	DC	N1-C2-O2	5.63	122.28	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	760	LEU	CA-CB-CG	5.47	127.88	115.30
2	P	113	DA	O5'-P-OP2	-5.40	100.84	105.70
1	T	18	DG	C5-C6-N1	-5.36	108.82	111.50
2	P	112	DT	O4'-C1'-N1	5.23	111.66	108.00

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	T	369	0	202	7	0
2	P	281	0	159	4	0
3	A	7343	0	7236	102	0
4	A	4	0	0	0	0
5	A	1	0	0	2	0
6	A	29	0	13	1	0
7	A	260	0	0	12	0
7	P	17	0	0	1	0
7	T	22	0	0	0	0
All	All	8326	0	7610	111	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 7.

All (111) 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:T:16:DG:H2"	1:T:17:DC:H5'	1.54	0.88
3:A:640:LYS:HG3	3:A:641:PHE:H	1.38	0.88
3:A:408:MET:HE3	3:A:651:LEU:HB3	1.61	0.83
3:A:737:THR:HG22	7:A:1062:HOH:O	1.78	0.83
3:A:640:LYS:HG3	3:A:641:PHE:N	1.96	0.80
3:A:6:LEU:HG	3:A:211:VAL:HG21	1.69	0.74
3:A:836:ARG:HD3	7:A:1222:HOH:O	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:41:CYS:HB2	3:A:42:PRO:HD2	1.71	0.73
3:A:408:MET:CE	3:A:651:LEU:HB3	2.19	0.73
3:A:514:LEU:HD13	3:A:533:LEU:HD11	1.72	0.70
3:A:373:LEU:HD13	3:A:470:VAL:HG11	1.74	0.68
3:A:456:CYS:O	3:A:674:MET:HG3	1.94	0.68
3:A:183:ILE:O	7:A:1052:HOH:O	2.12	0.67
3:A:839:ASN:HB2	3:A:840:PRO:HD2	1.78	0.65
3:A:881:GLU:HG3	7:A:1037:HOH:O	1.95	0.65
1:T:16:DG:H2"	1:T:17:DC:C5'	2.26	0.64
3:A:373:LEU:HD23	3:A:380:ILE:HG22	1.80	0.64
3:A:408:MET:HE2	3:A:688:ILE:HG12	1.82	0.62
3:A:707:ARG:NH2	3:A:731:GLU:OE1	2.32	0.62
3:A:505:ASN:O	3:A:505:ASN:CG	2.39	0.61
3:A:697:GLY:HA2	3:A:755:GLU:O	2.01	0.60
1:T:10:DC:H2"	1:T:11:DA:H5"	1.84	0.59
3:A:436:VAL:HG22	3:A:437:ALA:O	2.02	0.59
3:A:731:GLU:O	3:A:737:THR:HG21	2.04	0.57
5:A:1005:CL:CL	7:A:1061:HOH:O	2.54	0.57
3:A:618:LEU:HD13	3:A:626:TYR:O	2.05	0.57
3:A:641:PHE:HD2	3:A:646:HIS:ND1	2.03	0.57
3:A:232:ASN:OD1	5:A:1005:CL:CL	2.60	0.57
3:A:280:PHE:HD2	3:A:558:ASN:OD1	1.87	0.56
1:T:7:DA:H5"	3:A:705:LYS:HD3	1.88	0.55
3:A:385:SER:HB2	7:A:1239:HOH:O	2.06	0.55
2:P:102:DG:H2"	2:P:103:DC:OP2	2.07	0.55
3:A:223:ILE:HB	3:A:224:PRO:HD3	1.87	0.55
1:T:8:DA:H2'	1:T:9:DG:C8	2.41	0.54
1:T:2:DC:H5'	3:A:574:TRP:CD1	2.43	0.54
3:A:72:ILE:O	3:A:76:GLU:HG3	2.08	0.53
3:A:330:ARG:HA	3:A:333:GLN:HE21	1.73	0.53
3:A:641:PHE:HA	3:A:646:HIS:CD2	2.44	0.53
3:A:373:LEU:HD13	3:A:470:VAL:CG1	2.37	0.53
3:A:640:LYS:O	3:A:641:PHE:HB2	2.09	0.52
3:A:389:GLN:HB3	7:A:1243:HOH:O	2.09	0.52
3:A:738:PRO:HA	7:A:1073:HOH:O	2.09	0.52
3:A:52:ILE:HD12	3:A:428:GLU:HB3	1.90	0.51
3:A:203:ASN:ND2	3:A:241:ARG:HH12	2.07	0.51
3:A:818:ASN:CG	3:A:818:ASN:O	2.49	0.51
3:A:870:VAL:HG13	3:A:874:LYS:HD3	1.92	0.51
3:A:66:ARG:HE	3:A:70:GLN:HE21	1.57	0.51
3:A:153:ASN:HA	3:A:158:ASN:HD22	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:271:LEU:HD21	3:A:356:GLN:HA	1.92	0.50
3:A:896:SER:C	3:A:898:PHE:H	2.15	0.50
3:A:641:PHE:HA	3:A:646:HIS:CG	2.46	0.50
3:A:696:LYS:HB2	3:A:756:GLY:HA3	1.93	0.49
3:A:685:ARG:NH2	3:A:714:ASP:OD1	2.21	0.49
3:A:202:LEU:HD21	3:A:230:ILE:HD13	1.93	0.49
2:P:113:DA:H5'	3:A:734:LYS:HG2	1.95	0.49
1:T:11:DA:OP1	3:A:874:LYS:HE3	2.13	0.48
2:P:113:DA:H2"	2:P:114:DC:H5'	1.95	0.48
3:A:224:PRO:HA	3:A:263:ILE:HD13	1.95	0.48
3:A:696:LYS:HB2	3:A:756:GLY:CA	2.44	0.47
6:A:1006:TTP:O1B	6:A:1006:TTP:H3'	2.15	0.47
3:A:131:HIS:HD2	3:A:156:TYR:OH	1.97	0.47
3:A:810:THR:HG22	3:A:813:ARG:HH22	1.80	0.47
3:A:614:GLU:HG2	3:A:631:LYS:HE3	1.96	0.46
3:A:2:LYS:NZ	7:A:1106:HOH:O	2.48	0.46
3:A:253:ILE:HG22	3:A:253:ILE:O	2.15	0.46
3:A:76:GLU:HG2	7:A:1114:HOH:O	2.15	0.46
3:A:203:ASN:O	3:A:207:GLN:HG2	2.16	0.46
3:A:609:CYS:C	3:A:611:THR:H	2.19	0.45
3:A:126:PRO:HB3	3:A:224:PRO:HB2	1.98	0.45
3:A:507:ASN:ND2	3:A:531:LYS:O	2.50	0.45
3:A:413:THR:O	3:A:414:SER:C	2.54	0.45
3:A:836:ARG:HH11	3:A:836:ARG:HB3	1.80	0.45
3:A:11:ILE:HD12	3:A:16:PHE:CD2	2.52	0.45
3:A:288:TYR:HA	3:A:293:ILE:CD1	2.47	0.45
3:A:159:VAL:HG21	3:A:317:HIS:CD2	2.52	0.45
3:A:609:CYS:HB2	3:A:611:THR:HG22	1.99	0.45
3:A:415:LEU:HD22	3:A:623:ASP:HB3	1.99	0.44
3:A:1:MET:HE2	7:A:1257:HOH:O	2.17	0.44
3:A:696:LYS:HB2	3:A:756:GLY:N	2.33	0.43
3:A:772:ARG:HD3	7:A:1086:HOH:O	2.18	0.43
3:A:640:LYS:CG	3:A:641:PHE:H	2.21	0.43
3:A:663:ILE:HG21	3:A:683:MET:HB3	2.00	0.43
3:A:486:LYS:HE2	3:A:490:LEU:HD11	1.99	0.43
3:A:401:PRO:O	3:A:402:ASN:HB2	2.19	0.43
3:A:51:ASP:OD1	3:A:51:ASP:C	2.56	0.43
3:A:844:LYS:H	3:A:844:LYS:HG3	1.68	0.42
3:A:410:PHE:HB3	3:A:683:MET:HG2	2.01	0.42
3:A:286:PRO:HB3	3:A:782:VAL:HG21	2.01	0.42
2:P:106:DA:H8	7:P:133:HOH:O	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:507:ASN:HD22	3:A:507:ASN:N	2.18	0.42
3:A:604:TYR:OH	3:A:658:ARG:HB3	2.20	0.42
3:A:712:VAL:CG2	3:A:726:LYS:HB2	2.50	0.42
3:A:360:SER:HA	3:A:361:PRO:HD2	1.85	0.42
3:A:408:MET:CE	3:A:688:ILE:HG12	2.50	0.42
3:A:150:ASP:OD2	3:A:317:HIS:CE1	2.73	0.42
3:A:428:GLU:OE2	3:A:469:GLY:HA2	2.20	0.41
3:A:157:GLY:O	3:A:313:ARG:NH2	2.53	0.41
3:A:640:LYS:CG	3:A:641:PHE:N	2.75	0.41
3:A:896:SER:OG	3:A:898:PHE:HB2	2.19	0.41
3:A:810:THR:HG21	3:A:843:ASP:HB3	2.03	0.41
3:A:669:GLU:O	3:A:672:GLU:HB3	2.21	0.41
3:A:839:ASN:HB2	3:A:840:PRO:CD	2.50	0.41
3:A:171:GLN:C	3:A:173:GLN:H	2.23	0.41
3:A:253:ILE:CG2	3:A:253:ILE:O	2.70	0.40
3:A:14:SER:HB2	3:A:30:GLU:HG3	2.02	0.40
3:A:313:ARG:HG3	3:A:320:TYR:CD2	2.56	0.40
3:A:848:TRP:CE2	3:A:854:ILE:HG12	2.55	0.40
3:A:749:ILE:O	3:A:753:LEU:HG	2.22	0.40
3:A:639:SER:O	3:A:640:LYS:CB	2.69	0.40
3:A:412:LEU:HB2	3:A:623:ASP:HB2	2.03	0.40
3:A:292:TYR:OH	3:A:739:LYS:NZ	2.52	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
3	A	899/909 (99%)	836 (93%)	56 (6%)	7 (1%)	24 41

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	253	ILE
3	A	640	LYS
3	A	820	ASP
3	A	172	GLU
3	A	579	ASP
3	A	414	SER
3	A	900	MET

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	795/805 (99%)	736 (93%)	59 (7%)	17 31

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	6	LEU
3	A	25	ARG
3	A	44	SER
3	A	73	LYS
3	A	98	ASN
3	A	112	ASN
3	A	113	PHE
3	A	121	ASP
3	A	154	SER
3	A	180	SER
3	A	195	LYS
3	A	197	LEU
3	A	208	LYS
3	A	218	VAL
3	A	278	LYS
3	A	279	LYS
3	A	287	SER
3	A	302	LYS
3	A	303	LEU
3	A	309	ILE

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Mol	Chain	Res	Type
3	A	330	ARG
3	A	332	LEU
3	A	339	GLN
3	A	343	LEU
3	A	357	SER
3	A	373	LEU
3	A	413	THR
3	A	453	VAL
3	A	470	VAL
3	A	482	ARG
3	A	505	ASN
3	A	525	GLU
3	A	531	LYS
3	A	556	GLN
3	A	580	LEU
3	A	594	LEU
3	A	608	VAL
3	A	618	LEU
3	A	642	ARG
3	A	649	ASP
3	A	674	MET
3	A	685	ARG
3	A	726	LYS
3	A	728	MET
3	A	737	THR
3	A	790	LYS
3	A	818	ASN
3	A	819	ILE
3	A	835	LEU
3	A	836	ARG
3	A	844	LYS
3	A	854	ILE
3	A	860	ASP
3	A	862	VAL
3	A	880	LEU
3	A	894	LYS
3	A	897	LEU
3	A	899	ASP
3	A	900	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	70	GLN
3	A	98	ASN
3	A	112	ASN
3	A	131	HIS
3	A	158	ASN
3	A	203	ASN
3	A	228	ASN
3	A	333	GLN
3	A	505	ASN
3	A	507	ASN
3	A	546	GLN
3	A	564	ASN
3	A	678	GLN
3	A	761	GLN
3	A	823	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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	$\#\|Z\| > 2$
2	DOC	P	115	1,2	11,19,20	0.58	0	14,26,29	1.66	2 (14%)

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
2	DOC	P	115	1,2	-	0/3/18/19	0/2/2/2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	P	115	DOC	O4'-C1'-N1	2.90	112.74	107.72
2	P	115	DOC	C2-N3-C4	3.63	120.73	115.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 5 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	$\#\|Z\| > 2$
6	TTP	A	1006	4	21,30,30	0.54	0	31,47,47	2.08	7 (22%)

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
6	TTP	A	1006	4	-	0/18/34/34	0/2/2/2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	1006	TTP	C5-C4-N3	-5.10	119.45	125.14
6	A	1006	TTP	C2'-C1'-N1	-3.95	104.55	114.16
6	A	1006	TTP	PB-O3B-PG	-3.26	121.73	132.67
6	A	1006	TTP	PB-O3A-PA	-2.28	126.31	132.73
6	A	1006	TTP	O3A-PA-O5'	2.14	108.61	102.94
6	A	1006	TTP	O4'-C1'-N1	4.15	114.90	107.72
6	A	1006	TTP	C4-N3-C2	6.23	120.64	115.25

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
6	A	1006	TTP	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, 95^{th} 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	18/18 (100%)	-0.15	0 100 100	35, 46, 69, 88	0
2	P	13/14 (92%)	-0.03	0 100 100	38, 48, 62, 68	0
3	A	901/909 (99%)	0.11	24 (2%) 58 62	33, 53, 79, 114	0
All	All	932/941 (99%)	0.10	24 (2%) 59 63	33, 53, 79, 114	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	819	ILE	7.0
3	A	901	PHE	6.2
3	A	257	TYR	5.0
3	A	258	GLY	5.0
3	A	256	MET	4.5
3	A	895	ALA	4.3
3	A	44	SER	3.5
3	A	898	PHE	3.5
3	A	610	GLY	3.3
3	A	66	ARG	3.2
3	A	900	MET	3.2
3	A	252	VAL	3.0
3	A	508	LEU	3.0
3	A	120	PRO	2.8
3	A	530	ILE	2.5
3	A	46	ALA	2.5
3	A	636	VAL	2.5
3	A	254	GLU	2.4
3	A	639	SER	2.2
3	A	130	LYS	2.1
3	A	121	ASP	2.1
3	A	897	LEU	2.1
3	A	817	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
3	A	302	LYS	2.0

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^{th} 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
2	DOC	P	115	18/19	0.98	0.20	-	32,35,39,40	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^{th} 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
6	TTP	A	1006	29/29	0.97	0.08	-4.22	29,33,40,41	0
4	CA	A	1001	1/1	0.97	0.10	-4.85	56,56,56,56	0
5	CL	A	1005	1/1	0.83	0.35	-	108,108,108,108	0
4	CA	A	1002	1/1	0.94	0.13	-	101,101,101,101	0
4	CA	A	1004	1/1	0.90	0.06	-	85,85,85,85	0
4	CA	A	1003	1/1	0.84	0.20	-	94,94,94,94	0

6.5 Other polymers (i)

There are no such residues in this entry.